

WEEKLY REPORT FOR WEEK 38 (APRIL 21, 2019 – APRIL 27, 2019)

Report No. 53005-81-RPT-055

September 2019

Prepared for:

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Subcontract 53005, Release 81

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**Weekly Report for Week 38
(April 21, 2019 – April 27, 2019)**

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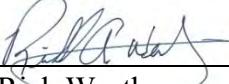
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Date: 09/19/2019

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Record of Revision

Revision	Date	Pages/Sections Changed	Brief Description
0	09/2019	All	Original Issue.

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Acronyms and Abbreviations

AOP	Abnormal Operating Procedure
COC	Chain of Custody
COPC	Chemical of Potential Concern
CSM	Central Shift Manager
CSO	Central Shift Office
IH	Industrial Hygiene
MDL	Method Detection Limit
ML	Mobile Laboratory
NDEA	N-nitrosodiethylamine
NDMA	N-nitrosodimethylamine
NEMA	N-nitrosomethylethylamine
NMOR	N-nitrosomorpholine
PTR-TOF	Proton Transfer Reaction – Time-of-Flight
PTR-MS	Proton Transfer Reaction – Mass Spectrometer
QA	Quality Assurance
QC	Quality Control
R&D	Research and Development
RL	Reporting Limit
SME	Subject Matter Expert
VOC	Volatile Organic Compound
WRPS	Washington River Protection Solutions, LLC

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1.0 INTRODUCTION

During the week of April 22, 2019, through April 28, 2019, the Mobile Laboratory (ML) performed area monitoring around the 200 East Area of the Hanford Site, as well as source characterization in support of the Washington River Protection Solutions, LLC (WRPS) Fugitive Emissions Team. The data team continued processing data collected from the previous week. The reporting team worked towards the completion of weekly reports for Weeks 33 through 37 and monthly reports.

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2.0 APRIL 22, 2019 – AREA MONITORING AND AOP-015 EVENT

2.1 Quality Assessment

Data from April 22, 2019, were assessed using Procedure 17124-DOE-HS-102, “Mobile Laboratory Data Processing – Analysis.” A Data Exchange Checklist was completed. The data were accepted by TerraGraphics with the following comments. Report No. 66409-RPT-004, *Mobile Laboratory Operational Procedure*, was adequately documented and all checks passed the acceptance limits.

2.2 Summary

On April 22, 2019, ML personnel arrived at the TerraGraphics warehouse in Pasco, WA at 04:50. The Quality Assurance/Quality Control (QA/QC) zero-air/span checks were performed on the LI-COR^{®1} CO₂ monitor, the Picarro NH₃ analyzer, and the Proton Transfer Reaction – Mass Spectrometer (PTR-MS) beginning at 05:08. The ML arrived on the Hanford Site and checked in with the Central Shift Manager (CSM) at 06:01. The ML began a site survey loop of A Farms at 06:12. The ML was parked at the southeast corner of A Farm prior to starting a site survey loop at 07:30. The ML Operators proceeded to park near the northwest side of A Farm and remained there until 09:18, when an Abnormal Operating Procedure (AOP)-015 occurred. The Operators reported to the Central Shift Office (CSO) for a pre-job briefing and headed to U Farm, the location of the AOP-015, at 09:37. The ML arrived at U Farm and completed set-up of the 208-foot heated line at 09:54. At 10:17, the Industrial Hygiene (IH) Technicians moved the 208-foot line as close to the change trailer as possible. At 14:02, the AOP-015 was exited and Operators began the process to disconnect the 208-foot line. At 14:40, operators checked out with the CSM and the ML departed the site. After stopping to pick up a Swagelok^{®2} order, the ML arrived back at the TerraGraphics warehouse at 15:50.

¹ LI-COR is a registered trademark of LI-COR, Inc., Lincoln, Nebraska.

² Swagelok is a registered trademark of Swagelok Company, Solon, Ohio.

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Figure 2-1. Location of the Mobile Laboratory for the Duration of the Monitoring Period.

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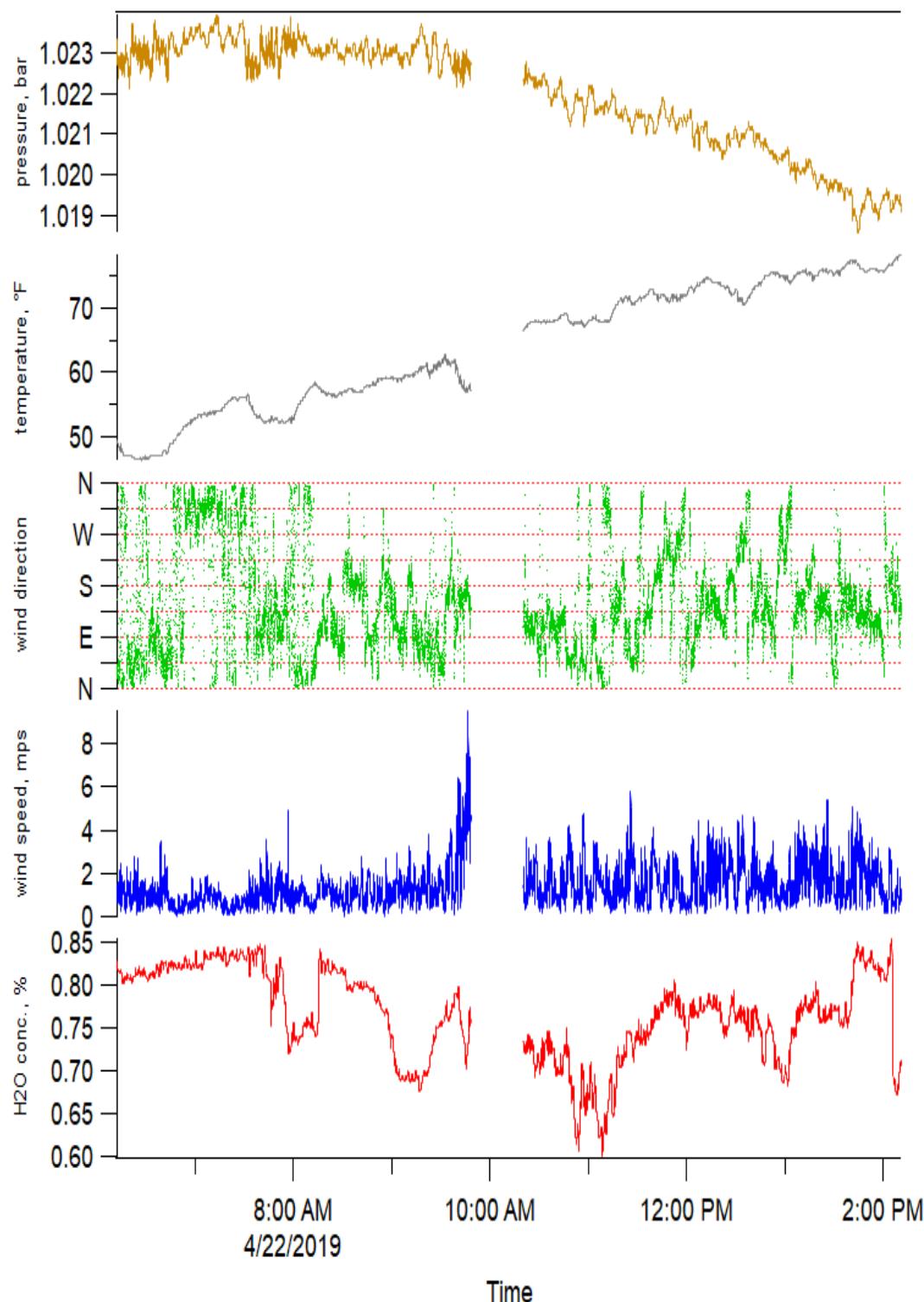


Figure 2-2. Weather Data for the Duration of the Monitoring Period.

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2.3 Samples Collected

Continuous air monitoring was performed using the following instrumentation:

- Proton Transfer Reaction – Time-of-Flight (PTR-TOF) 6000 X2,
- LI-COR CO₂ Monitor,
- Picarro Ammonia Monitor, and
- Airmar[®]³ Weather Station.

Confirmatory air samples were not collected during this period.

2.4 Area Monitoring

The ML performed area monitoring from approximately 06:12 to 09:18. The section below displays the chemical of potential concern (COPC) statistical results during that monitoring period.

³ Airmar is a registered trademark of Airmar Technology Corporation, Milford, New Hampshire.

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Table 2-1. Chemical of Potential Concern Statistical Information for the Area Monitoring on April 22, 2019. (2 Sheets)

COPC #	COPC Name	OEL (ppb)	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel St. Dev. (%)	Max. (ppb)	Median (ppb)
1	ammonia	25000	6.225	15.117†	2.501	16.546	19.967	14.926†
2	formaldehyde	300	1.721	<1.721	0.160	16.728	2.069	<1.721
3	methanol	200000	5.758	9.093†	1.309	14.401	31.024	8.944†
4	acetonitrile	20000	0.085	0.115†	0.035	30.372	0.690	0.113†
5	acetaldehyde	25000	1.027	1.868†	0.286	15.290	4.132	1.848†
6	ethylamine	5000	0.069	<0.069	0.022	60.293	0.149	<0.069
7	1,3-butadiene	1000	0.183	0.243†	0.076	31.323	0.764	0.237†
8	propanenitrile	6000	0.107	<0.107	0.029	41.530	0.226	<0.107
9	2-propenal	100	0.340	<0.340	0.059	44.631	0.800	<0.340
10	1-butanol + butenes	20000	0.214	<0.214	0.051	38.302	0.684	<0.214
11	methyl isocyanate	20	0.069	<0.069	0.029	51.972	0.190	<0.069
12	methyl nitrite	100	0.098	0.122†	0.046	37.793	1.081	0.120†
13	furan	1	0.062	<0.062	0.020	51.288	0.142	<0.062
14	butanenitrile	8000	0.039	<0.039	0.020	59.178	0.133	<0.039
15	but-3-en-2-one + 2,3-dihydrofuran + 2,5-dihydrofuran	200, 1, 1	0.041	0.059†	0.024	41.597	N/A*	N/A*
16	butanal	25000	0.061	0.181†	0.044	24.374	0.507	0.179†
17	NDMA**	0.3	0.082	<0.082	0.024	170.407	0.181	<0.082
18	benzene	500	0.236	<0.236	0.062	35.619	1.087	<0.236
19	2,4-pentadienenitrile + pyridine	300, 1000	0.085	<0.085	0.021	41.818	0.145	<0.085
20	2-methylene butanenitrile	300	0.036	<0.036	0.013	66.048	0.086	<0.036
21	2-methylfuran	1	0.043	0.047†	0.025	53.309	0.199	0.044†
22	pentanenitrile	6000	0.036	<0.036	0.015	71.669	0.102	<0.036
23	3-methyl-3-butene-2-one + 2-methyl-2-butenal	20, 30	0.043	0.051†	0.027	52.213	0.207	0.048†
24	NEMA**	0.3	0.058	<0.058	0.023	147.645	0.181	<0.058
25	2,5-dimethylfuran	1	0.032	<0.032	0.020	68.511	0.156	<0.032
26	hexanenitrile	6000	0.031	<0.031	0.012	84.308	0.080	<0.031
27	2-hexanone (MBK)	5000	0.036	<0.036	0.018	75.365	0.115	<0.036
28	NDEA**	0.1	0.034	<0.034	0.014	144.283	0.096	<0.034
29	butyl nitrite + 2-nitro-2-methylpropane	100, 300	0.058	<0.058	0.016	47.429	0.111	<0.058

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Table 2-1. Chemical of Potential Concern Statistical Information for the Area Monitoring on April 22, 2019. (2 Sheets)

COPC #	COPC Name	OEL (ppb)	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel St. Dev. (%)	Max. (ppb)	Median (ppb)
30	2,4-dimethylpyridine	500	0.036	<0.036	0.014	89.652	0.186	<0.036
31	2-propylfuran + 2-ethyl-5-methylfuran	1	0.027	<0.027	0.015	90.089	0.108	<0.027
32	heptanenitrile	6000	0.027	<0.027	0.010	88.437	0.071	<0.027
33	4-methyl-2-hexanone	500	0.033	<0.033	0.014	81.695	0.097	<0.033
34	NMOR**	0.6	0.021	<0.021	0.010	171.498	0.073	<0.021
35	butyl nitrate	2500	0.022	<0.022	0.009	117.865	0.060	<0.022
36	2-ethyl-2-hexenal + 4-(1-methylpropyl)-2,3-dihydrofuran + 3-(1,1-dimethylethyl)-2,3-dihydrofuran	100, 1, 1	0.028	<0.028	0.013	82.677	0.087	<0.028
37	6-methyl-2-heptanone	8000	0.028	<0.028	0.012	78.887	0.094	<0.028
38	2-pentylfuran	1	0.026	<0.026	0.013	78.273	0.081	<0.026
39	biphenyl	200	0.022	<0.022	0.010	122.375	0.070	<0.022
40	2-heptylfuran	1	0.067	<0.067	0.017	38.417	0.127	<0.067
41	1,4-butanediol dinitrate	50	0.036	<0.036	0.012	81.392	0.075	<0.036
42	2-octylfuran	1	0.020	<0.02	0.010	199.378	0.069	<0.02
43	1,2,3-propanetriol 1,3-dinitrate	50	0.011	<0.011	0.005	354.338	0.062	<0.011
44	PCB	1000	0.034	<0.034	0.011	65.590	0.090	<0.034
45	6-(2-furanyl)-6-methyl-2-heptanone	1	0.025	<0.025	0.009	92.834	0.059	<0.025
46	furfural acetophenone	1	0.064	<0.064	0.017	39.118	0.117	<0.064
N/A*	The maximum peak value for but-3-en-2-one + 2,3 dihydrofuran + 2,5 dihydrofuran was 0.263 ppb and the median value was 0.056† ppb. The PTR-MS results for but-3-en-2-one + 2,3 dihydrofuran + 2,5 dihydrofuran are not compared to OEL concentrations because: 1) the result is suspect due to a known biogenic interferant (methacrolein) that is expected to be in concentrations that occasionally exceed the dihydrofuran OEL, and 2) this combination of COPCs have OEL concentrations that differ by a factor of 200, which provide widely variant bases for these numbers.							
**	Nitrosamine results are suspect due to isobaric interferants causing positive bias that have been encountered during previous background [53005-81-RPT-007, PTR-MS Mobile Laboratory Vapor Monitoring Background Study, (3/18/2018 – 4/20/2018), and Fiscal Year 2017 Mobile Laboratory Vapor Monitoring at the Hanford Site: Monitoring During Waste Disturbing Activities and Background Study, RJ Lee Group, Inc.].							
<	COPC Averages below the MDL.							
†	COPC Averages between the RL and the MDL.							
	COPC Averages >100% of the OEL.							
	COPC Averages 50-100% of the OEL.							
	COPC Averages 10-50% of the OEL.							

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Figure 2-3 through Figure 2-51 display 46 COPC signals, overlaid with the same signal smoothed using a 1-minute moving average (in cases where a moving average assists with data visualization), and CO₂, for the monitoring period April 22, 2019. If within range of the plot's left axis, a green horizontal line representing 50% of the COPC's OEL, a blue horizontal line representing the COPC's OEL, a horizontal purple line representing the Reporting Limit (RL), and a pink horizontal line representing the Method Detection Limit (MDL) are shown.

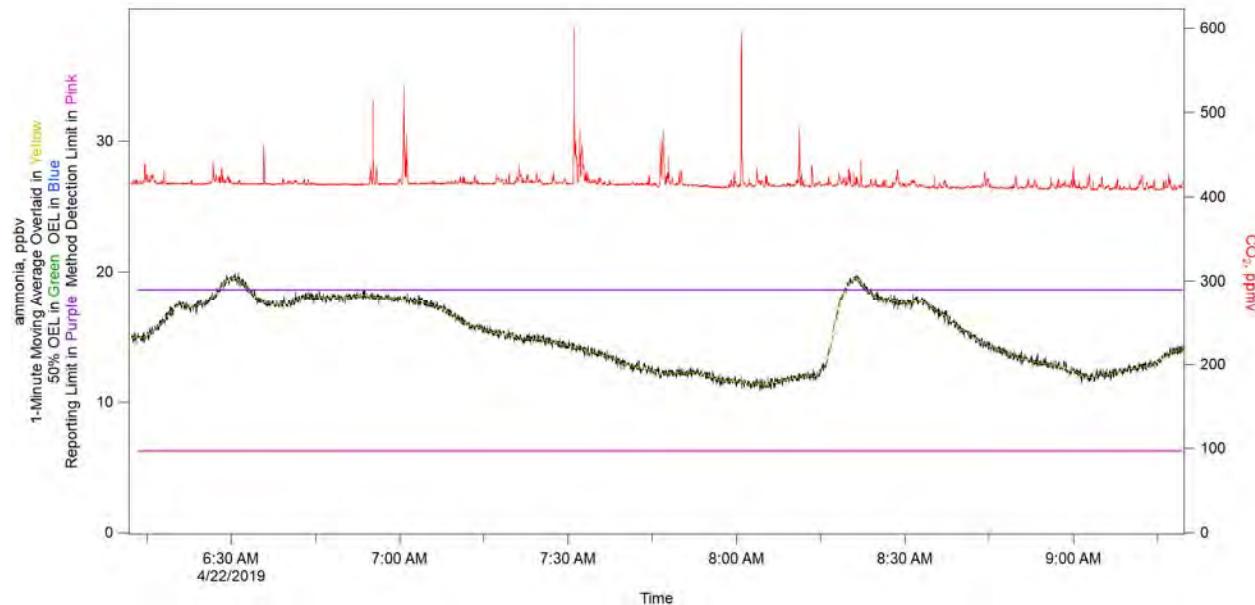


Figure 2-3. Ammonia.

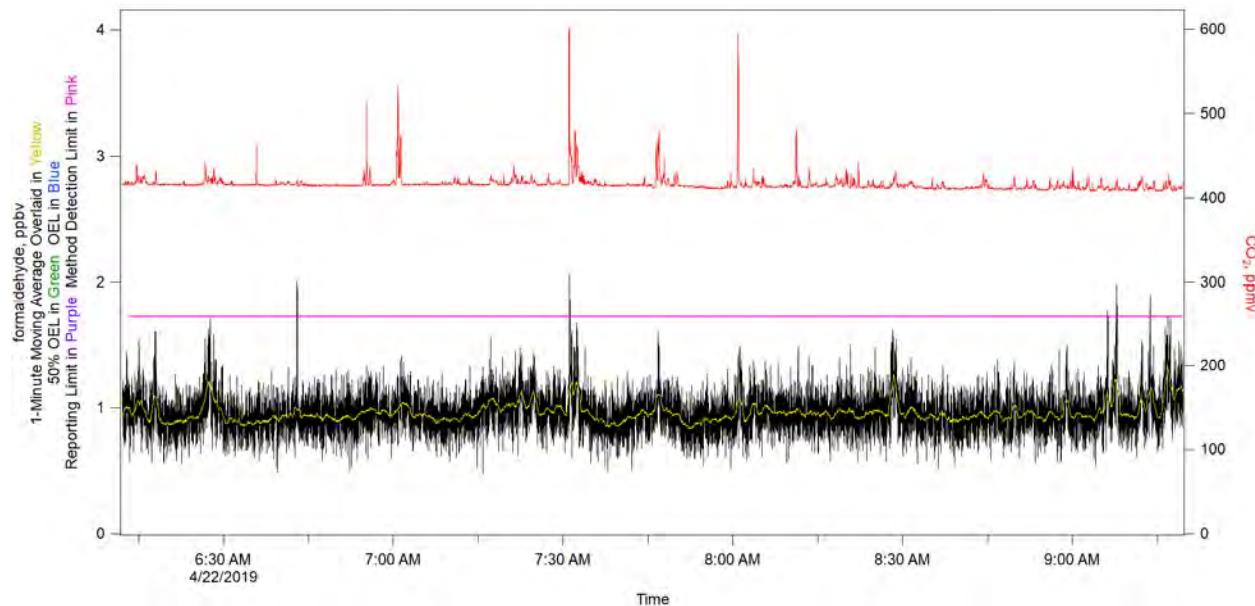


Figure 2-4. Formaldehyde.

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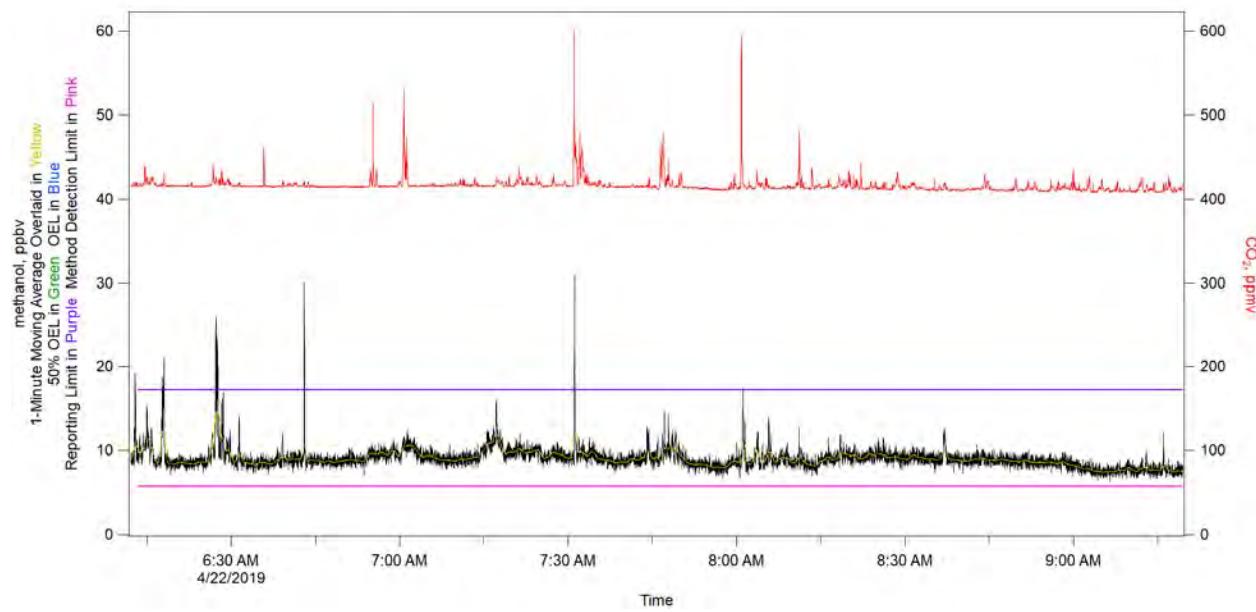


Figure 2-5. Methanol.

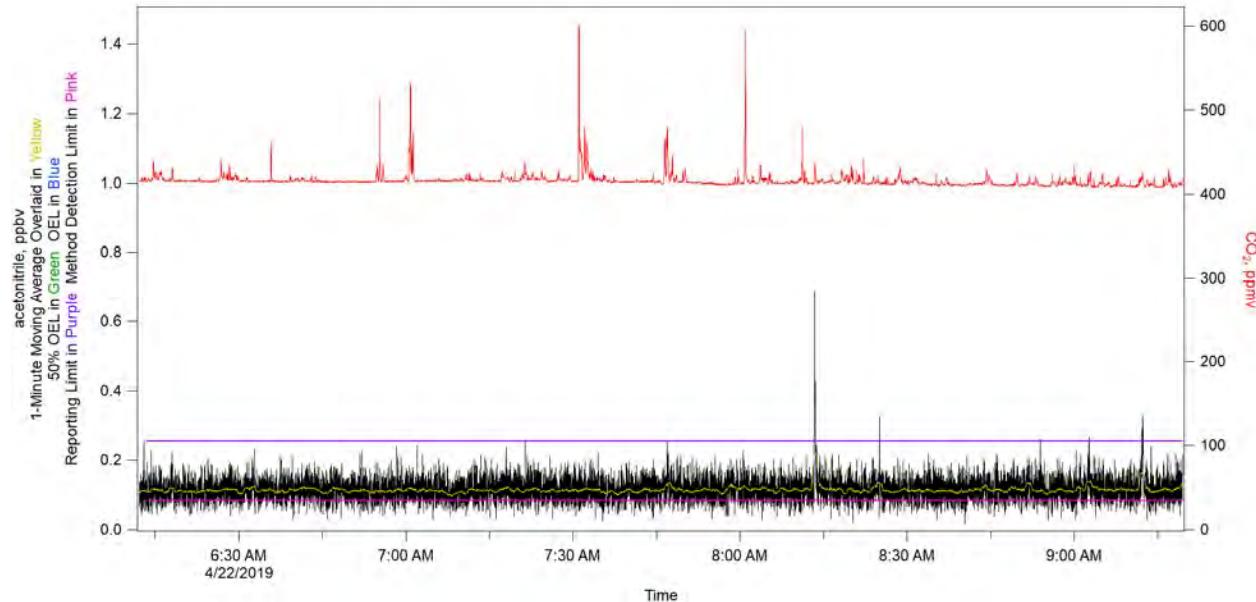


Figure 2-6. Acetonitrile.

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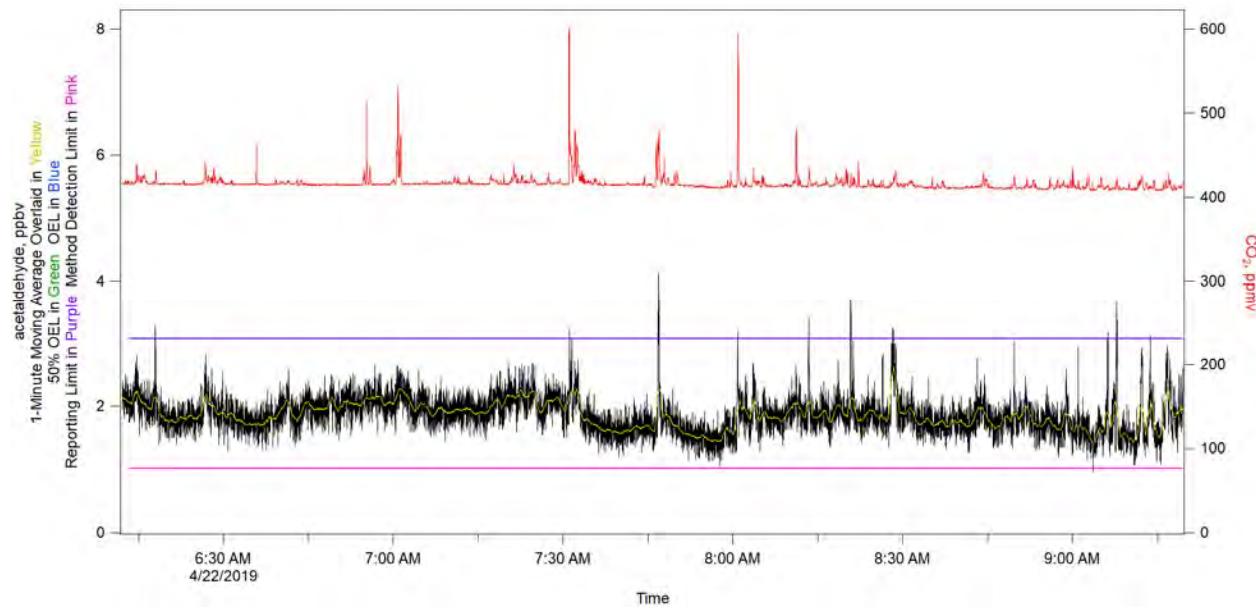


Figure 2-7. Acetaldehyde.

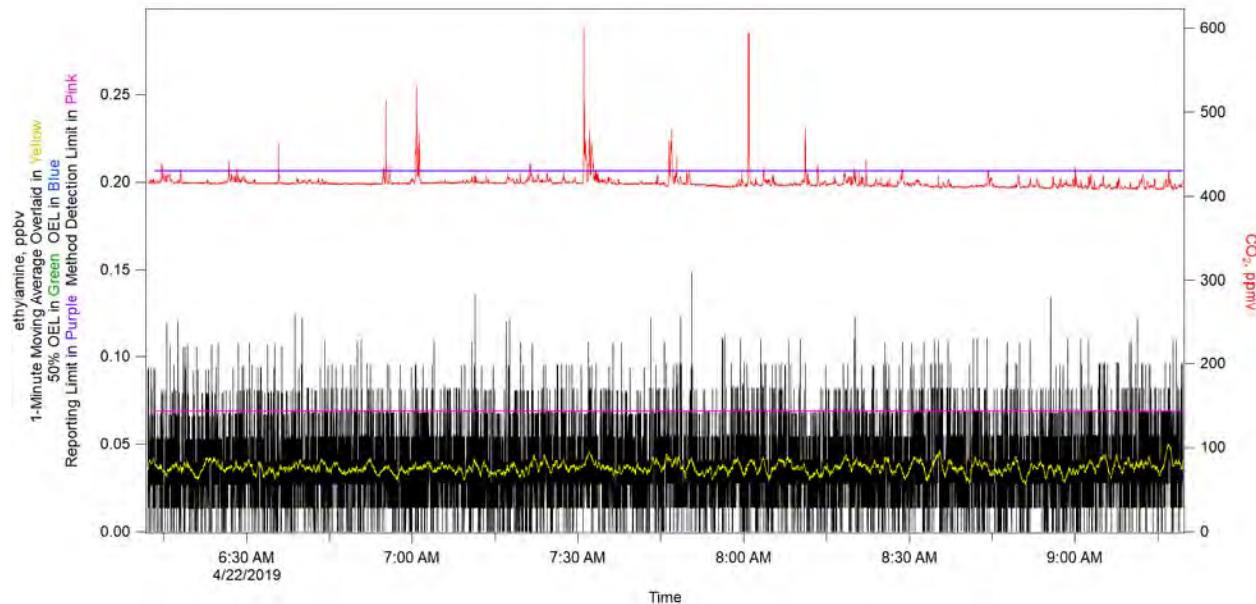


Figure 2-8. Ethylamine.

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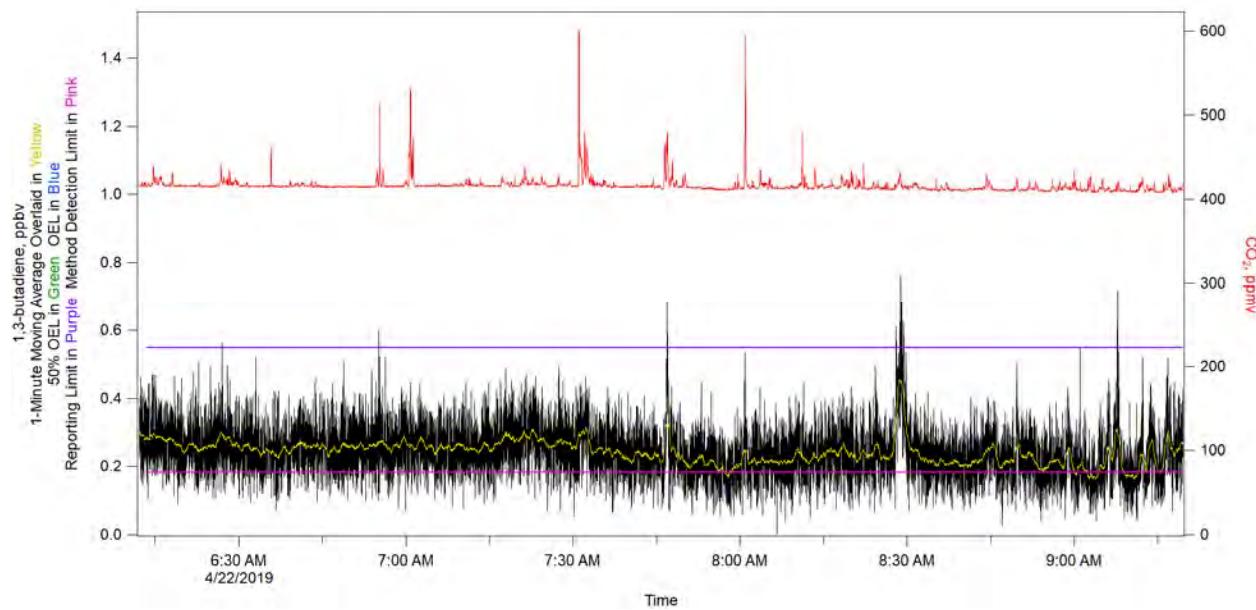


Figure 2-9. 1,3-butadiene.

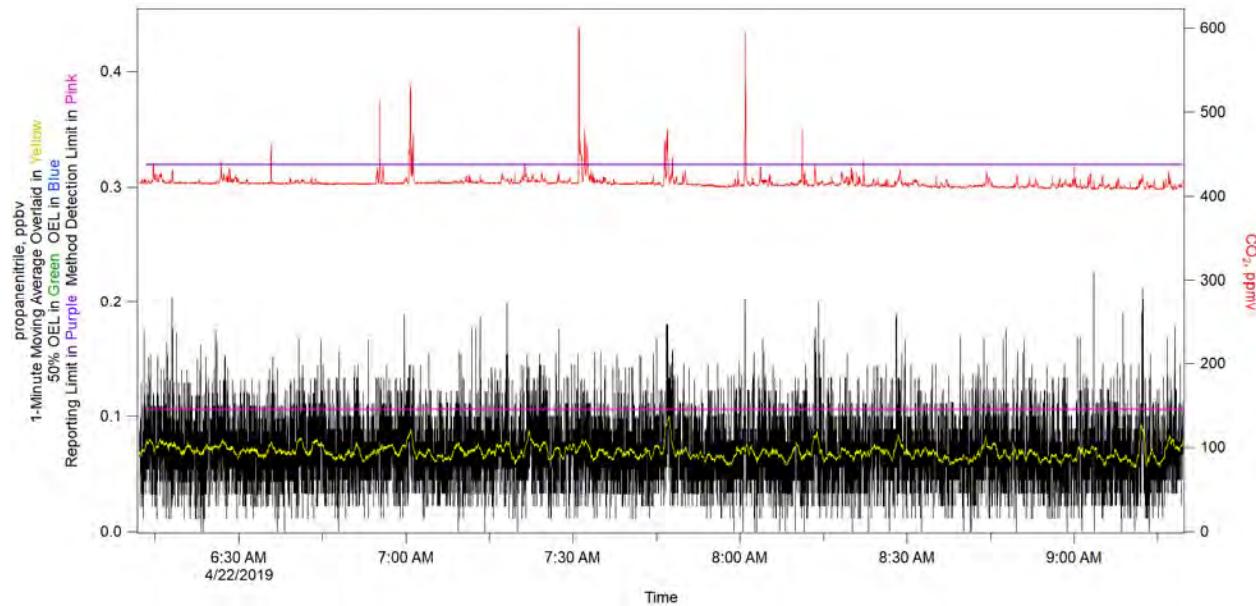


Figure 2-10. Propanenitrile.

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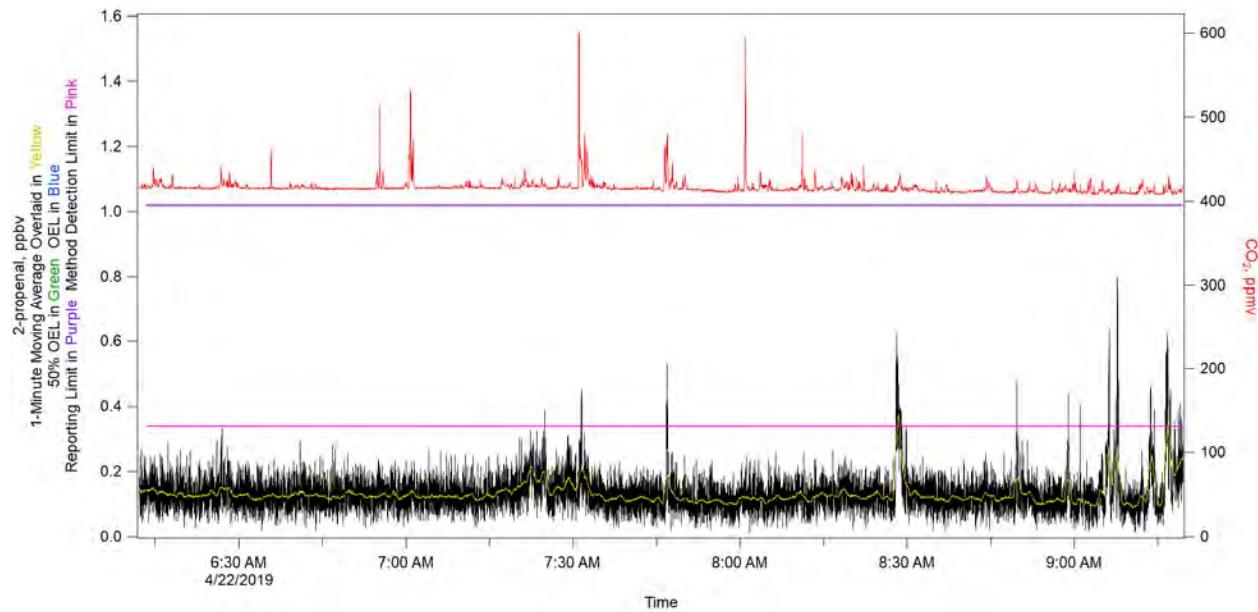


Figure 2-11. 2-propenal.

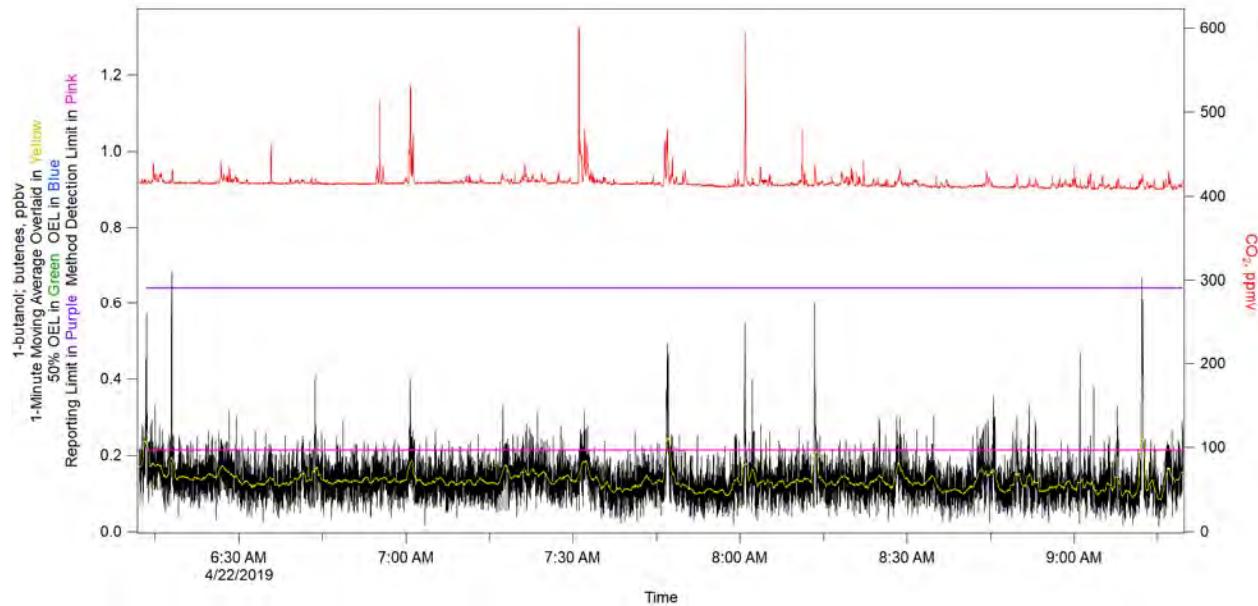


Figure 2-12. 1-butanol; Butenes.

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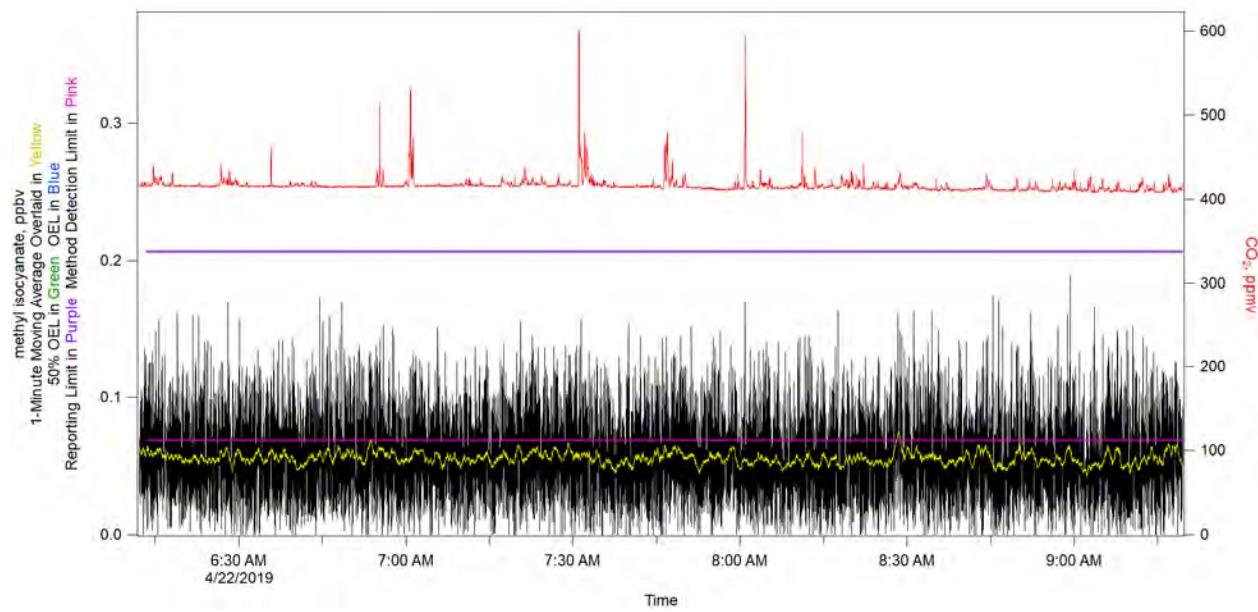


Figure 2-13. Methyl Isocyanate.

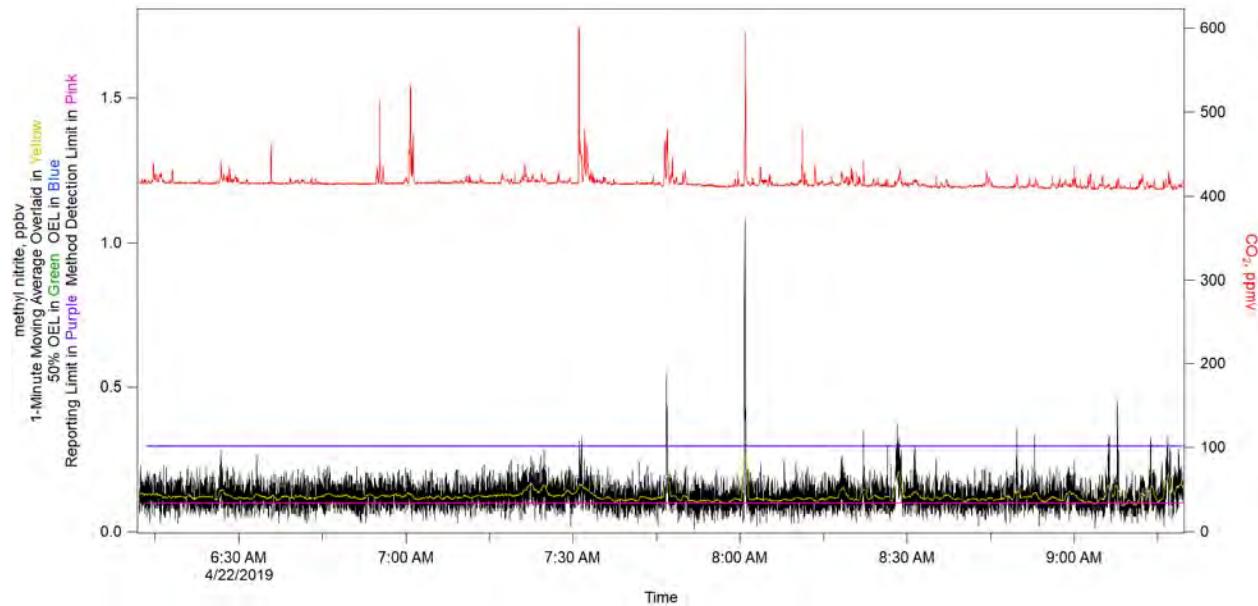


Figure 2-14. Methyl Nitrite.

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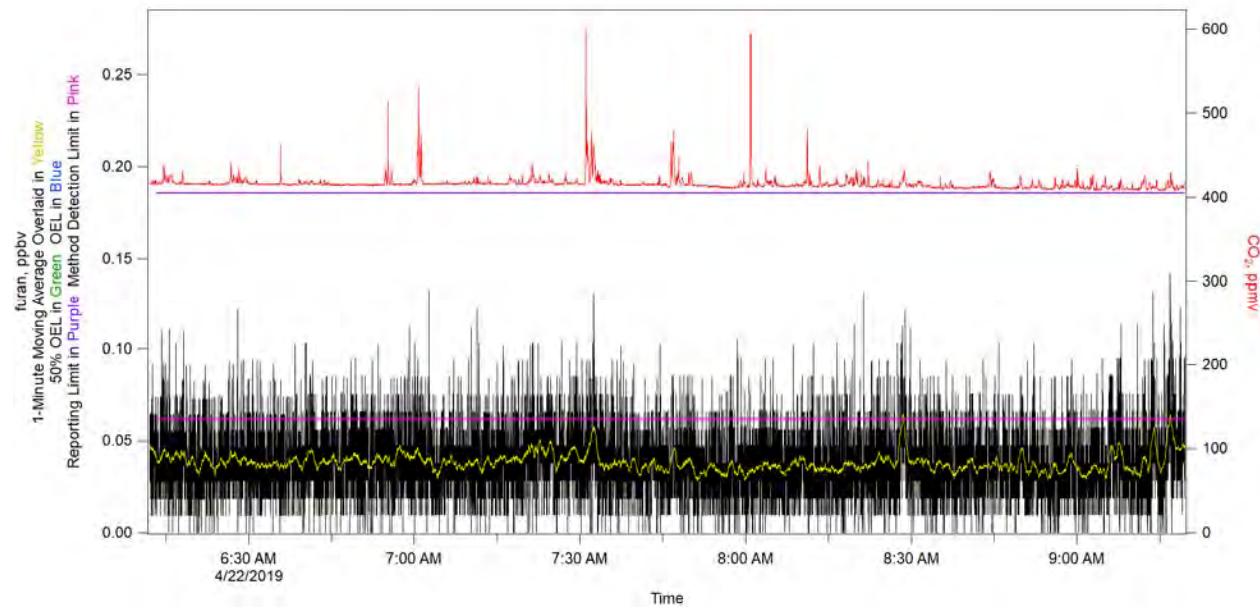


Figure 2-15. Furan.

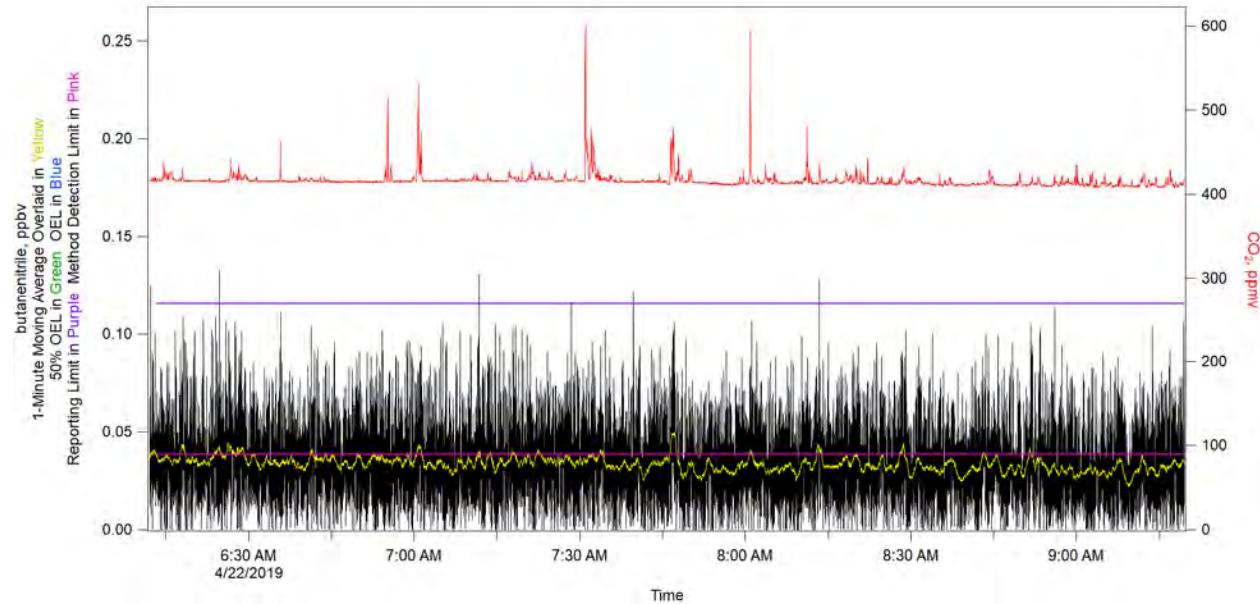


Figure 2-16. Butanenitrile.

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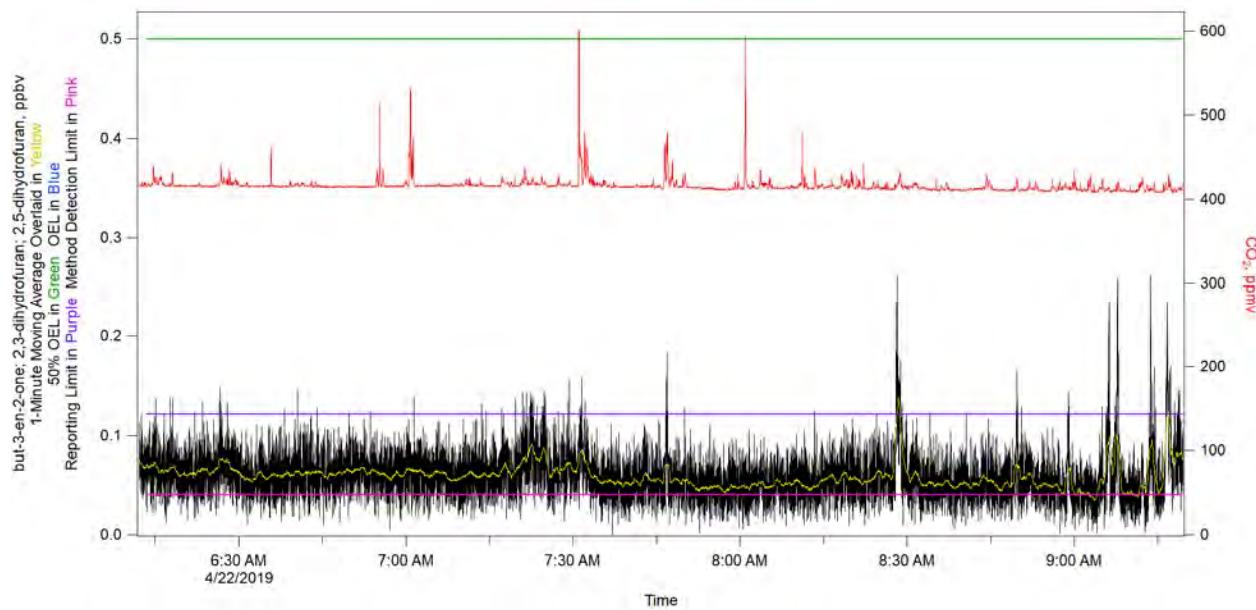


Figure 2-17. But-3-en-2-one; 2,3-dihydrofuran; 2,5-dihydrofuran.

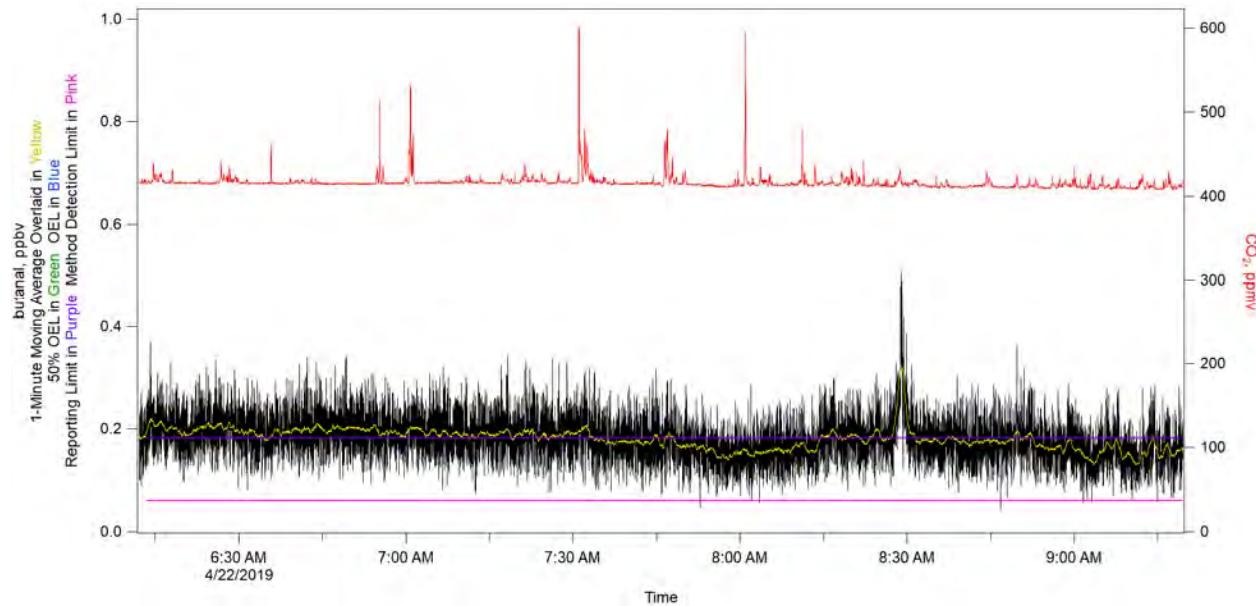


Figure 2-18. Butanal.

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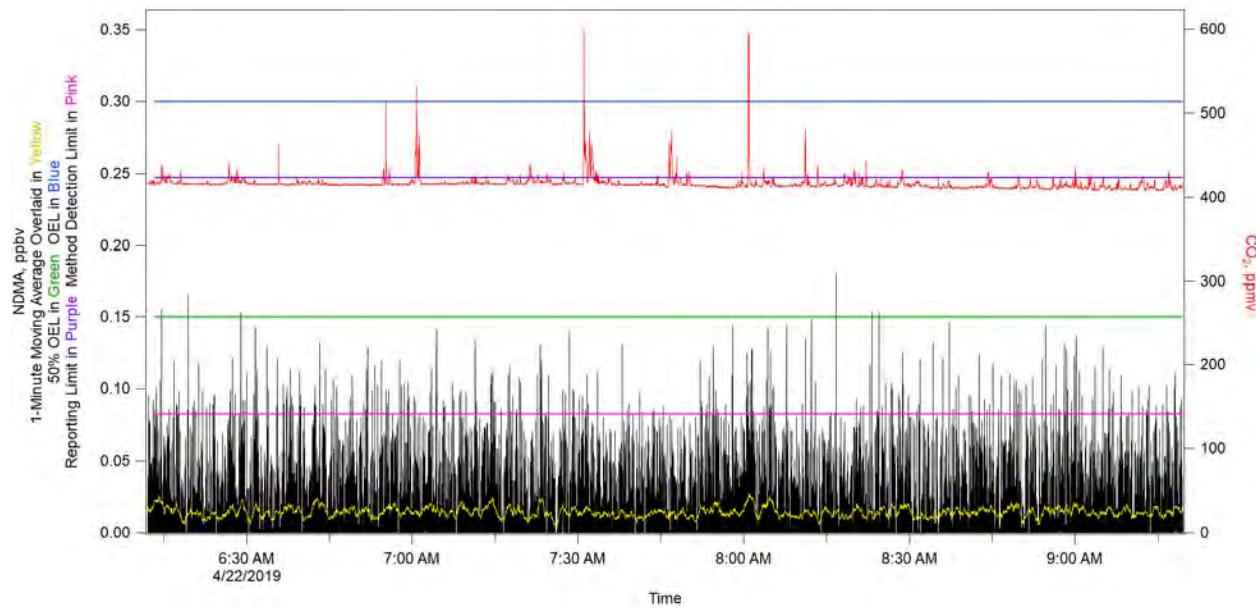


Figure 2-19. N-nitrosodimethylamine (NDMA).

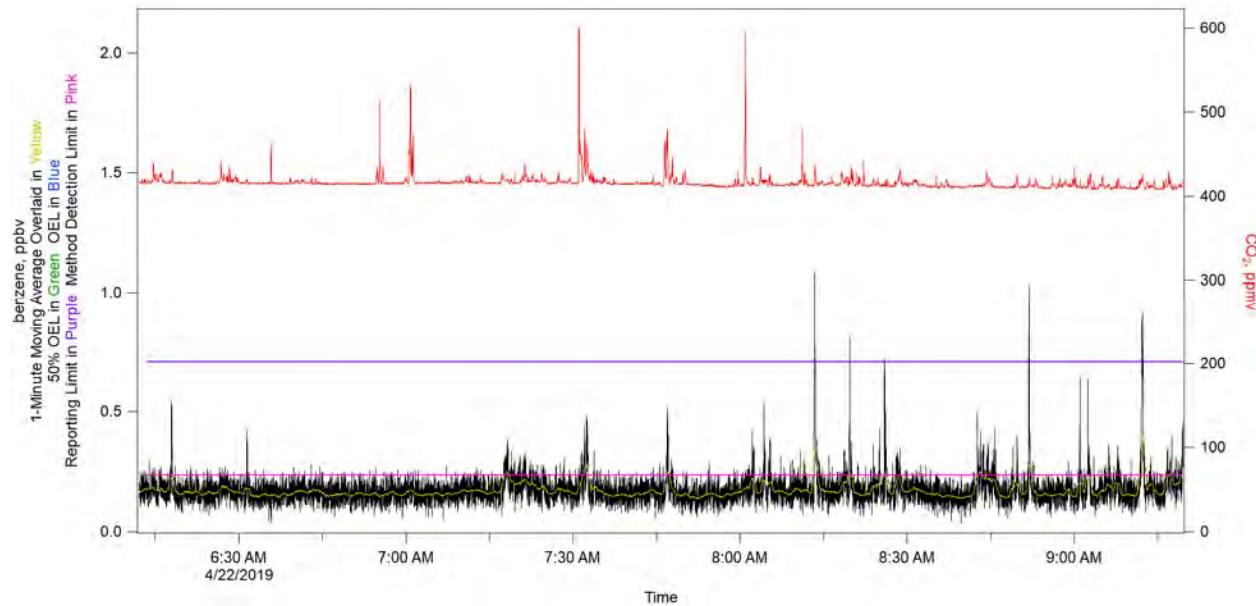


Figure 2-20. Benzene.

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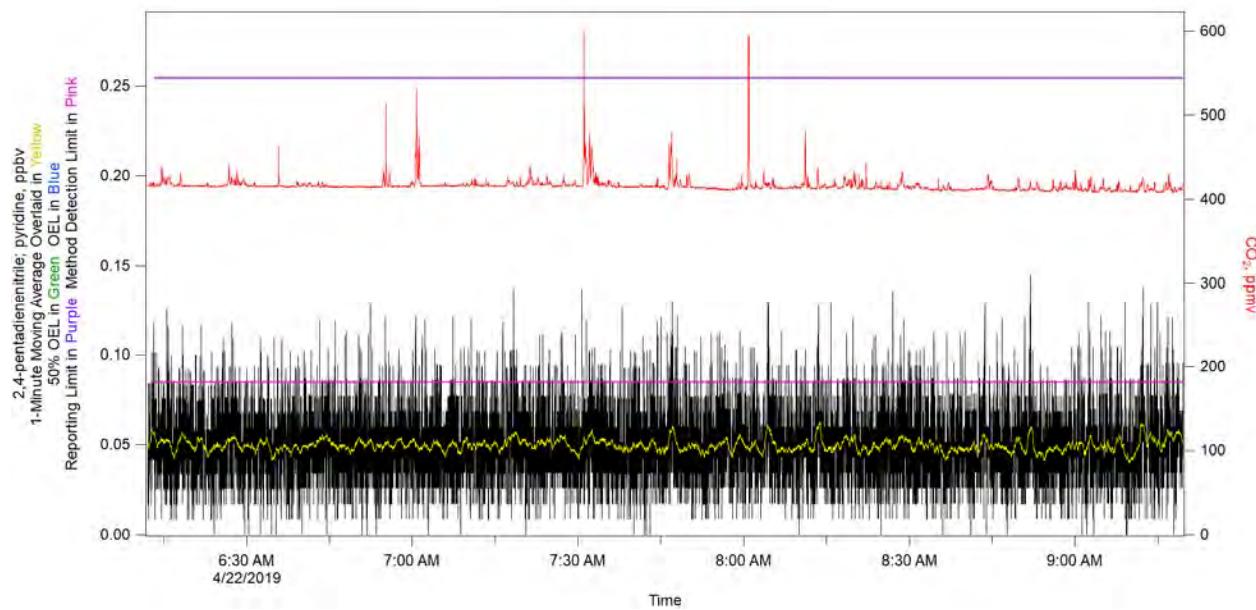


Figure 2-21. 2,4-pentadienenitrile; Pyridine.

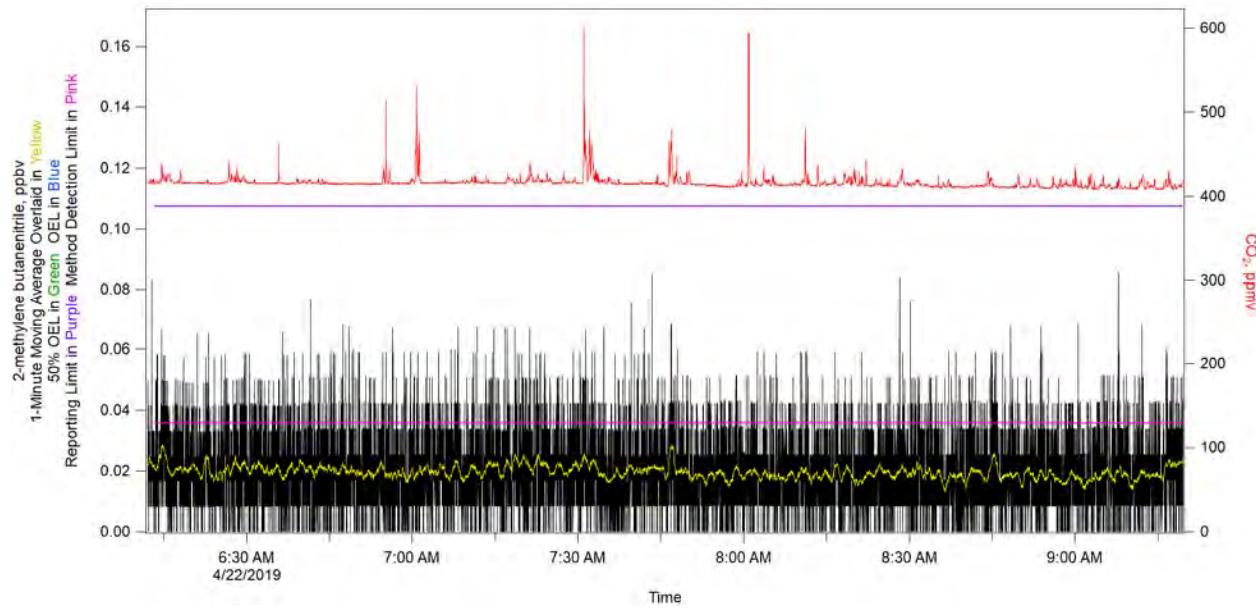


Figure 2-22. 2-methylene Butanenitrile.

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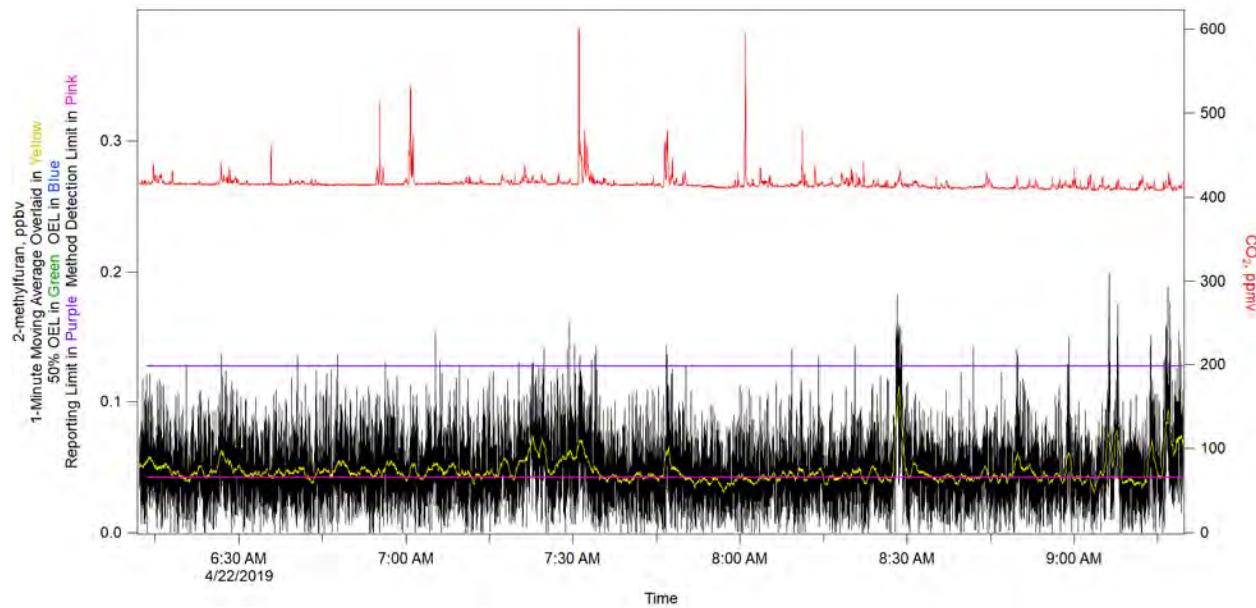


Figure 2-23. 2-methylfuran.

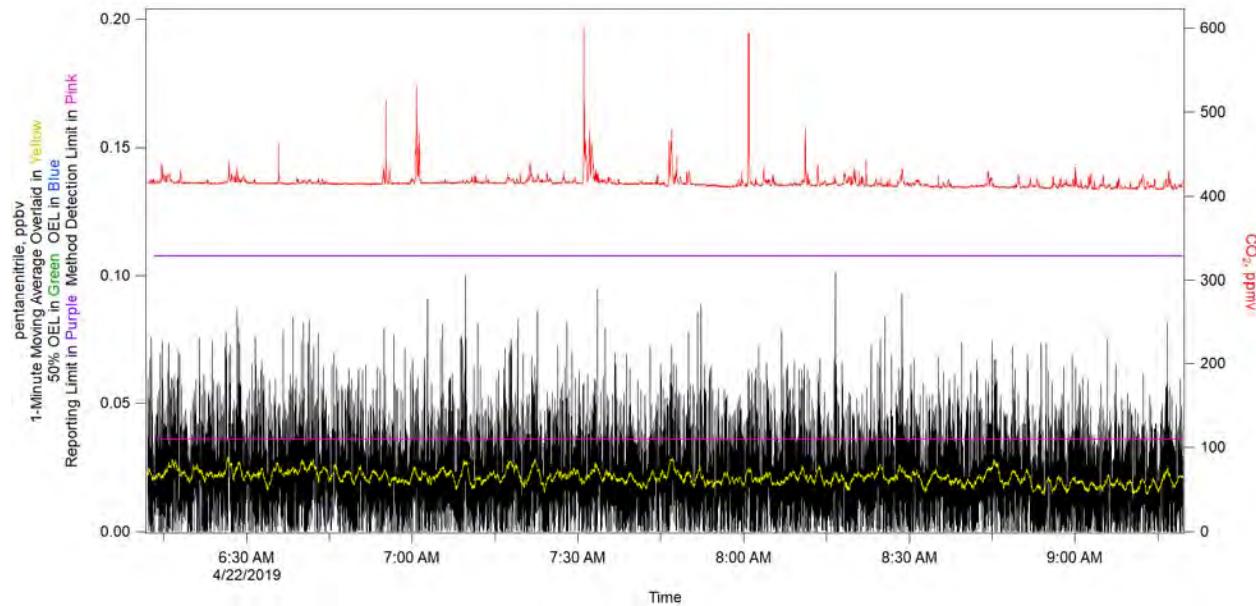


Figure 2-24. Pentanenitrile.

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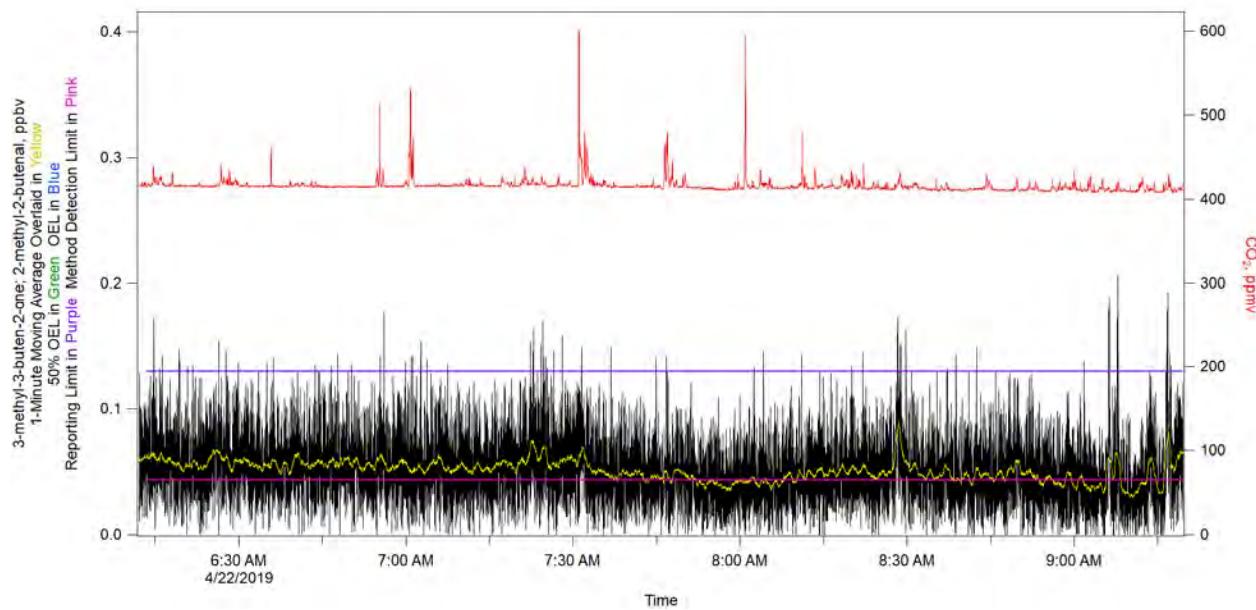


Figure 2-25. 3-methyl-3-buten-2-one; 2-methyl-2-butenal.

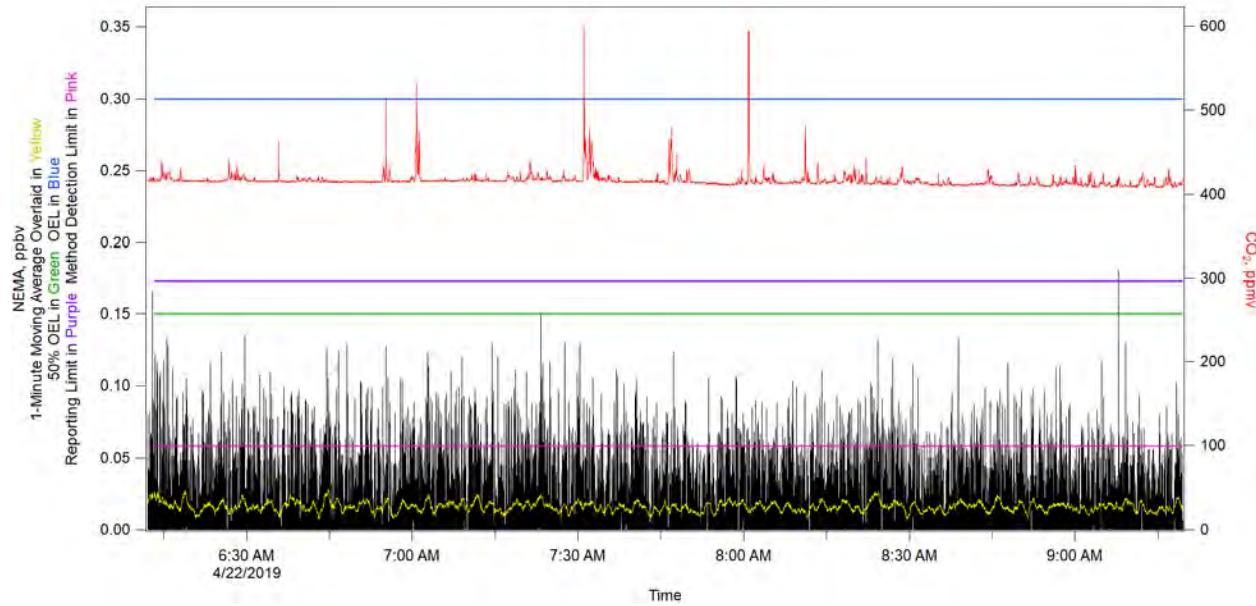


Figure 2-26. N-nitrosomethylethylamine (NEMA).

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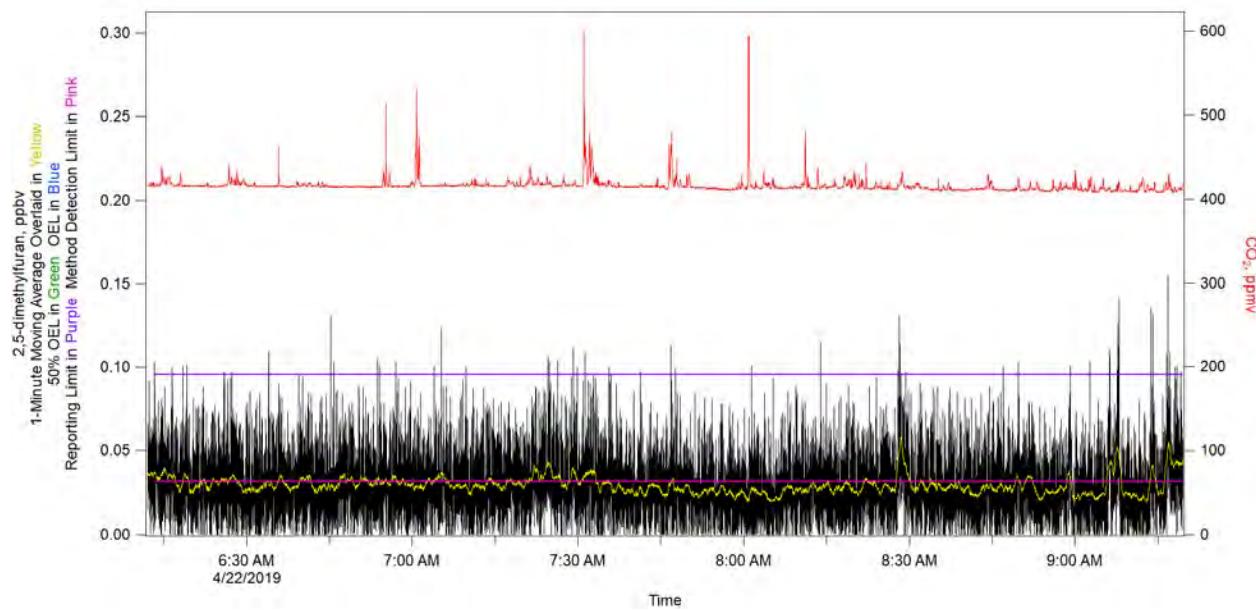


Figure 2-27. 2,5-dimethylfuran.

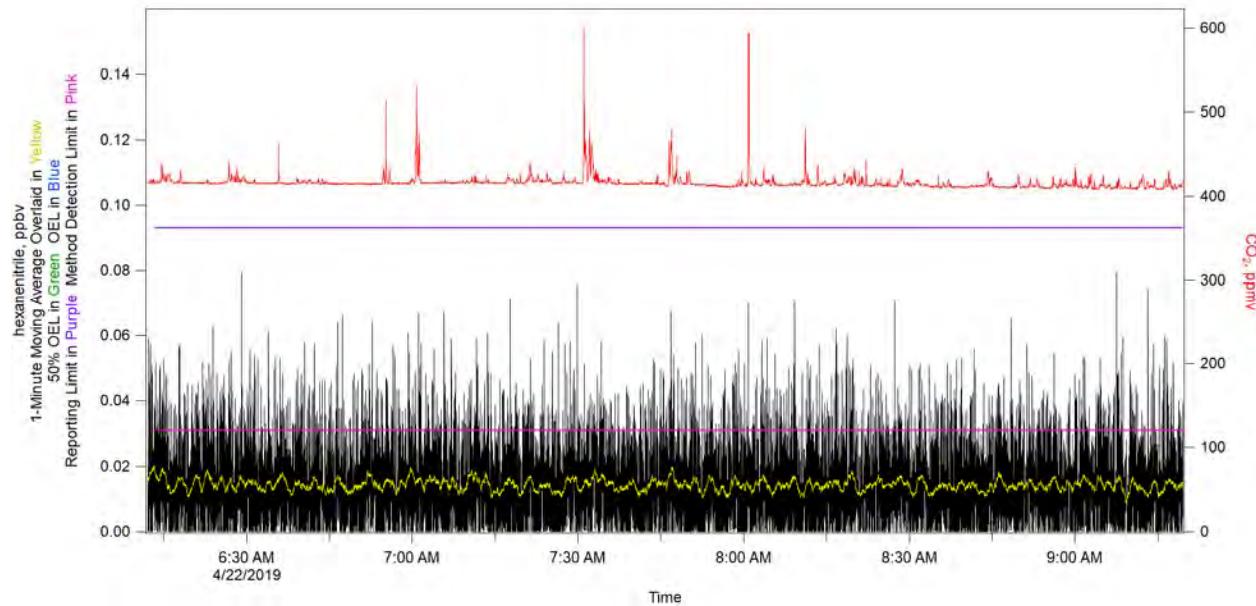


Figure 2-28. Hexanenitrile.

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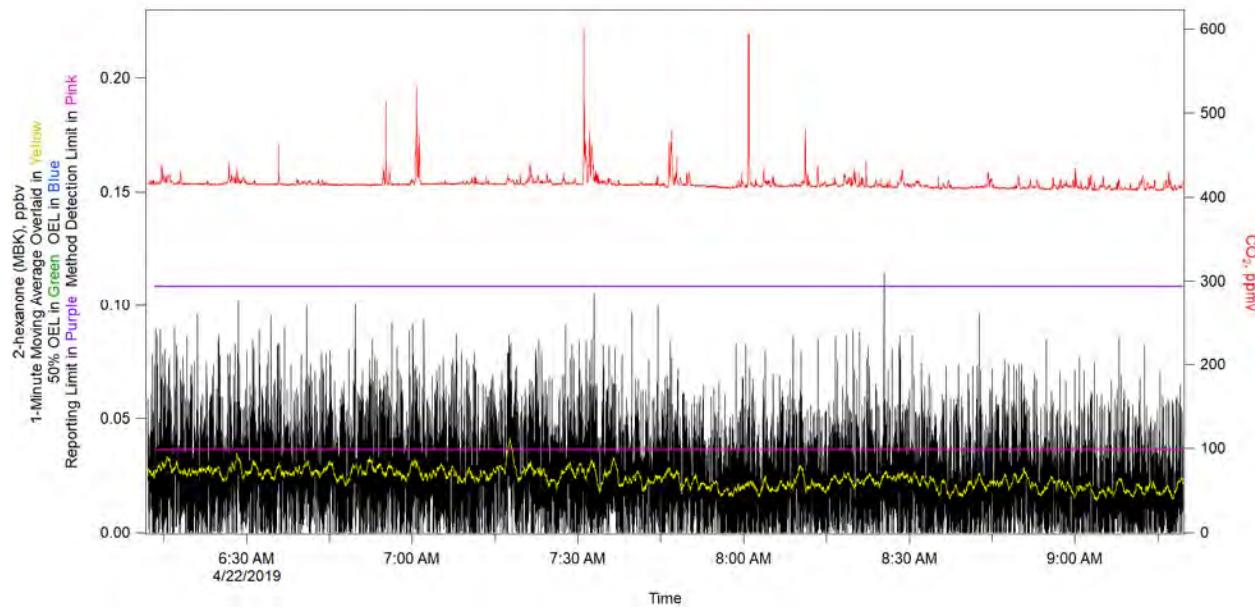


Figure 2-29. 2-hexanone (MBK).

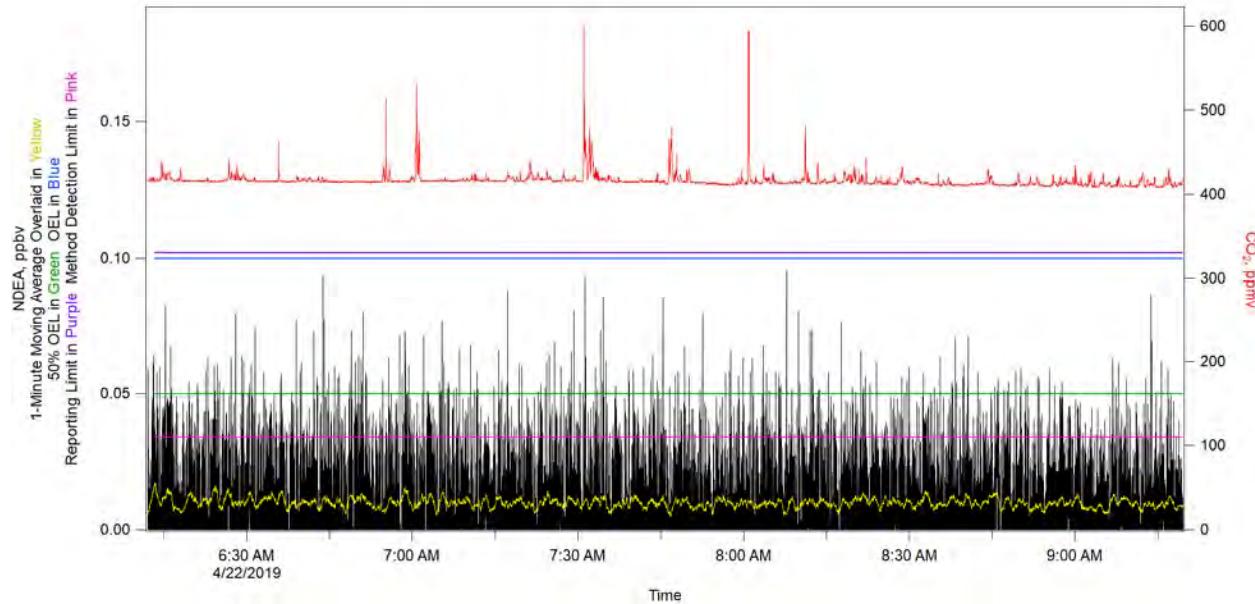


Figure 2-30. N-nitrosodiethylamine (NDEA).

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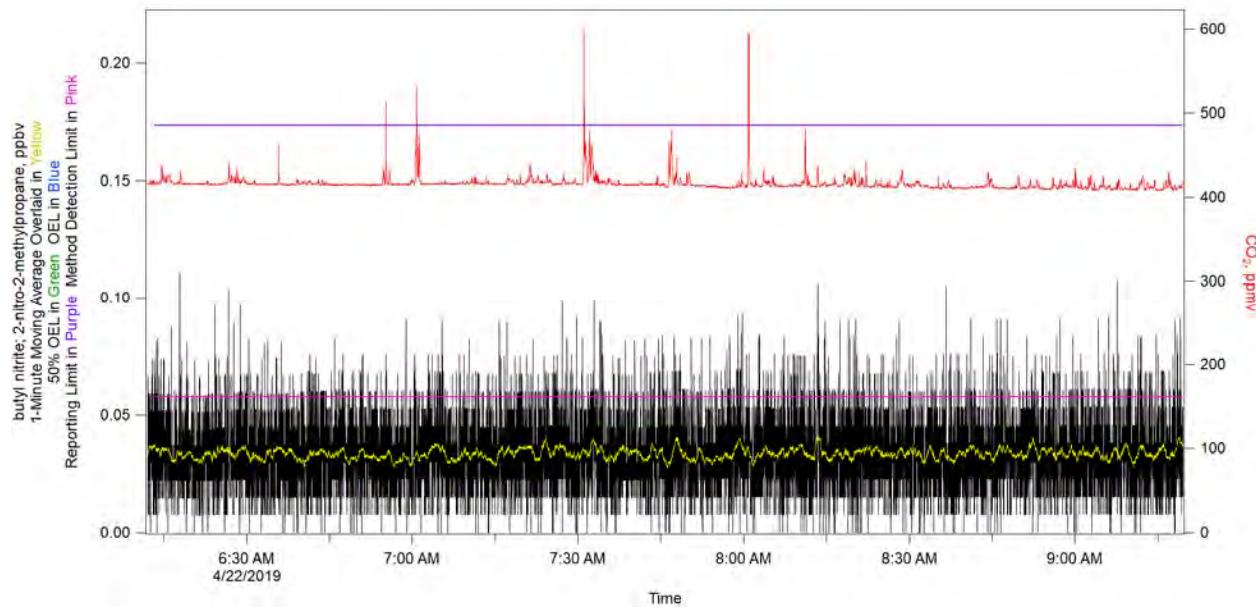


Figure 2-31. Butyl Nitrite; 2-nitro-2-methylpropane.

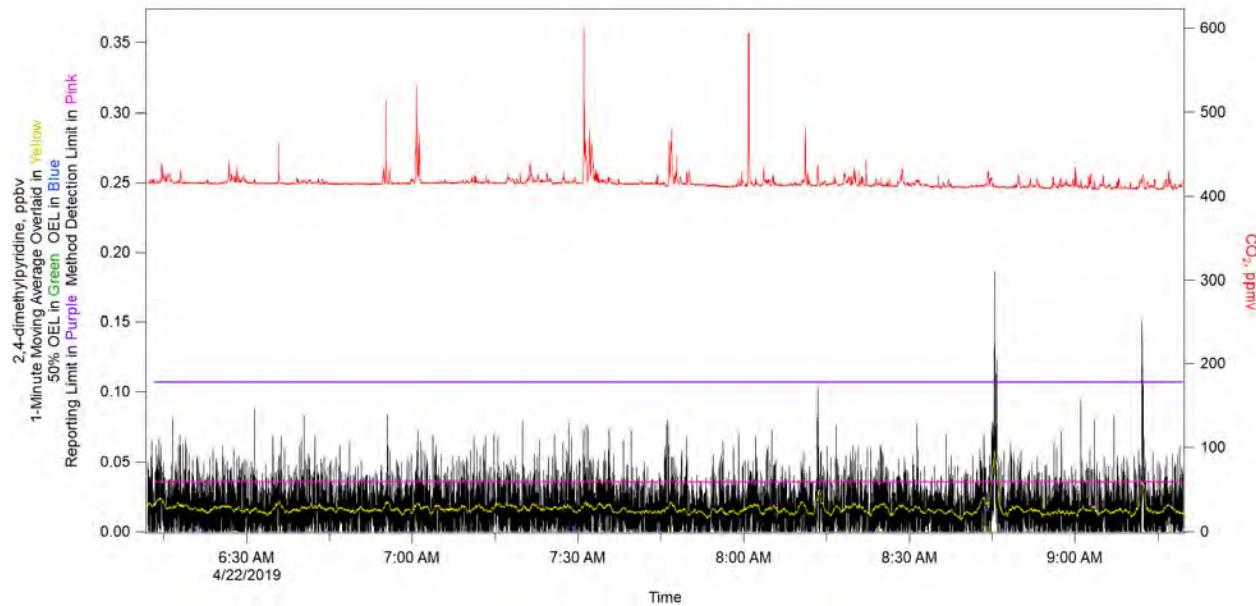


Figure 2-32. 2,4-dimethylpyridine.

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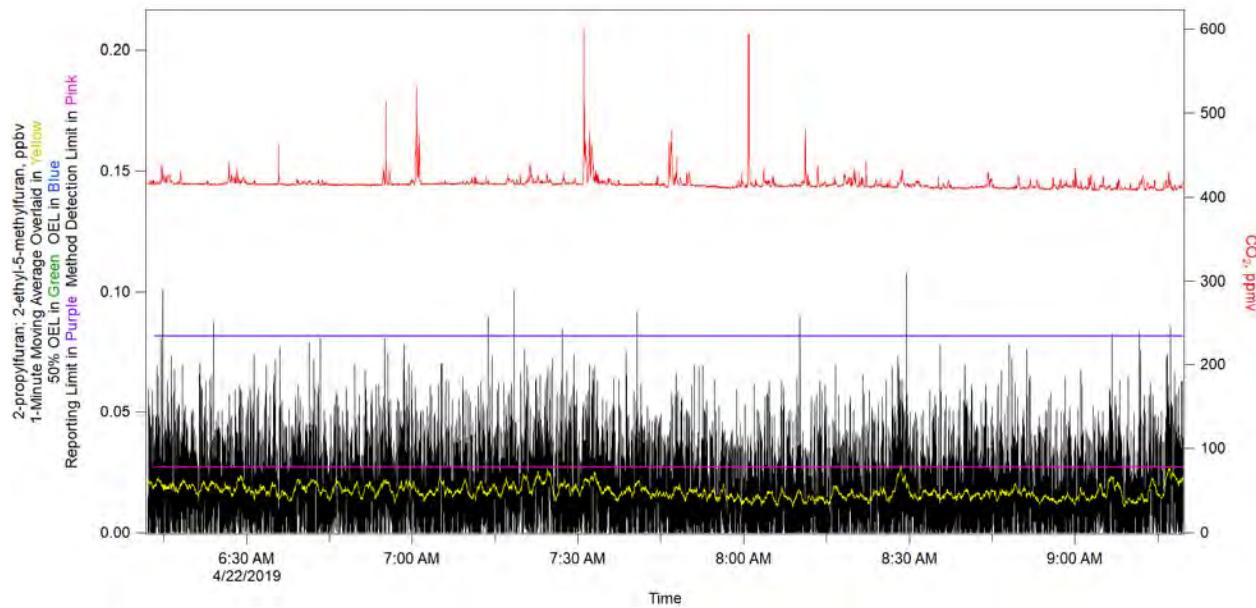


Figure 2-33. 2-propylfuran; 2-ethyl-5-methylfuran.

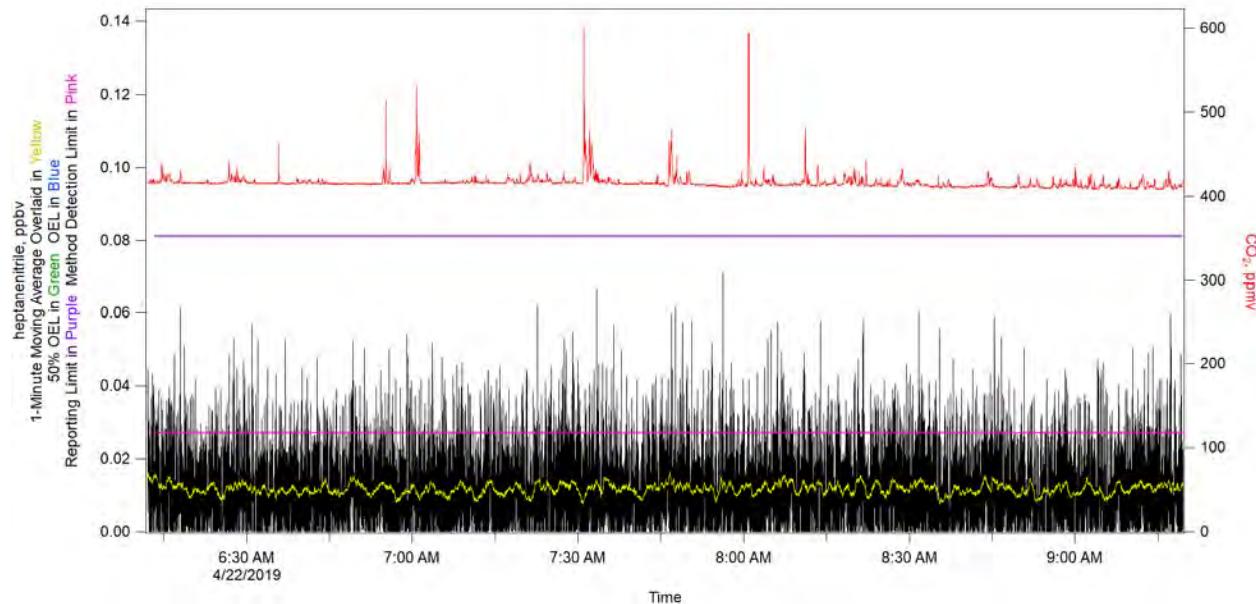


Figure 2-34. Heptanenitrile.

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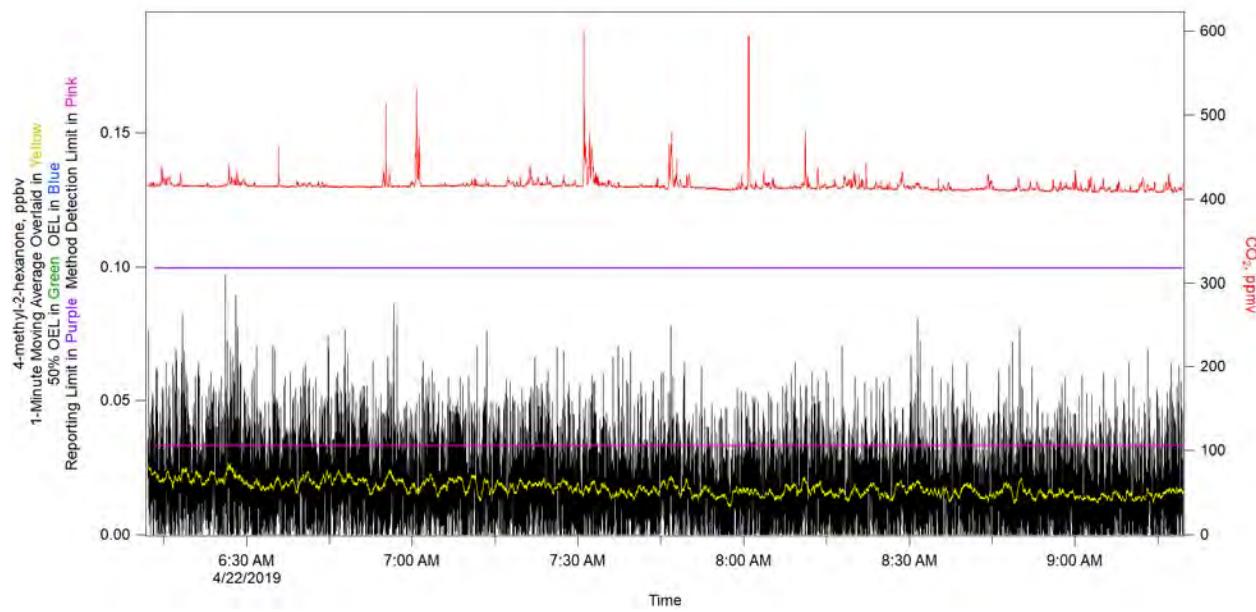


Figure 2-35. 4-methyl-2-hexanone.

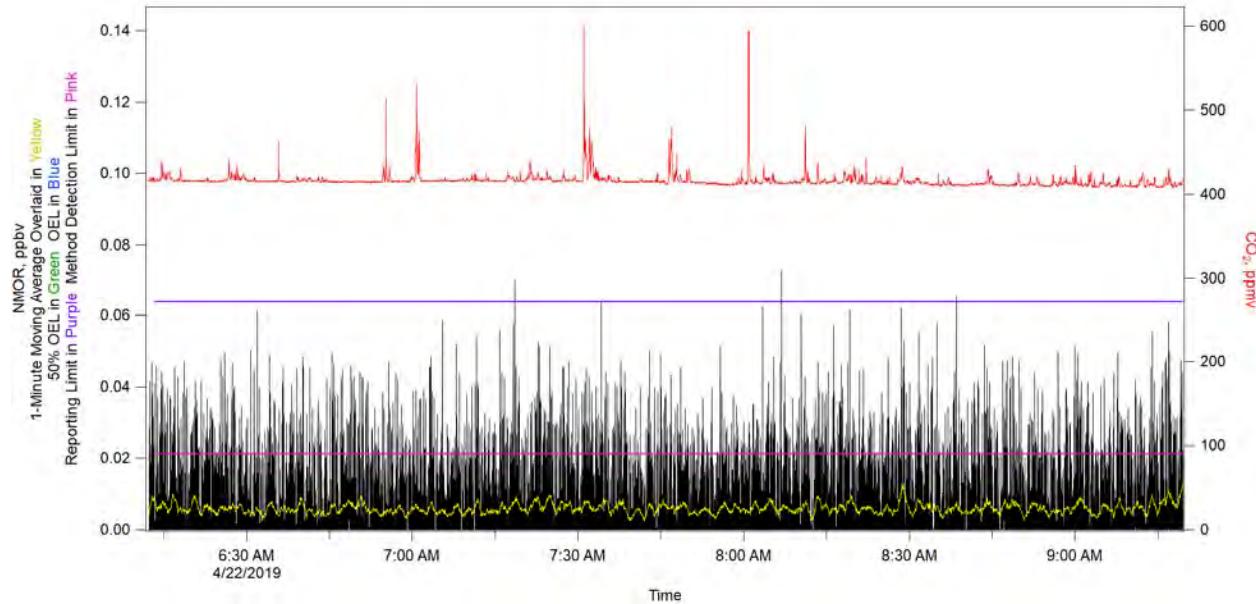


Figure 2-36. N-nitrosomorpholine (NMOR).

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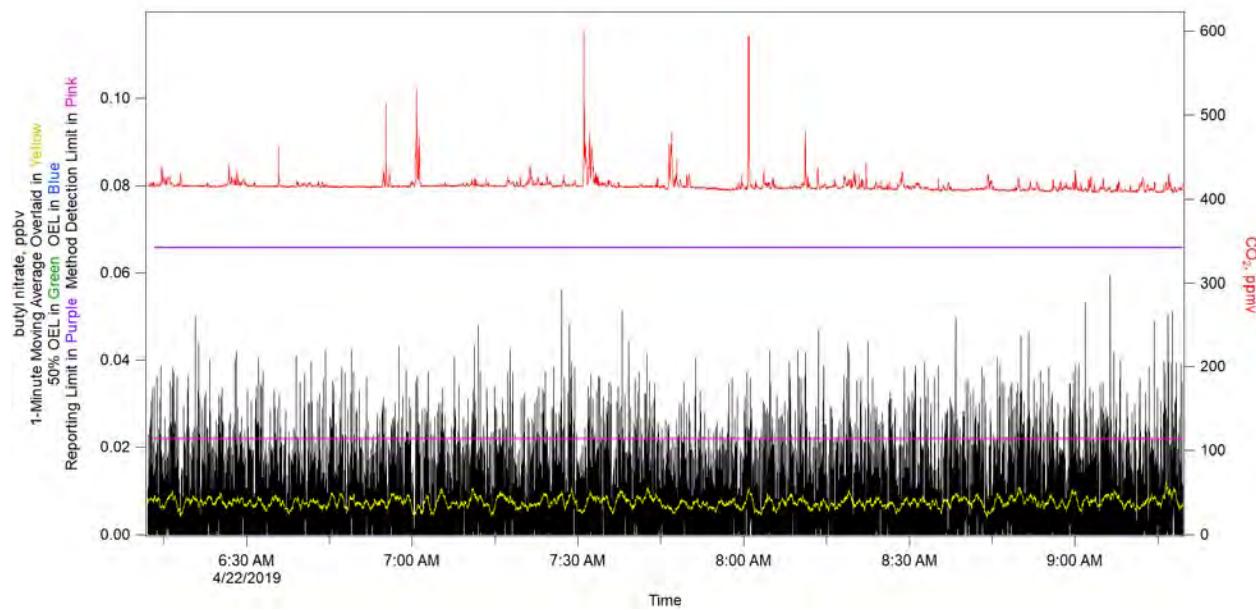
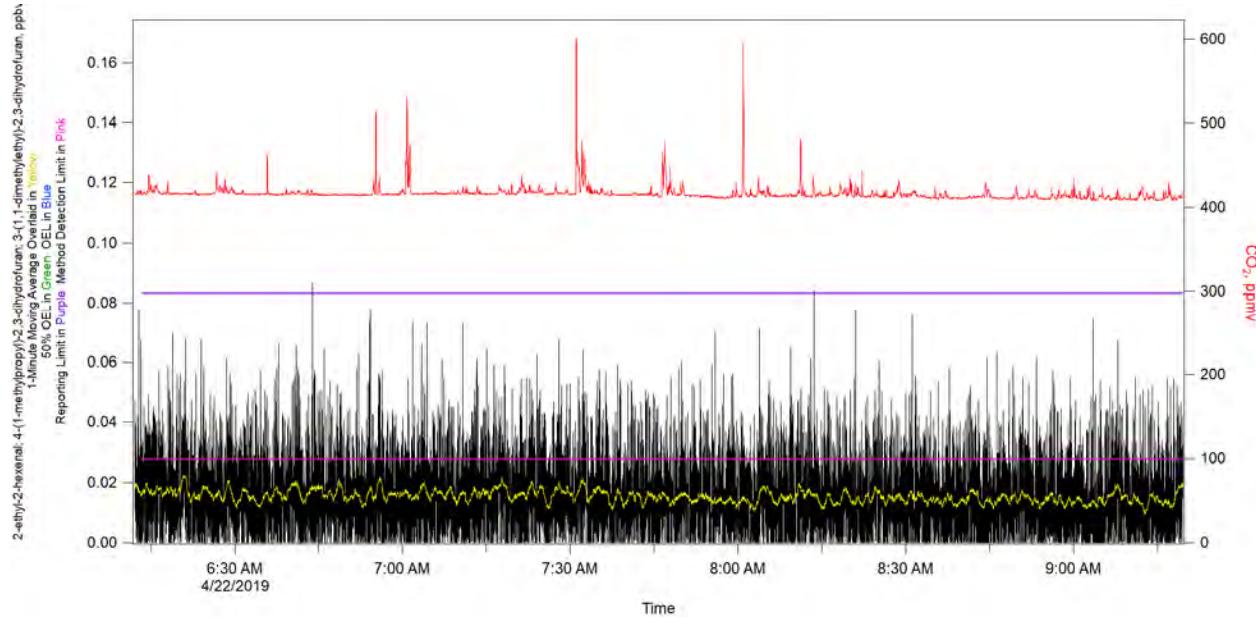


Figure 2-37. Butyl Nitrate.



**Figure 2-38. 2-ethyl-2-hexenal; 4-(1-methylpropyl)-2,3-dihydrofuran
3-(1,1-dimethylethyl)-2,3-dihydrofuran.**

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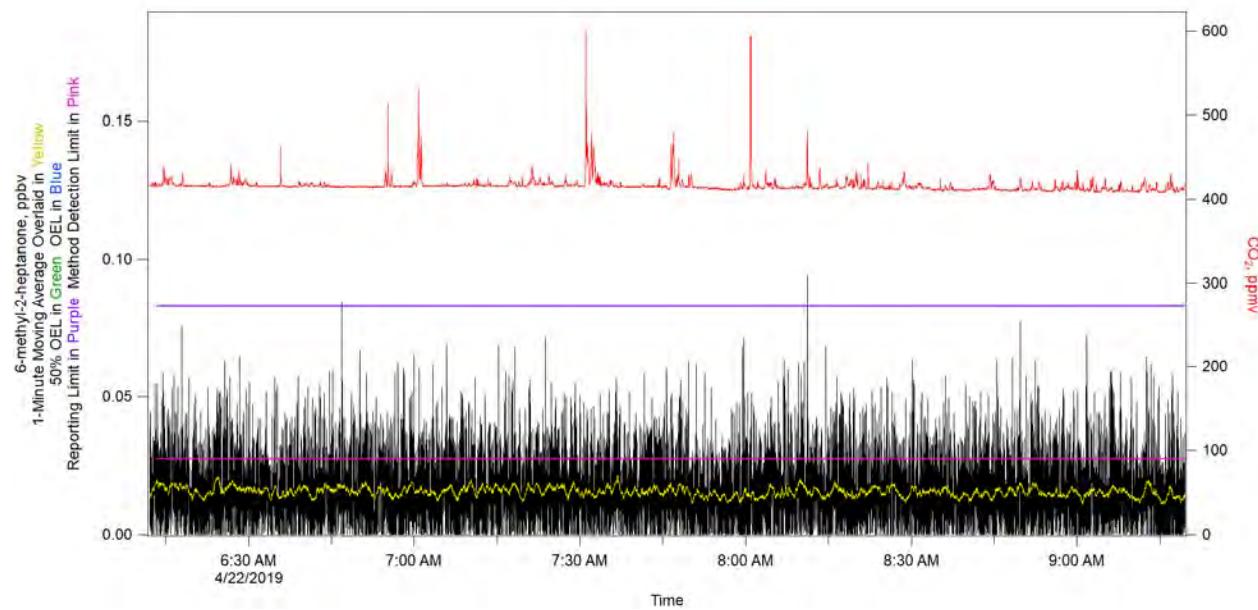


Figure 2-39. 6-methyl-2-heptanone.

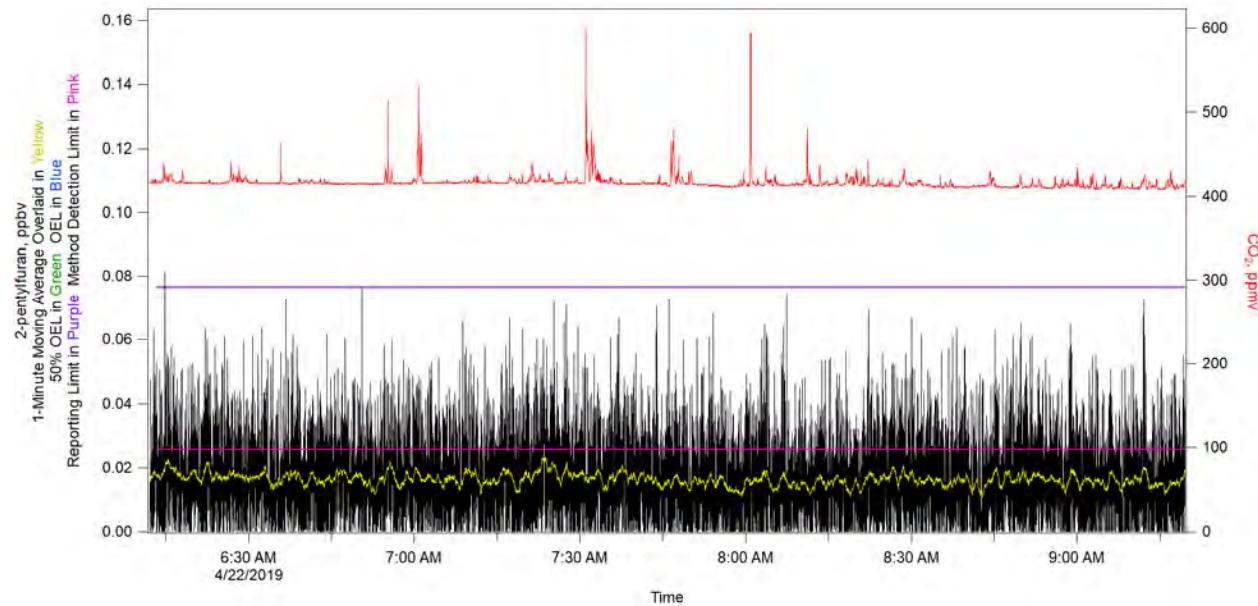


Figure 2-40. 2-pentylfuran.

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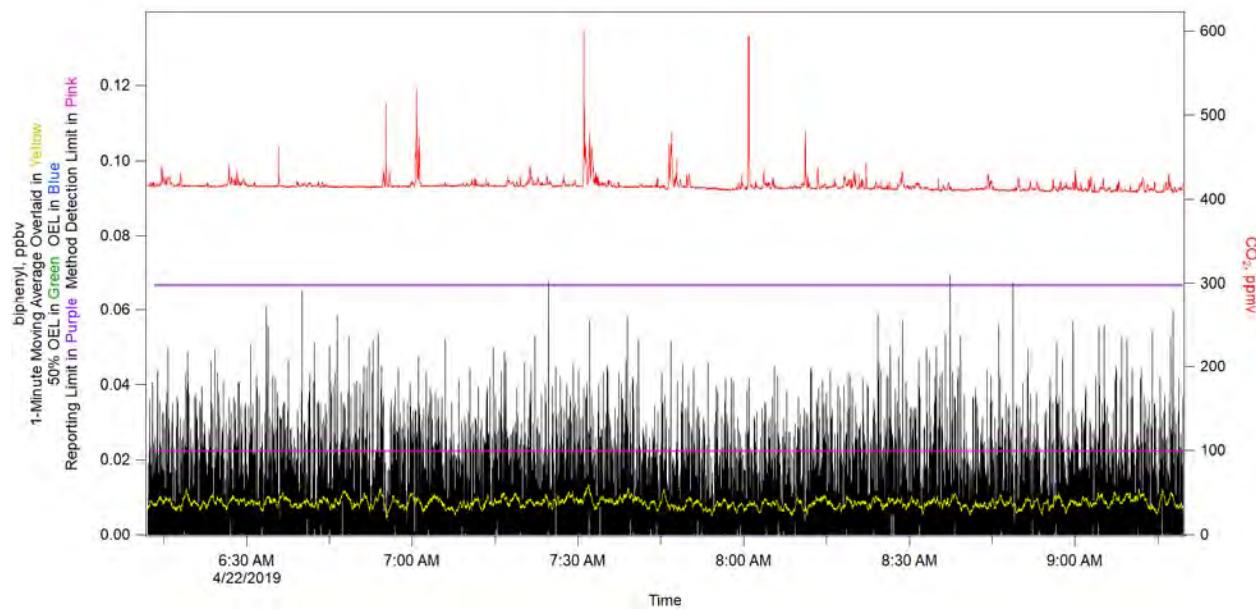


Figure 2-41. Biphenyl.

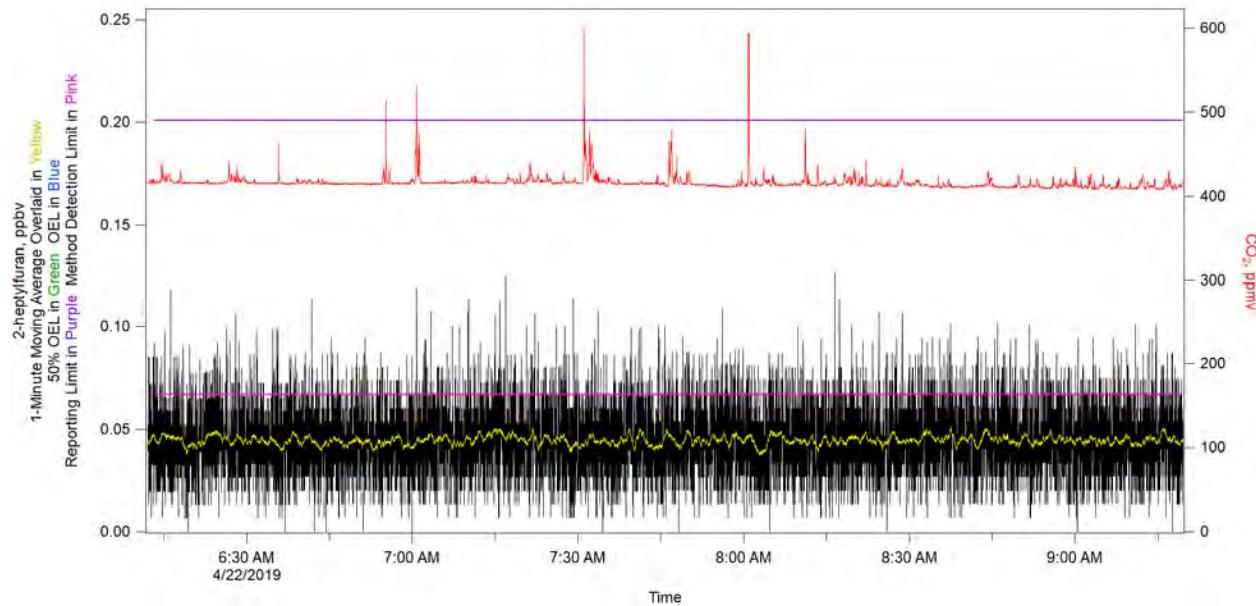


Figure 2-42. 2-heptylfuran.

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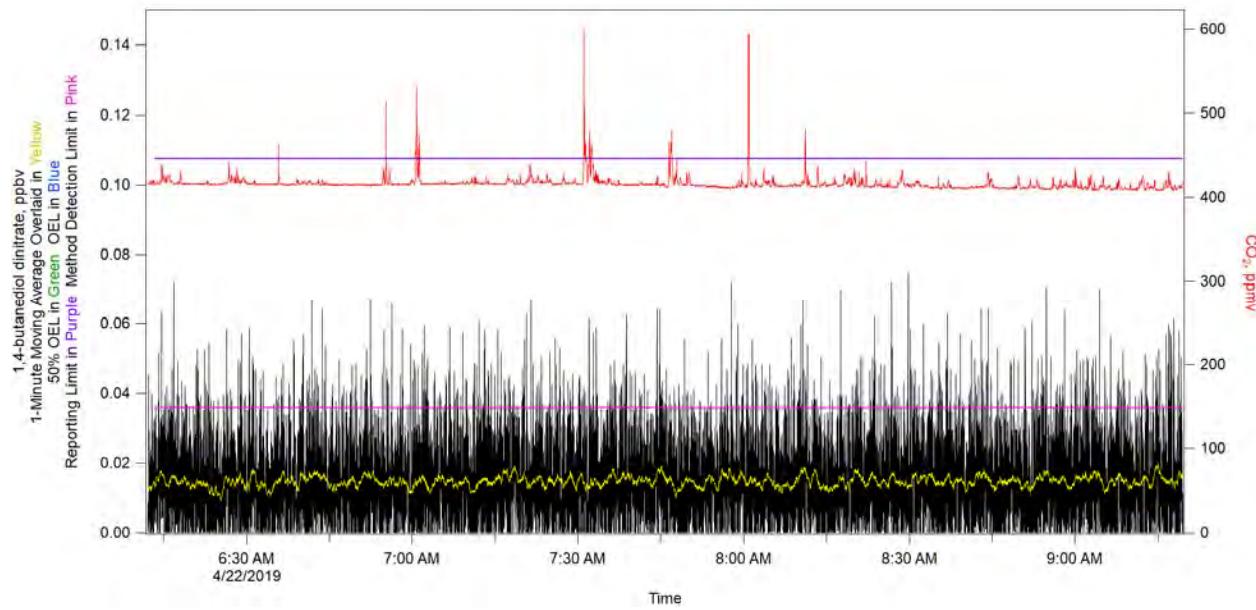


Figure 2-43. 1,4-butanediol Dinitrate.

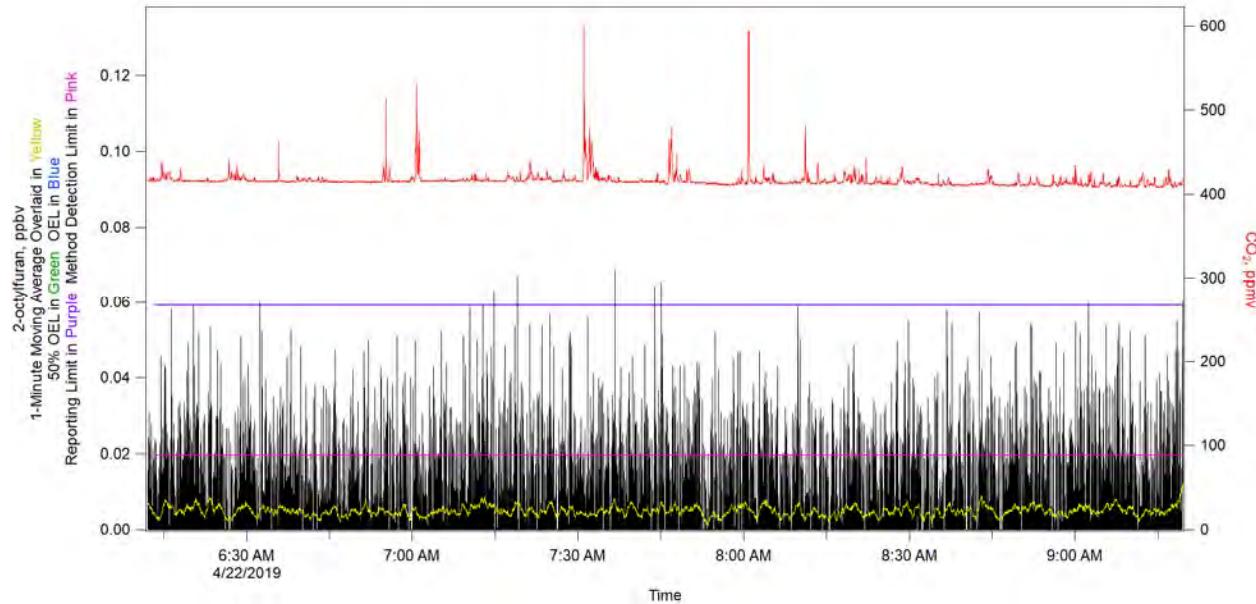


Figure 2-44. 2-octylfuran.

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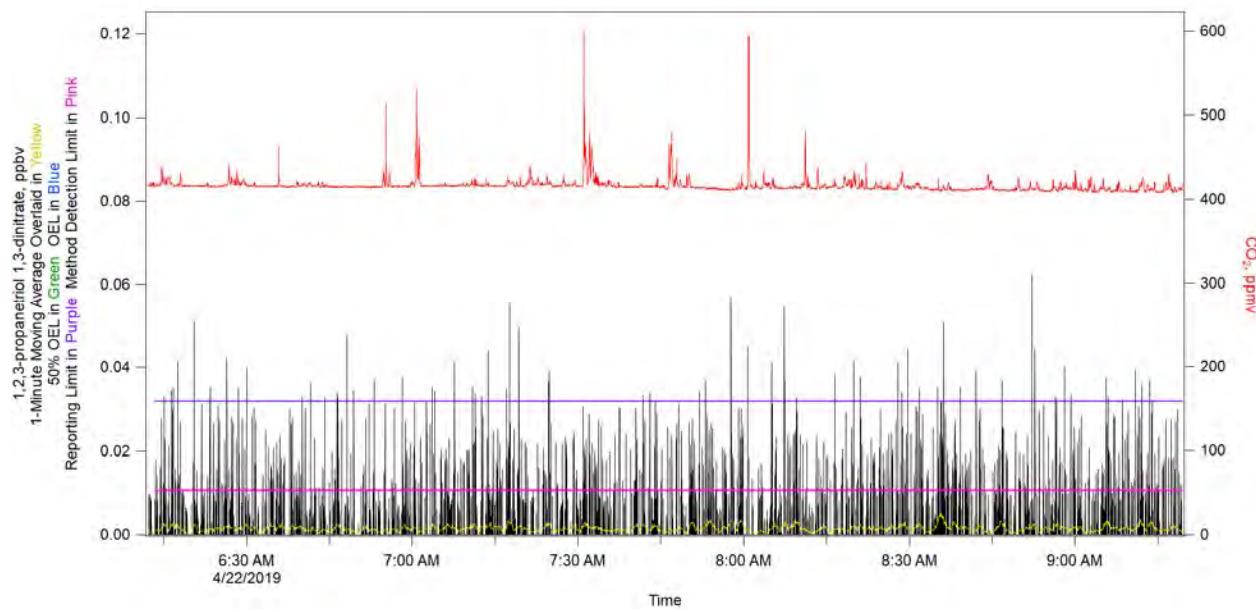


Figure 2-45. 1,2,3-propanetriol 1,3-dinitrate.

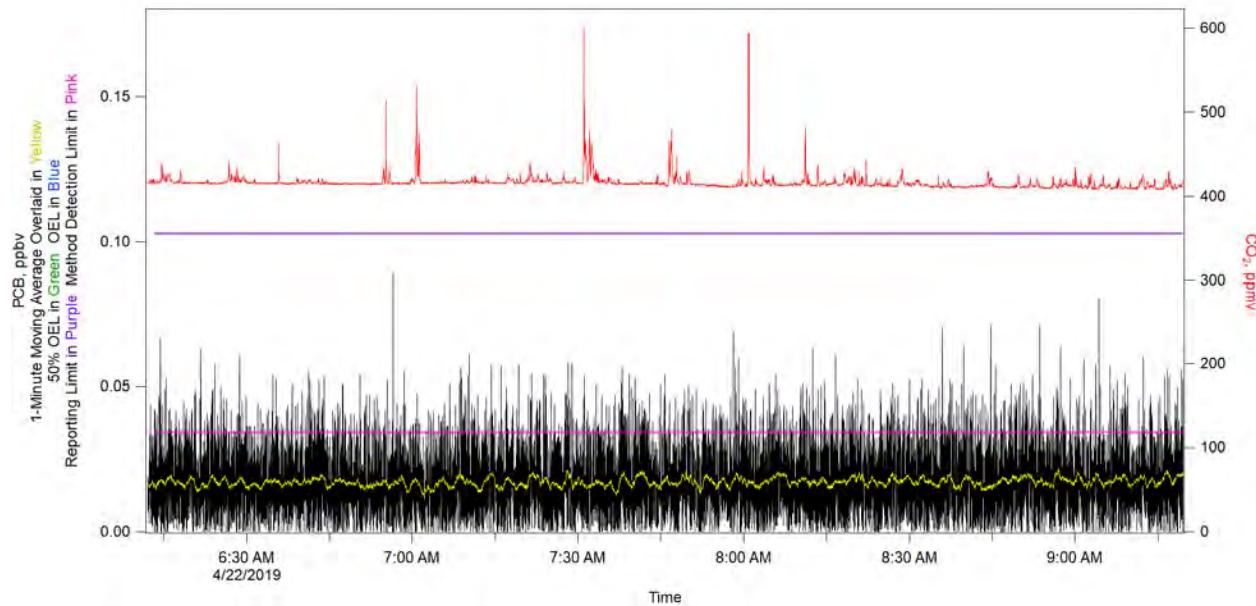


Figure 2-46. PCB.

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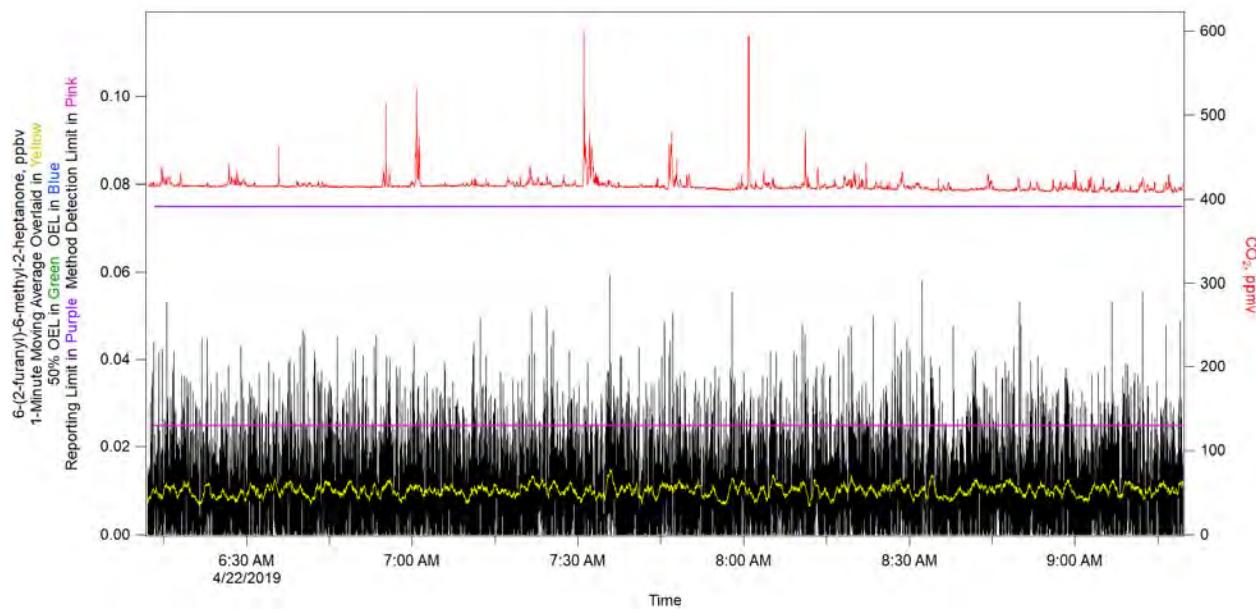


Figure 2-47. 6-(2-furanyl)-6-methyl-2-heptanone.

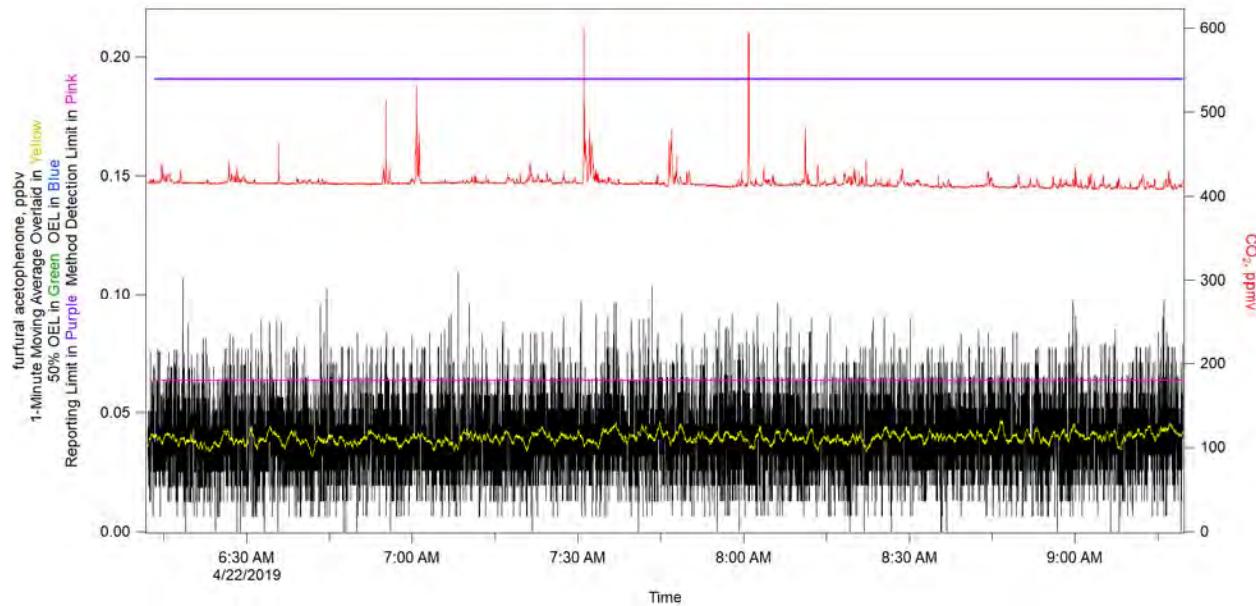


Figure 2-48. Furfural Acetophenone.

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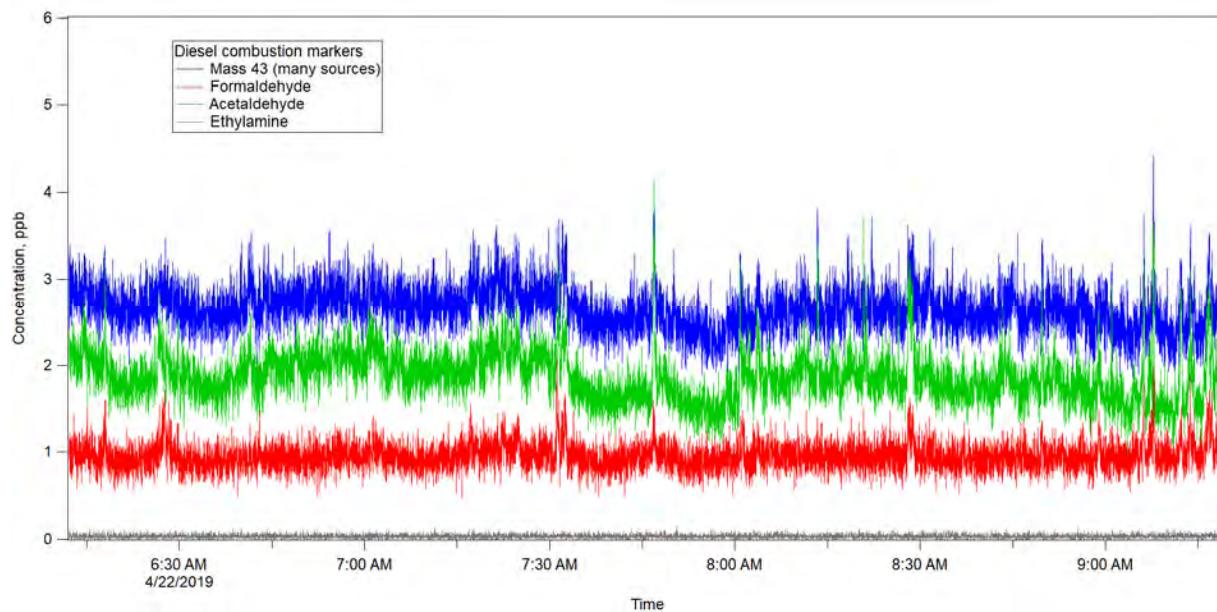


Figure 2-49. Diesel Combustion Markers.

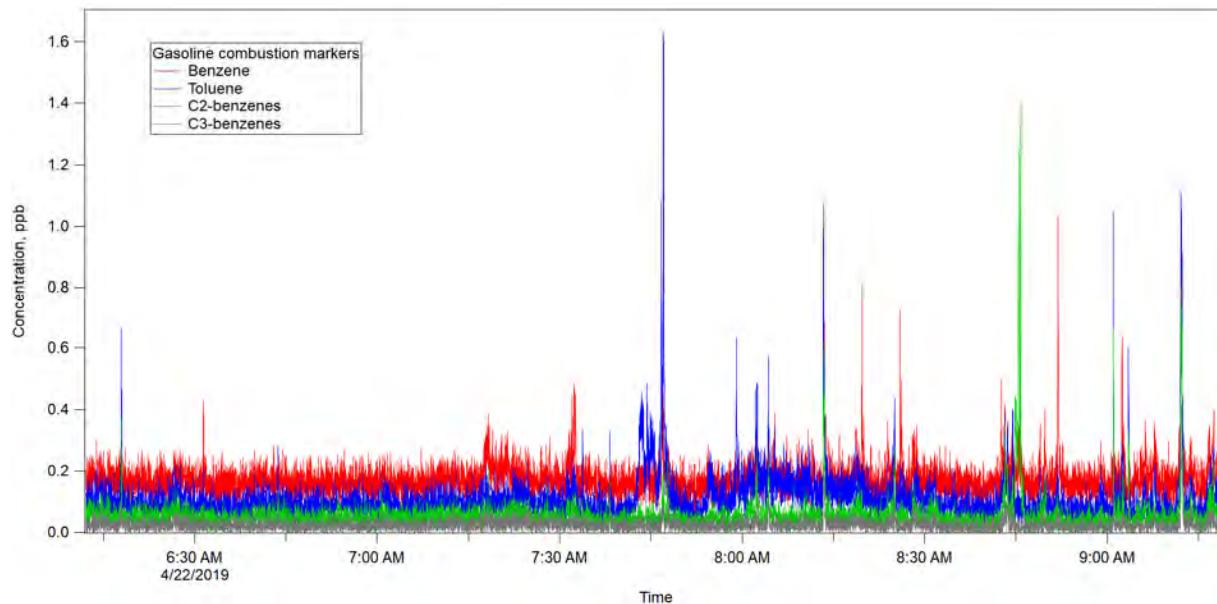


Figure 2-50. Gasoline Combustion Markers.

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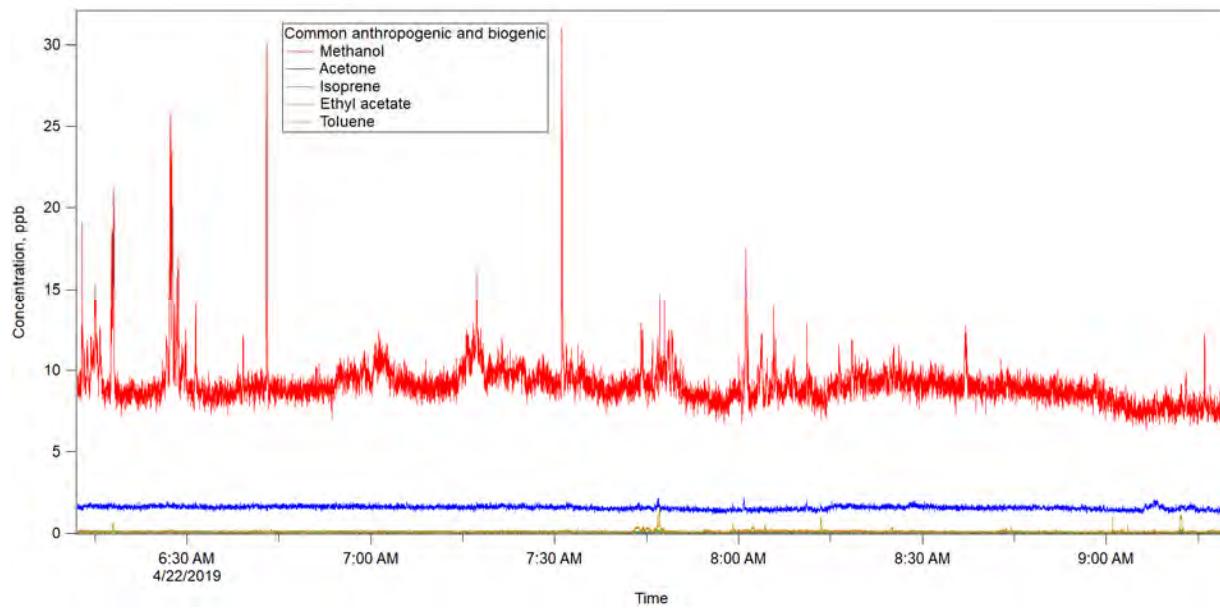


Figure 2-51. Plant and Human Markers.

2.5 AOP-015 Event

The ML Operators monitored the AOP-015 event near the U Farm Change Trailer from approximately 10:17 to 14:19. The section below provides the COPC statistical results from that monitoring period. 53005-81-COM-0519-005, Subcontract 53005, Release 81 – Transmittal of Revision 1 of Special Communication Report for (AOP)-015 Event on April 22, 2019, can be found in Appendix A.

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Table 2-2. Chemical of Potential Concern Statistical Information for the AOP-015 Monitoring on April 22, 2019. (2 Sheets)

COPC #	COPC Name	OEL (ppb)	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel St. Dev. (%)	Max. (ppb)	Median (ppb)
1	ammonia	25000	6.225	8.602†	0.401	4.658	10.190	8.579†
2	formaldehyde	300	1.721	2.225†	0.291	13.097	3.324	2.220†
3	methanol	200000	5.758	20.292	2.157	10.629	26.112	20.378
4	acetonitrile	20000	0.085	0.121†	0.033	27.005	0.260	0.119†
5	acetaldehyde	25000	1.027	2.438†	0.302	12.387	3.625	2.425†
6	ethylamine	5000	0.069	<0.069	0.024	58.427	0.166	<0.069
7	1,3-butadiene	1000	0.183	0.285†	0.078	27.319	0.714	0.282†
8	propanenitrile	6000	0.107	<0.107	0.030	38.903	0.226	<0.107
9	2-propenal	100	0.340	<0.34	0.049	30.420	0.398	<0.34
10	1-butanol + butenes	20000	0.214	<0.214	0.050	31.860	0.410	<0.214
11	methyl isocyanate	20	0.069	0.083†	0.035	42.416	0.266	0.080†
12	methyl nitrite	100	0.098	0.217†	0.050	22.983	0.433	0.217†
13	furan	1	0.062	<0.062	0.022	44.090	0.155	<0.062
14	butanenitrile	8000	0.039	<0.039	0.021	58.294	0.147	<0.039
15	but-3-en-2-one; 2,3-dihydrofuran + 2,5-dihydrofuran	200, 1, 1	0.041	0.051†	0.019	38.469	N/A*	N/A*
16	butanal	25000	0.061	0.264	0.054	20.465	0.518	0.262
17	NDMA**	0.3	0.082	<0.082	0.030	198.986	0.269	<0.082
18	benzene	500	0.236	0.413†	0.075	18.184	0.699	0.412†
19	2,4-pentadienenitrile + pyridine	300, 1000	0.085	<0.085	0.024	36.674	0.192	<0.085
20	2-methylene butanenitrile	300	0.036	<0.036	0.014	59.873	0.103	<0.036
21	2-methylfuran	1	0.043	0.046†	0.024	52.210	0.178	0.044†
22	Pantanenitrile	6000	0.036	<0.036	0.016	67.023	0.119	<0.036
23	3-methyl-3-butene-2-one + 2-methyl-2-butenal	20, 30	0.043	<0.043	0.024	58.595	0.161	<0.043
24	NEMA**	0.3	0.058	<0.058	0.029	154.500	0.181	<0.058
25	2,5-dimethylfuran	1	0.032	<0.032	0.020	72.220	0.125	<0.032
26	hexanenitrile	6000	0.031	<0.031	0.014	77.287	0.098	<0.031
27	2-hexanone (MBK)	5000	0.036	<0.036	0.021	66.718	0.149	<0.036
28	NDEA**	0.1	0.034	<0.034	0.016	134.378	0.107	<0.034
29	butyl nitrite + 2-nitro-2-methylpropane	100, 300	0.058	<0.058	0.020	39.035	0.132	<0.058
30	2,4-dimethylpyridine	500	0.036	<0.036	0.016	76.457	0.118	<0.036

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Table 2-2. Chemical of Potential Concern Statistical Information for the AOP-015 Monitoring on April 22, 2019. (2 Sheets)

COPC #	COPC Name	OEL (ppb)	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel St. Dev. (%)	Max. (ppb)	Median (ppb)
31	2-propylfuran + 2-ethyl-5-methylfuran	1	0.027	<0.027	0.014	113.359	0.102	<0.027
32	heptanenitrile	6000	0.027	<0.027	0.012	81.009	0.084	<0.027
33	4-methyl-2-hexanone	500	0.033	<0.033	0.015	81.918	0.092	<0.033
34	NMOR**	0.6	0.021	<0.021	0.012	200.988	0.093	<0.021
35	butyl nitrate	2500	0.022	<0.022	0.010	110.338	0.067	<0.022
36	2-ethyl-2-hexenal + 4-(1-methylpropyl)-2,3-dihydrofuran + 3-(1,1-dimethylethyl)-2,3-dihydrofuran	100, 1, 1	0.028	<0.028	0.014	80.121	0.100	<0.028
37	6-methyl-2-heptanone	8000	0.028	<0.028	0.013	81.765	0.089	<0.028
38	2-pentylfuran	1	0.026	<0.026	0.013	81.122	0.078	<0.026
39	biphenyl	200	0.022	<0.022	0.014	106.903	0.079	<0.022
40	2-heptylfuran	1	0.067	<0.067	0.020	33.915	0.150	<0.067
41	1,4-butanediol dinitrate	50	0.036	<0.036	0.016	64.364	0.101	<0.036
42	2-octylfuran	1	0.020	<0.02	0.010	204.239	0.093	<0.02
43	1,2,3-propanetriol 1,3-dinitrate	50	0.011	<0.011	0.010	264.422	0.080	<0.011
44	PCB	1000	0.034	<0.034	0.016	53.239	0.109	<0.034
45	6-(2-furanyl)-6-methyl-2-heptanone	1	0.025	<0.025	0.010	93.233	0.080	<0.025
46	furfural acetophenone	1	0.064	<0.064	0.019	36.634	0.135	<0.064
N/A*	The maximum peak value for but-3-en-2-one + 2,3 dihydrofuran + 2,5 dihydrofuran was 0.143 ppb and the median value was 0.049† ppb. The PTR-MS results for but-3-en-2-one + 2,3 dihydrofuran + 2,5 dihydrofuran are not compared to OEL concentrations because: 1) the result is suspect due to a known biogenic interferant (methacrolein) that is expected to be in concentrations that occasionally exceed the dihydrofuran OEL, and 2) this combination of COPCs have OEL concentrations that differ by a factor of 200, which provide widely variant bases for these numbers.							
**	Nitrosamine results are suspect due to isobaric interferants causing positive bias that have been encountered during previous background [53005-81-RPT-007, PTR-MS Mobile Laboratory Vapor Monitoring Background Study, (3/18/2018 – 4/20/2018), and Fiscal Year 2017 Mobile Laboratory Vapor Monitoring at the Hanford Site: Monitoring During Waste Disturbing Activities and Background Study, RJ Lee Group, Inc.].							
<	COPC Averages below the MDL.							
†	COPC Averages between the RL and the MDL.							
	COPC Averages >100% of the OEL.							
	COPC Averages 50-100% of the OEL.							
	COPC Averages 10-50% of the OEL.							

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Figure 2-52 through Figure 2-100 display 46 COPC signals, overlaid with the same signal smoothed using a 1-minute moving average (in cases where a moving average assists with data visualization), and CO₂, for the monitoring period April 22, 2019. If within range of the plot's left axis, a green horizontal line representing 50% of the COPC's OEL, a blue horizontal line representing the COPC's OEL, a horizontal purple line representing the RL, and a pink horizontal line representing the MDL are shown.

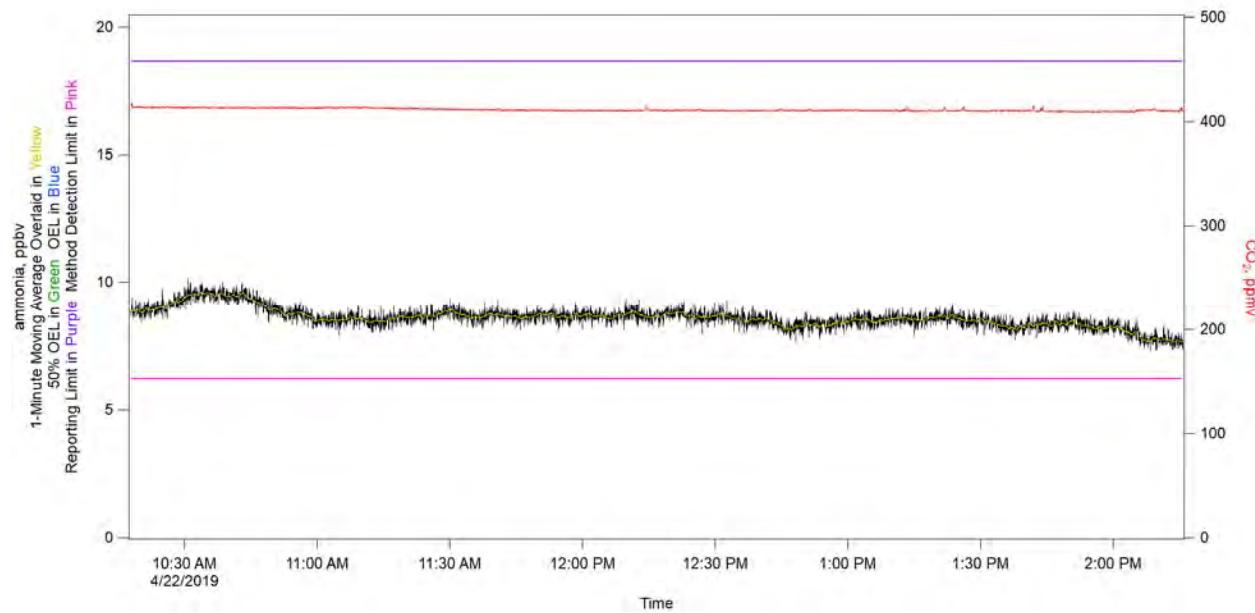


Figure 2-52. Ammonia.

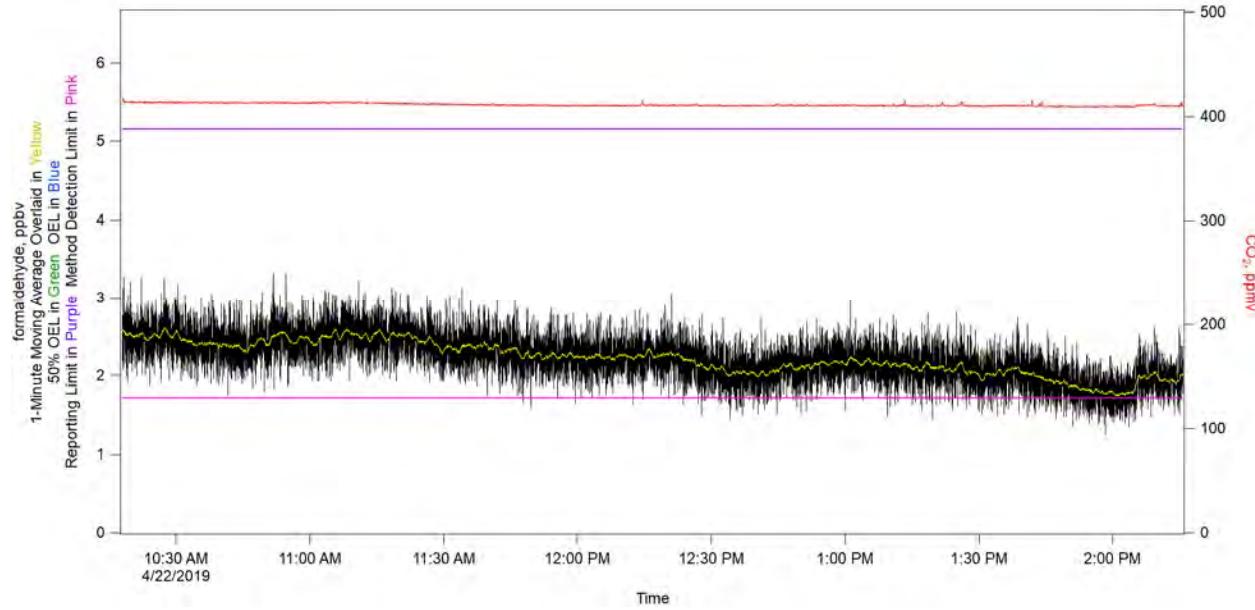


Figure 2-53. Formaldehyde.

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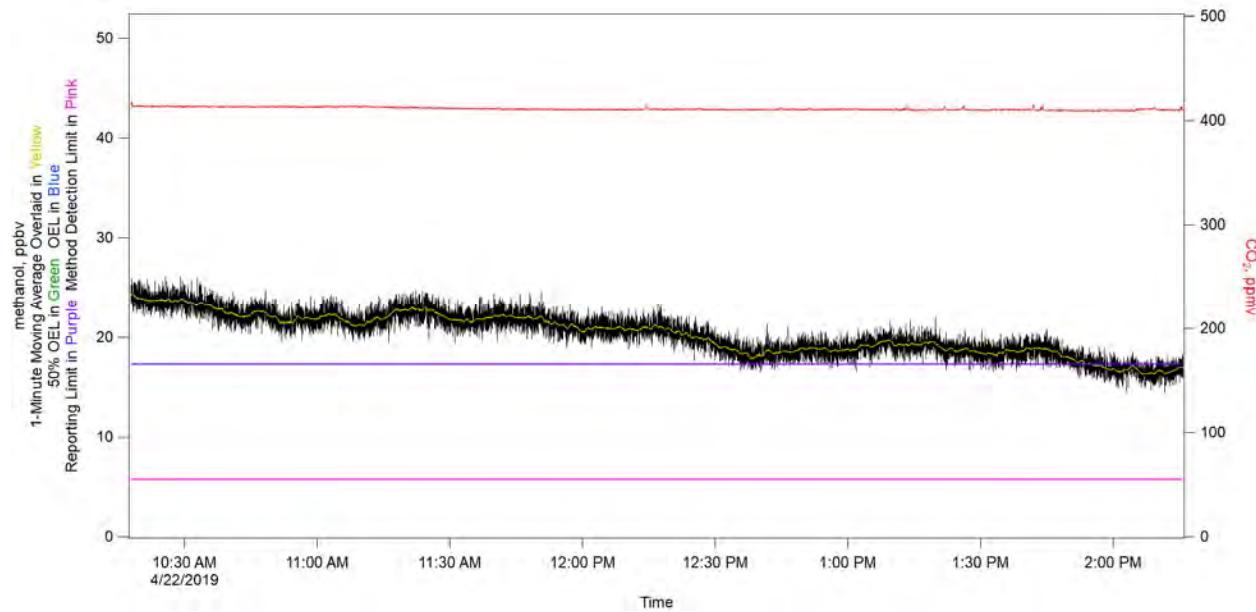


Figure 2-54. Methanol.

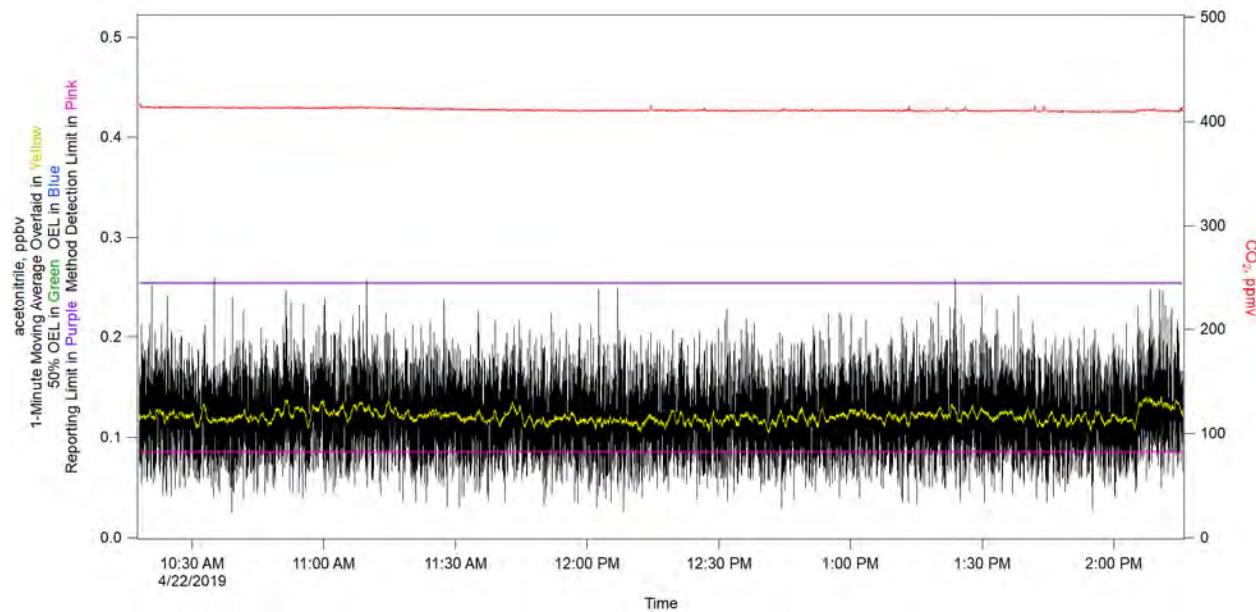


Figure 2-55. Acetonitrile.

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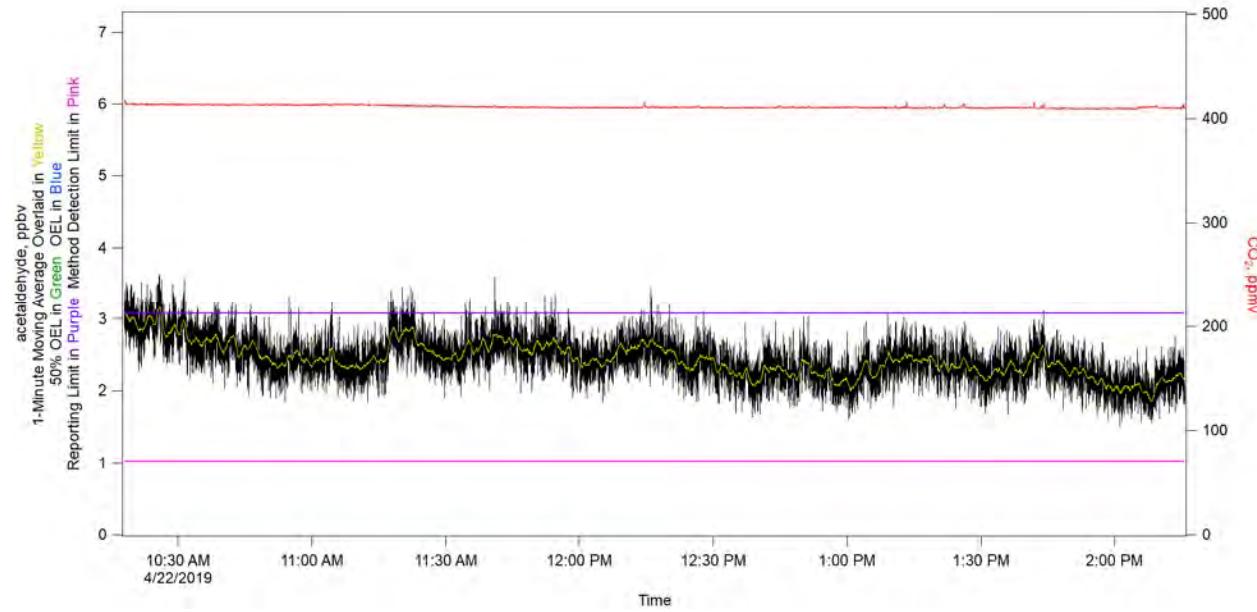


Figure 2-56. Acetaldehyde.

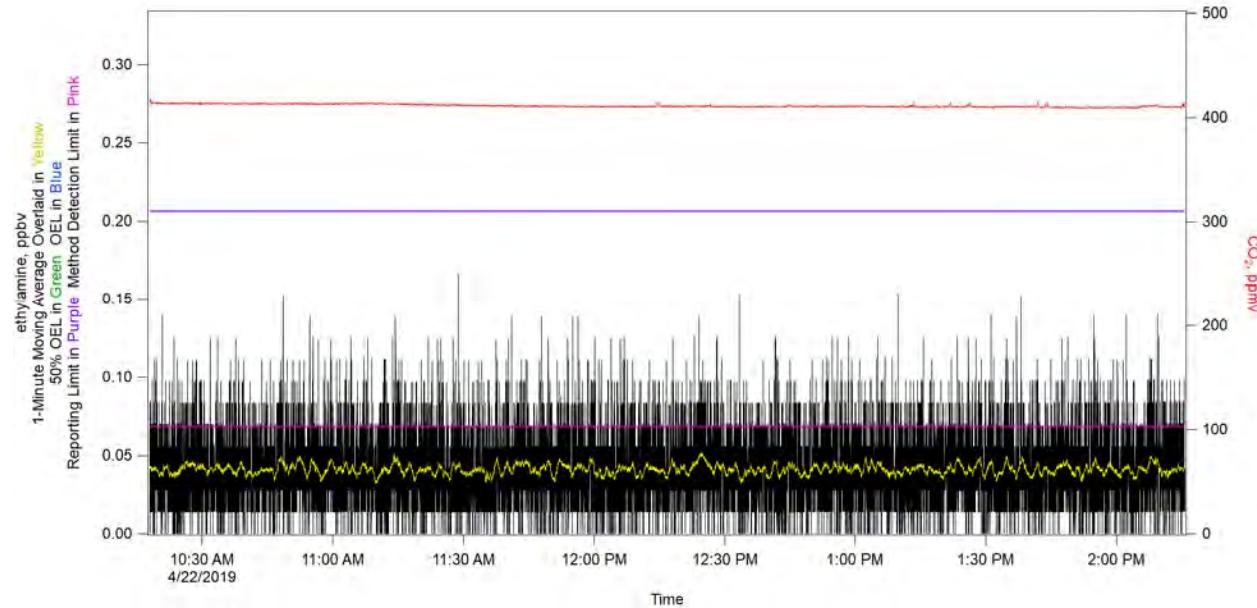


Figure 2-57. Ethylamine.

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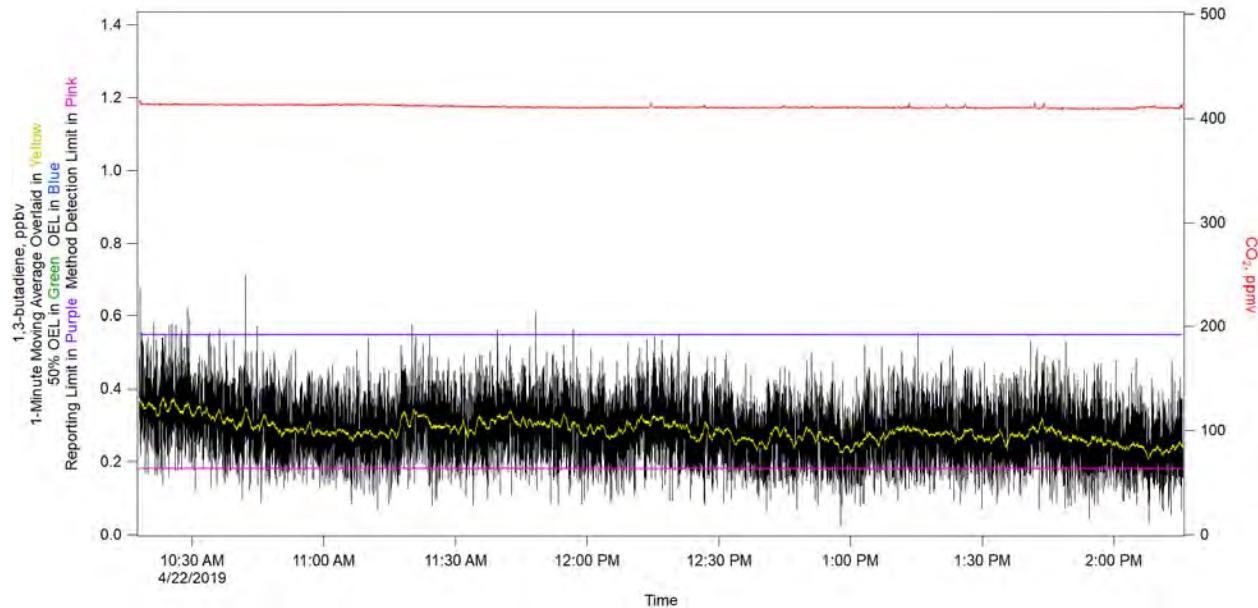


Figure 2-58. 1,3-butadiene.

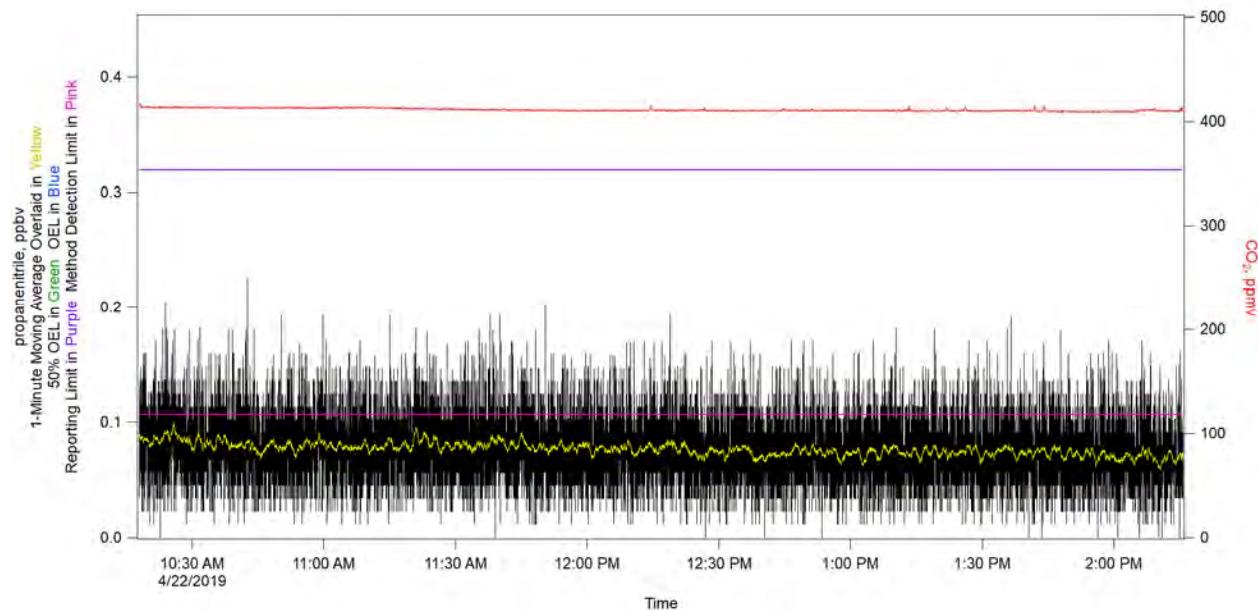


Figure 2-59. Propanenitrile.

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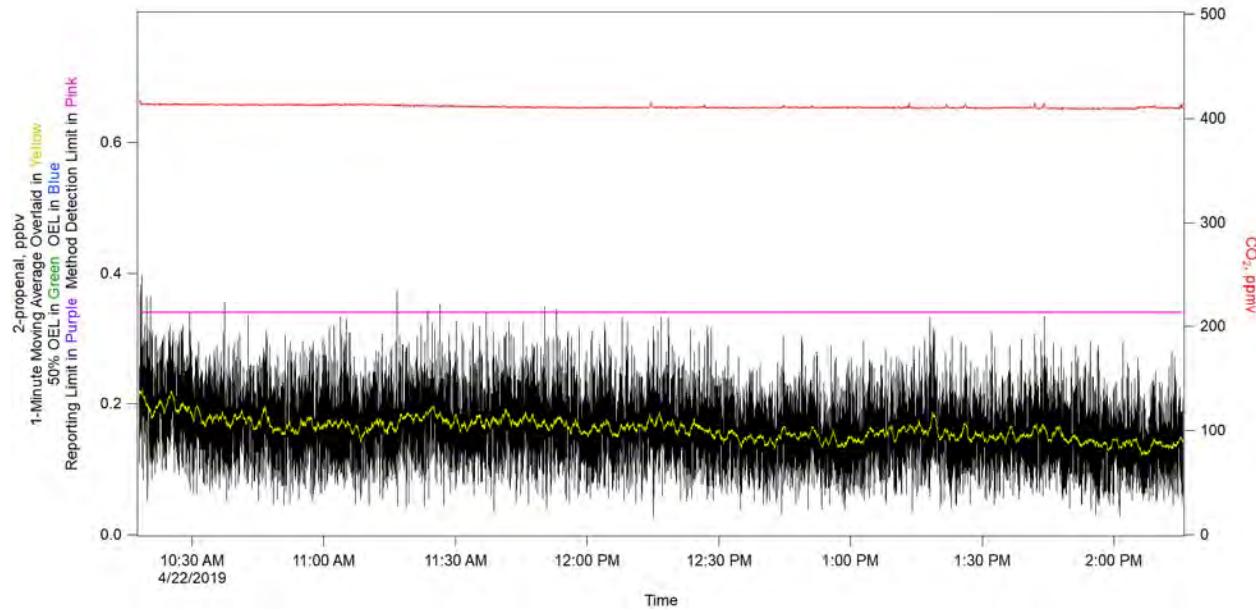


Figure 2-60. 2-propenal.

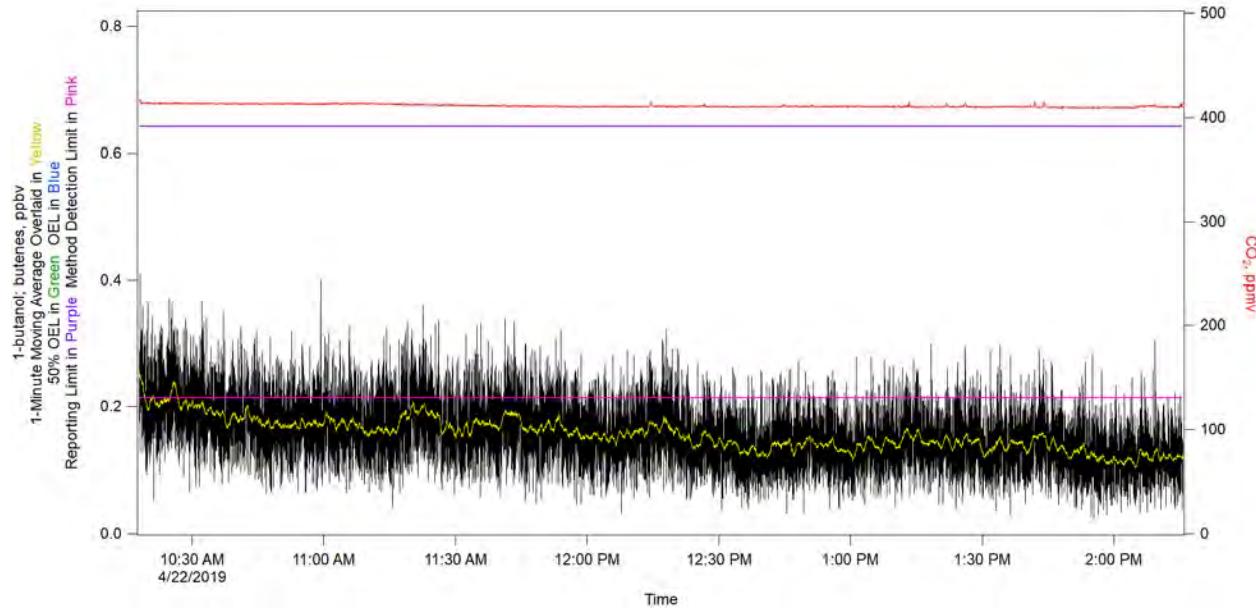


Figure 2-61. 1-butanol; Butenes.

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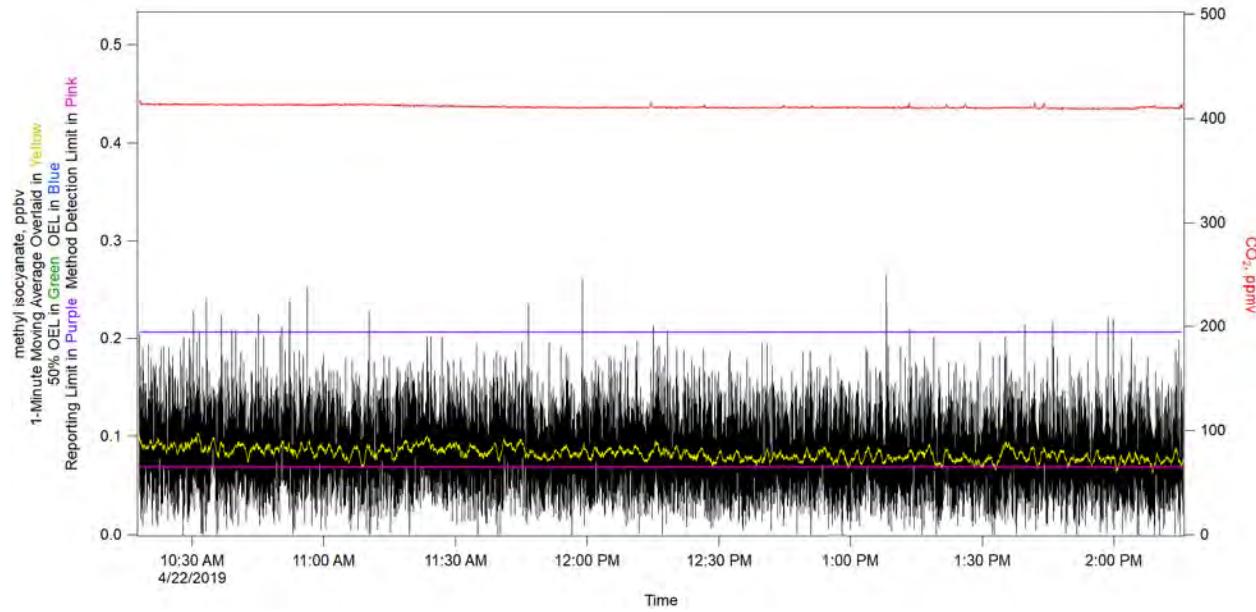


Figure 2-62. Methyl Isocyanate.

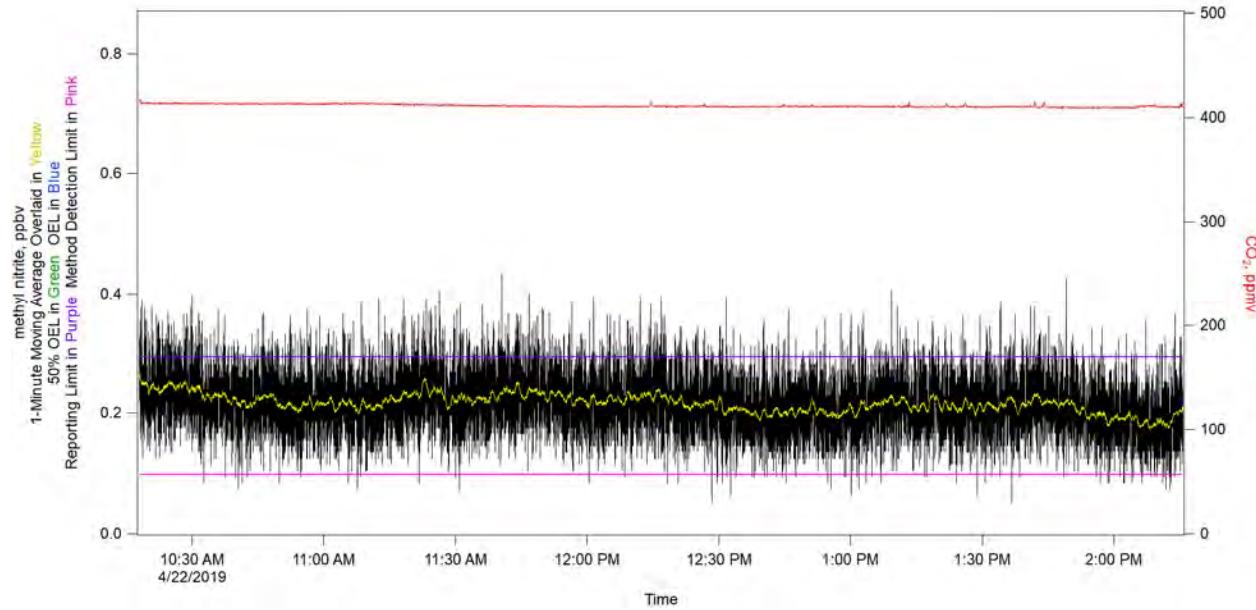


Figure 2-63. Methyl Nitrite.

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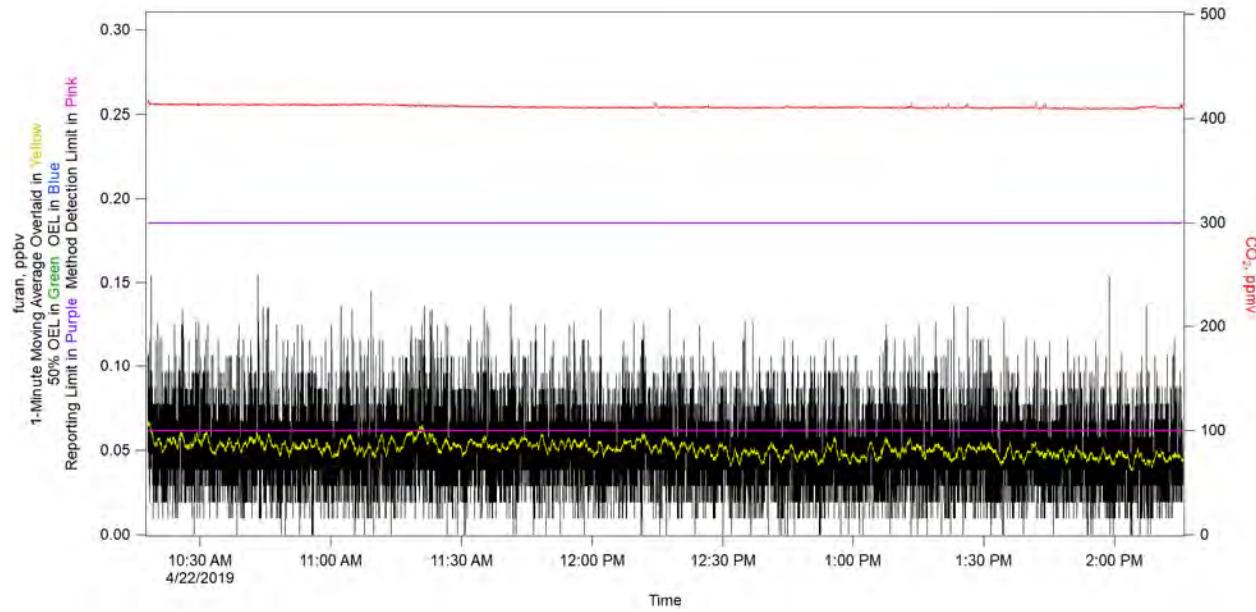


Figure 2-64. Furan.

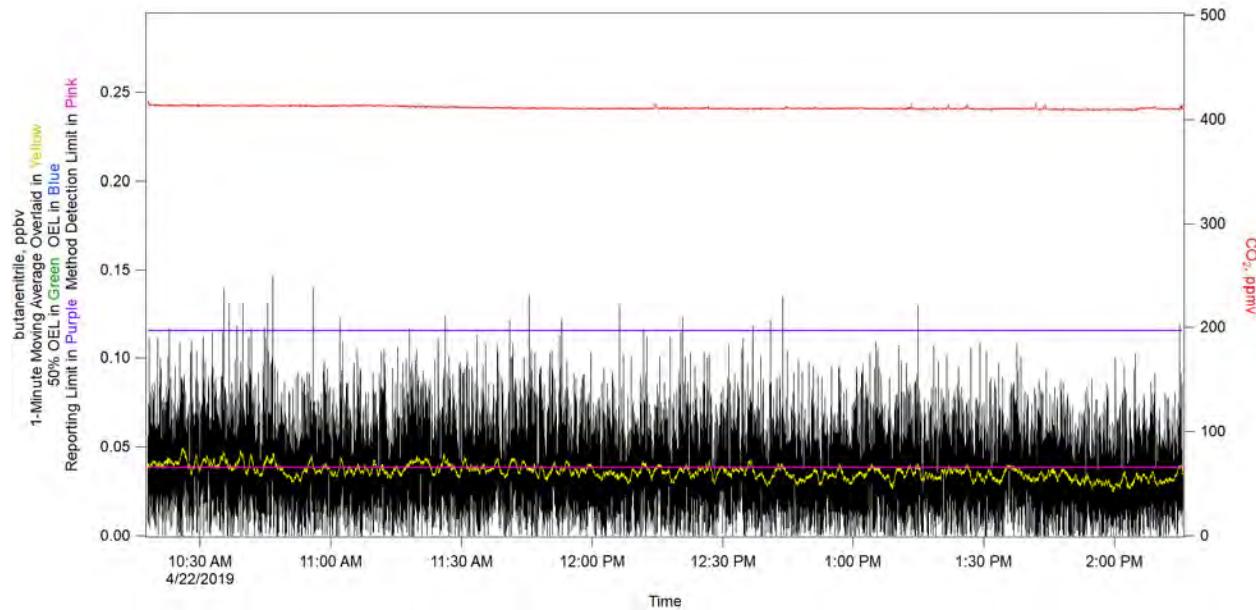


Figure 2-65. Butanenitrile.

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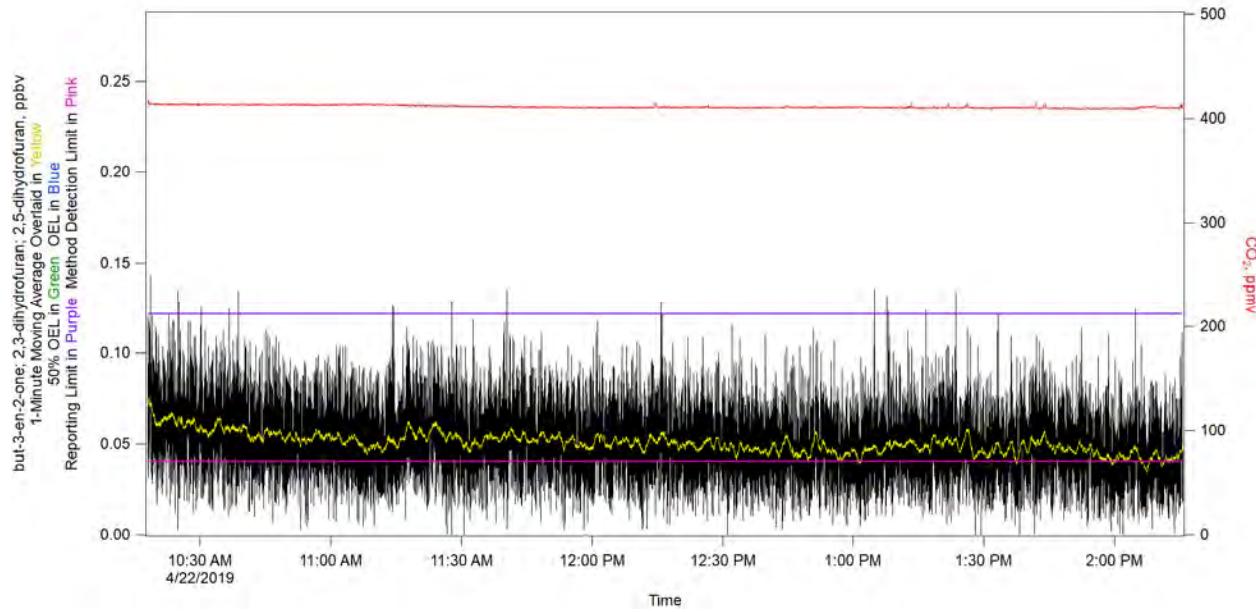


Figure 2-66. But-3-en-2-one; 2,3-dihydrofuran; 2,5-dihydrofuran.

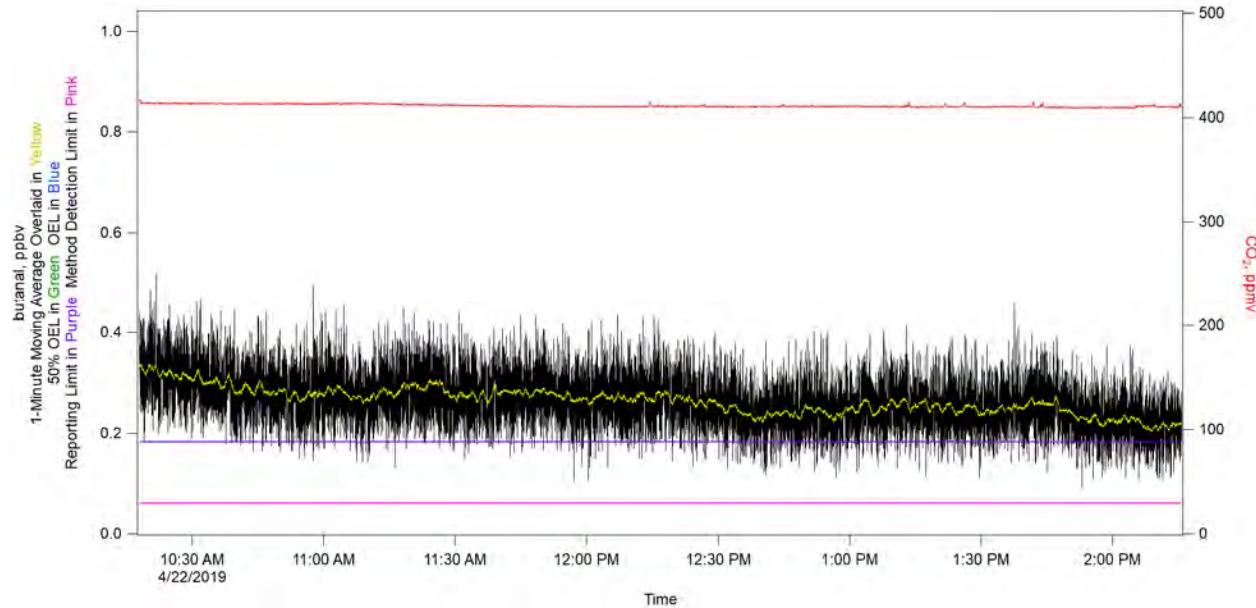


Figure 2-67. Butanal.

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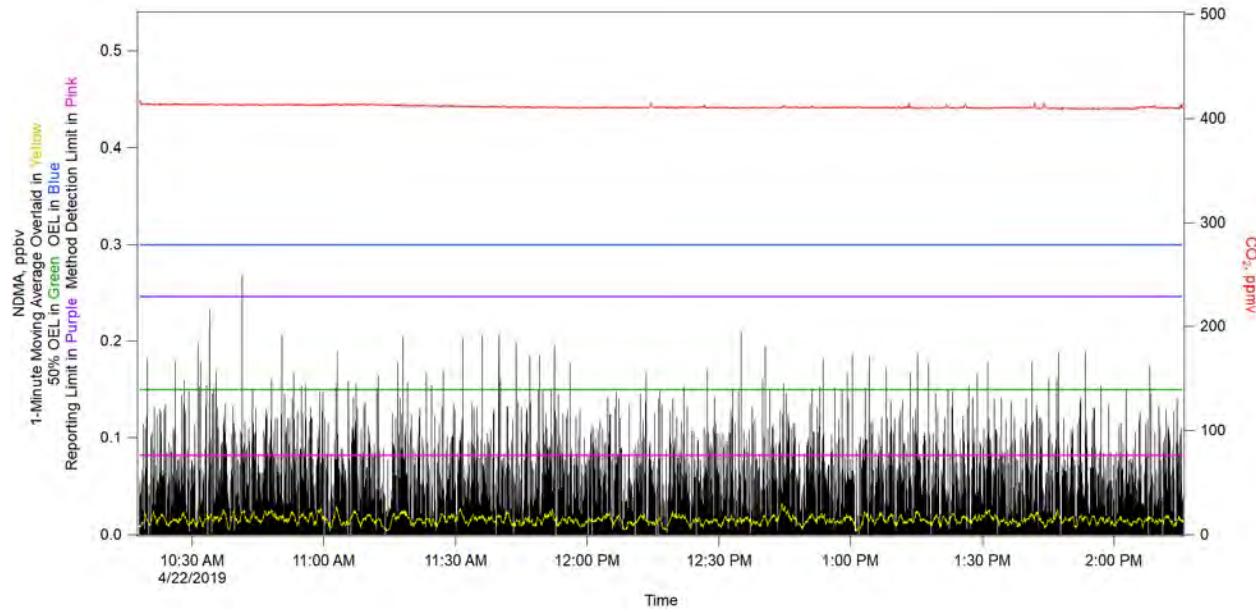


Figure 2-68. N-nitrosodimethylamine (NDMA).

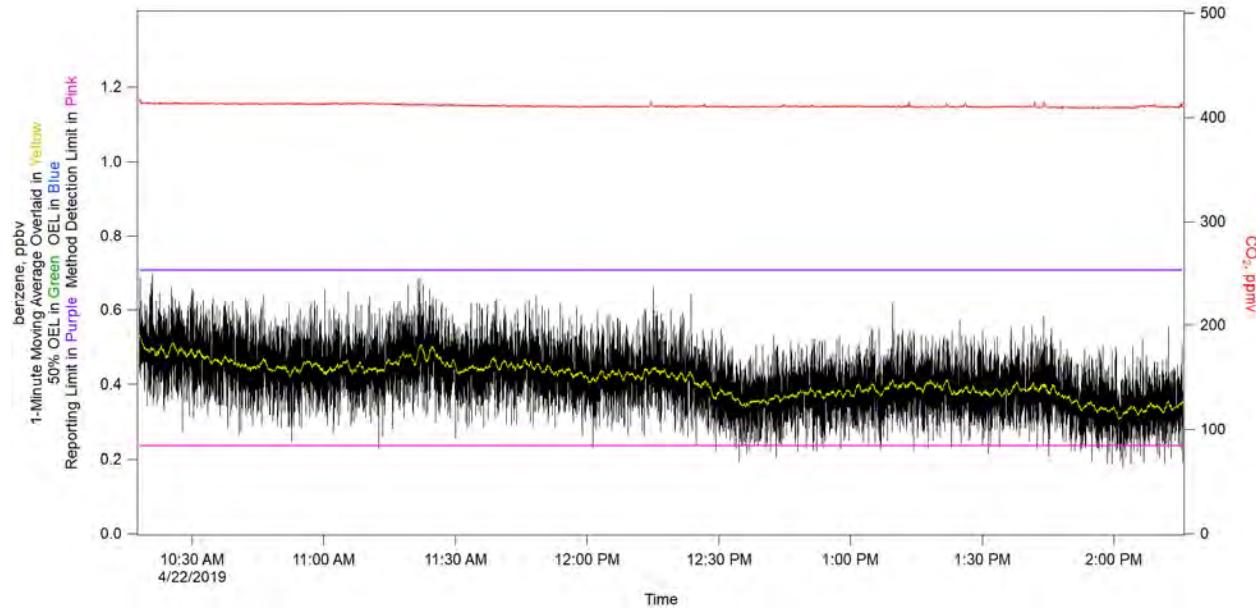


Figure 2-69. Benzene.

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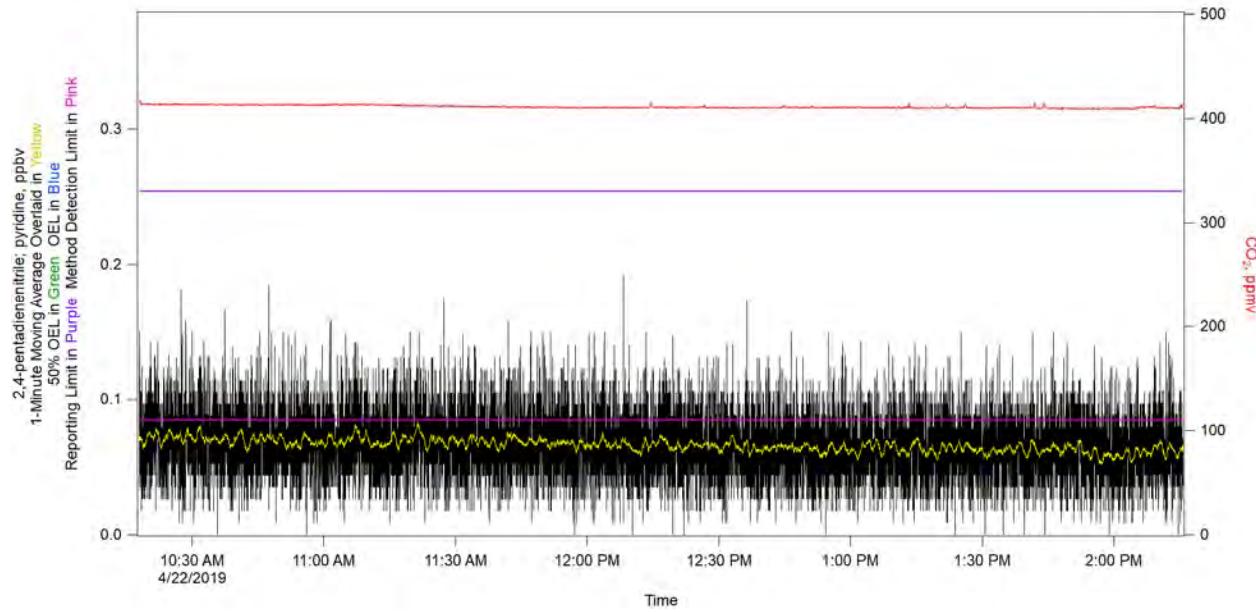


Figure 2-70. 2,4-pentadienenitrile; Pyridine.

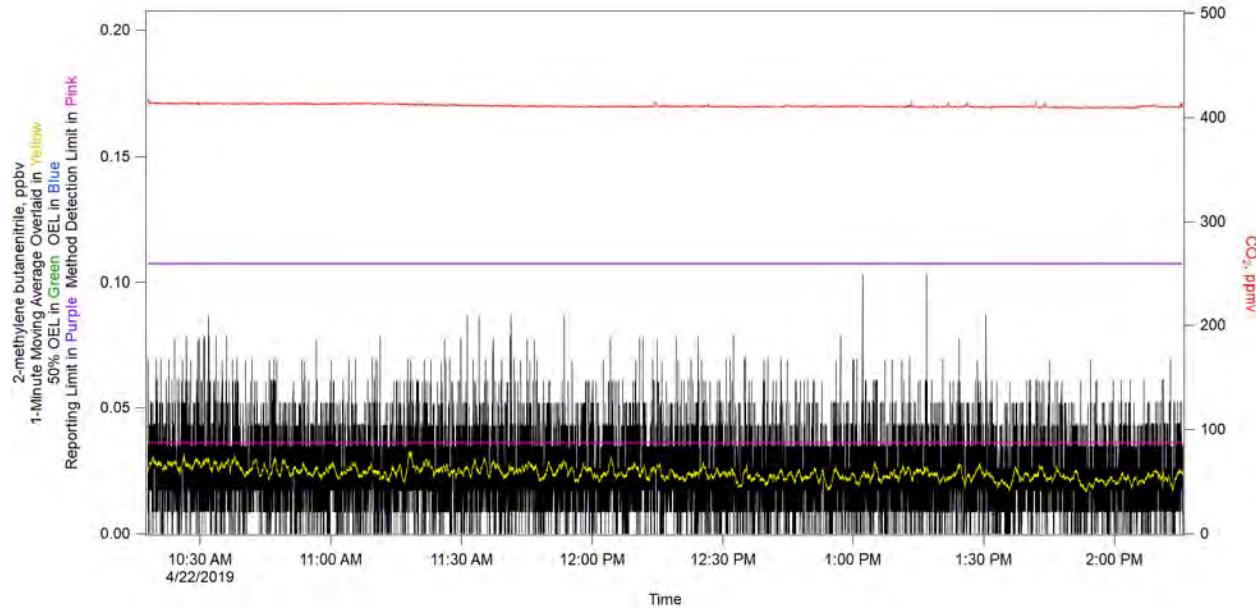


Figure 2-71. 2-methylene Butanenitrile.

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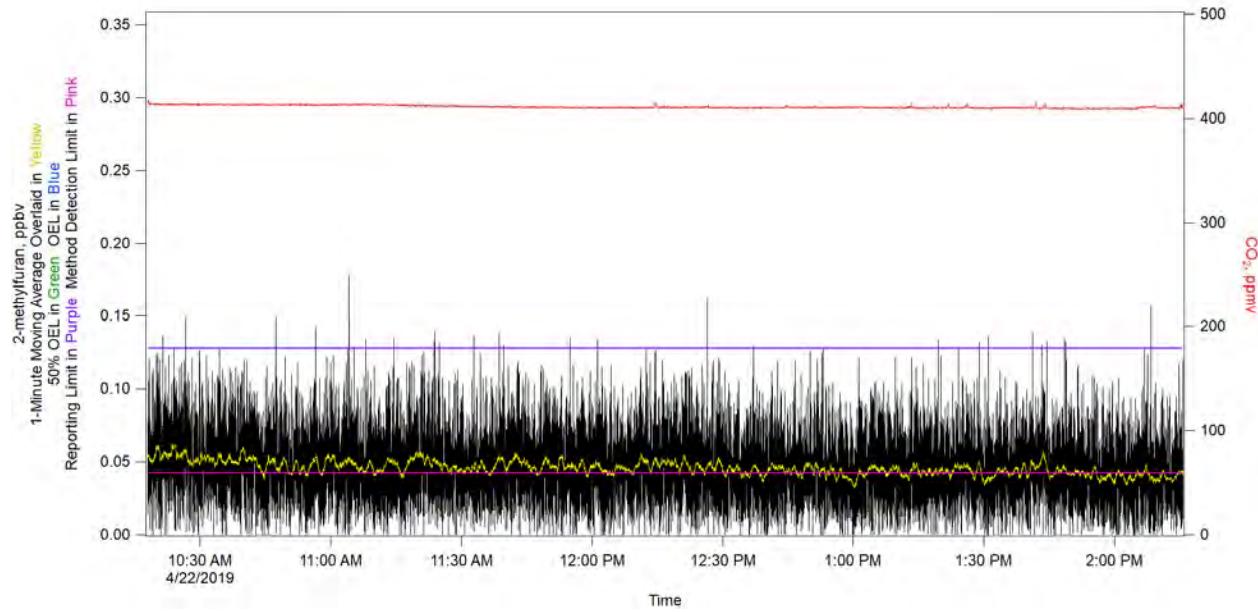


Figure 2-72. 2-methylfuran.

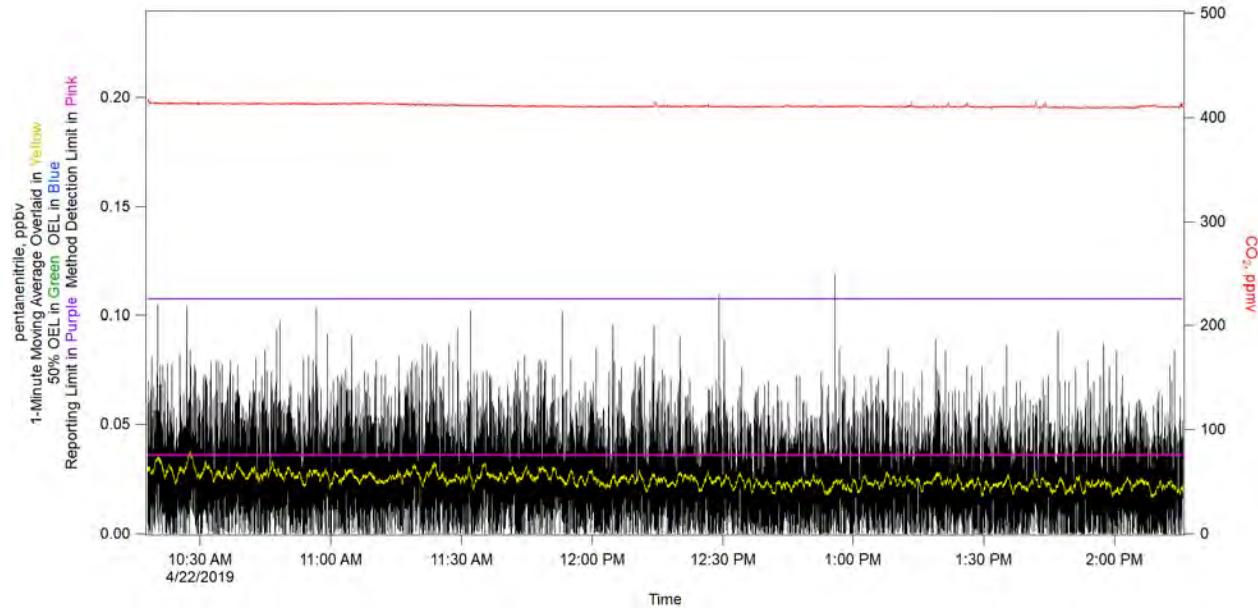


Figure 2-73. Pentanenitrile.

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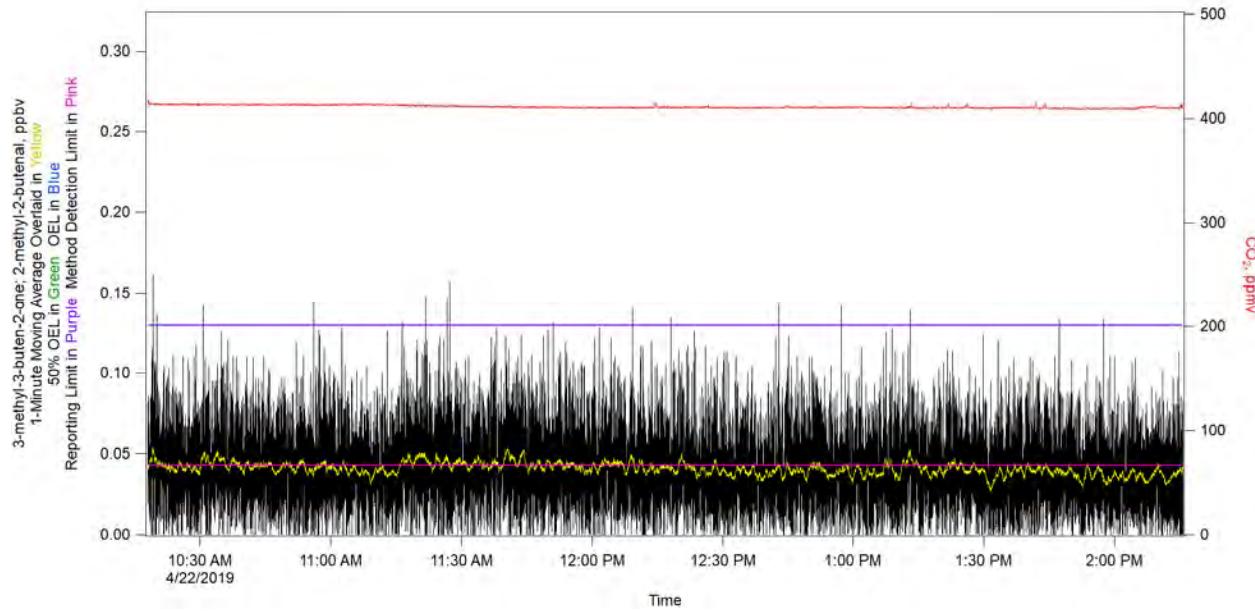


Figure 2-74. 3-methyl-3-buten-2-one; 2-methyl-2-butenal.

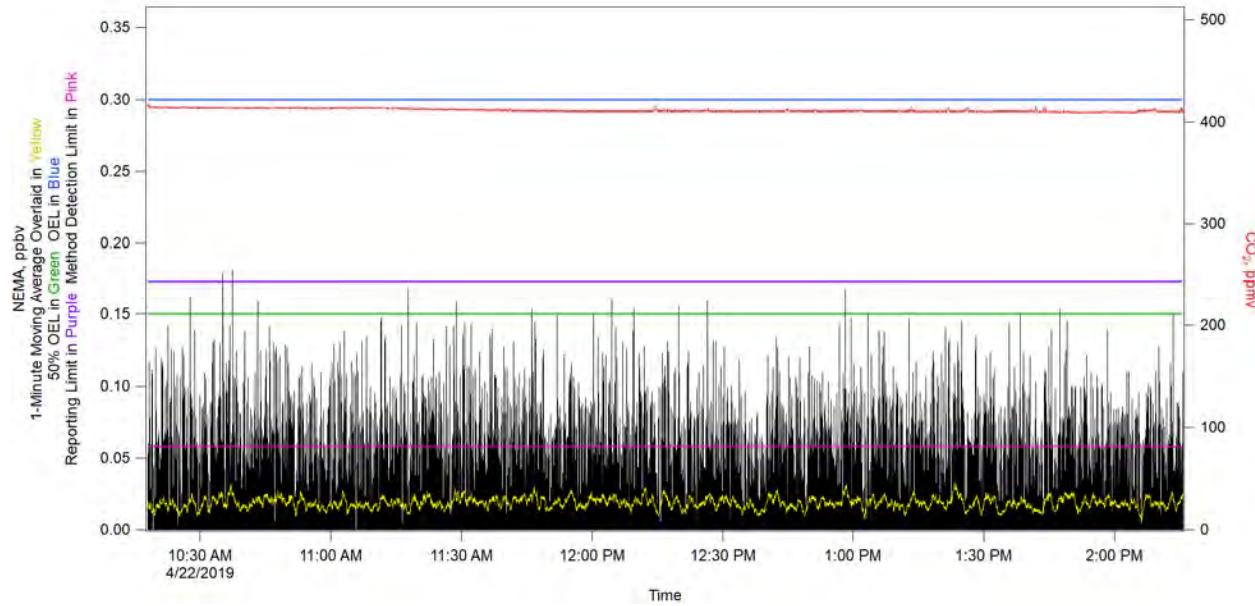


Figure 2-75. N-nitrosomethylethylamine (NEMA).

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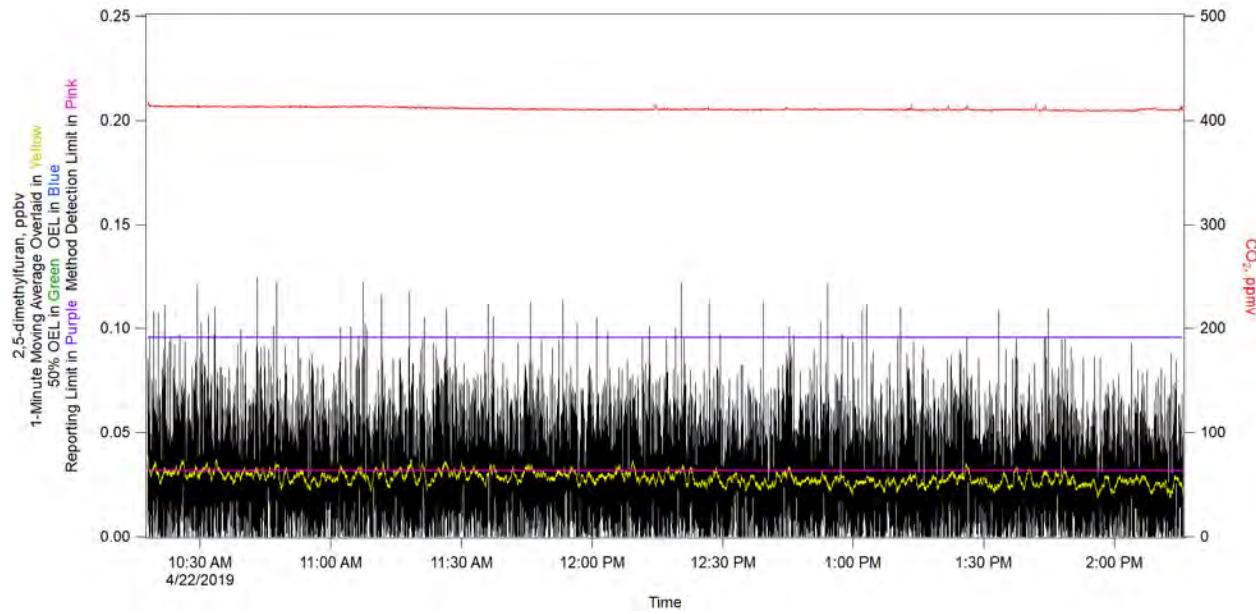


Figure 2-76. 2,5-dimethylfuran.

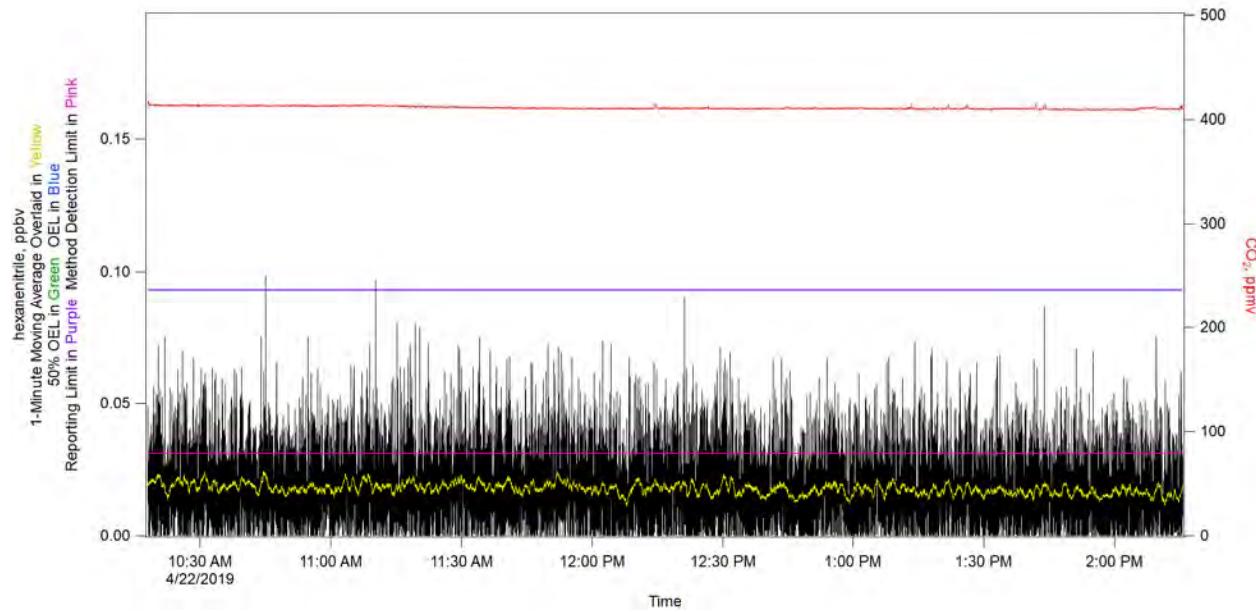


Figure 2-77. Hexanenitrile.

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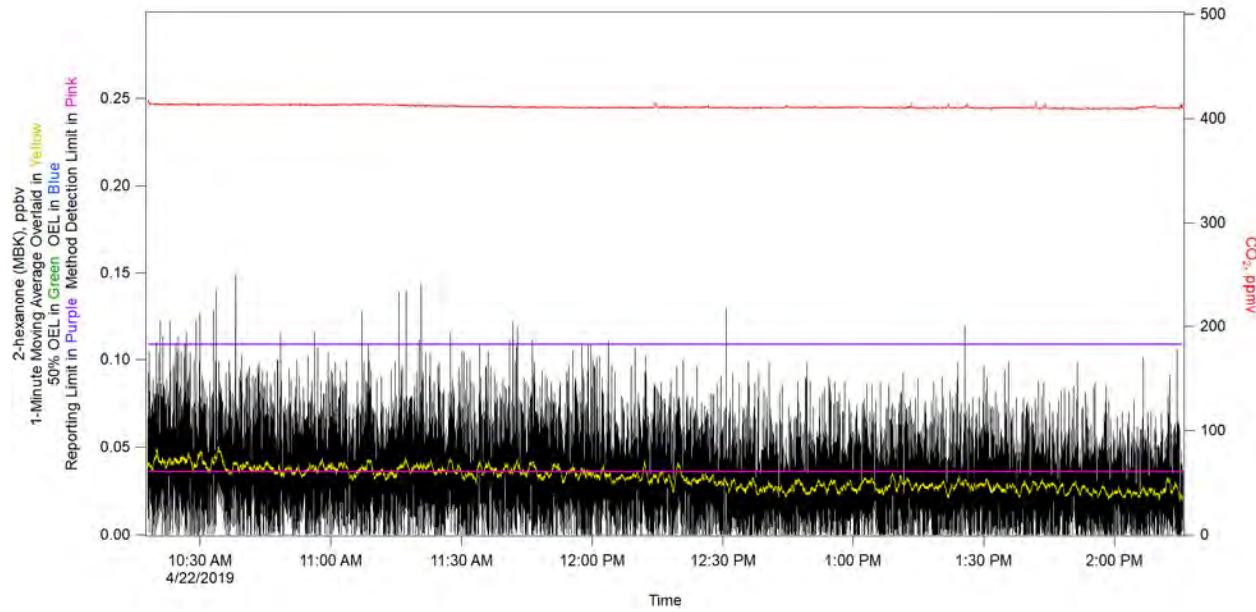


Figure 2-78. 2-hexanone (MBK).

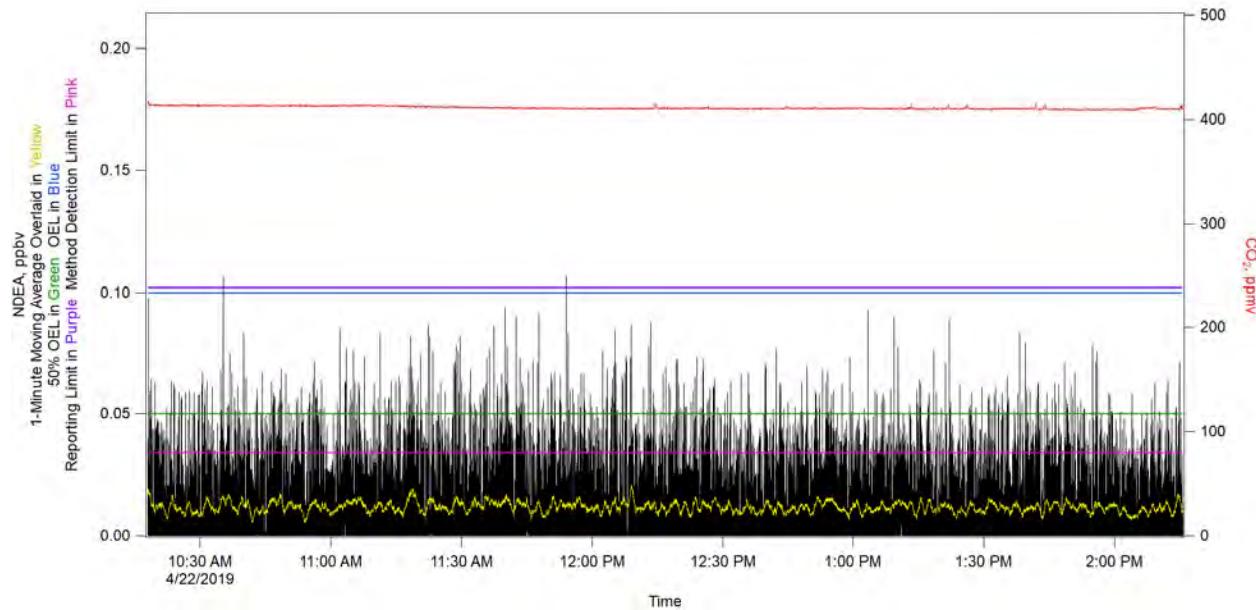


Figure 2-79. N-nitrosodiethylamine (NDEA).

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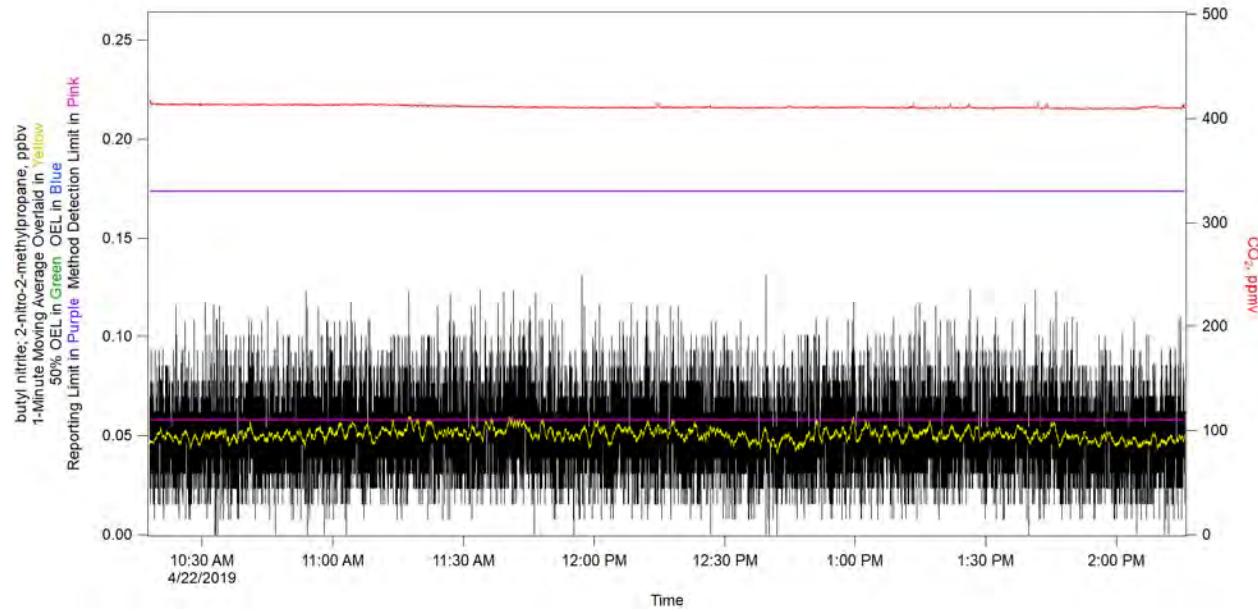


Figure 2-80. Butyl Nitrite; 2-nitro-2-methylpropane.

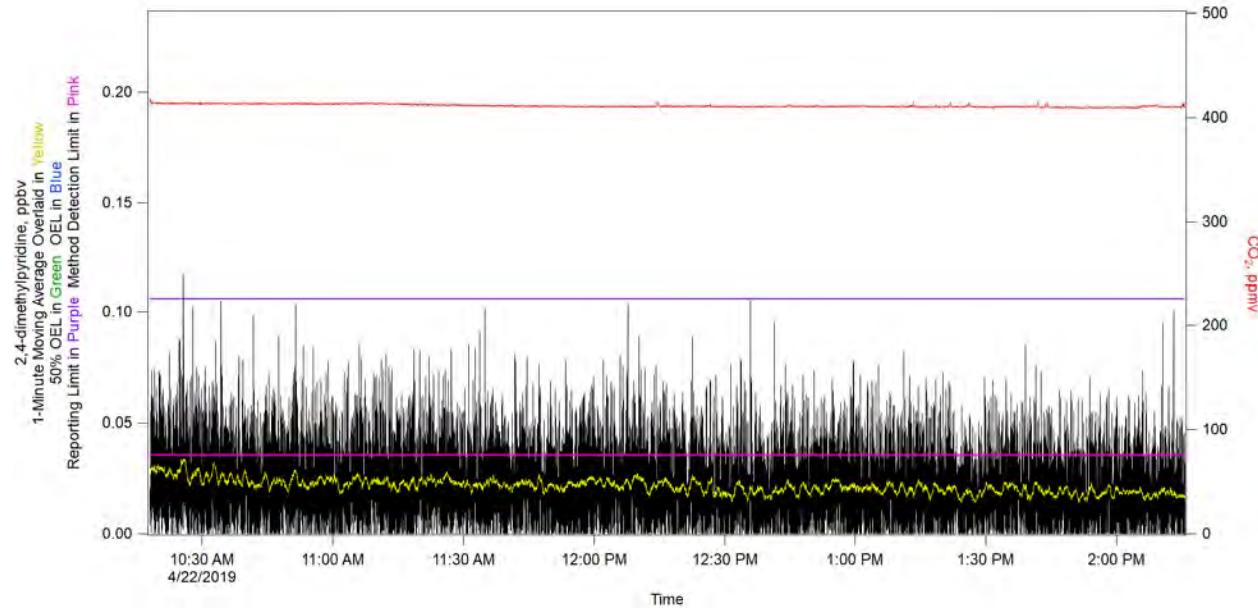


Figure 2-81. 2,4-dimethylpyridine.

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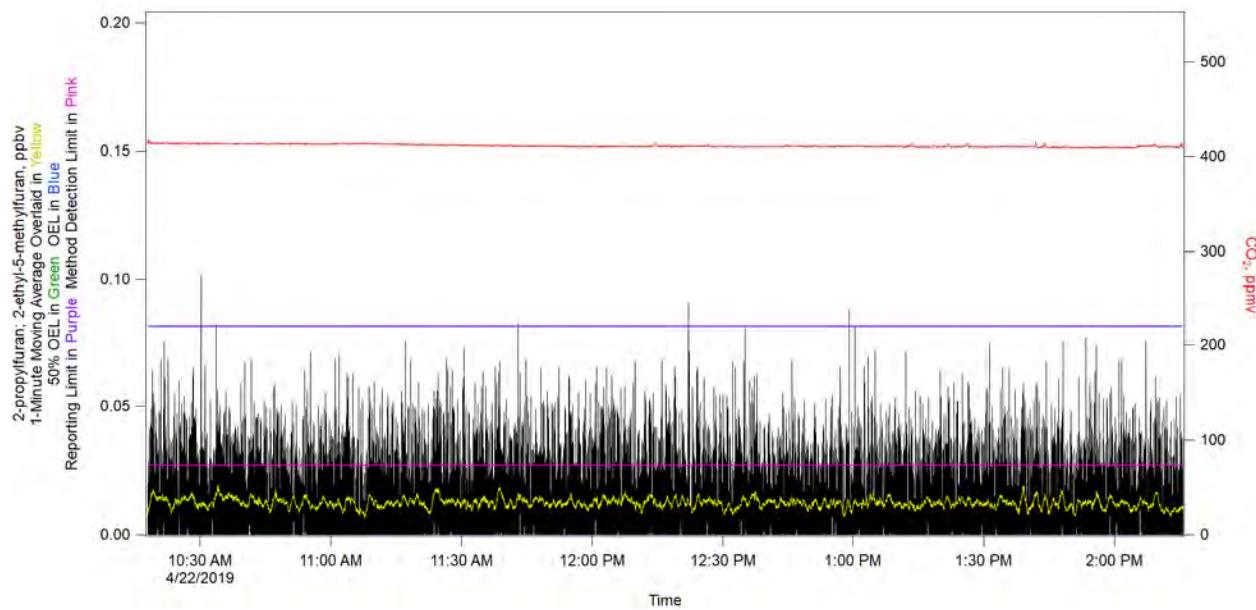


Figure 2-82. 2-propylfuran; 2-ethyl-5-methylfuran.

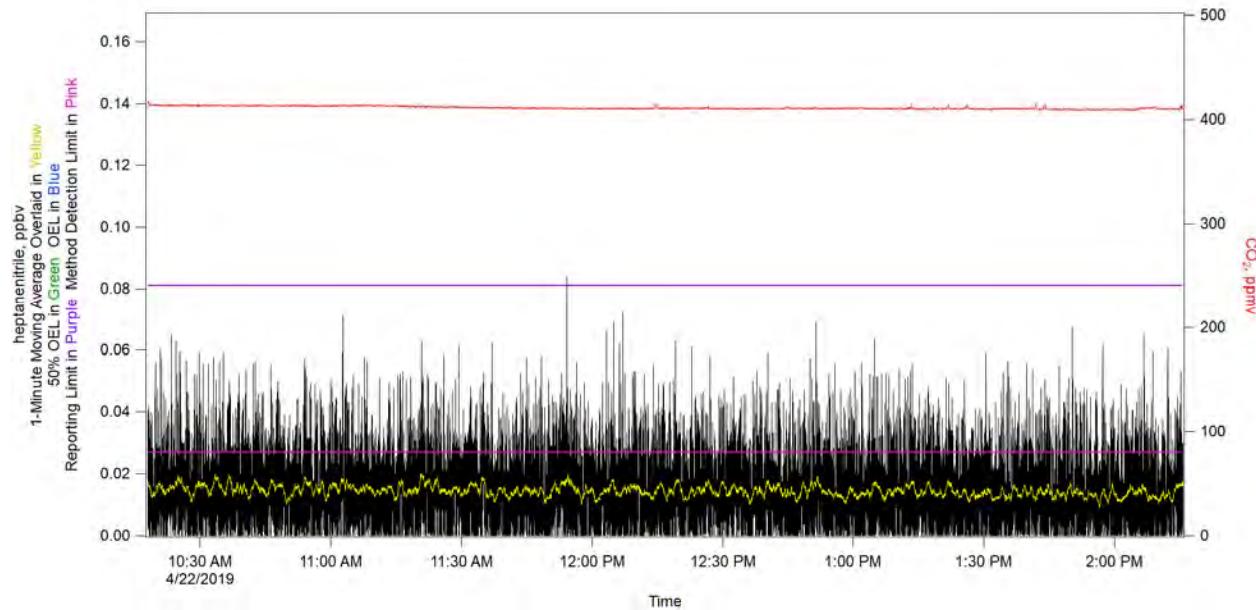


Figure 2-83. Heptanenitrile.

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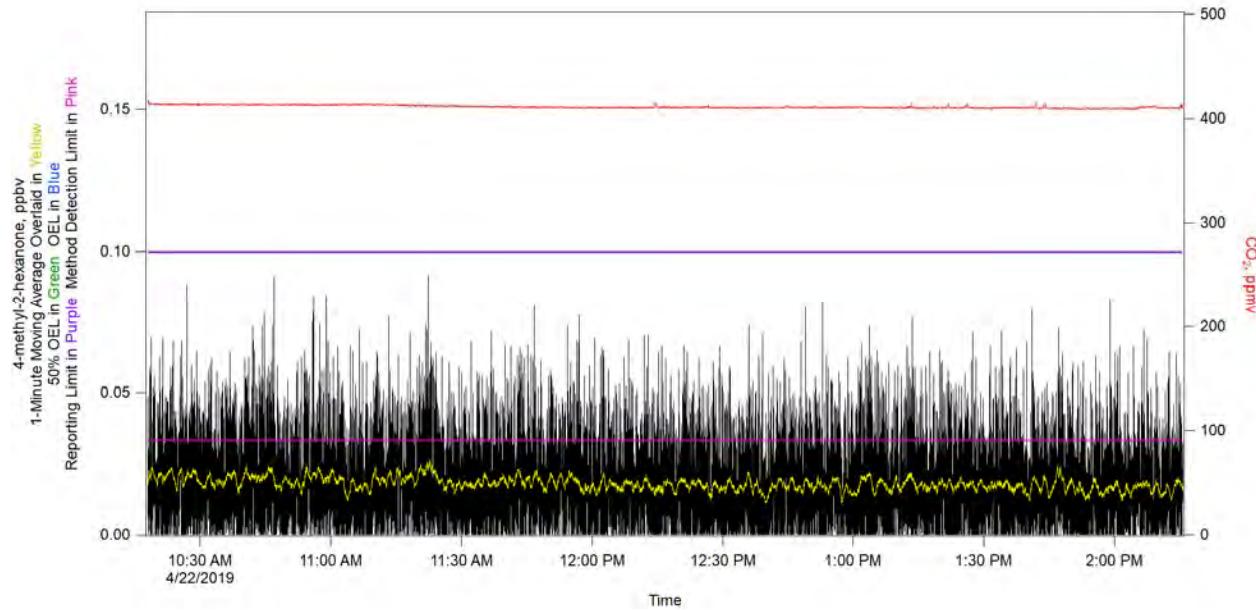


Figure 2-84. 4-methyl-2-hexanone.

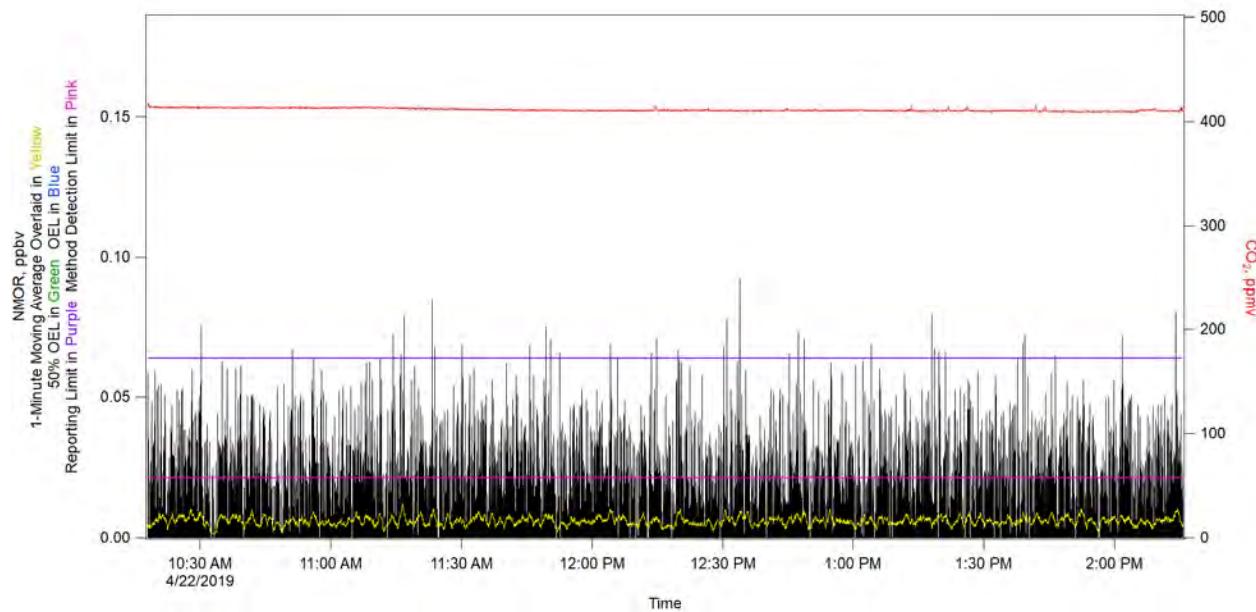


Figure 2-85. N-nitrosomorpholine (NMOR).

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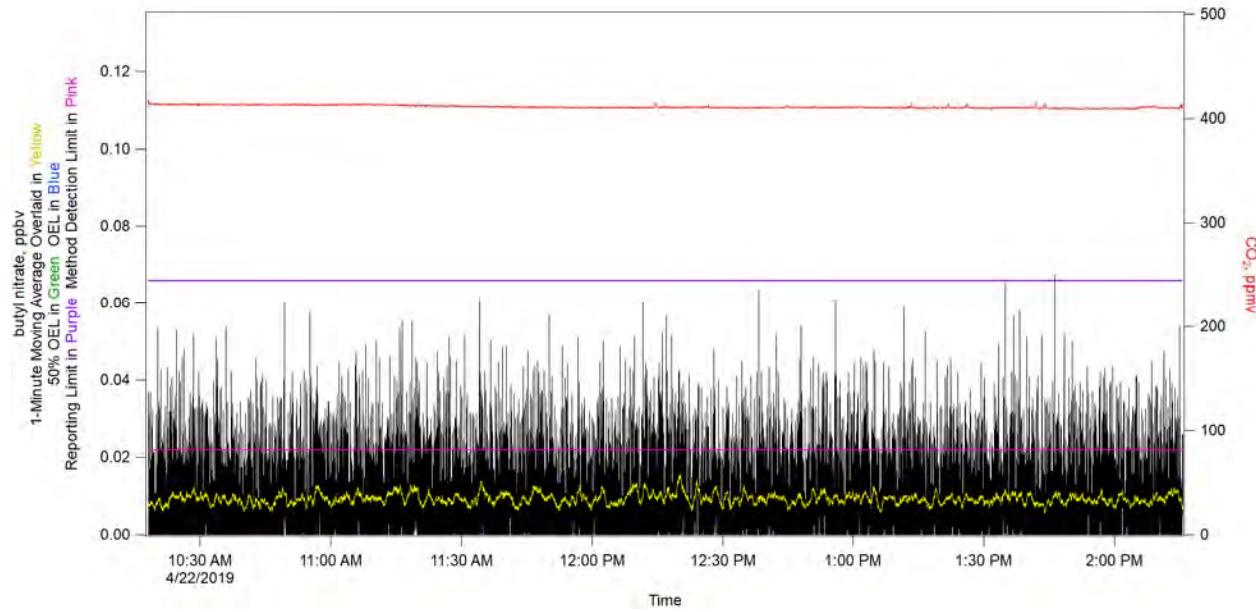
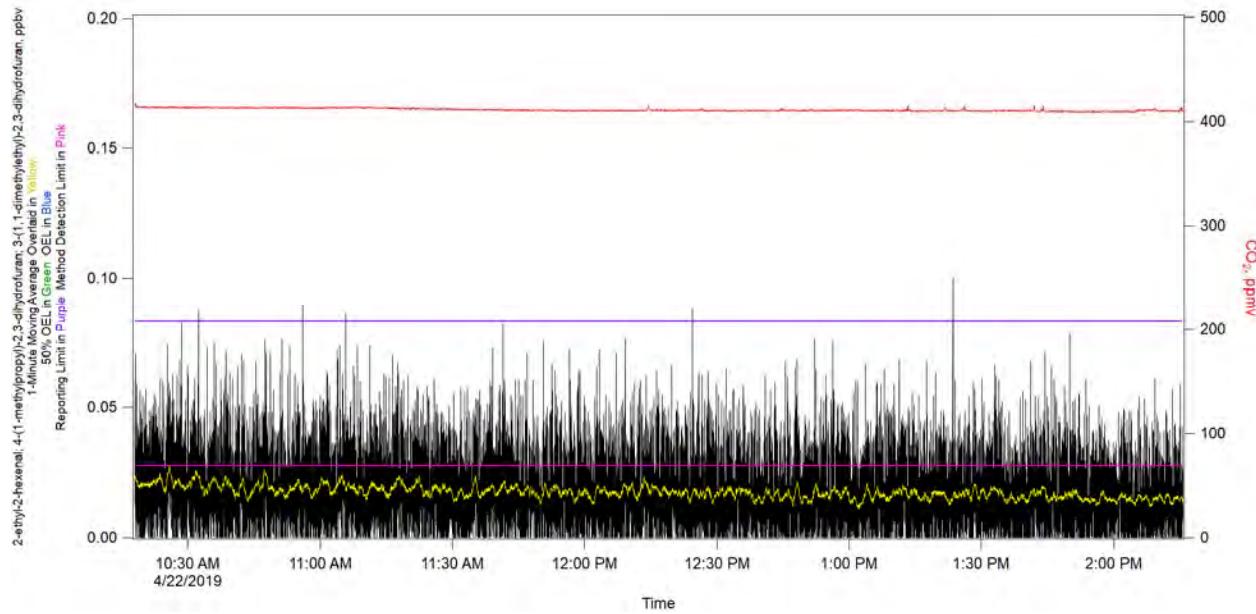


Figure 2-86. Butyl Nitrate.



**Figure 2-87. 2-ethyl-2-hexenal; 4-(1-methylpropyl)-2,3-dihydrofuran
3-(1,1-dimethylethyl)-2,3-dihydrofuran.**

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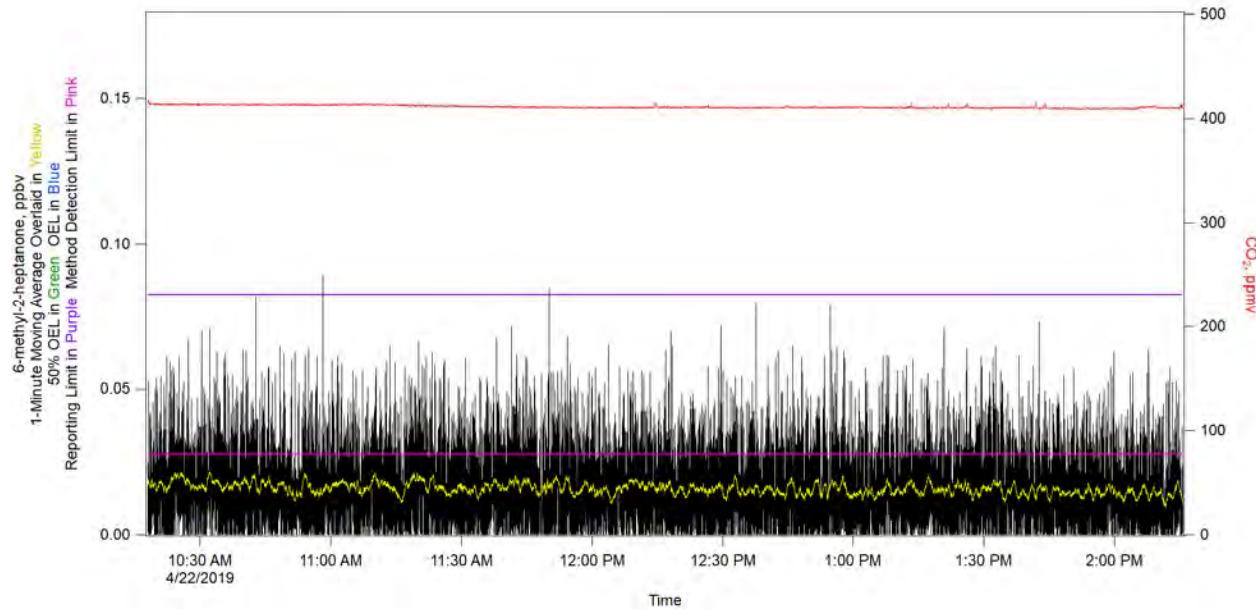


Figure 2-88. 6-methyl-2-heptanone.

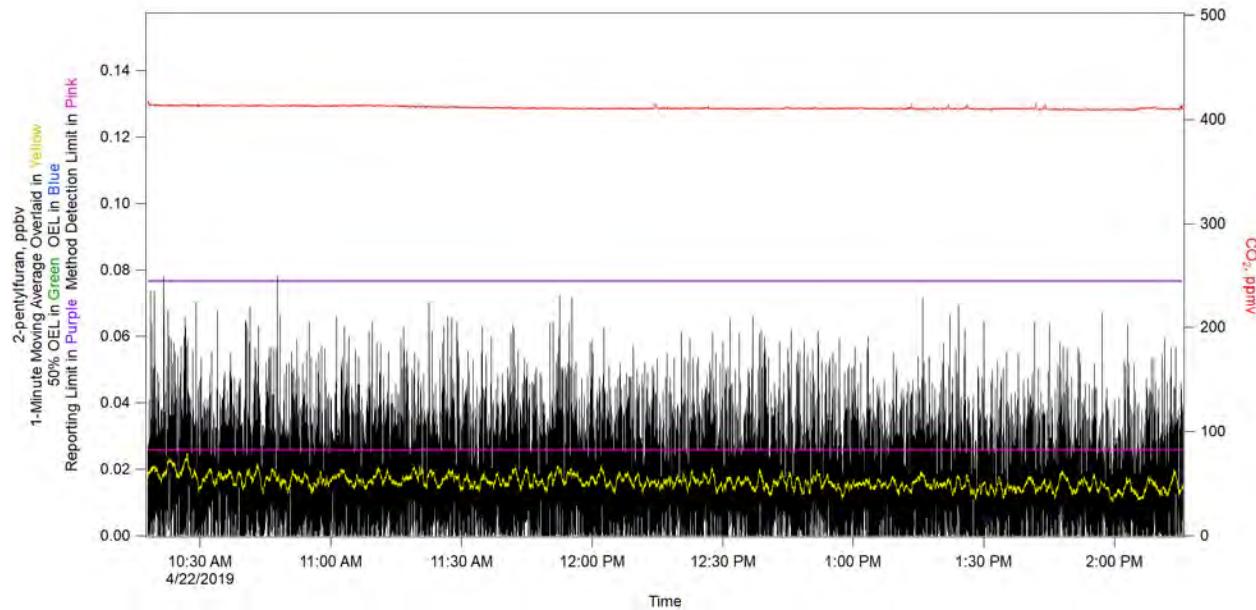


Figure 2-89. 2-pentylfuran.

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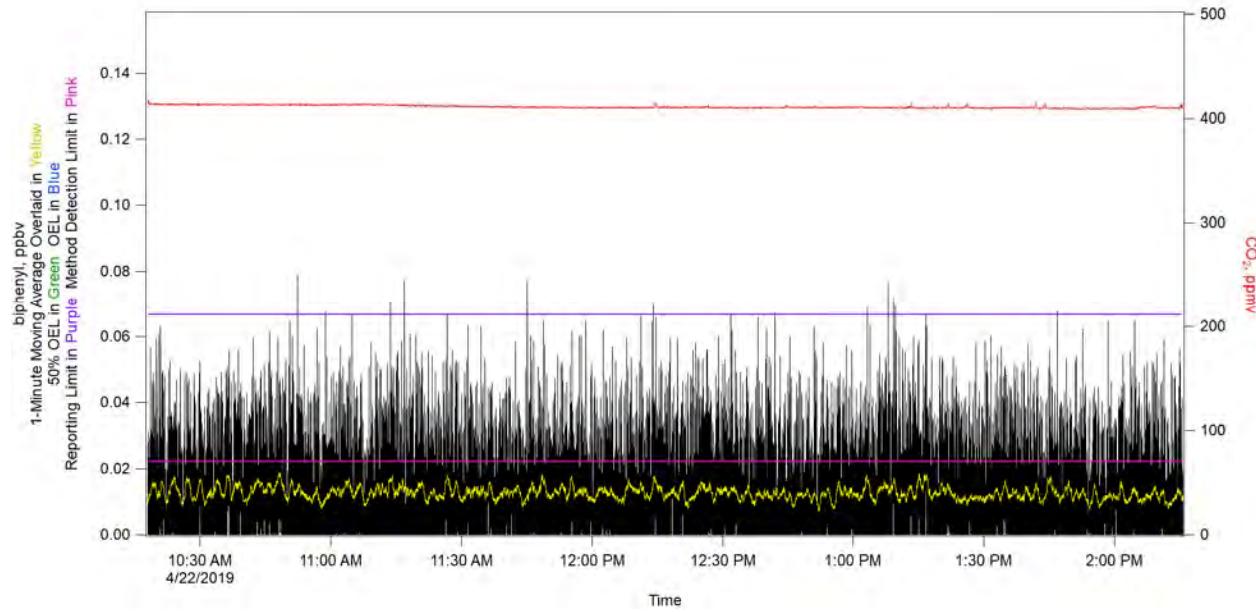


Figure 2-90. Biphenyl.

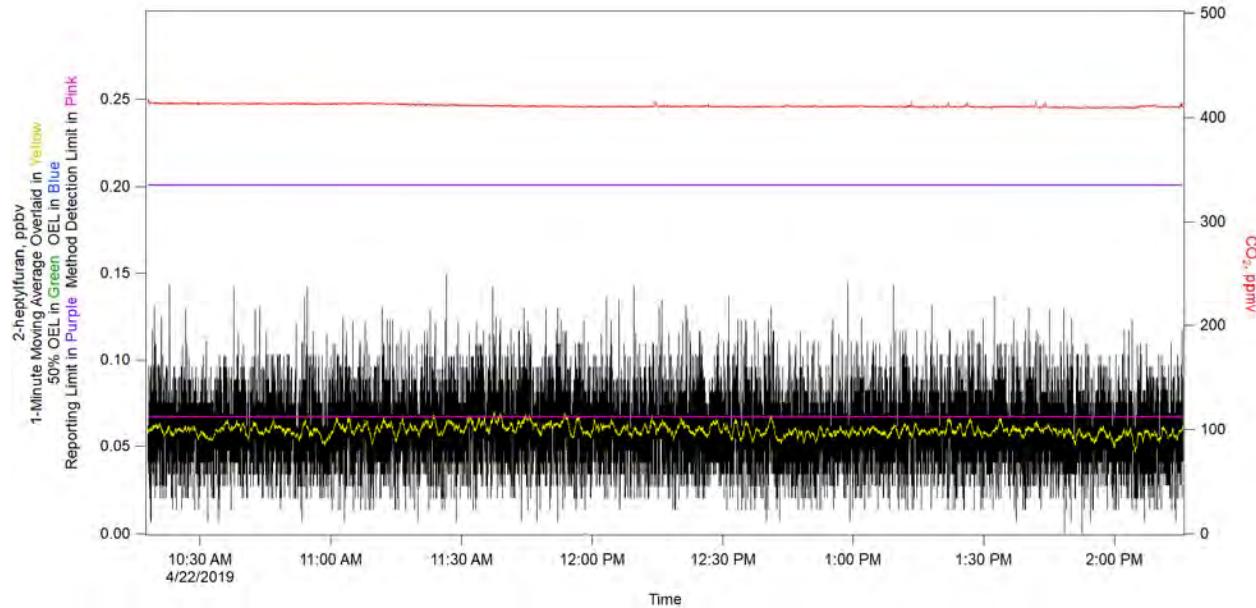


Figure 2-91. 2-heptylfuran.

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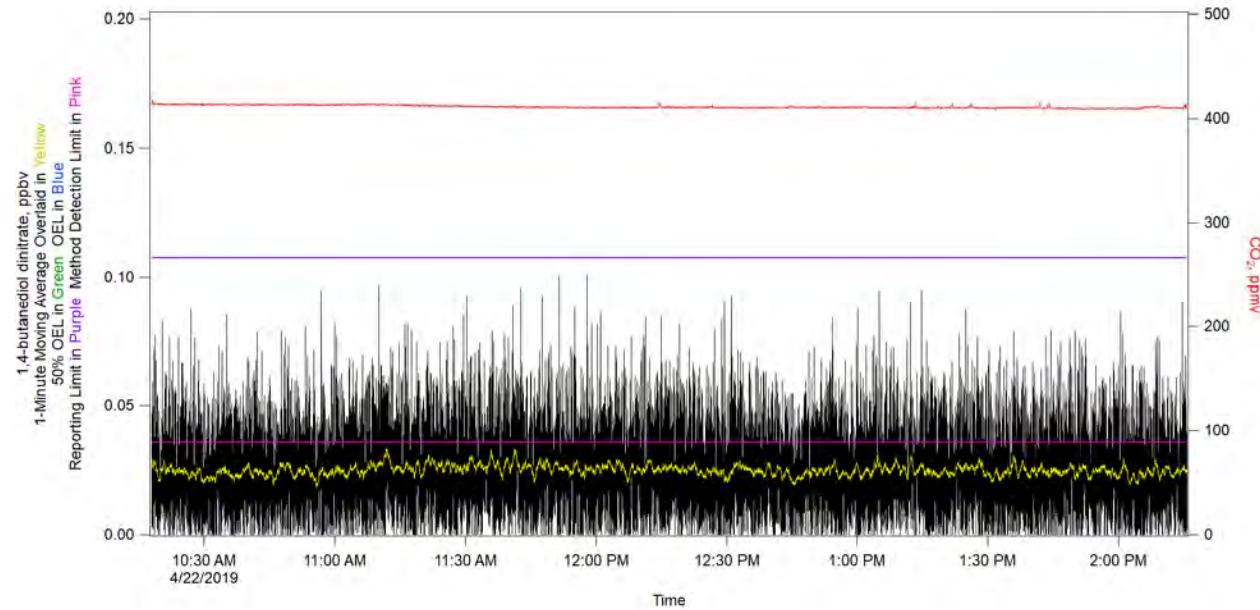


Figure 2-92. 1,4-butanediol Dinitrate.

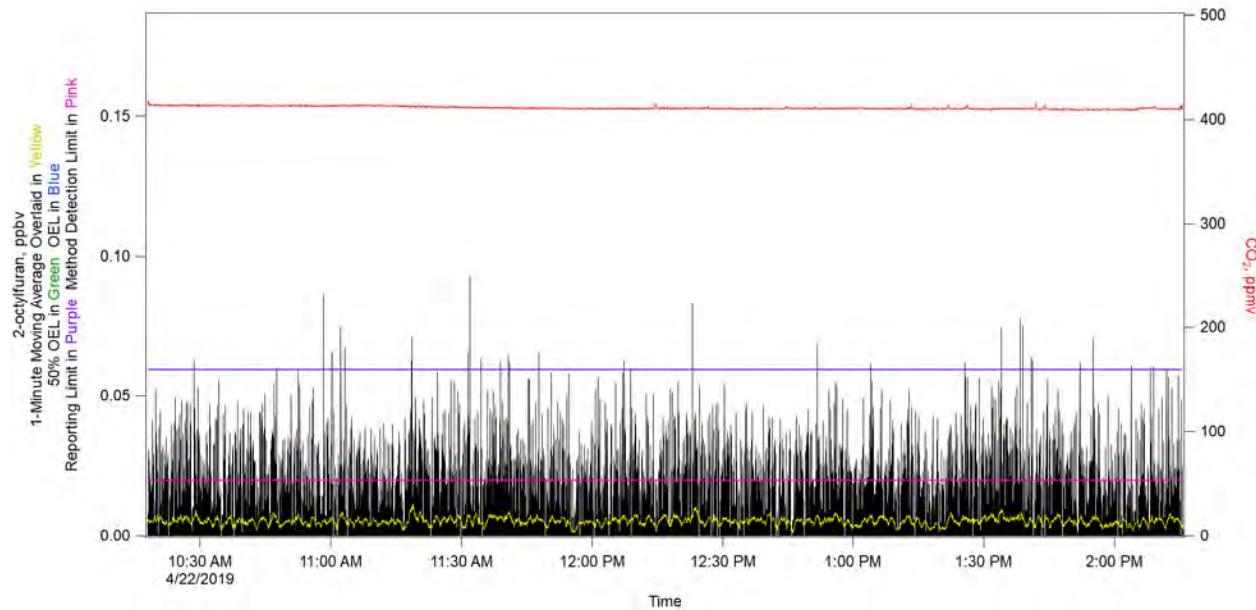


Figure 2-93. 2-octylfuran.

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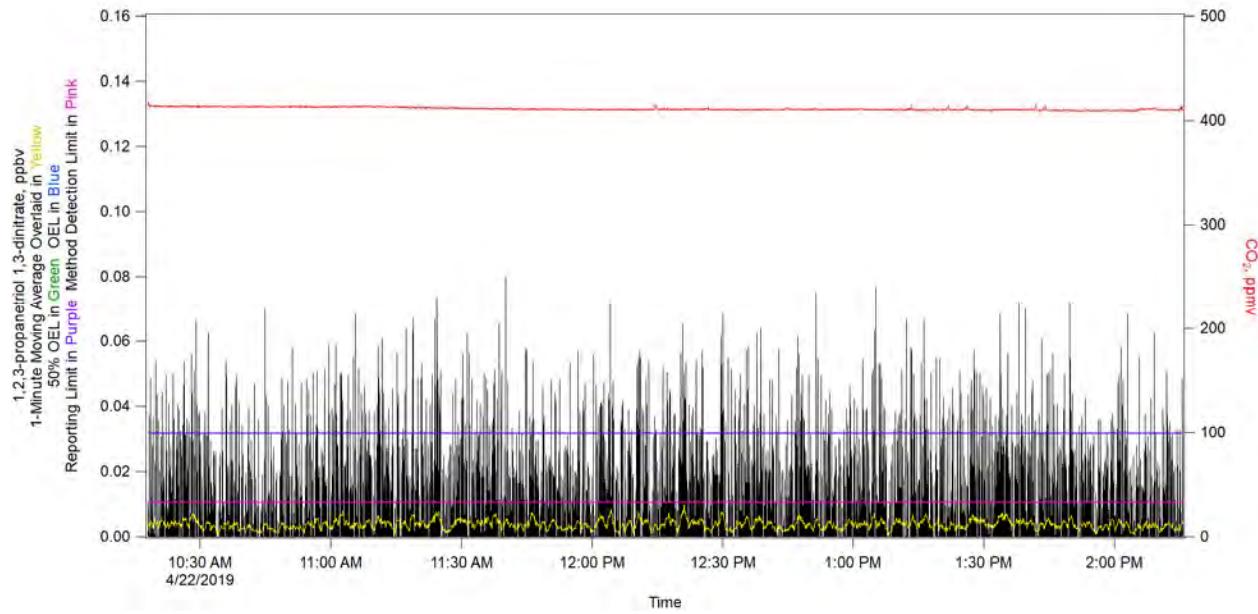


Figure 2-94. 1,2,3-propanetriol 1,3-dinitrate.

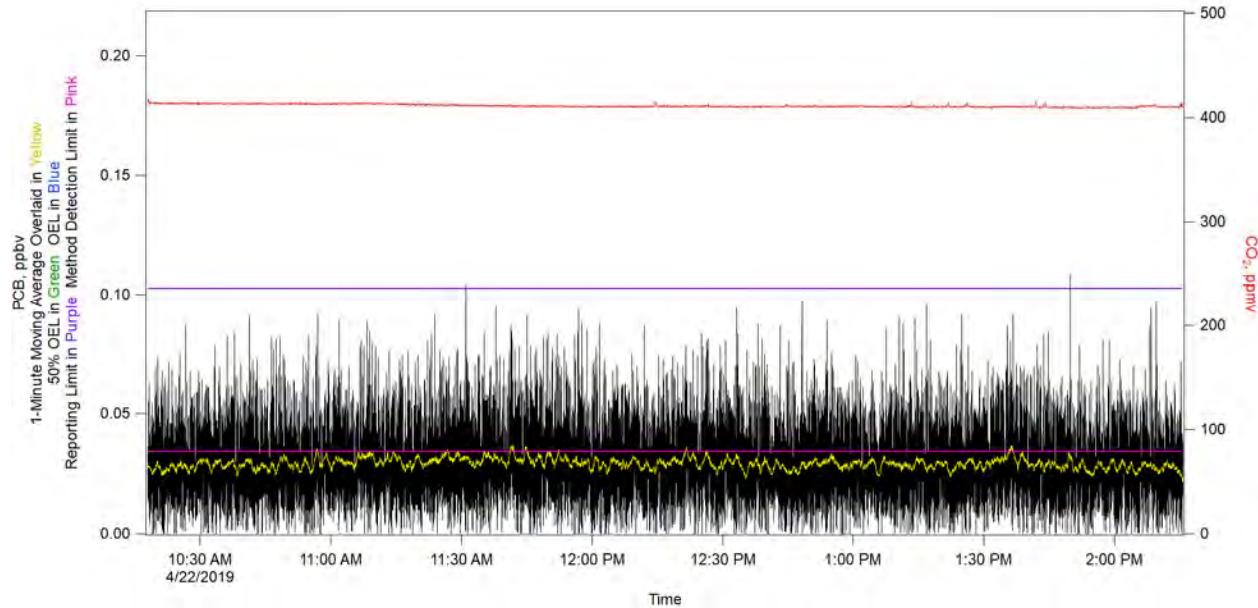


Figure 2-95. PCB.

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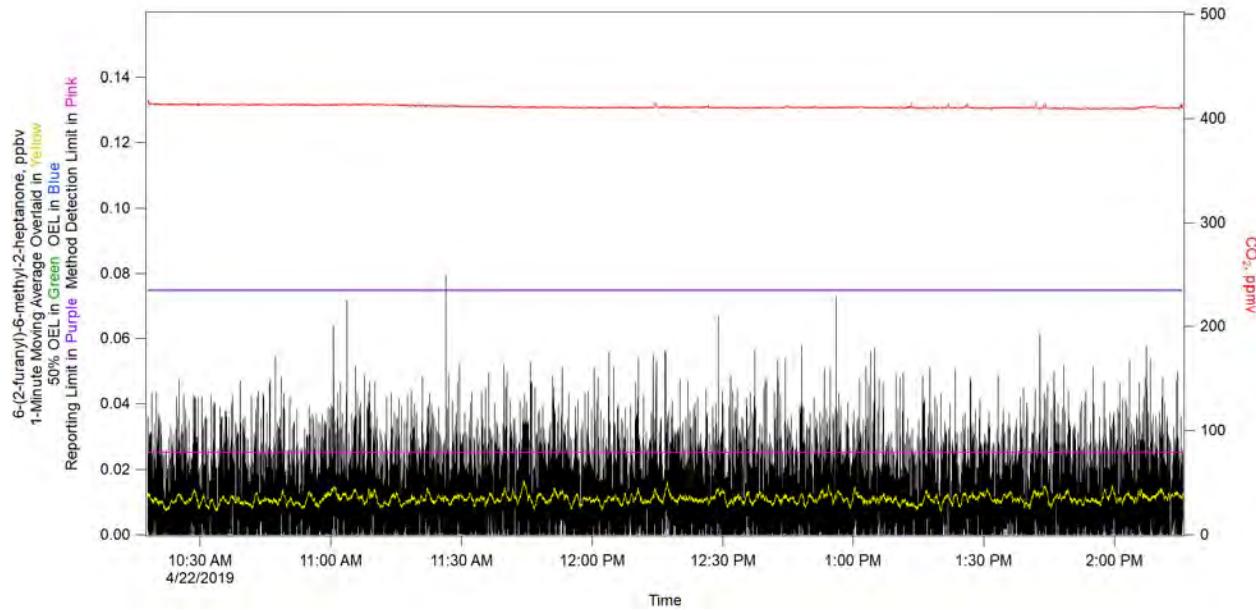


Figure 2-96. 6-(2-furanyl)-6-methyl-2-heptanone.

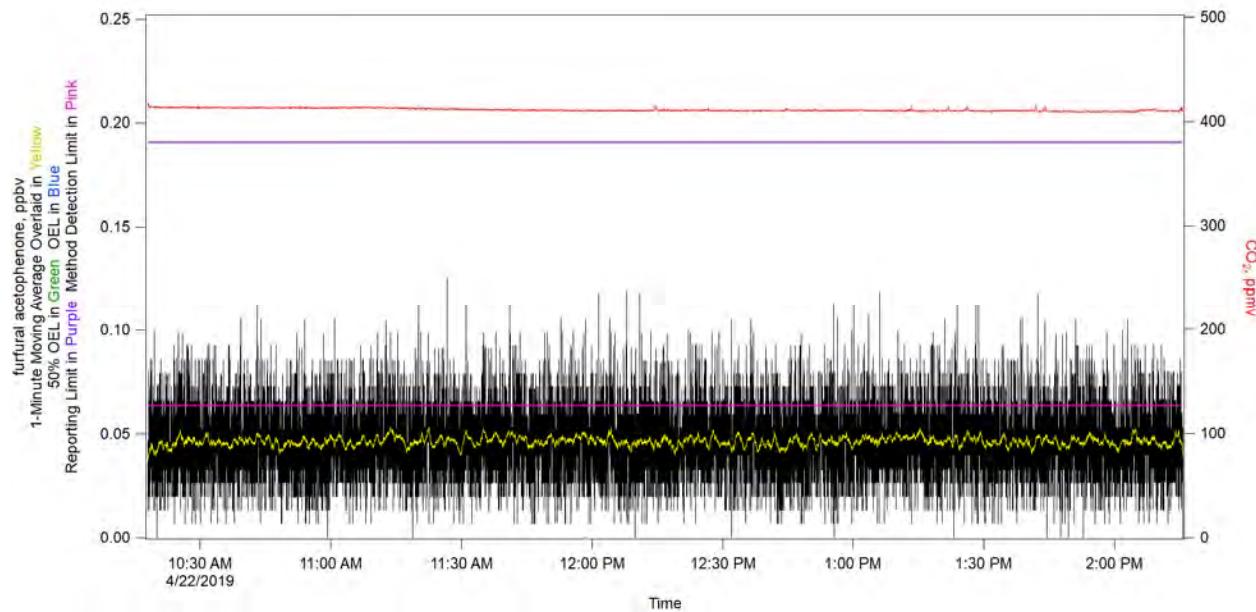


Figure 2-97. Furfural Acetophenone.

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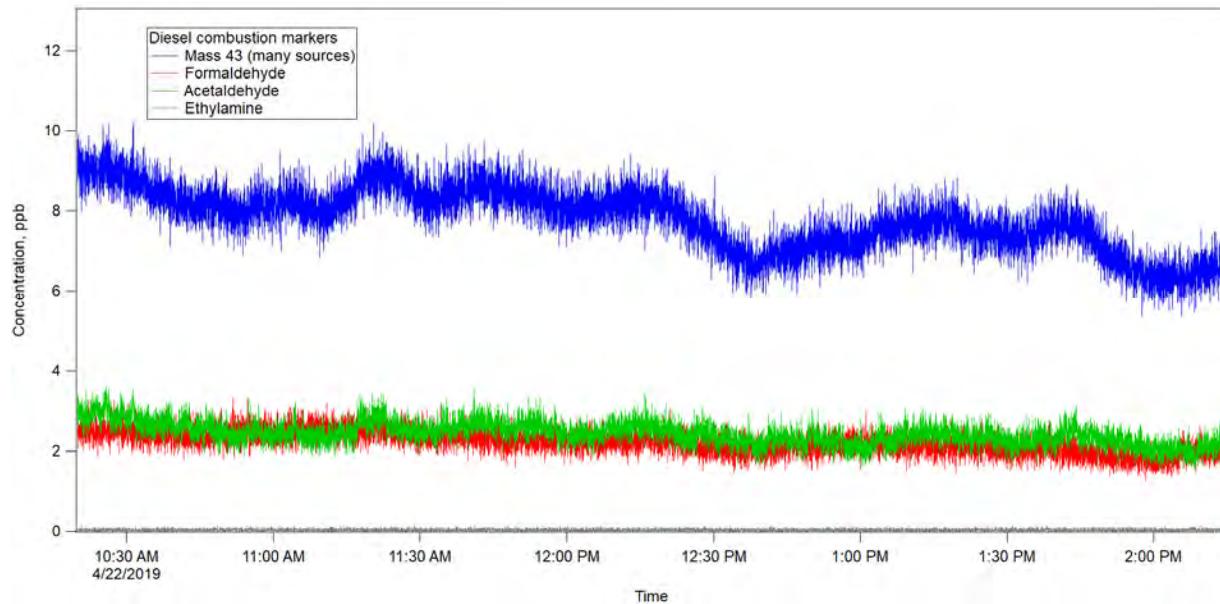


Figure 2-98. Diesel Combustion Markers.

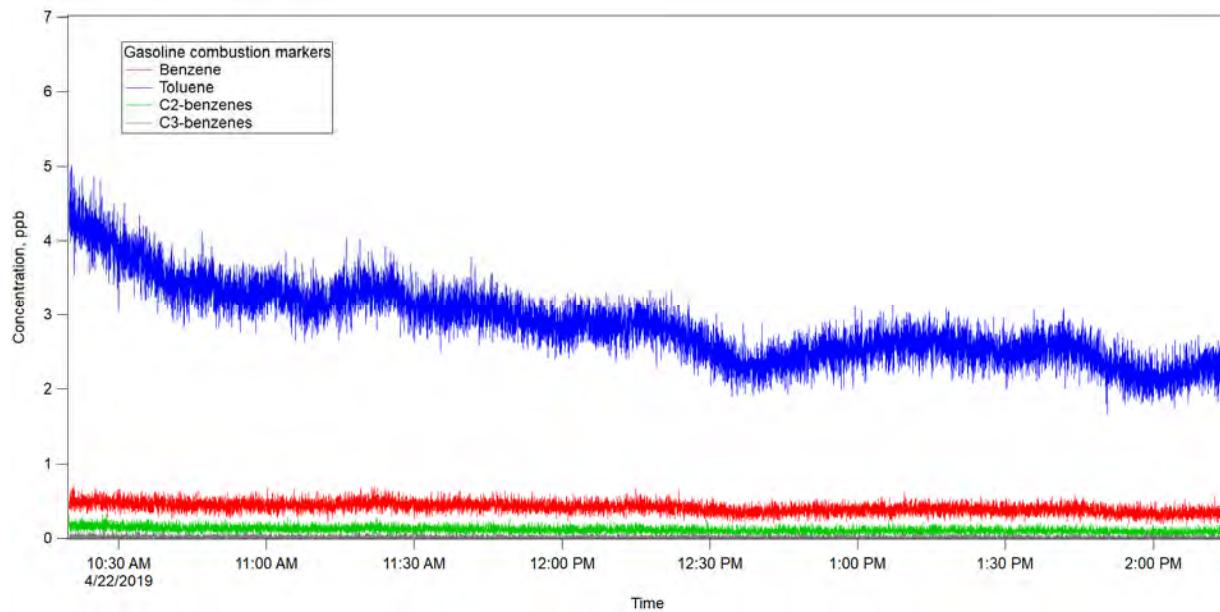


Figure 2-99. Gasoline Combustion Markers.

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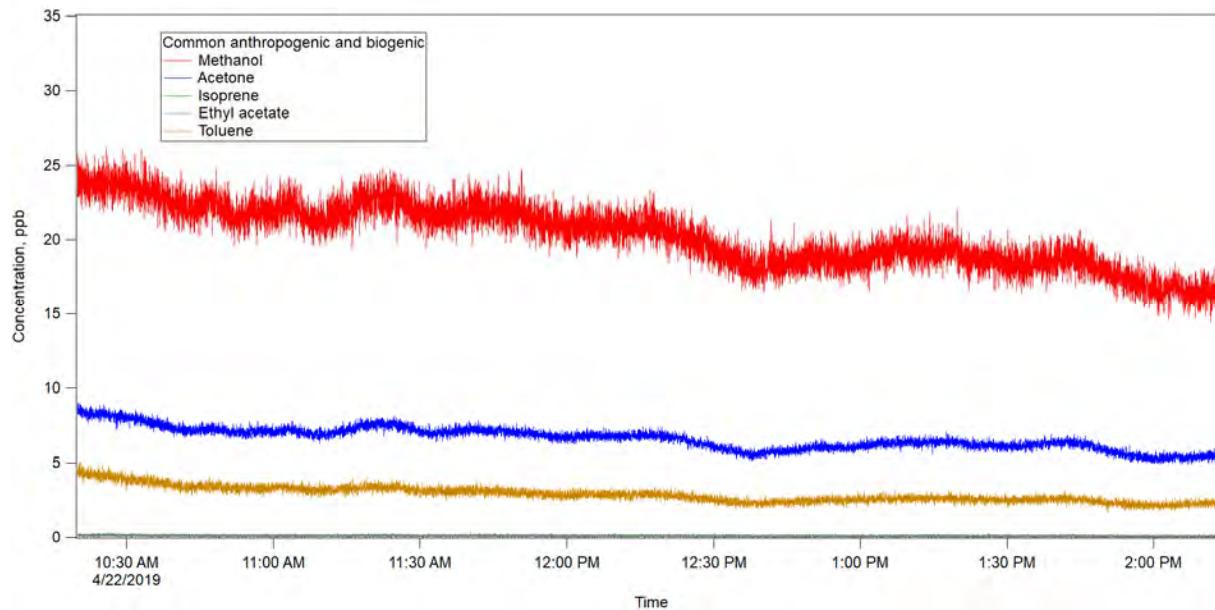


Figure 2-100. Plant and Human Markers.

2.6 Odor Compounds

The ML monitored the AOP-015 event near the U Farm Change Trailer from approximately 10:17 to 14:19. This section provides the odor compound statistical results from that monitoring period.

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Table 2-3. Odor Statistical Information for AOP-015 Monitoring on April 22, 2019.

Odor #	Odor Compound Name	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel St. Dev. (%)	Max. (ppb)	Median (ppb)
1	methyl mercaptan	0.147	<0.147	0.044	30.407	0.366	<0.147
2	Dimethylsulfide + ethanethiol	0.088	0.193†	0.045	23.122	0.402	0.193†
3	allyl mercaptan	0.033	<0.033	0.013	201.445	0.126	<0.033
4	1-propanethiol + isopropyl mercaptan	0.041	<0.041	0.023	139.480	0.182	<0.041
5	2-butene-1-thiol	0.062	<0.062	0.012	263.193	0.147	<0.062
6	diethyl sulfide + 2-methylpropane-2-thiol	0.281	0.373†	0.089	23.980	0.824	0.365†
7	thiopropanal sulfuroxide	0.024	0.126	0.051	40.564	0.353	0.123
8	dimethyl disulfide	0.027	0.053†	0.030	57.077	0.203	0.051†
9	1-pantanethiol + 2,2-dimethylpropane-1-thiol	0.046	<0.046	0.019	121.401	0.121	<0.046
10	benzenethiol	0.028	0.107	0.057	52.758	0.320	0.105
11	diallyl sulfide	0.022	<0.022	0.013	193.852	0.130	<0.022
12	methyl propyl disulfide	0.016	0.026†	0.032	124.046	0.187	0.011†
13	methylbenzenethiol	0.027	<0.027	0.017	114.929	0.111	<0.027
14	dimethyl trisulfide	0.020	<0.020	0.010	157.053	0.069	<0.020
15	(1-oxoethyl) thiophene	0.040	<0.040	0.024	170.062	0.164	<0.040
16	(1-oxopropyl) thiophene	0.029	0.031†	0.024	75.714	0.138	0.028†
17	dipropyl disulfide	0.024	<0.024	0.013	102.344	0.087	<0.024
18	methyl propyl trisulfide	0.019	<0.019	0.013	160.881	0.089	<0.019
19	dimethyl tetrasulfide	0.021	<0.021	0.009	92.339	0.057	<0.021
20	dipropyl trisulfide	0.022	<0.022	0.013	177.118	0.096	<0.022
21	diphenyl sulfide	0.026	<0.026	0.014	137.747	0.099	<0.026
<	COPC Average/Median Below the MDL.						
†	Average/Median Between the RL and the MDL.						

Figure 2-101 through Figure 2-121 display potential odor-causing compounds, overlaid with the same signal smoothed using a one-minute moving average, and CO₂, for the monitoring period of April 22, 2019. If within range of the plot's left axis, a horizontal purple line representing the RL, and a pink horizontal line representing the MDL are shown.

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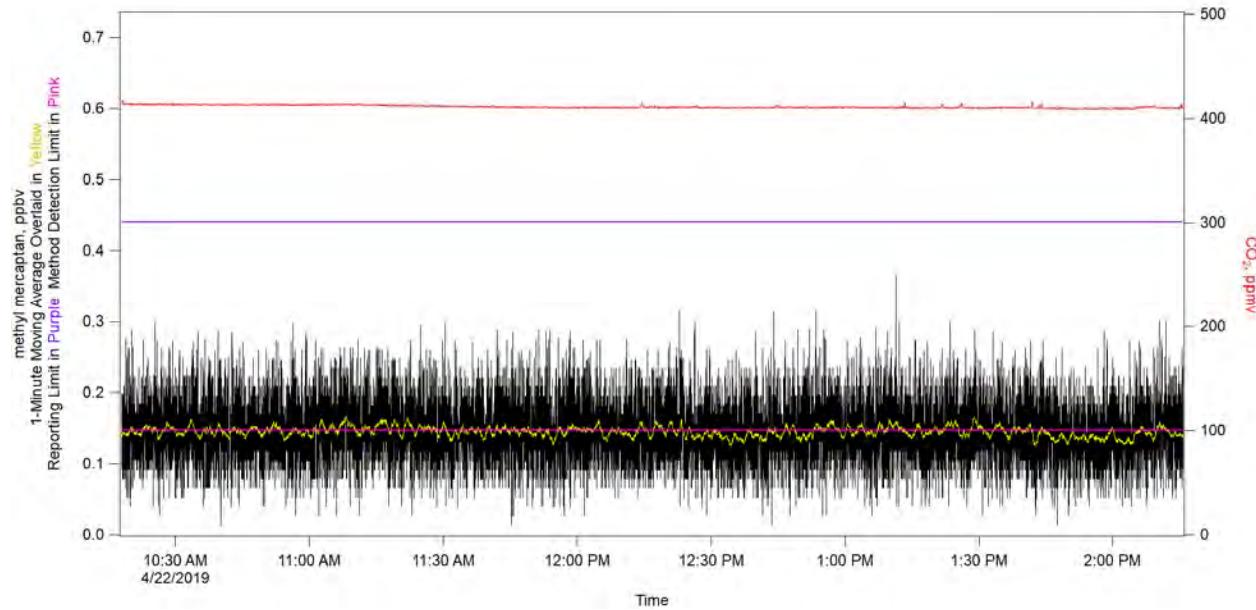


Figure 2-101. Methyl Mercaptan.

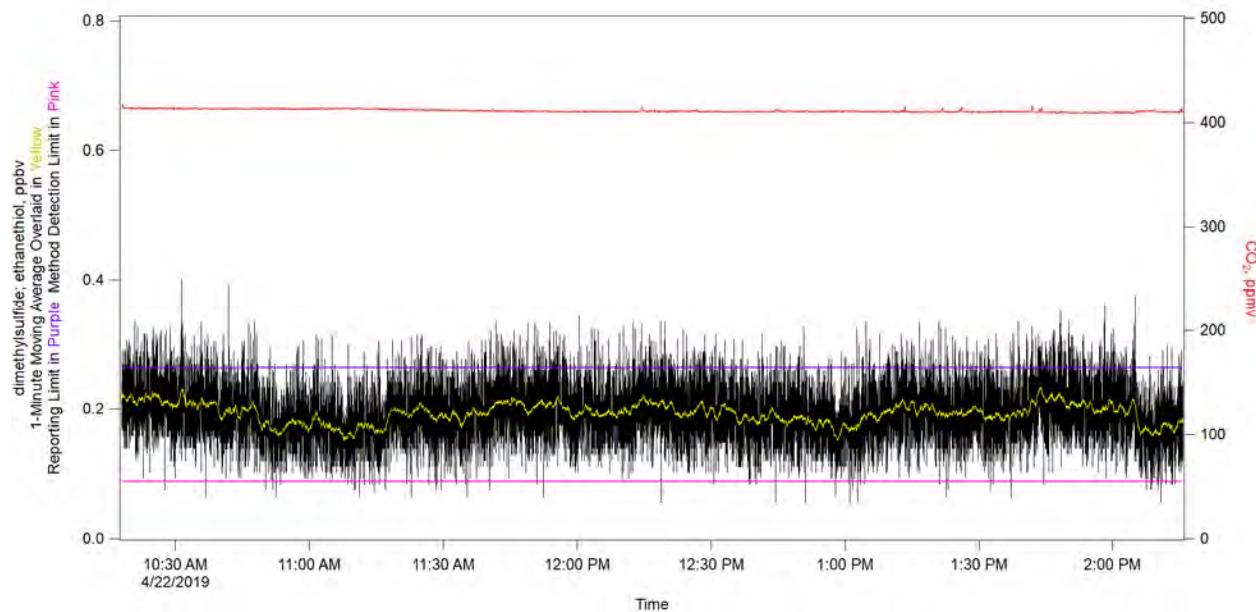


Figure 2-102. Dimethyl Sulfide; Ethanethiol.

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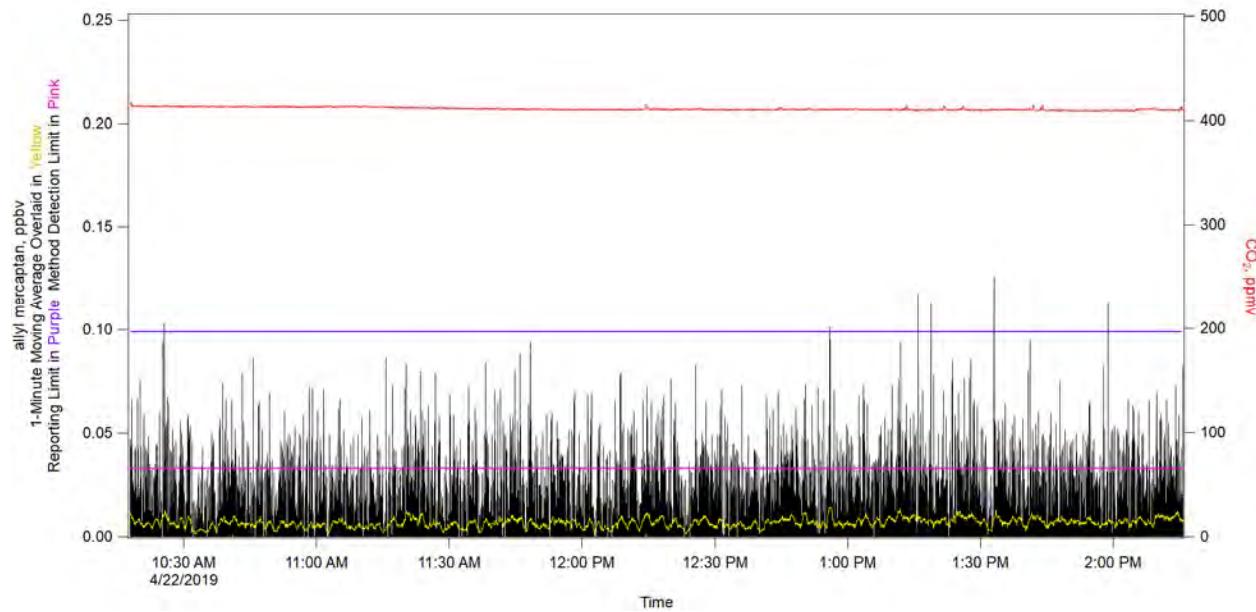


Figure 2-103. Allyl Mercaptan.

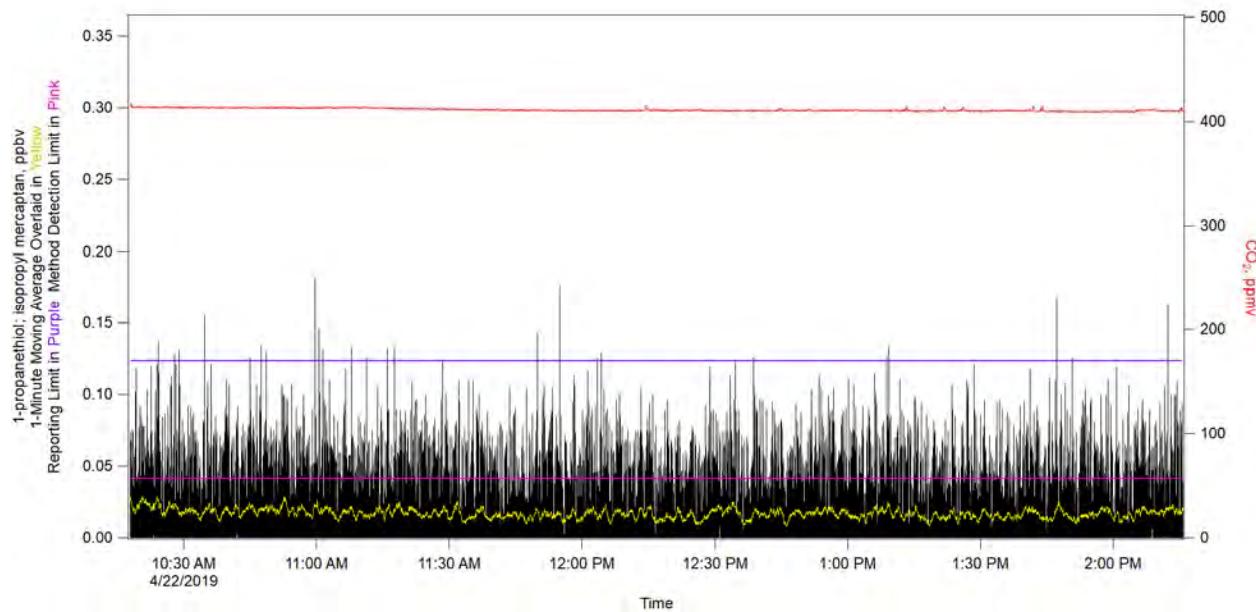


Figure 2-104. 1-propanethiol; Isopropyl Mercaptan.

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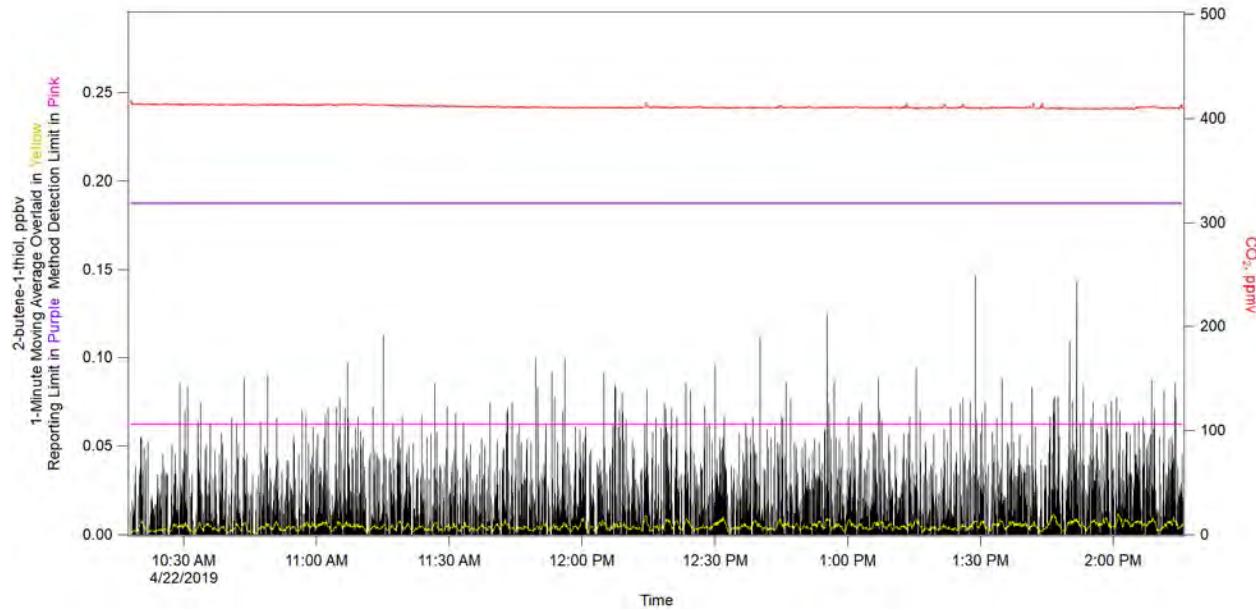


Figure 2-105. 2-butene-1-thiol.

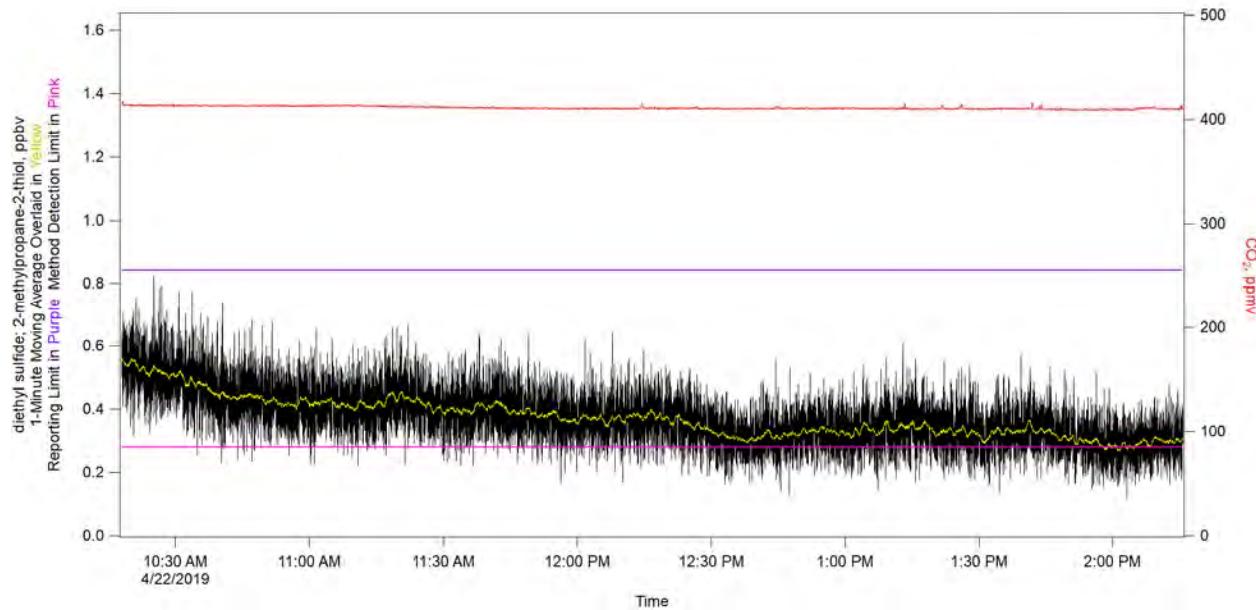


Figure 2-106. Diethyl Sulfide; 2-methylpropane-2-thiol.

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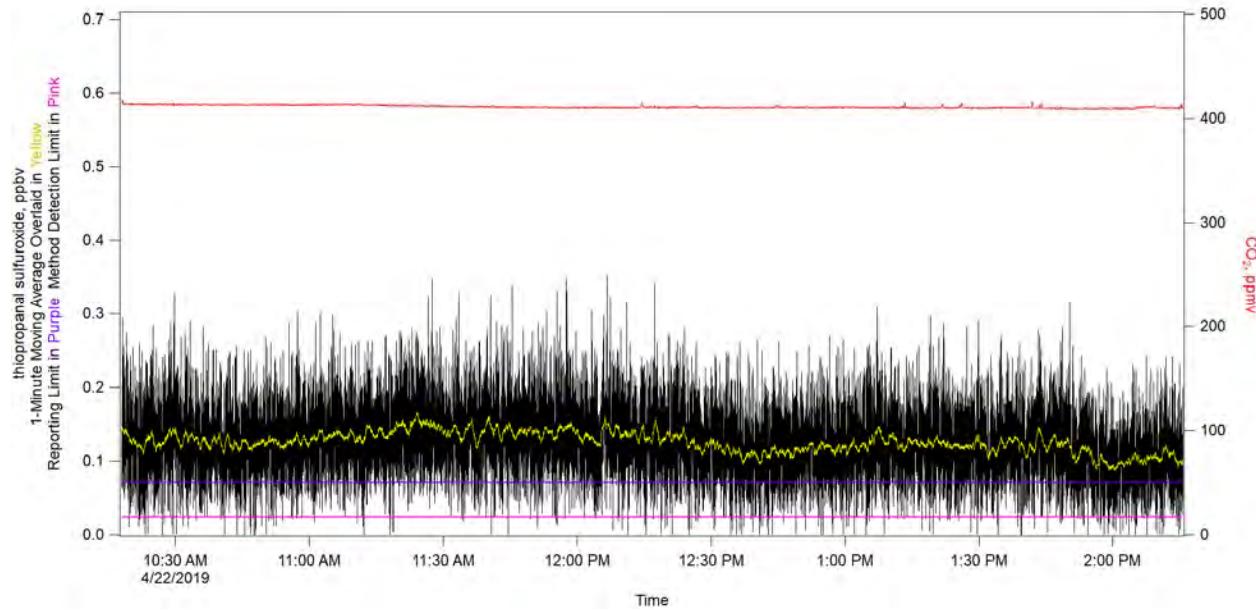


Figure 2-107. Thiopropanal Sulfuroxide.

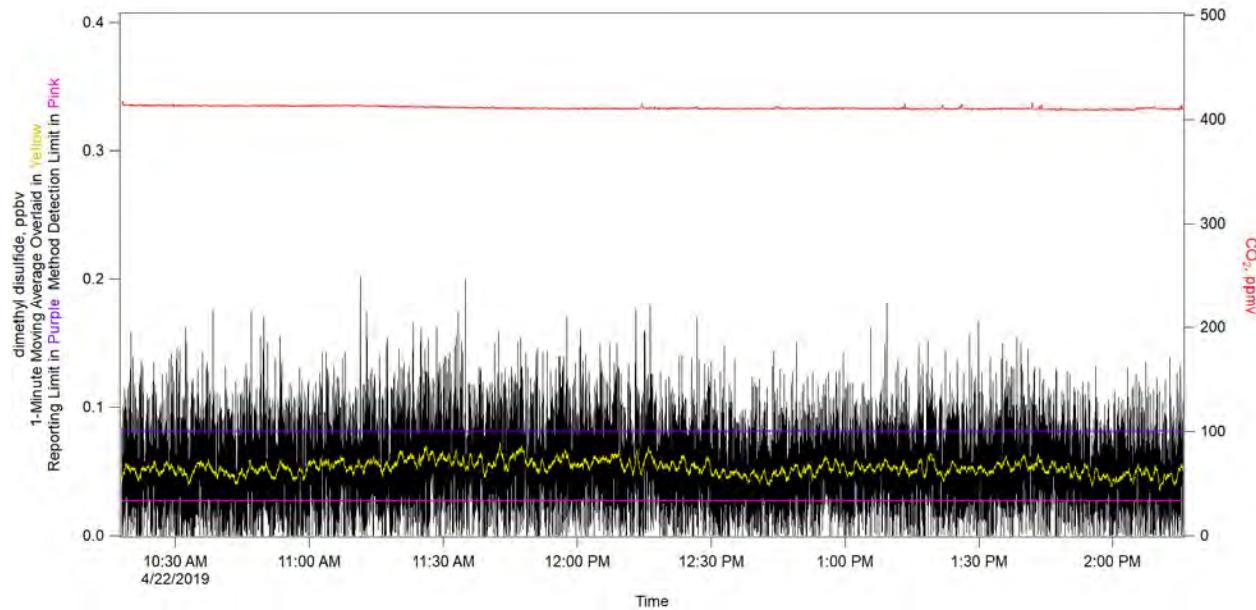


Figure 2-108. Dimethyl Disulfide.

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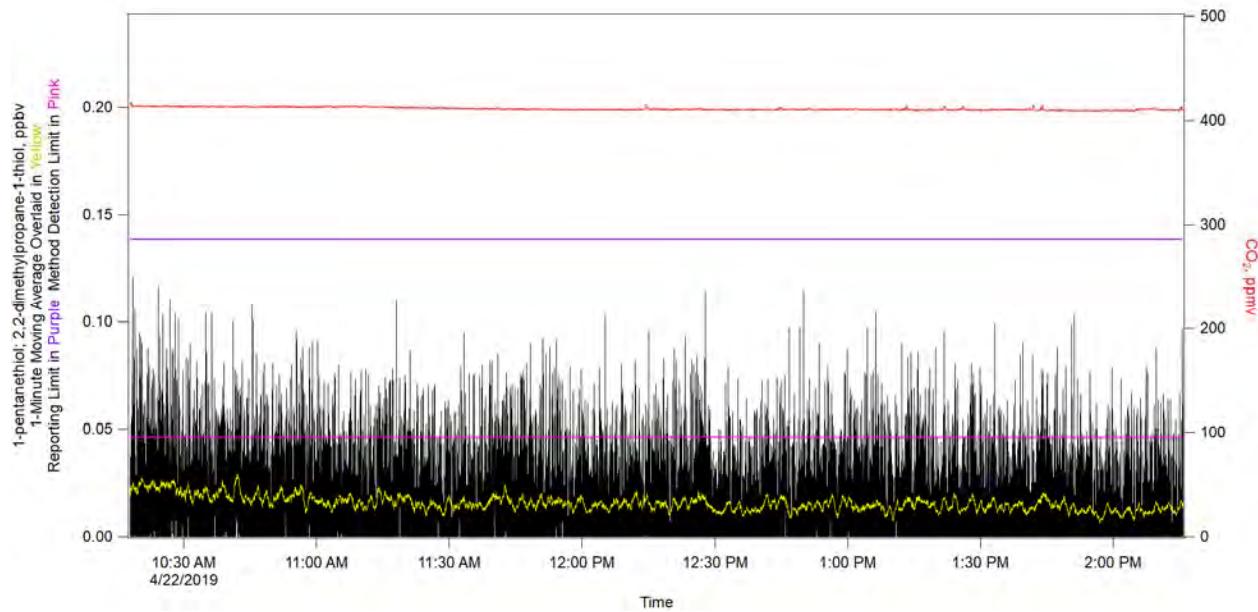


Figure 2-109. 1-pentanethiol; 2,2-dimethylpropane-1-thiol.

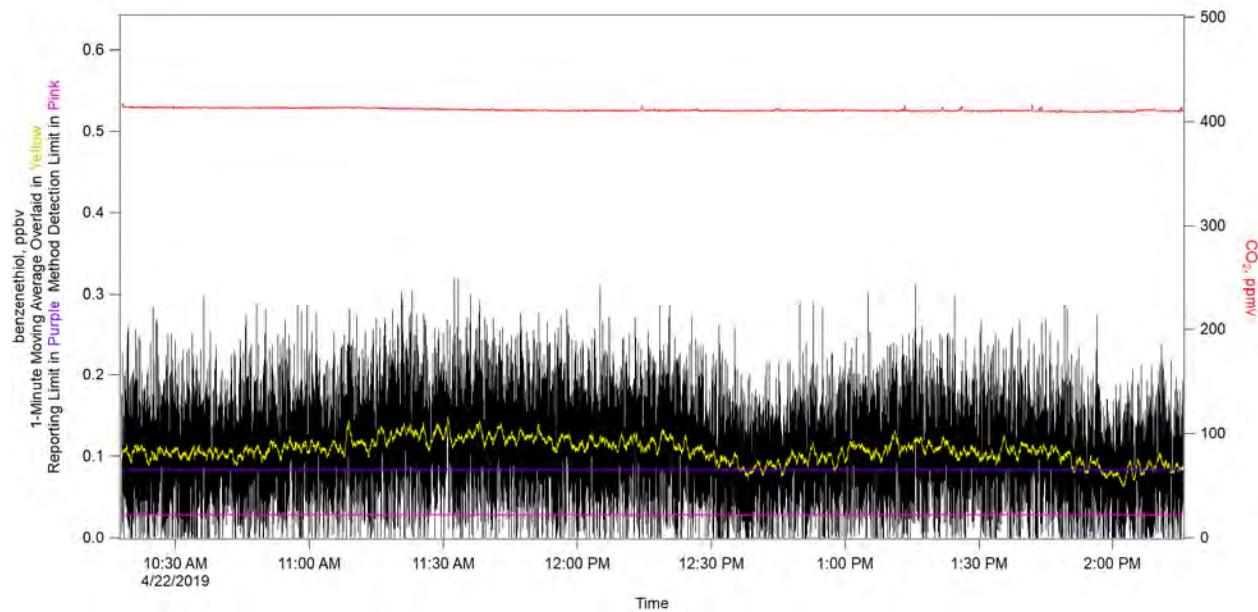


Figure 2-110. Benzenethiol.

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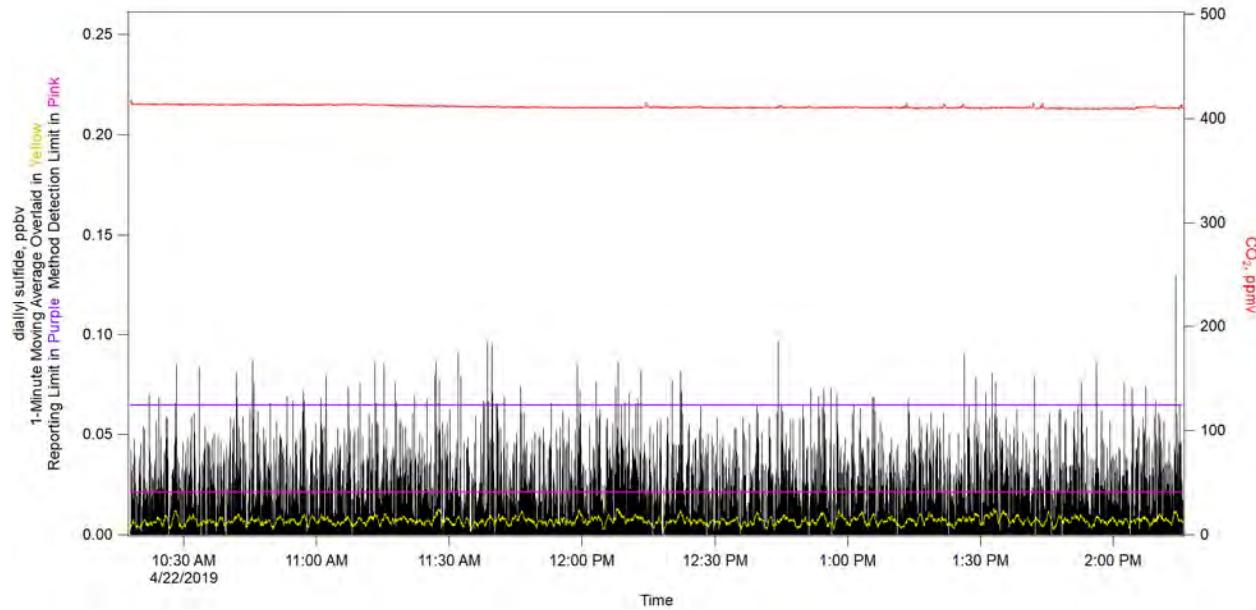


Figure 2-111. Diallyl Sulfide.

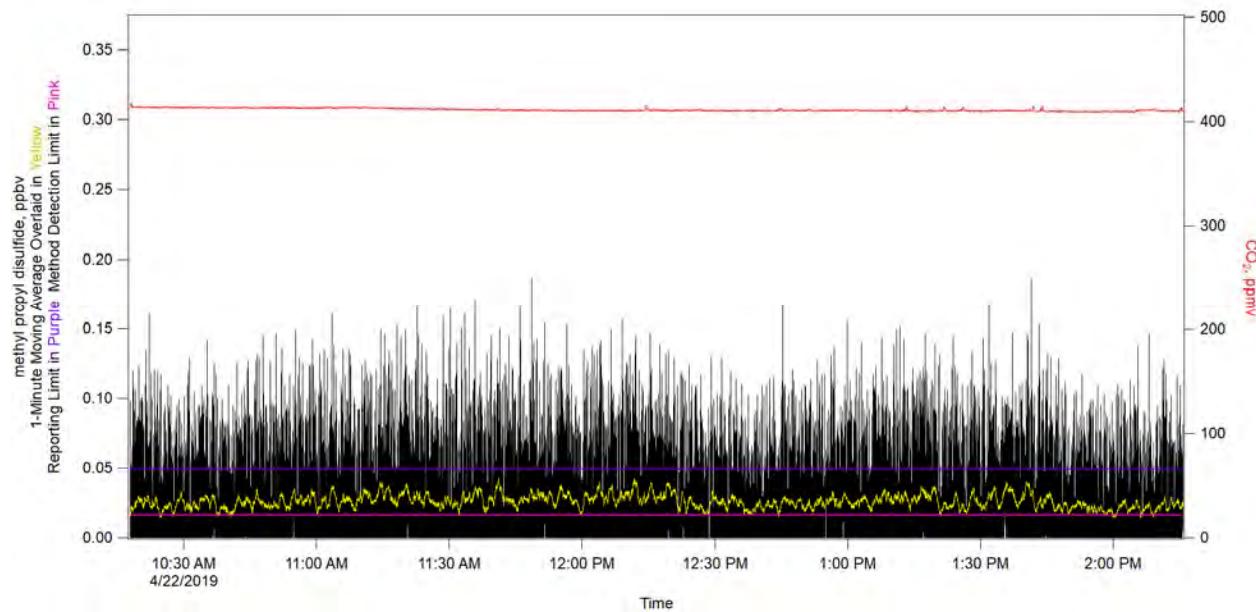


Figure 2-112. Methyl Propyl Disulfide.

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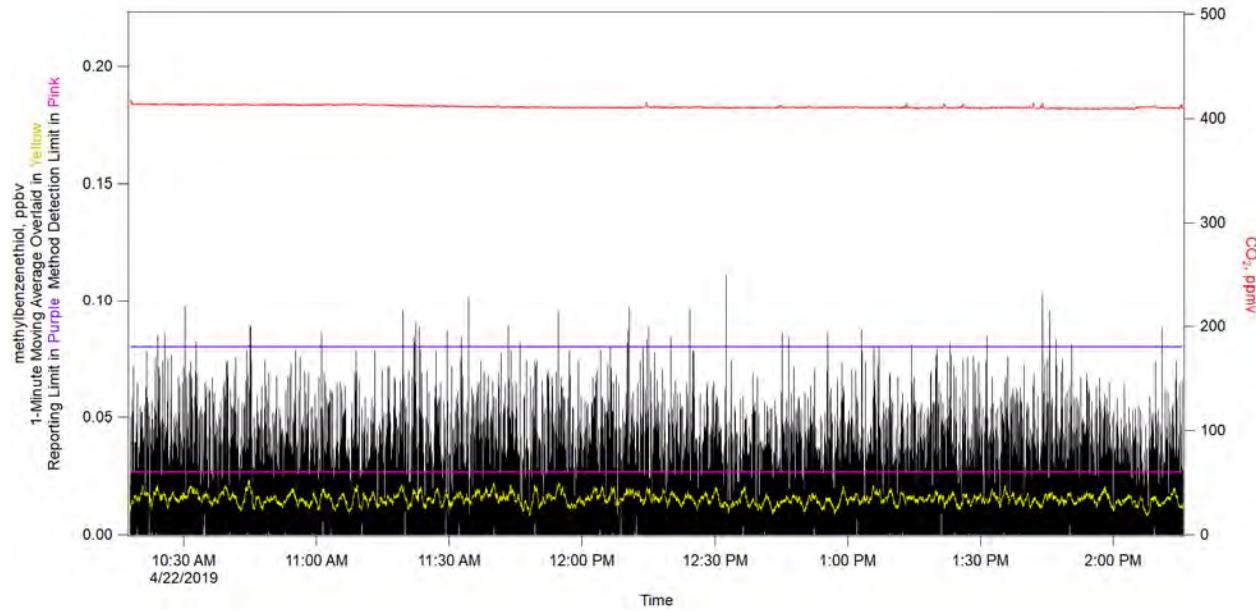


Figure 2-113. Methylbenzenethiol.

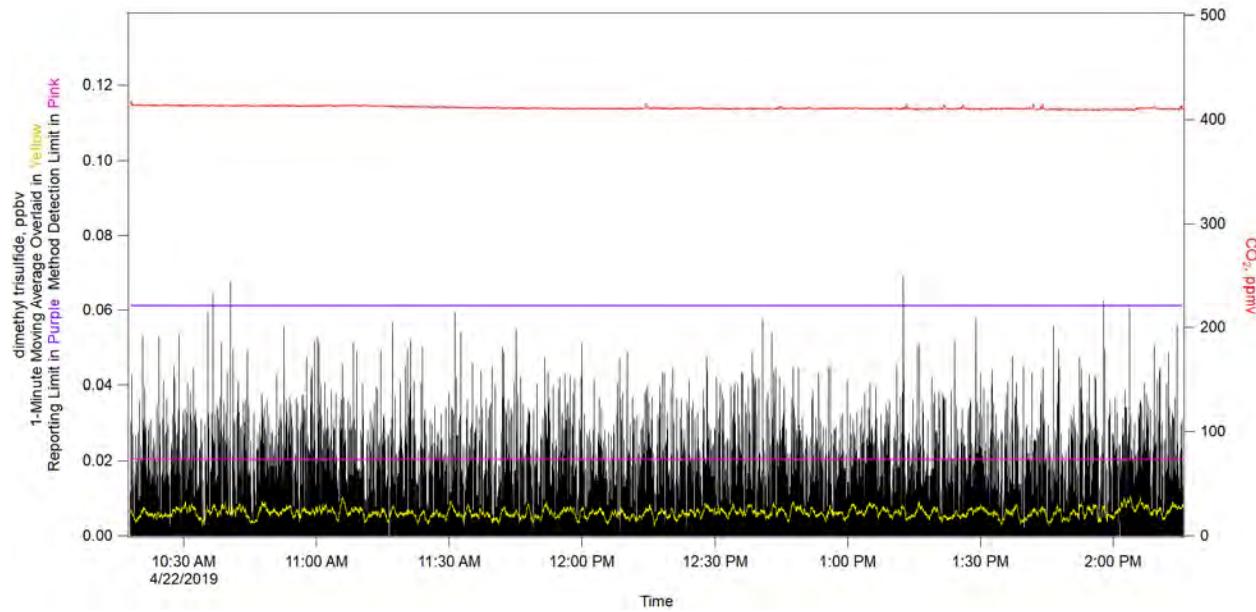


Figure 2-114. Dimethyl Trisulfide.

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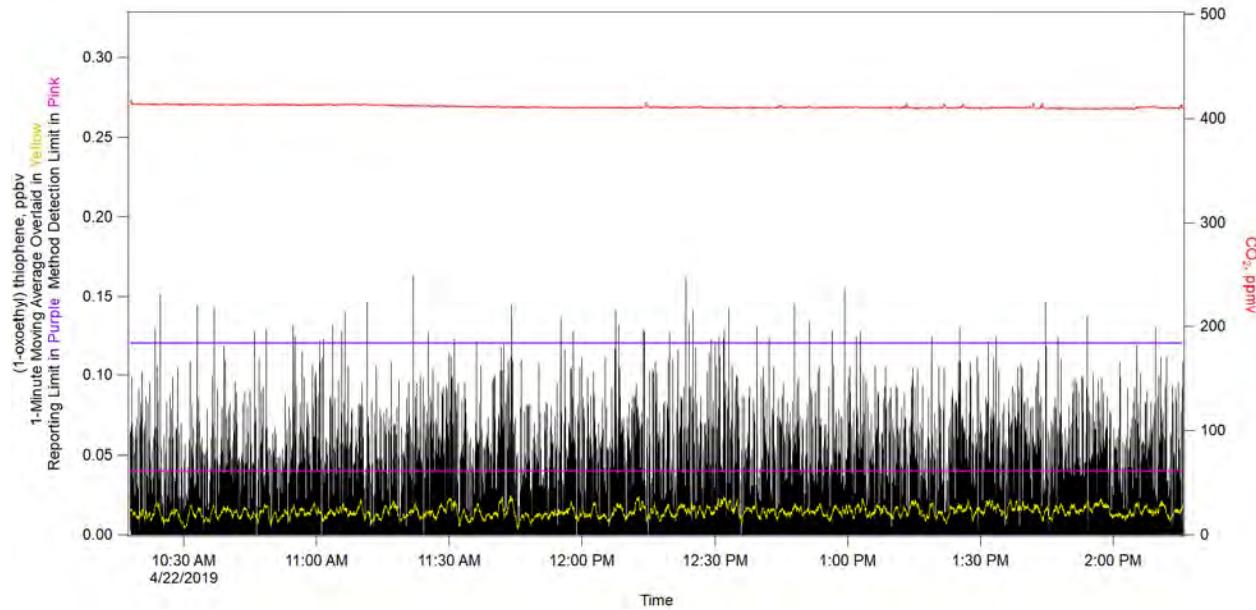


Figure 2-115. (1-oxoethyl) Thiophene.

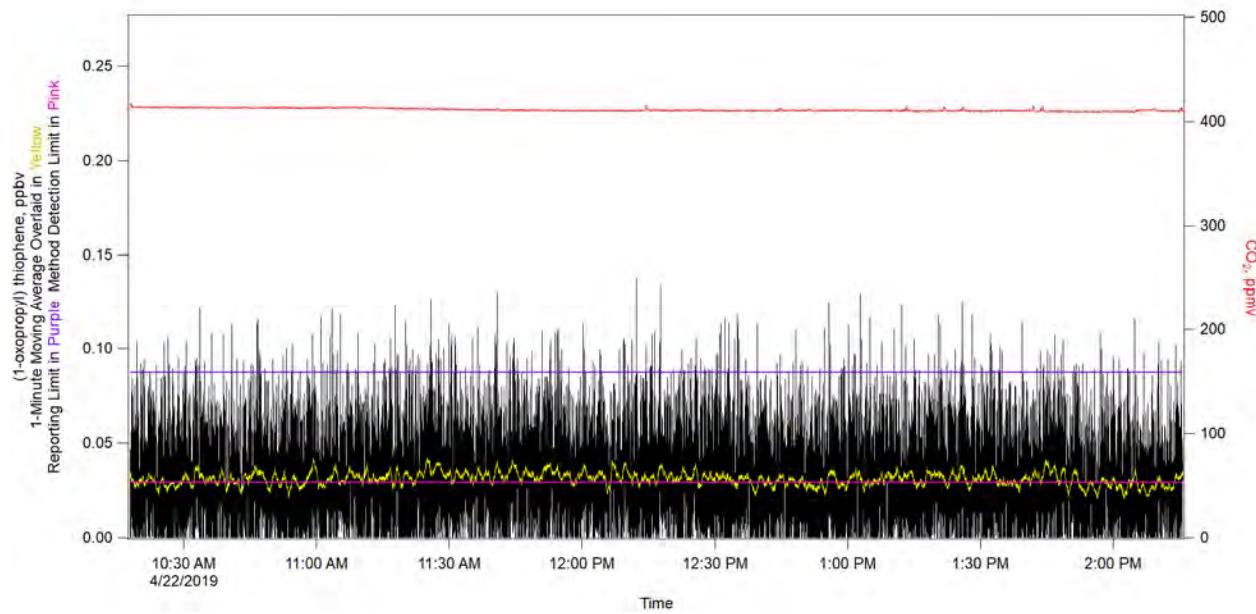


Figure 2-116. (1-oxopropyl) Thiophene.

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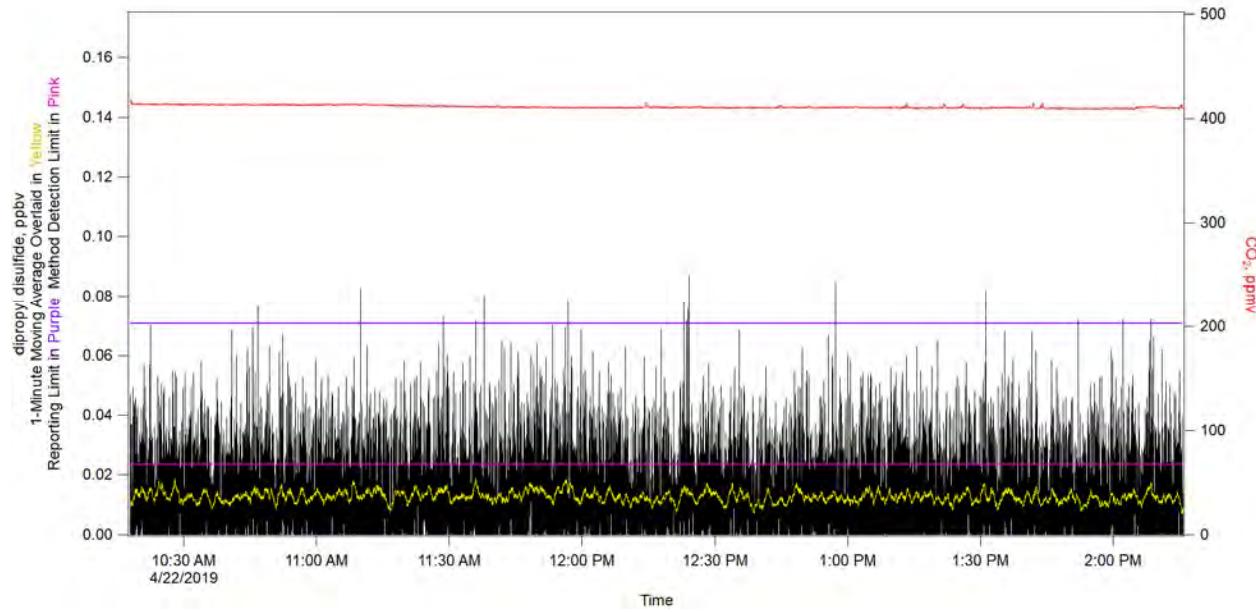


Figure 2-117. Dipropyl Disulfide.

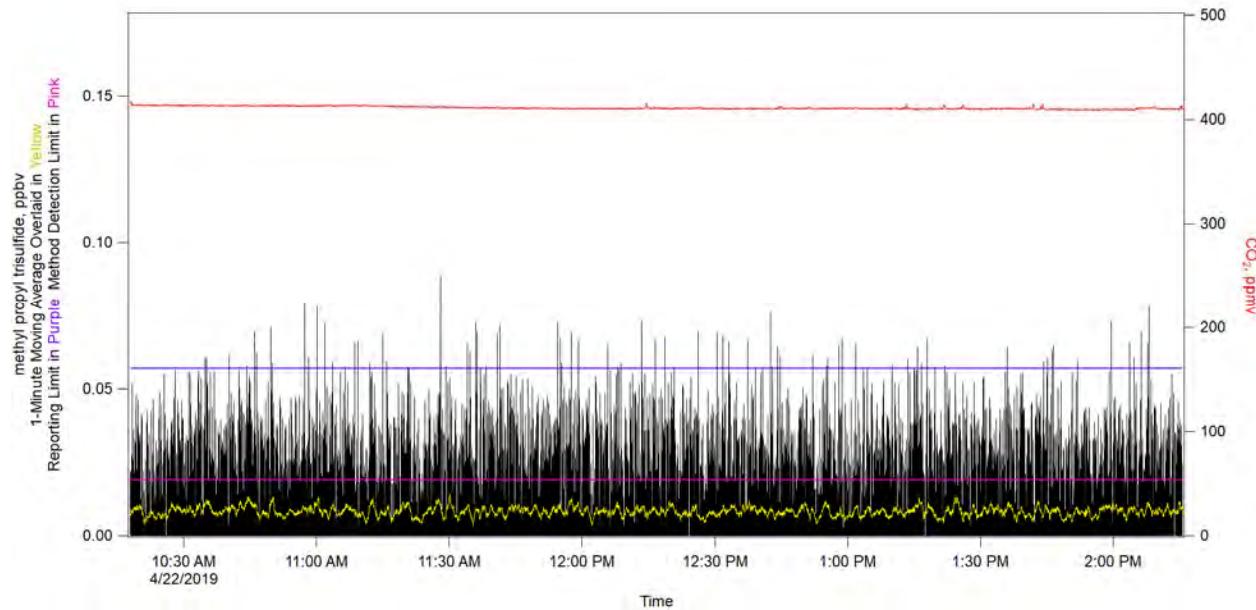


Figure 2-118. Methyl Propyl Trisulfide.

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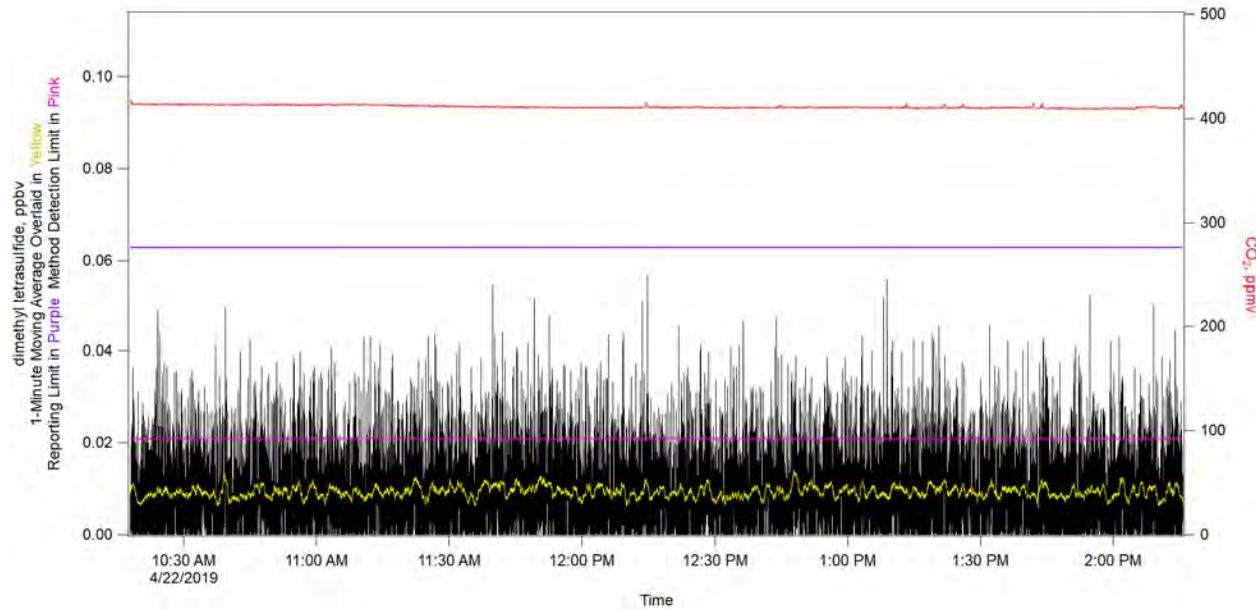


Figure 2-119. Dimethyl Tetrasulfide.

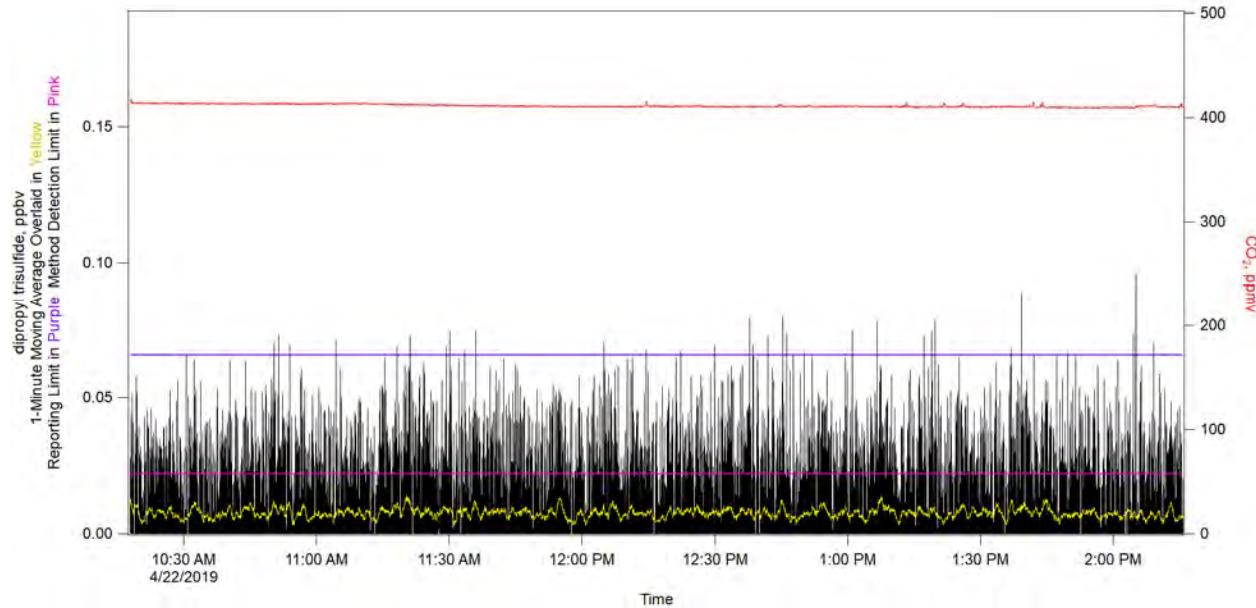


Figure 2-120. Dipropyl Trisulfide.

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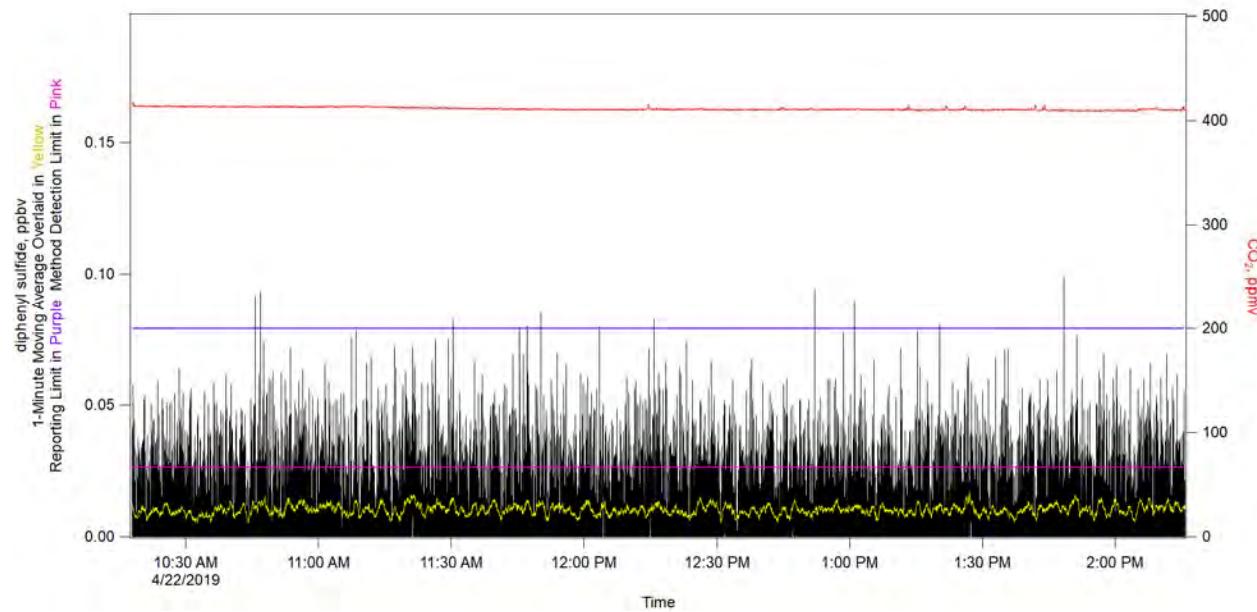


Figure 2-121. Diphenyl Sulfide.

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3.0 APRIL 23, 2019 – AREA MONITORING

3.1 Quality Assessment

Data from April 23, 2019, were assessed using Procedure 17124-DOE-HS-102. A Data Exchange Checklist was completed. The data were accepted by TerraGraphics with the following comments. Report No. 66409-RPT-004 was adequately documented and all checks passed the acceptance limits.

3.2 Summary

On April 23, 2019, the ML personnel arrived at the TerraGraphics warehouse at 04:50. The QA/QC zero-air/span checks were performed on the LI-COR CO₂ monitor, the Picarro NH₃ analyzer, and the PTR-MS beginning at 05:01. The ML arrived on the Hanford Site and checked in with the CSM at 06:20. The ML began a site survey loop of A Farms at 06:26. The ML relocated to the parking lot of AZ Farm, downwind of AZ Farm, at 06:58. Due to varying wind direction, the ML was moved to the northeast side of A Farm. The SME remotely logged in to the PTR-MS to tune the ion source at 08:03. The ML began a site survey loop at 08:40 until parking on the north side of 244-CR Waste Disposal Vault. The ML Operators performed another site survey loop at 10:28. The ML then proceeded to park east of AP Farm, downwind of a portable restroom, and remained there for approximately two hours. The ML Operators performed a site survey loop at 13:30. The Operators checked out with the CSM at 14:00, started a multipoint calibration using the Research and Development (R&D) gas standard, and the ML departed the site. The multipoint calibration was performed at request of the SME to confirm the transmission efficiency. The ML arrived back at the TerraGraphics warehouse at 14:50.

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Figure 3-1. Location of the Mobile Laboratory for the Duration of the Monitoring Period.

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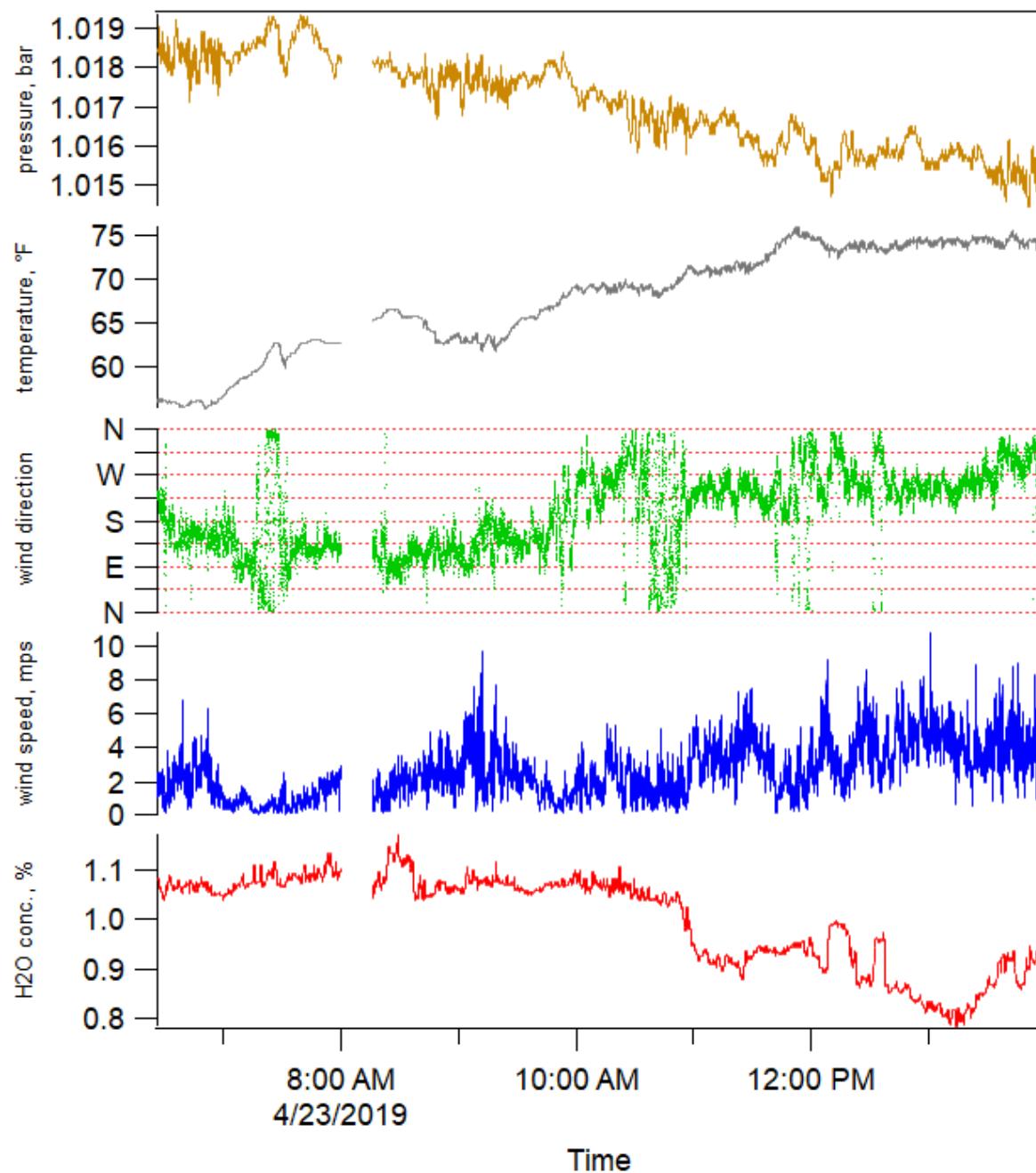


Figure 3-2. Weather Data for the Duration of the Monitoring Period.

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3.3 Samples Collected

Continuous air monitoring was performed using the following instrumentation:

- PTR-TOF 6000 X2,
- LI-COR CO₂ Monitor,
- Picarro Ammonia Monitor, and
- Airmar Weather Station.

Confirmatory air samples were not collected during this period.

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3.4 Area Monitoring

Table 3-1. Chemical of Potential Concern Statistical Information for the Monitoring Period of April 23, 2019. (2 Sheets)

COPC #	COPC Name	OEL (ppb)	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel St. Dev. (%)	Max. (ppb)	Median (ppb)
1	ammonia	25000	6.225	11.483†	3.526	30.710	22.783	10.694†
2	formaldehyde	300	1.721	<1.721	0.177	28.311	13.358	<1.721
3	methanol	200000	5.758	6.482†	1.008	15.553	28.017	6.286†
4	acetonitrile	20000	0.085	0.141†	0.036	25.717	0.469	0.141†
5	acetaldehyde	25000	1.027	1.367†	0.428	31.303	17.290	1.289†
6	ethylamine	5000	0.069	<0.069	0.020	70.042	0.148	<0.069
7	1,3-butadiene	1000	0.183	0.242†	0.098	40.591	2.969	0.235†
8	propanenitrile	6000	0.107	<0.107	0.040	72.910	3.053	<0.107
9	2-propenal	100	0.340	<0.340	0.052	58.819	1.113	<0.340
10	1-butanol + butenes	20000	0.214	<0.214	0.048	52.175	1.677	<0.214
11	methyl isocyanate	20	0.069	<0.069	0.028	61.255	1.126	<0.069
12	methyl nitrite	100	0.098	<0.098	0.040	42.378	0.552	<0.098
13	furan	1	0.062	<0.062	0.018	59.305	0.161	<0.062
14	butanenitrile	8000	0.039	<0.039	0.018	73.796	0.550	<0.039
15	but-3-en-2-one + 2,3-dihydrofuran + 2,5-dihydrofuran	200, 1, 1	0.041	<0.041	0.022	57.346	N/A*	N/A*
16	butanal	25000	0.061	0.107†	0.036	33.830	1.273	0.106†
17	NDMA**	0.3	0.082	<0.082	0.022	169.944	0.166	<0.082
18	benzene	500	0.236	<0.236	0.045	37.209	1.064	<0.236
19	2,4-pentadienenitrile + pyridine	300, 1000	0.085	<0.085	0.019	46.390	0.129	<0.085
20	2-methylene butanenitrile	300	0.036	<0.036	0.012	75.842	0.078	<0.036
21	2-methylfuran	1	0.043	<0.043	0.023	65.586	0.235	<0.043
22	pentanenitrile	6000	0.036	<0.036	0.020	125.650	1.596	<0.036
23	3-methyl-3-buten-2-one + 2-methyl-2-butenal	20, 30	0.043	<0.043	0.022	67.202	0.196	<0.043
24	NEMA**	0.3	0.058	<0.058	0.018	143.833	0.135	<0.058
25	2,5-dimethylfuran	1	0.032	<0.032	0.018	78.526	0.143	<0.032
26	hexanenitrile	6000	0.031	<0.031	0.011	93.948	0.078	<0.031
27	2-hexanone (MBK)	5000	0.036	<0.036	0.014	95.360	0.132	<0.036
28	NDEA**	0.1	0.034	<0.034	0.012	147.938	0.088	<0.034

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Table 3-1. Chemical of Potential Concern Statistical Information for the Monitoring Period of April 23, 2019. (2 Sheets)

COPC #	COPC Name	OEL (ppb)	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel St. Dev. (%)	Max. (ppb)	Median (ppb)
29	butyl nitrite + 2-nitro-2-methylpropane	100, 300	0.058	<0.058	0.015	51.200	0.101	<0.058
30	2,4-dimethylpyridine	500	0.036	<0.036	0.011	104.254	0.106	<0.036
31	2-propylfuran + 2-ethyl-5-methylfuran	1	0.027	<0.027	0.013	99.993	0.124	<0.027
32	heptanenitrile	6000	0.027	<0.027	0.009	96.807	0.069	<0.027
33	4-methyl-2-hexanone	500	0.033	<0.033	0.012	94.084	0.087	<0.033
34	NMOR**	0.6	0.021	<0.021	0.009	177.343	0.090	<0.021
35	butyl nitrate	2500	0.022	<0.022	0.008	124.356	0.056	<0.022
36	2-ethyl-2-hexenal + 4-(1-methylpropyl)-2,3-dihydrofuran + 3-(1,1-dimethylethyl)-2,3-dihydrofuran	100, 1, 1	0.028	<0.028	0.012	90.495	0.087	<0.028
37	6-methyl-2-heptanone	8000	0.028	<0.028	0.010	89.781	0.086	<0.028
38	2-pentylfuran	1	0.026	<0.026	0.012	84.944	0.086	<0.026
39	biphenyl	200	0.022	<0.022	0.010	125.006	0.072	<0.022
40	2-heptylfuran	1	0.067	<0.067	0.016	42.649	0.113	<0.067
41	1,4-butanediol dinitrate	50	0.036	<0.036	0.011	86.983	0.071	<0.036
42	2-octylfuran	1	0.020	<0.02	0.009	203.821	0.075	<0.02
43	1,2,3-propanetriol 1,3-dinitrate	50	0.011	<0.011	0.004	378.421	0.071	<0.011
44	PCB	1000	0.034	<0.034	0.010	71.473	0.075	<0.034
45	6-(2-furanyl)-6-methyl-2-heptanone	1	0.025	<0.025	0.009	100.951	0.075	<0.025
46	furfural acetophenone	1	0.064	<0.064	0.015	42.996	0.112	<0.064
N/A*	The maximum peak value for but-3-en-2-one + 2,3 dihydrofuran + 2,5 dihydrofuran was 0.365 ppb and the median value was <0.041 ppb. The PTR-MS results for but-3-en-2-one + 2,3 dihydrofuran + 2,5 dihydrofuran are not compared to OEL concentrations because: 1) the result is suspect due to a known biogenic interferant (methacrolein) that is expected to be in concentrations that occasionally exceed the dihydrofuran OEL, and 2) this combination of COPCs have OEL concentrations that differ by a factor of 200, which provide widely variant bases for these numbers.							
**	Nitrosamine results are suspect due to isobaric interferants causing positive bias that have been encountered during previous background [53005-81-RPT-007, PTR-MS Mobile Laboratory Vapor Monitoring Background Study, (3/18/2018 – 4/20/2018), and Fiscal Year 2017 Mobile Laboratory Vapor Monitoring at the Hanford Site: Monitoring During Waste Disturbing Activities and Background Study, RJ Lee Group, Inc.].							
<	COPC Averages below the MDL.							
†	COPC Averages between the RL and the MDL.							
	COPC Averages >100% of the OEL.							
	COPC Averages 50-100% of the OEL.							
	COPC Averages 10-50% of the OEL.							

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Figure 3-3 through Figure 3-51 display 46 COPC signals, overlaid with the same signal smoothed using a 1-minute moving average (in cases where a moving average assists with data visualization), and CO₂, for the monitoring period April 23, 2019. If within range of the plot's left axis, a green horizontal line representing 50% of the COPC's OEL, a blue horizontal line representing the COPC's OEL, a horizontal purple line representing the RL, and a pink horizontal line representing the MDL are shown. The tuning of the PTR-MS ion source from 08:03 to 08:16 performed by Dr. Matthew Erickson, the subject matter expert (SME), was removed from the following plots.

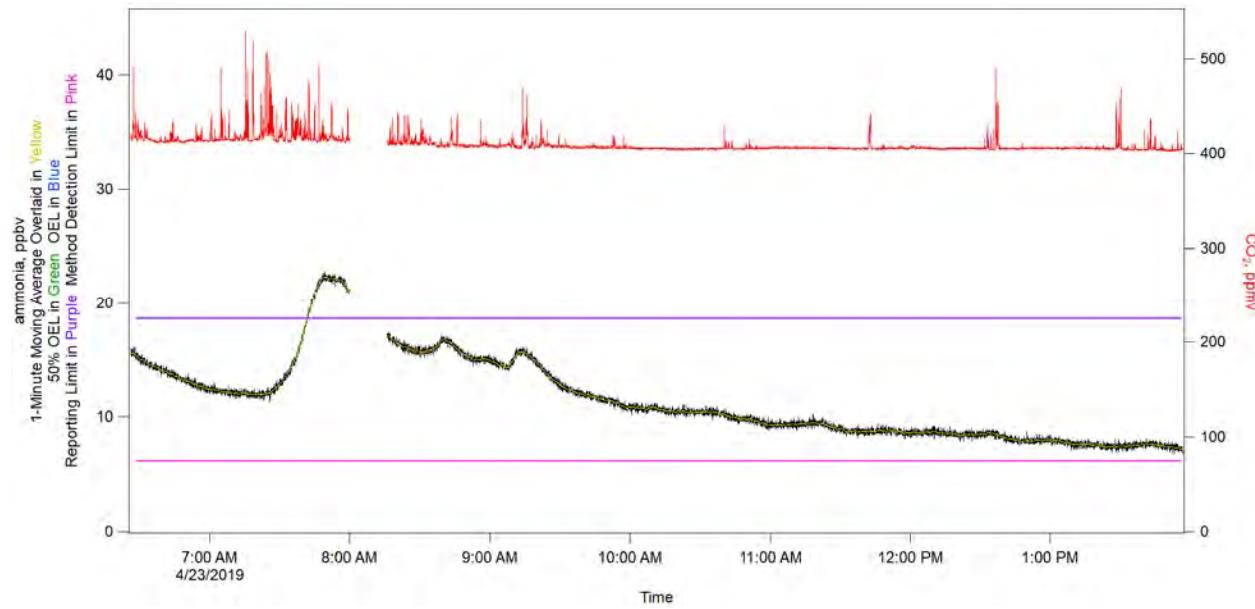


Figure 3-3. Ammonia



Figure 3-4. Formaldehyde.

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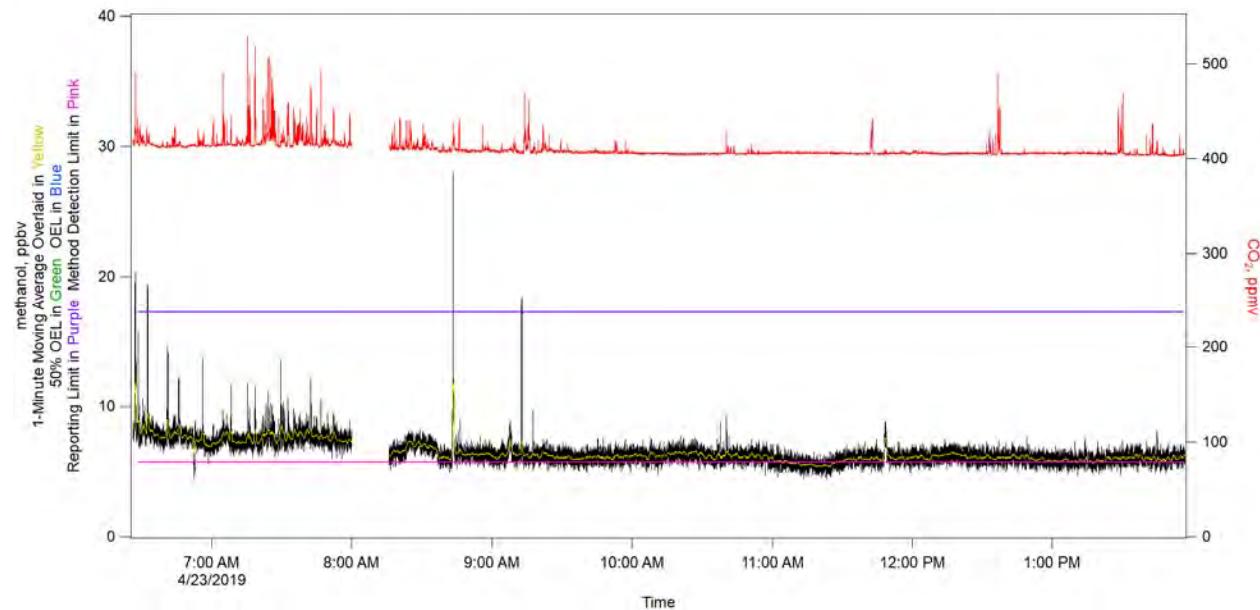


Figure 3-5. Methanol.

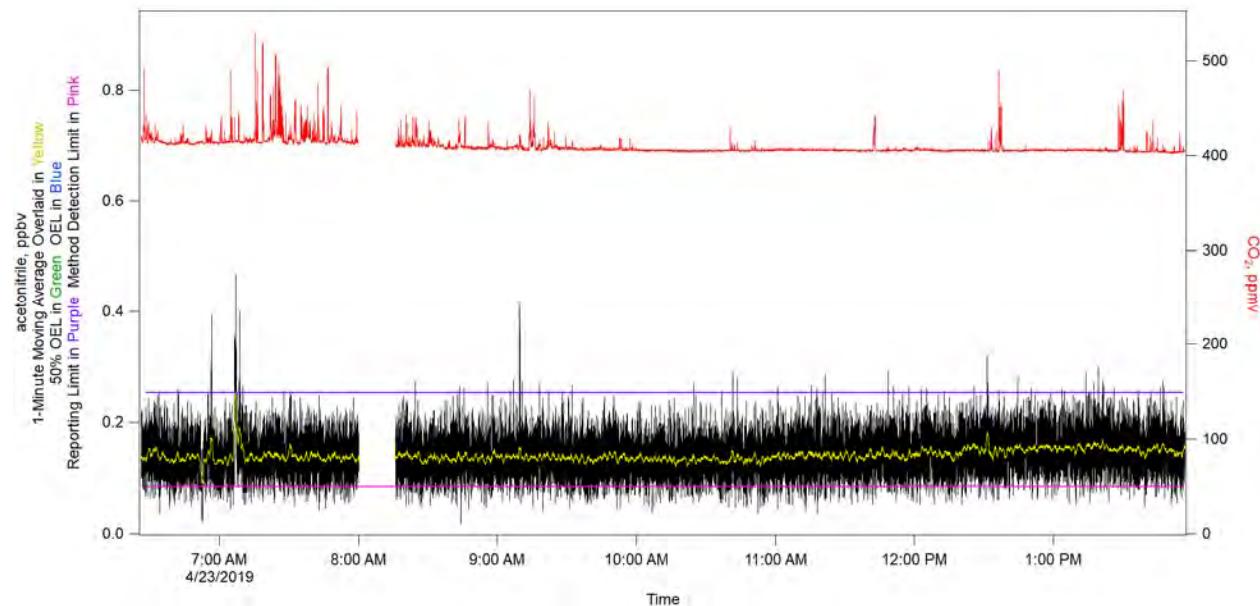


Figure 3-6. Acetonitrile.

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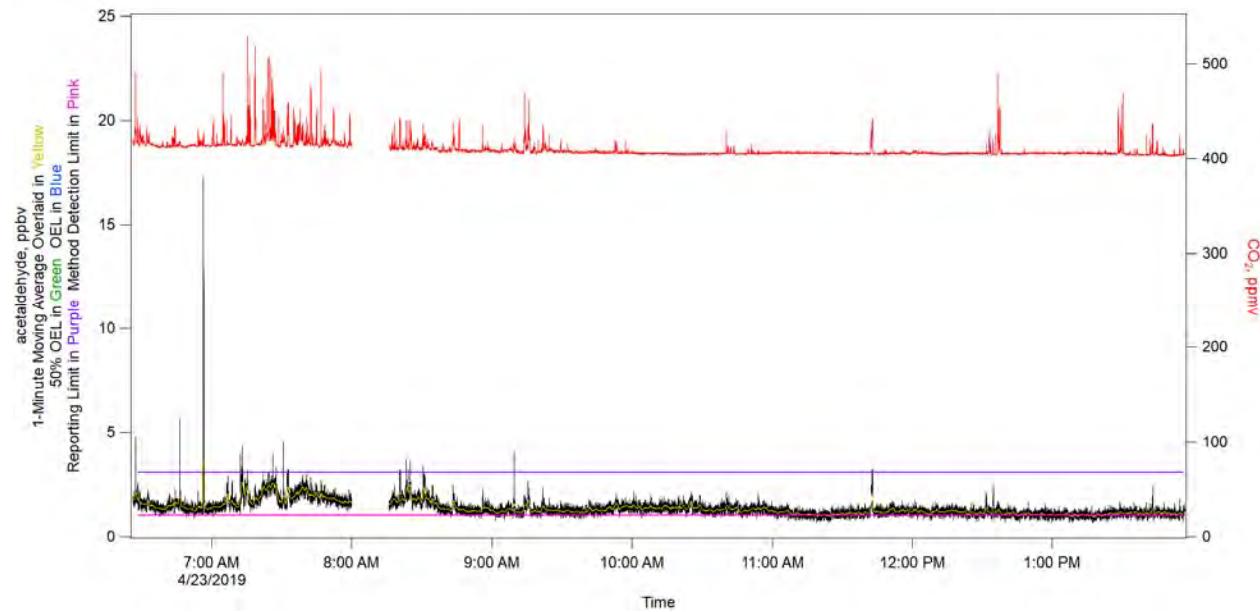


Figure 3-7. Acetaldehyde.

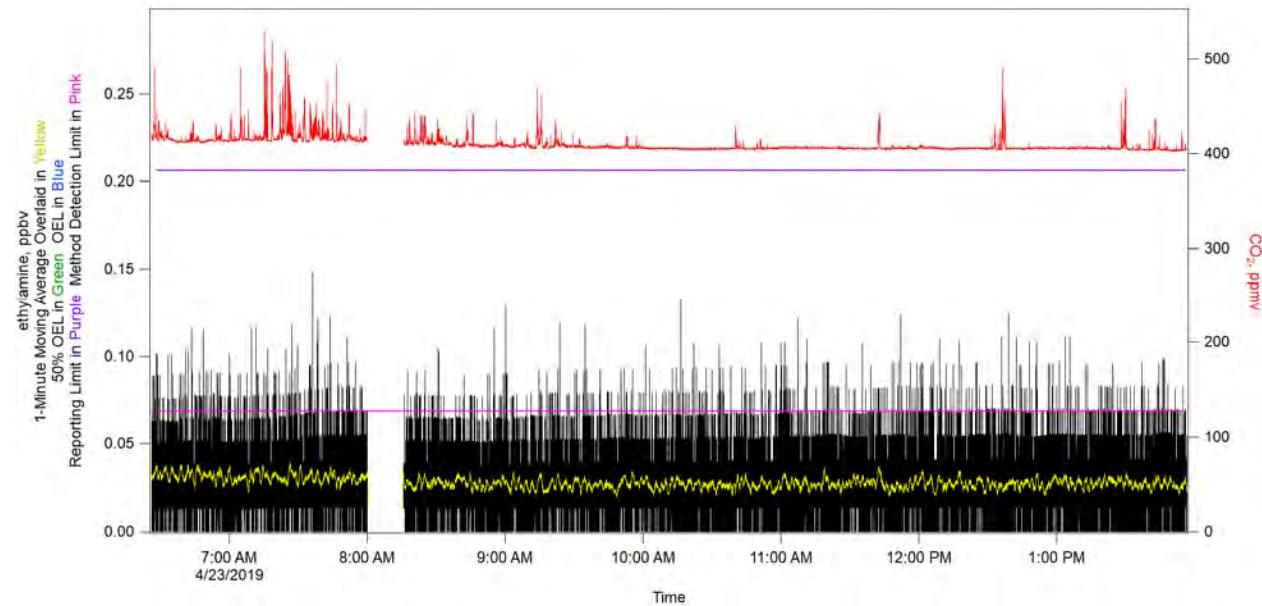


Figure 3-8. Ethylamine.

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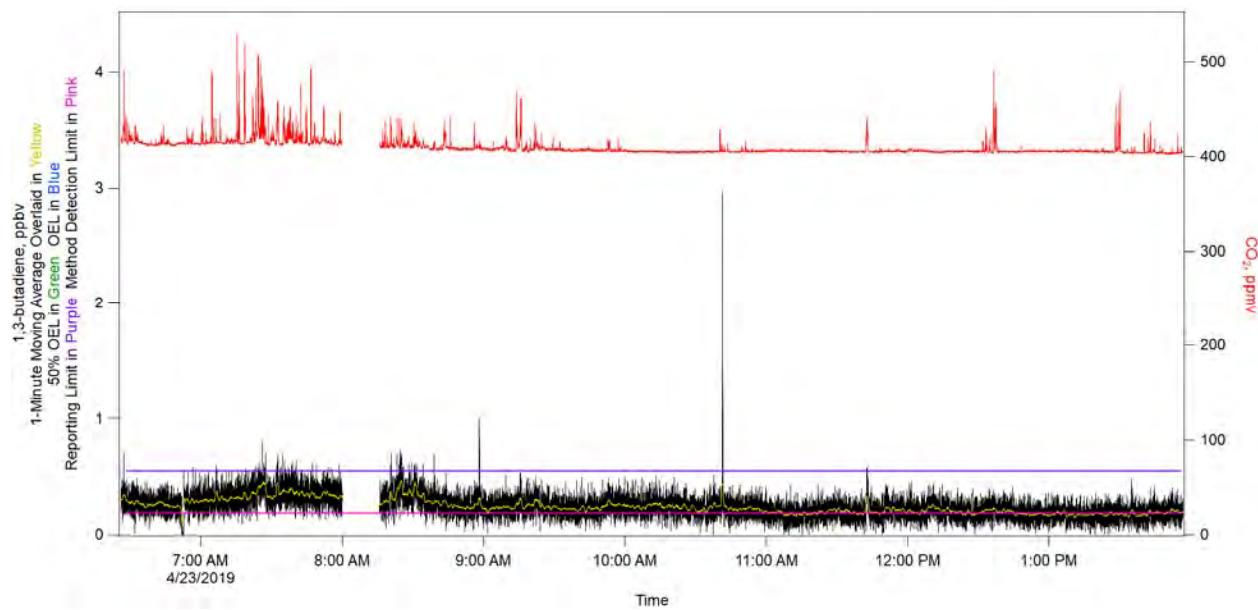


Figure 3-9. 1,3-butadiene.

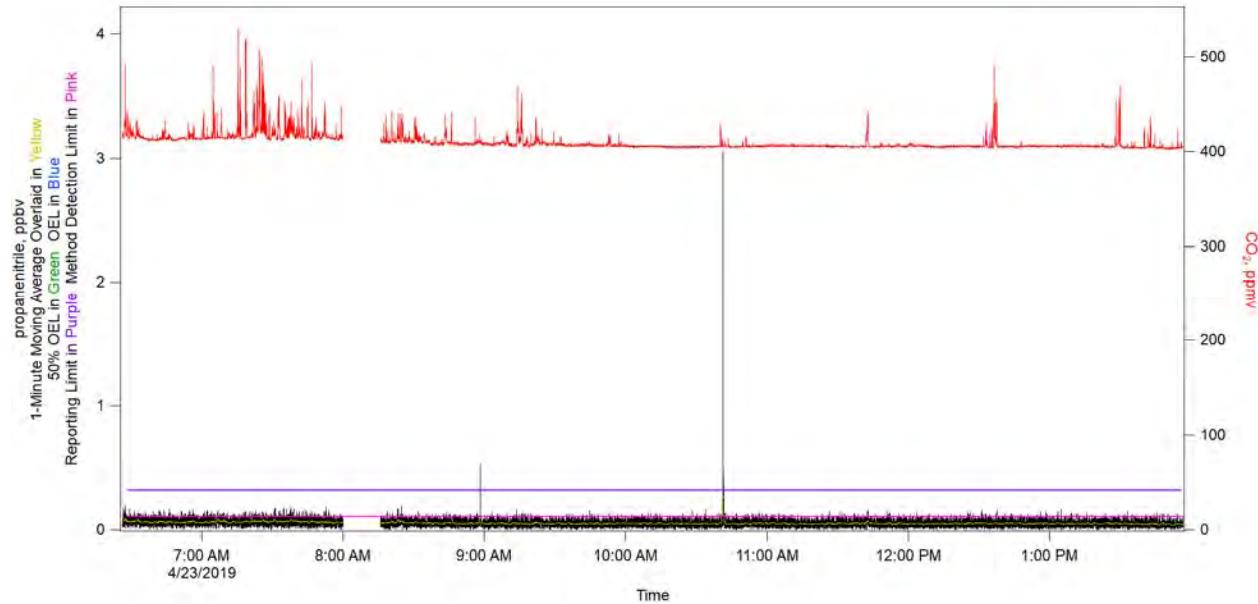


Figure 3-10. Propanenitrile.

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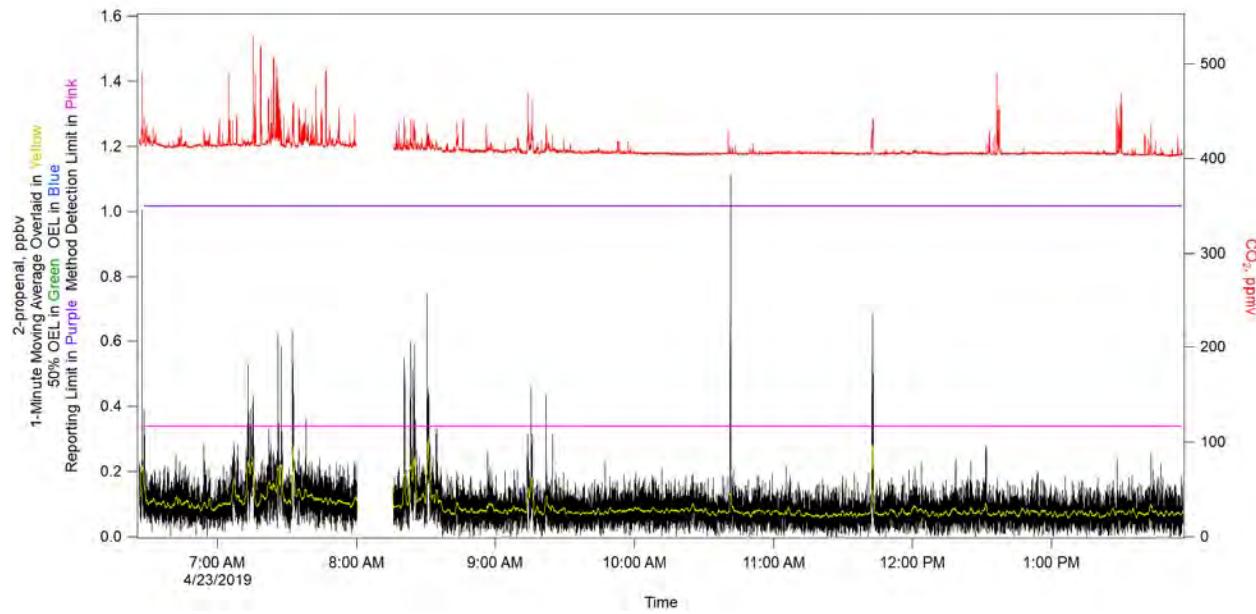


Figure 3-11. 2-propenal.

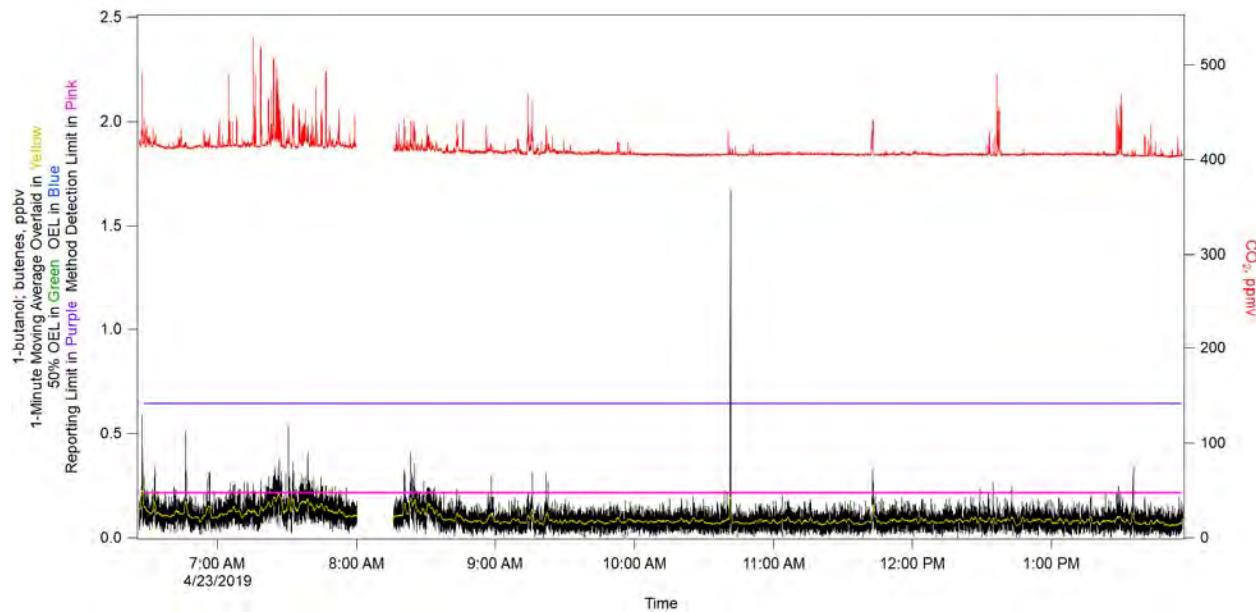


Figure 3-12. 1-butanol; Butenes.

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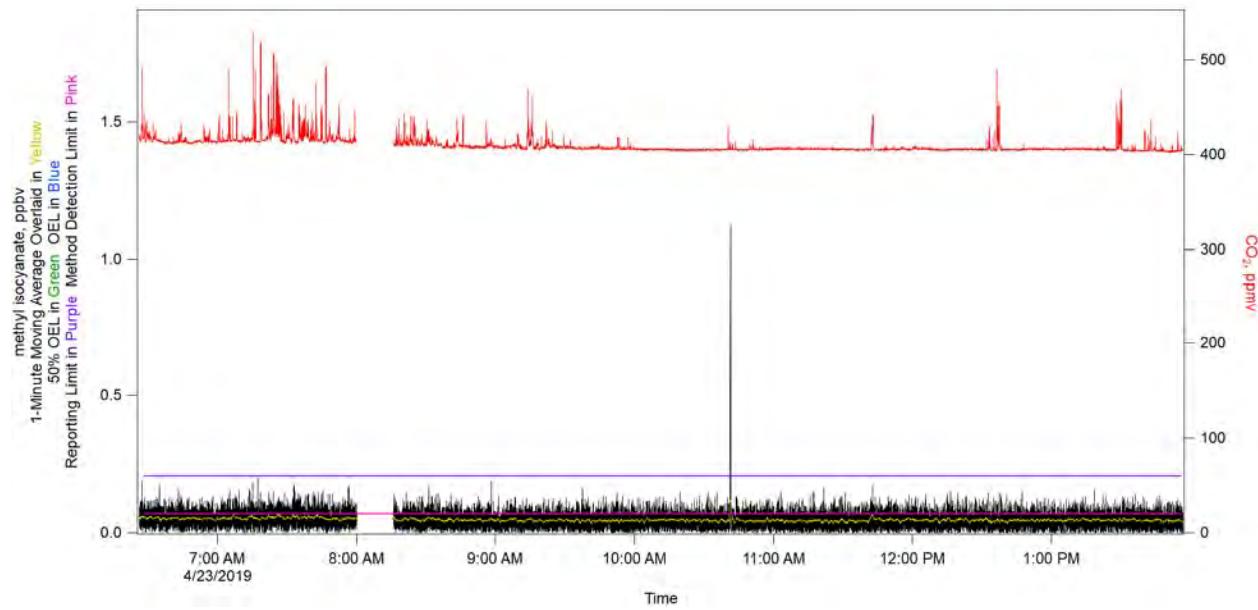


Figure 3-13. Methyl Isocyanate.

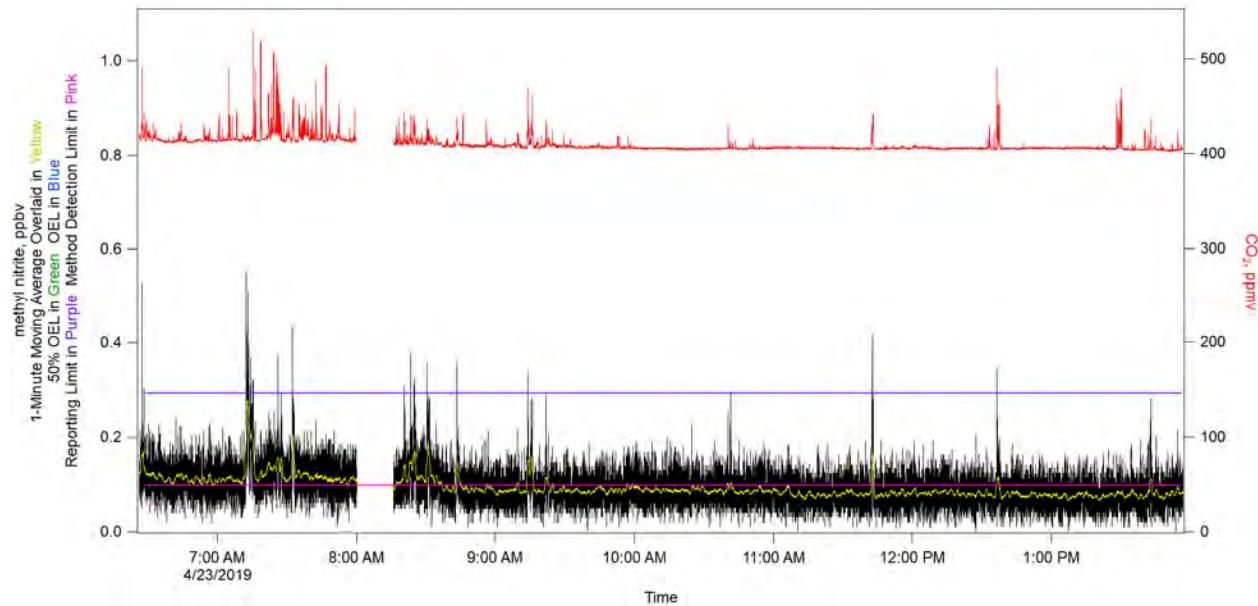


Figure 3-14. Methyl Nitrite.

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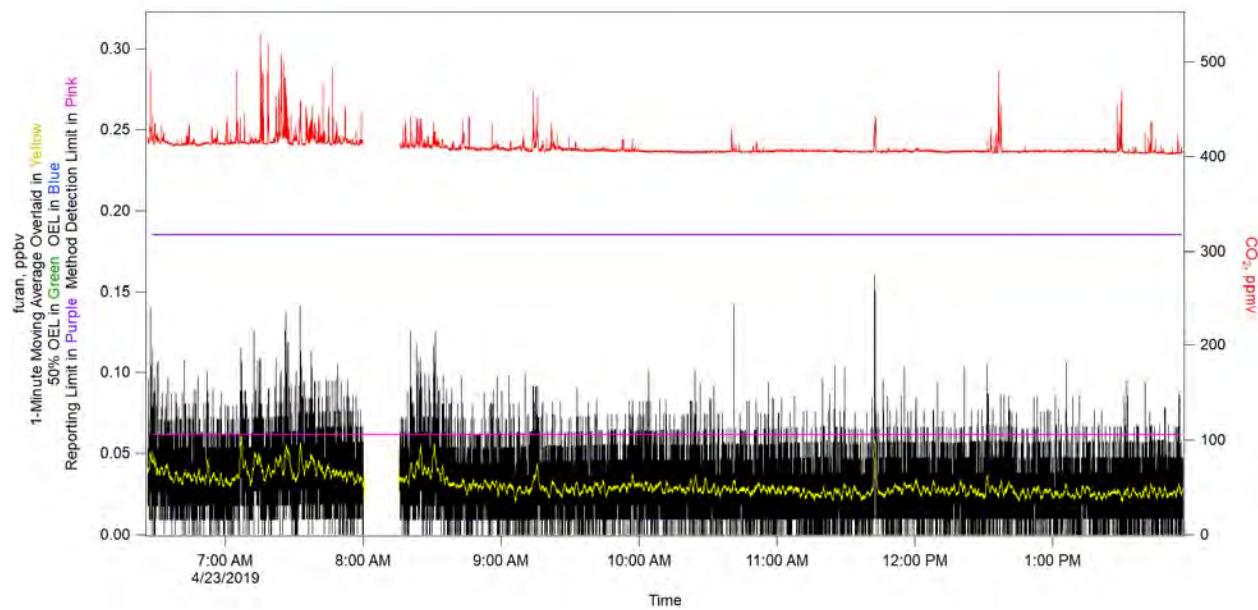


Figure 3-15. Furan.

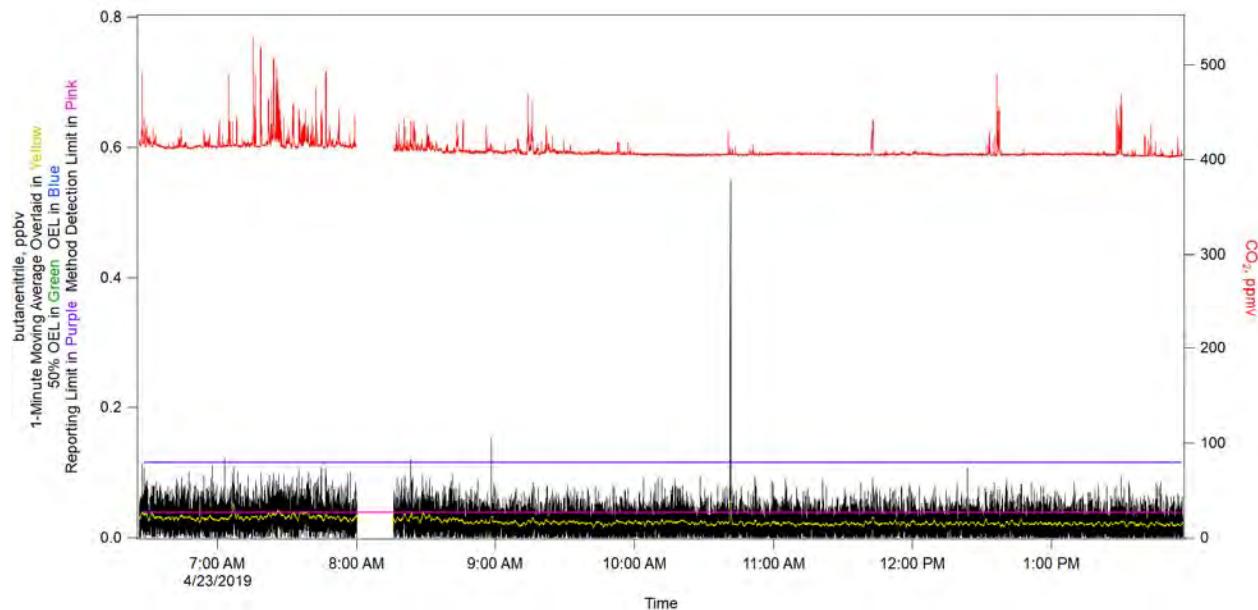


Figure 3-16. Butanenitrile.

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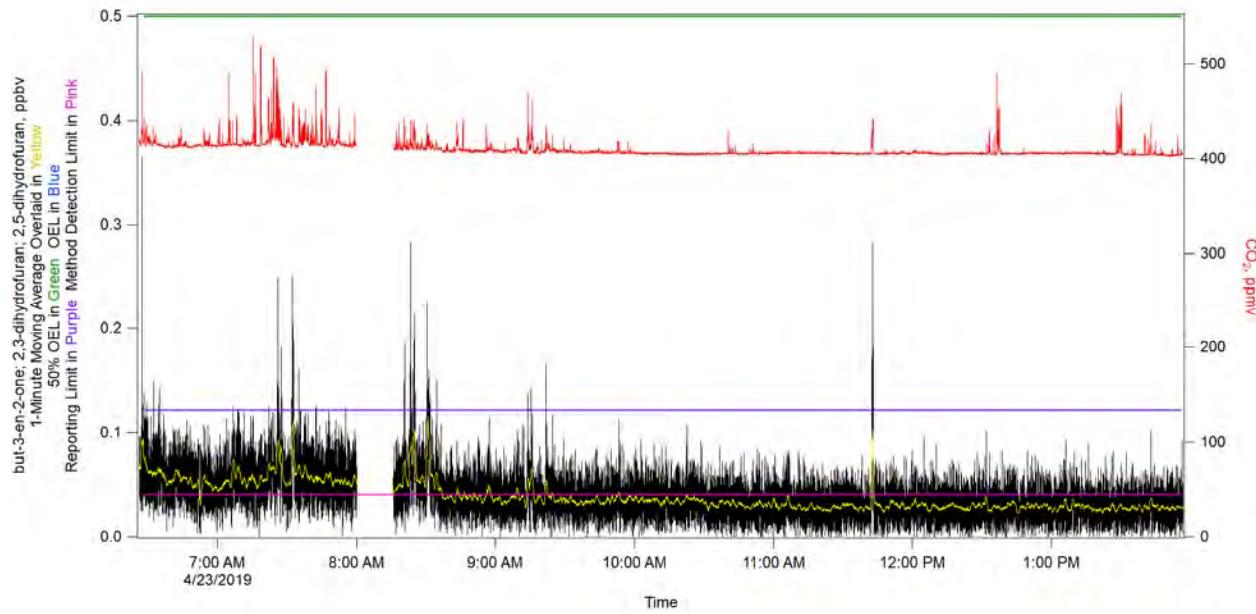


Figure 3-17. But-3-en-2-one; 2,3-dihydrofuran; 2,5-dihydrofuran.

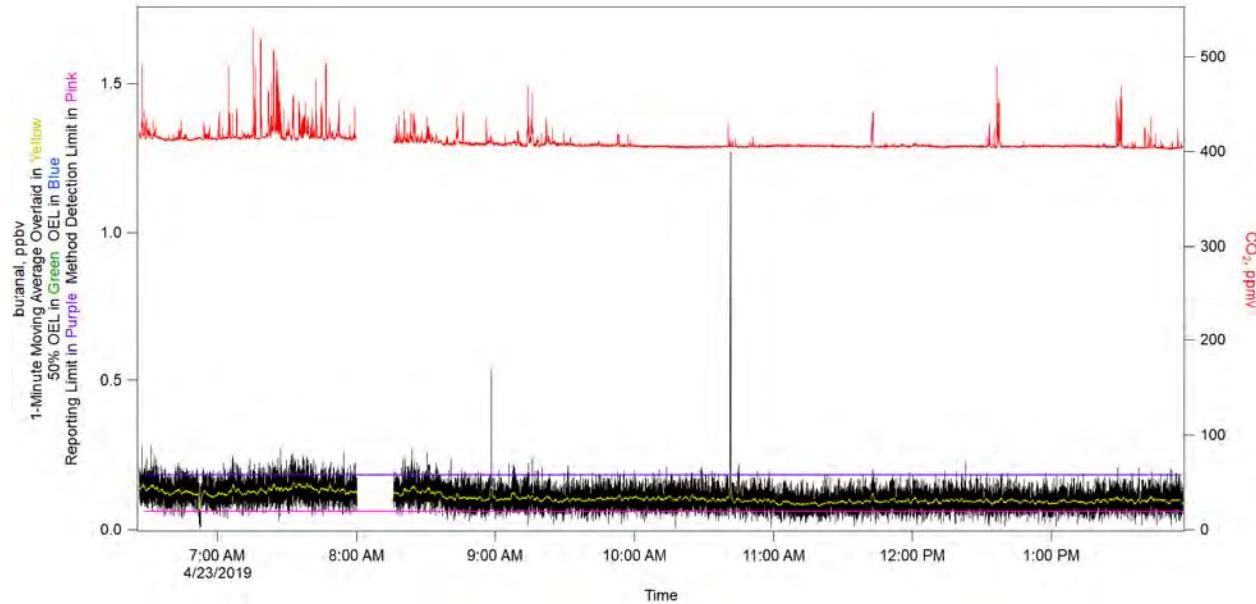


Figure 3-18. Butanal.

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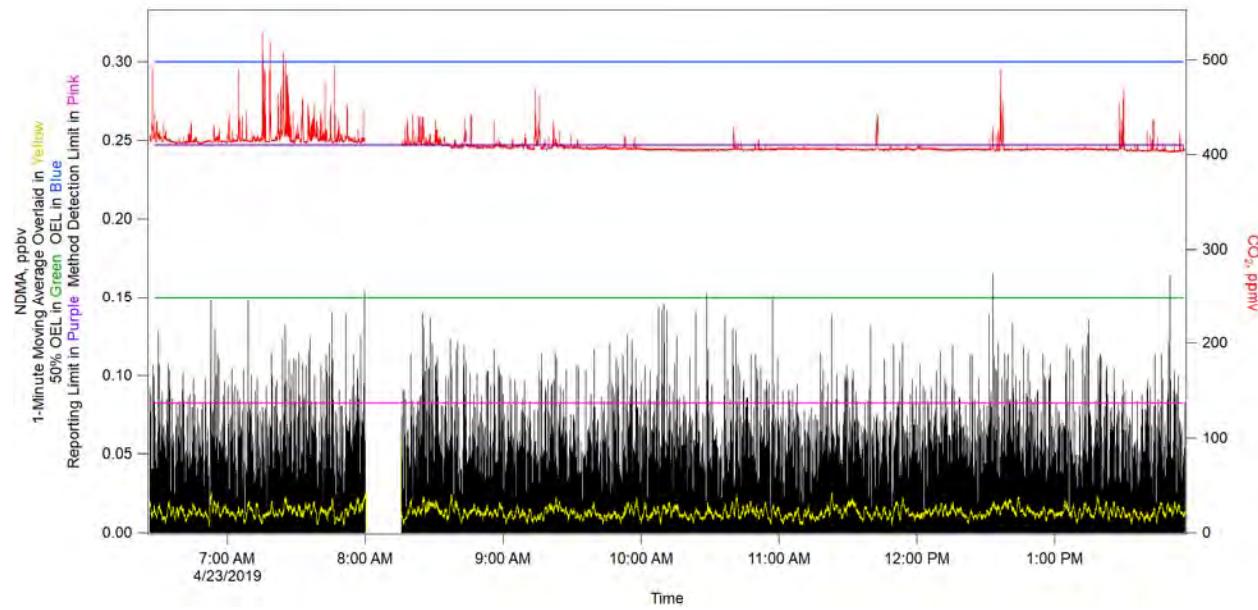


Figure 3-19. N-nitrosodimethylamine (NDMA).

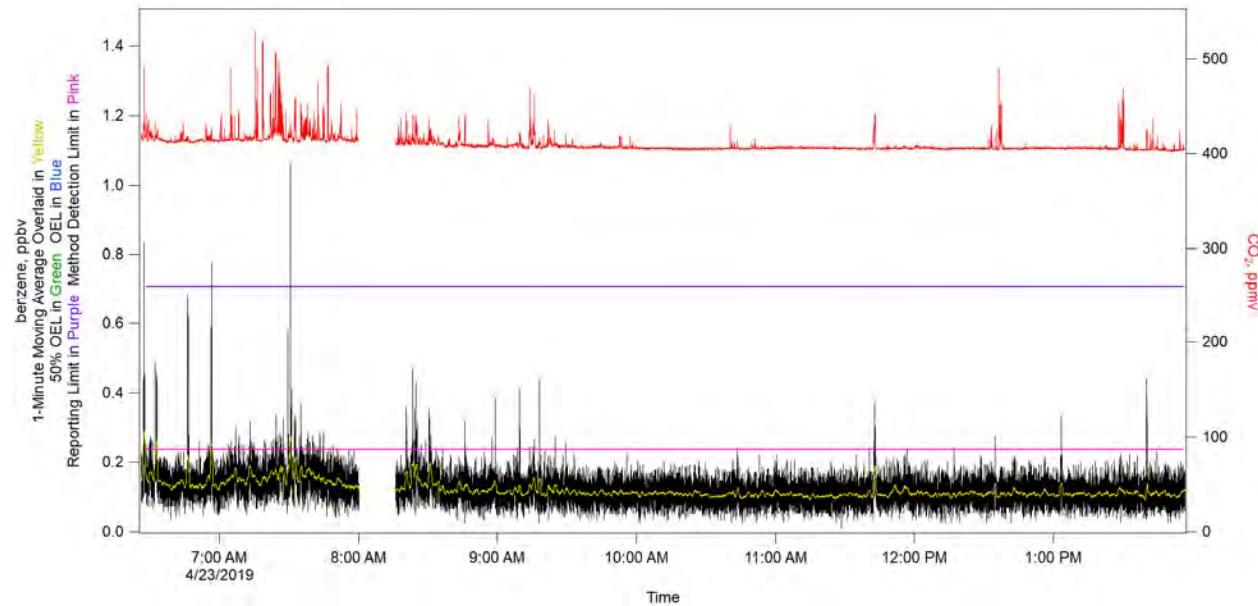


Figure 3-20. Benzene.

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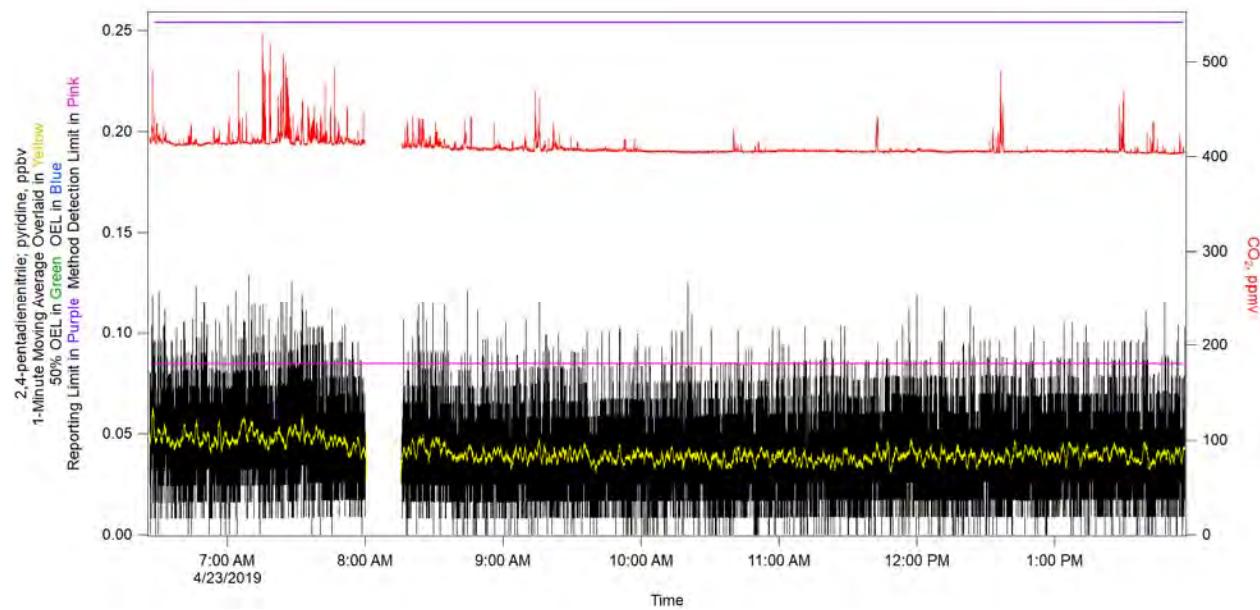


Figure 3-21. 2,4-pentadienenitrile; Pyridine.

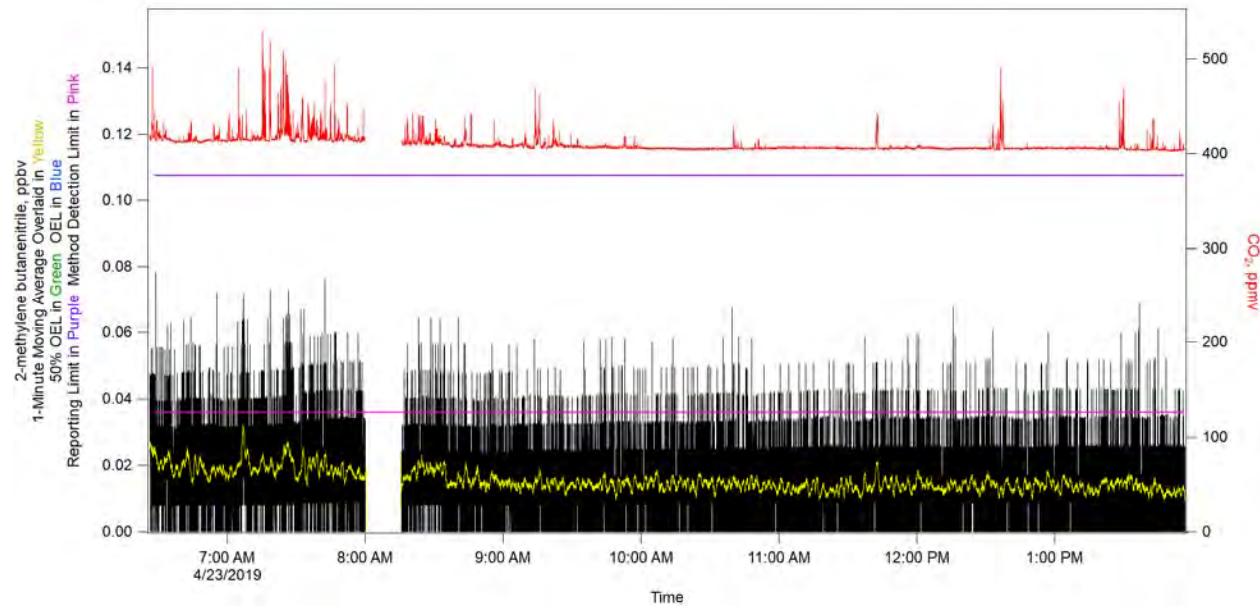


Figure 3-22. 2-methylene Butanenitrile.

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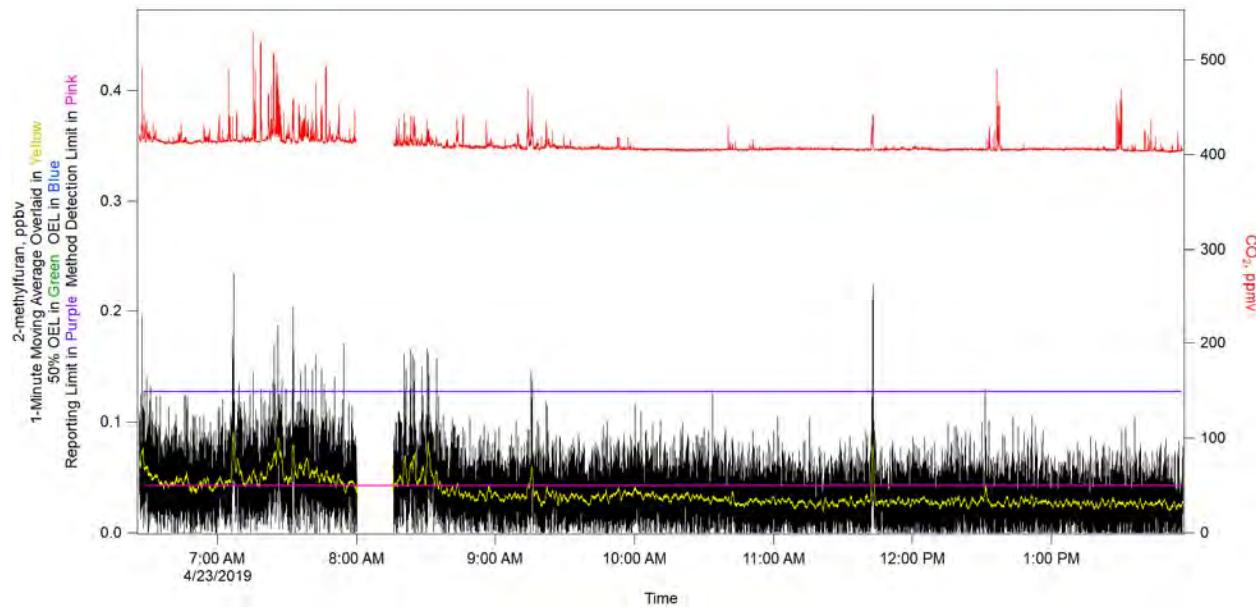


Figure 3-23. 2-methylfuran.

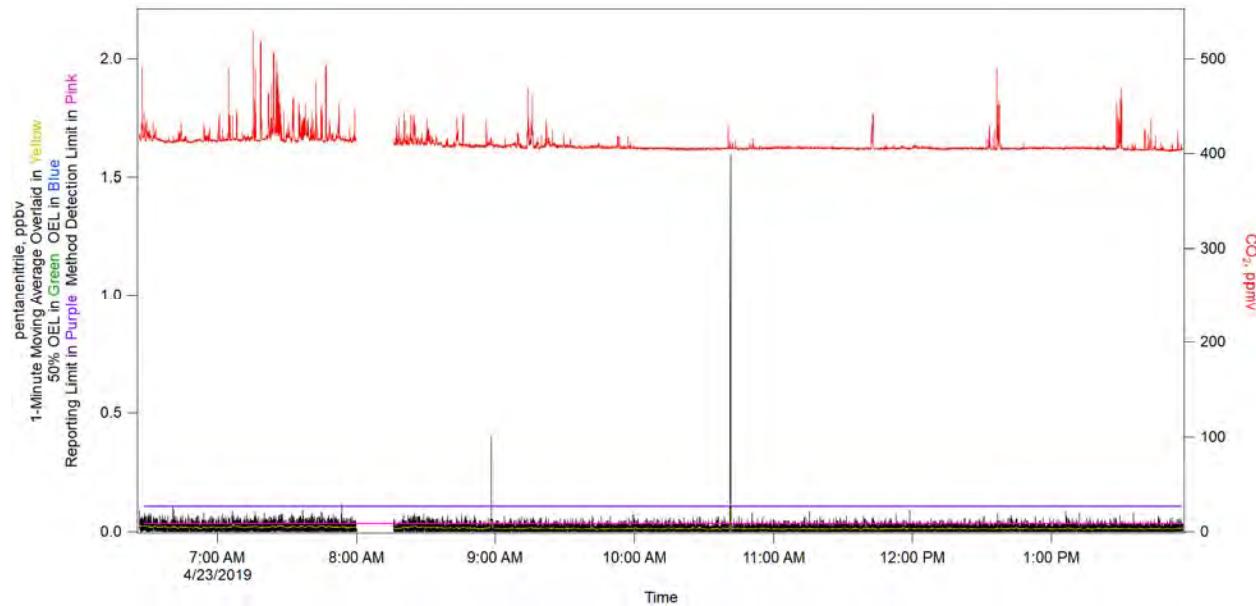


Figure 3-24. Pentanenitrile.

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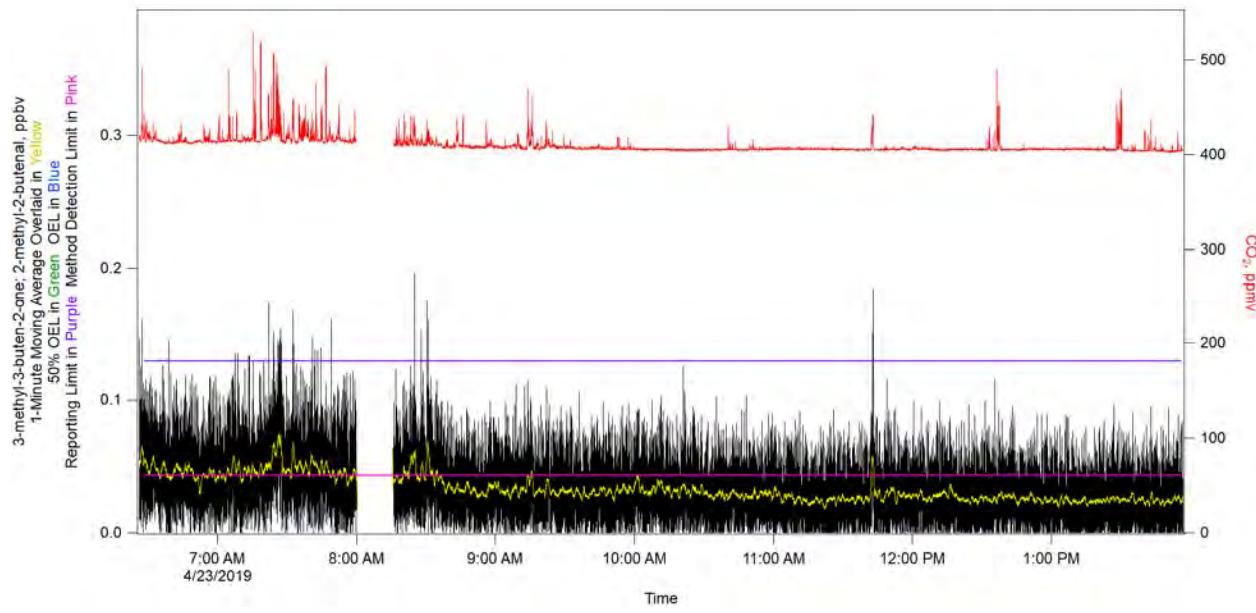


Figure 3-25. 3-methyl-3-buten-2-one; 2-methyl-2-butenal.

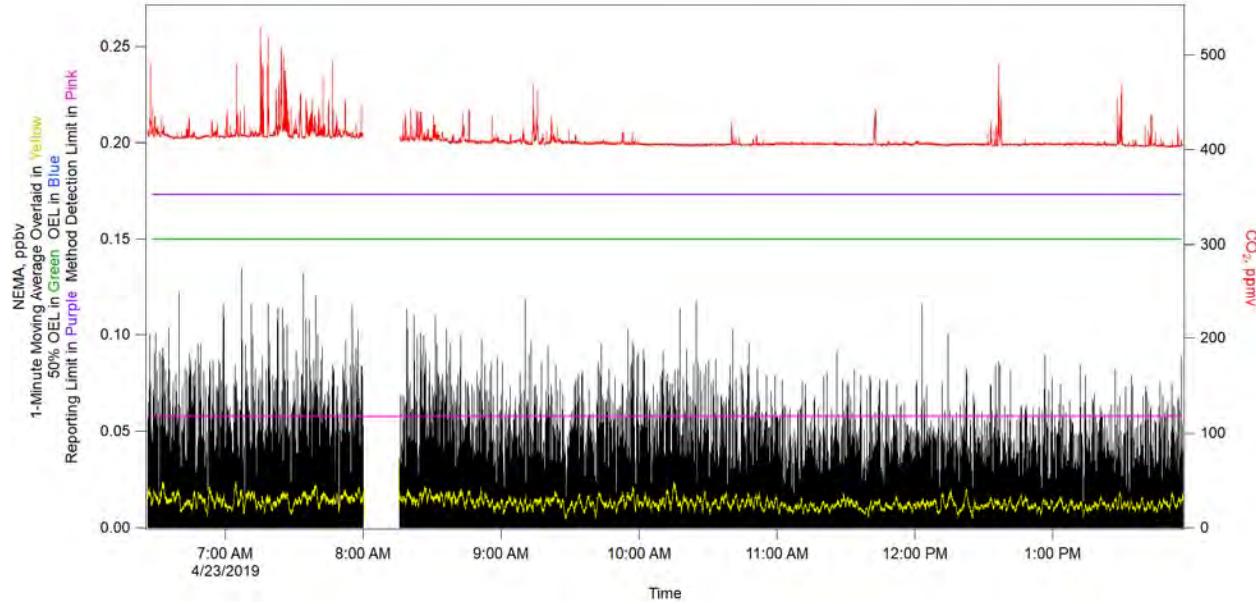


Figure 3-26. N-nitrosomethylethylamine (NEMA).

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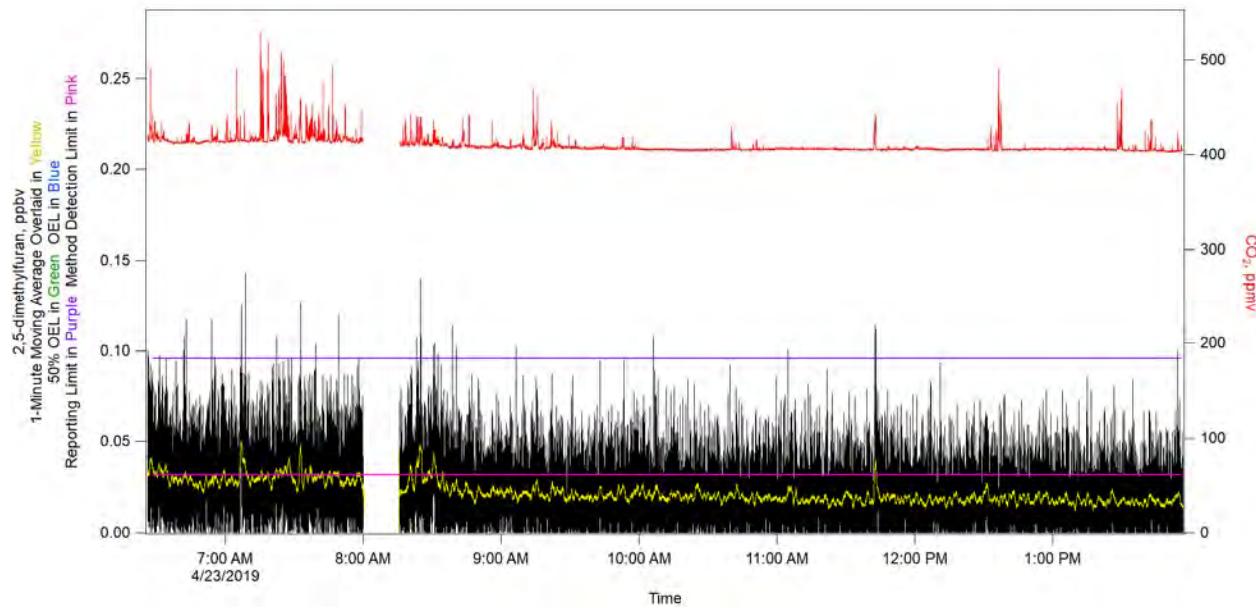


Figure 3-27. 2,5-dimethylfuran.

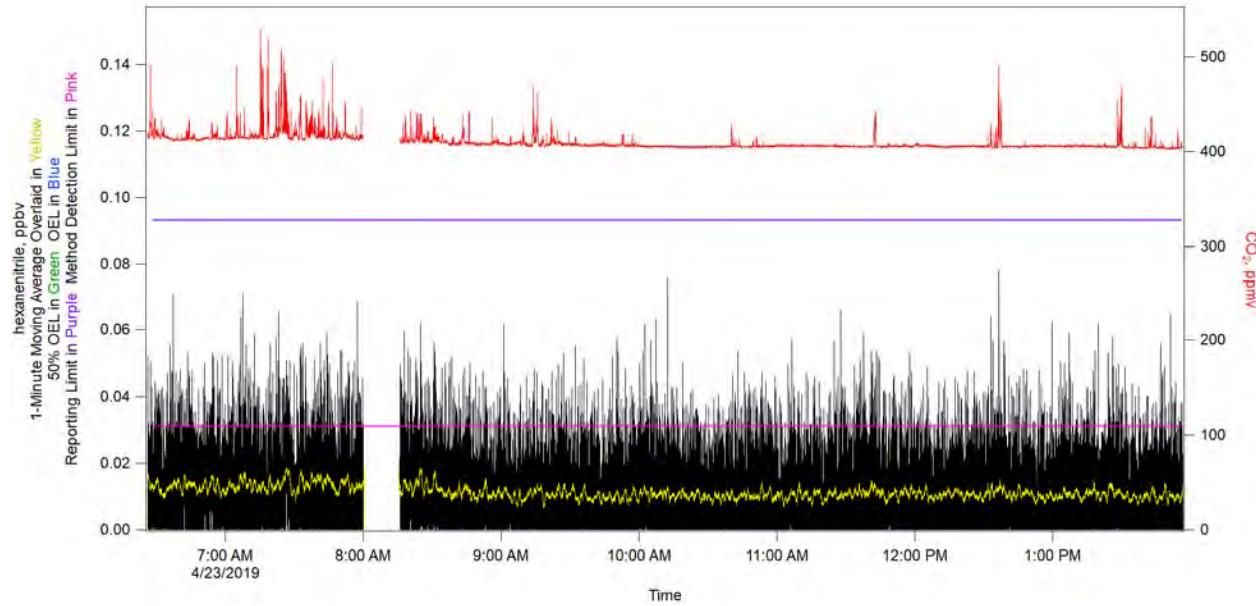


Figure 3-28. Hexanenitrile.

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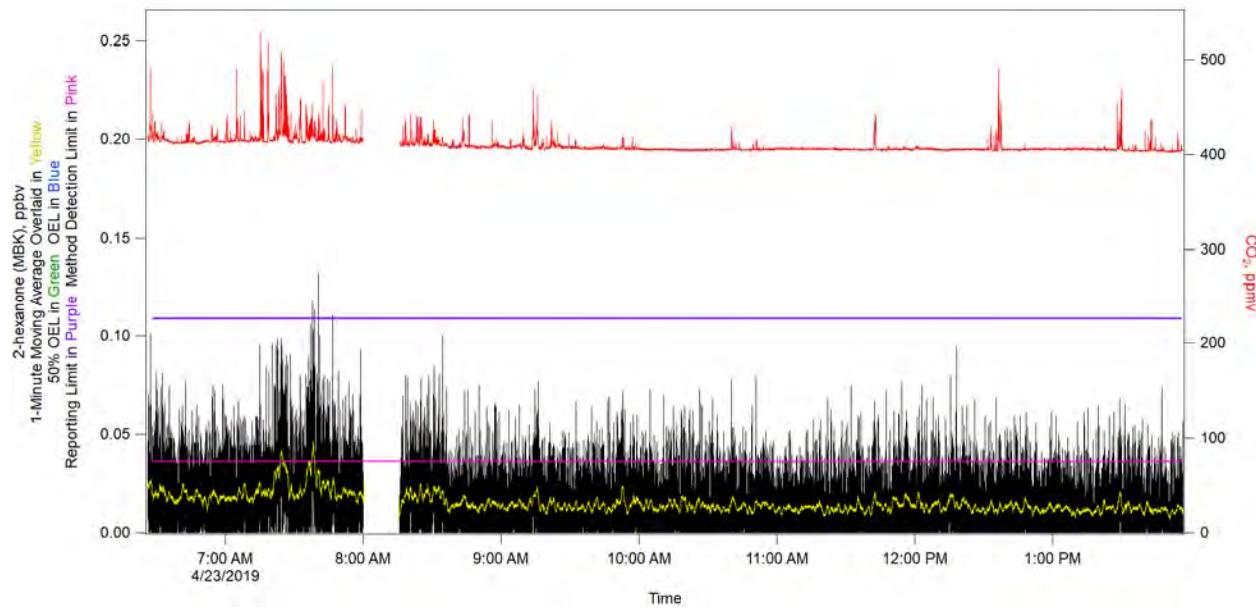


Figure 3-29. 2-hexanone (MBK).

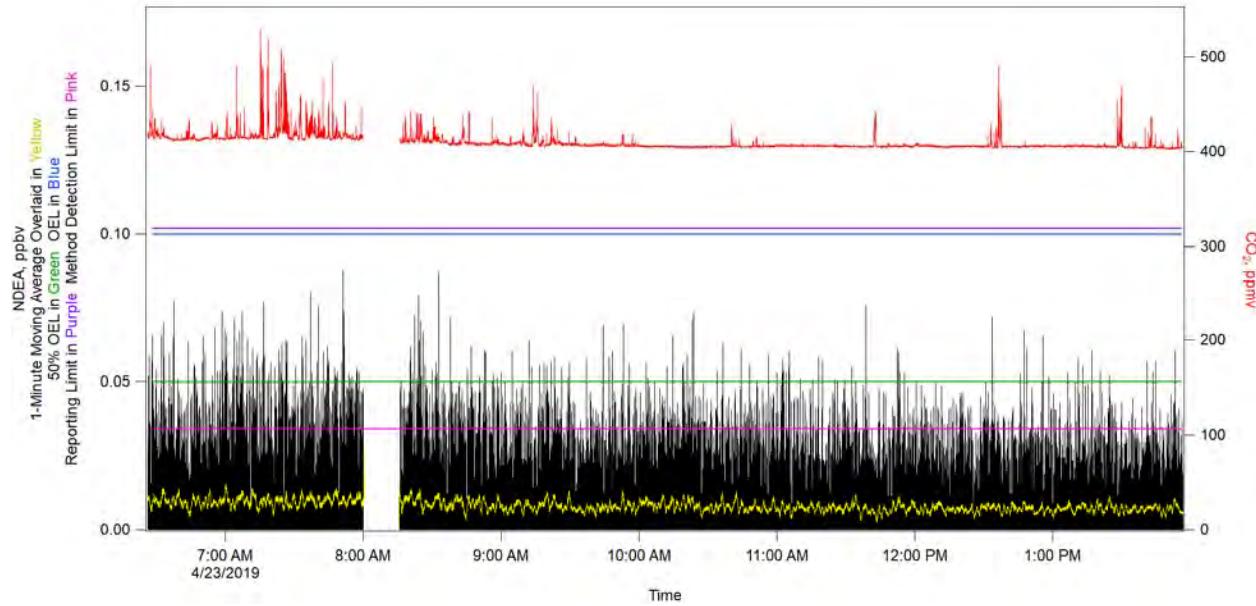


Figure 3-30. N-nitrosodiethylamine (NDEA).

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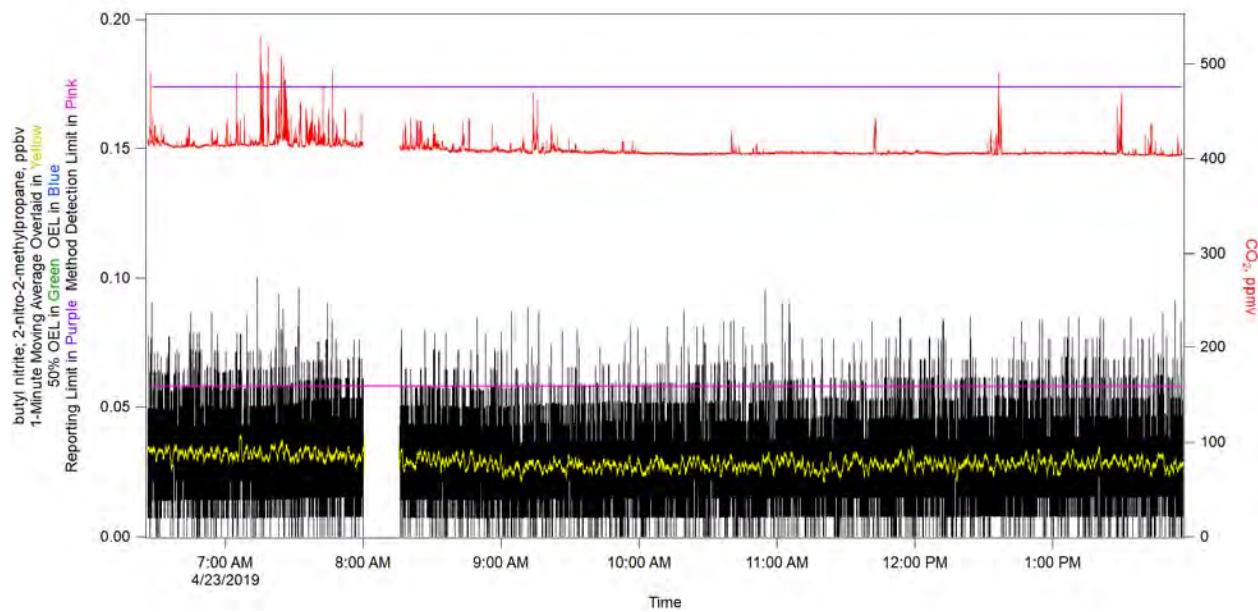


Figure 3-31. Butyl Nitrite; 2-nitro-2-methylpropane.

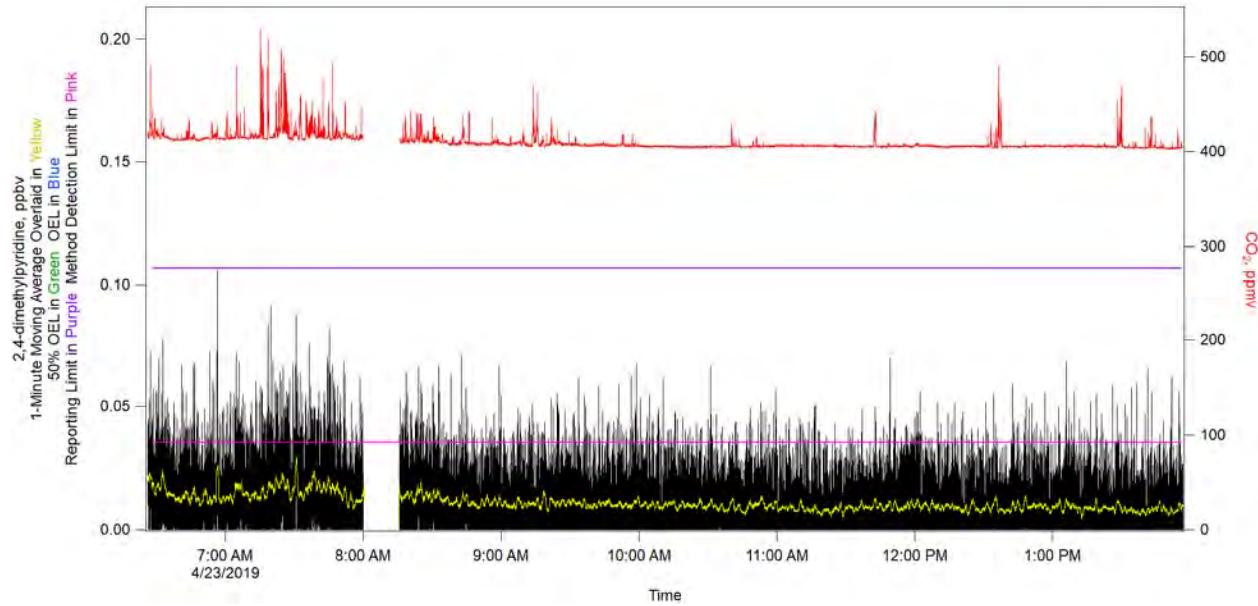


Figure 3-32. 2,4-dimethylpyridine.

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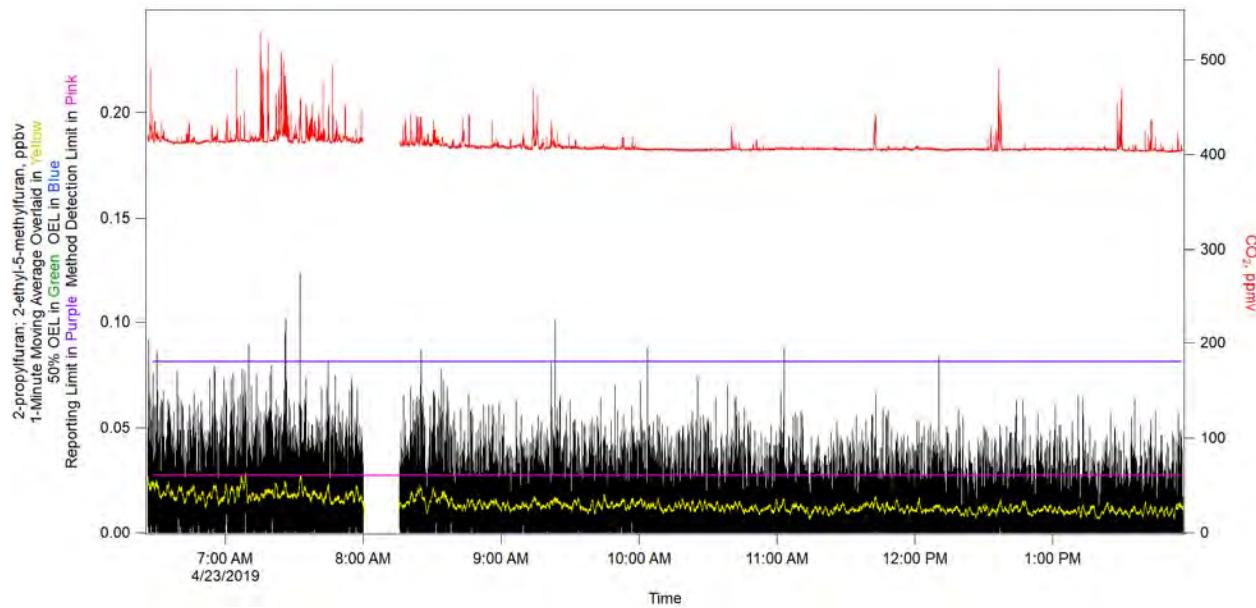


Figure 3-33. 2-propylfuran; 2-ethyl-5-methylfuran.

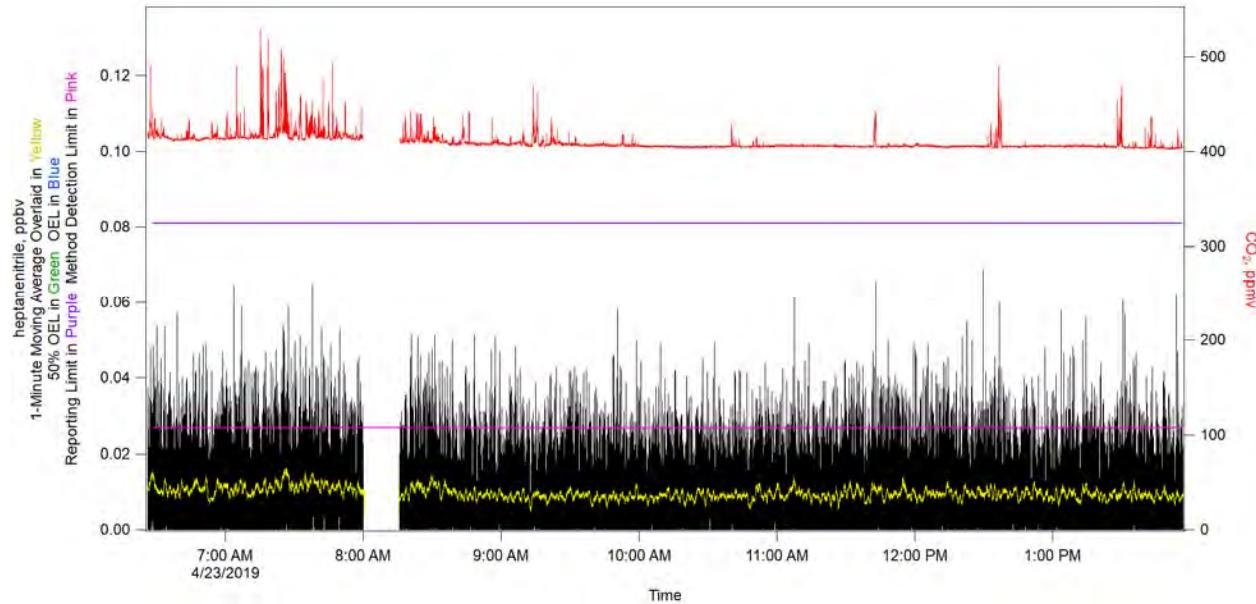


Figure 3-34. Heptanenitrile.

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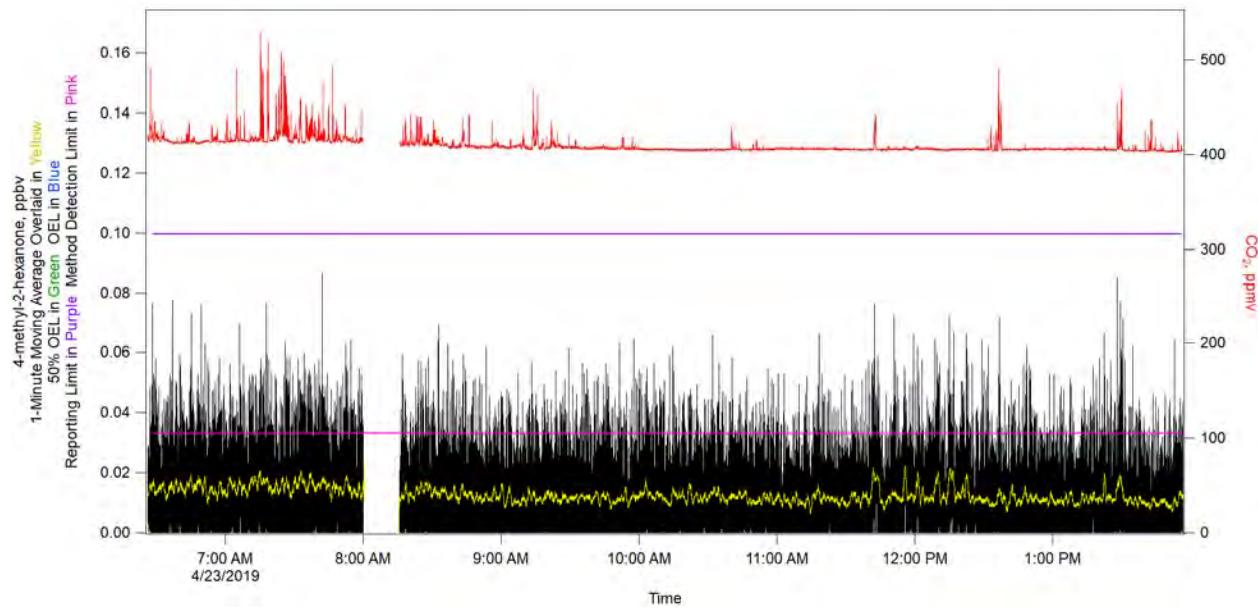


Figure 3-35. 4-methyl-2-hexanone.

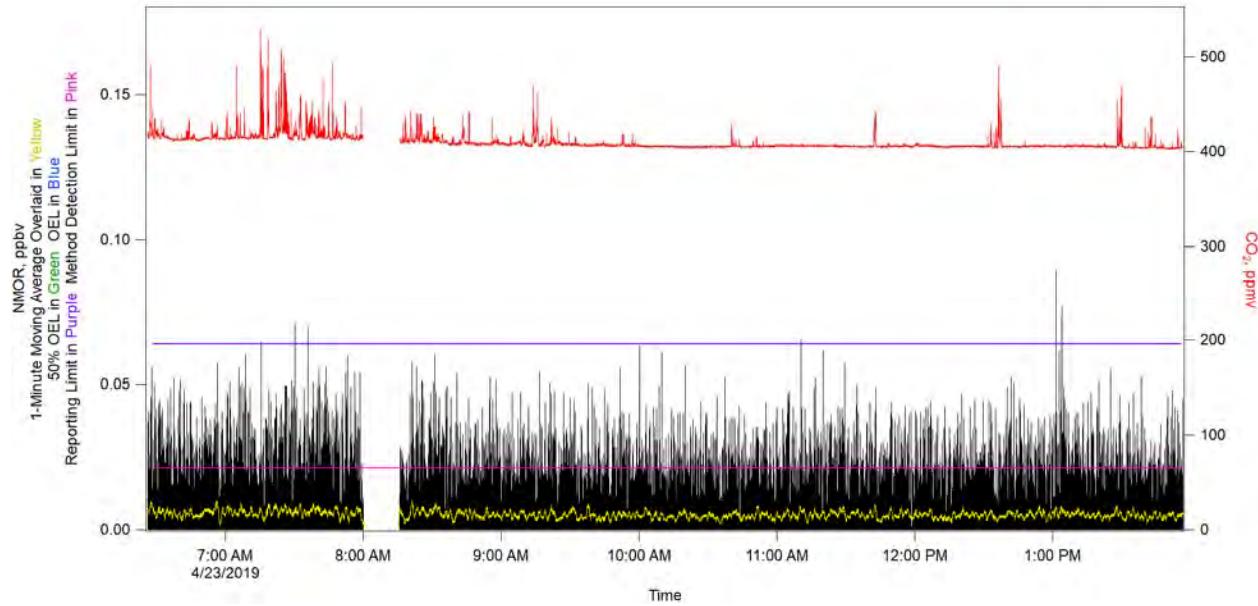


Figure 3-36. N-nitrosomorpholine (NMOR).

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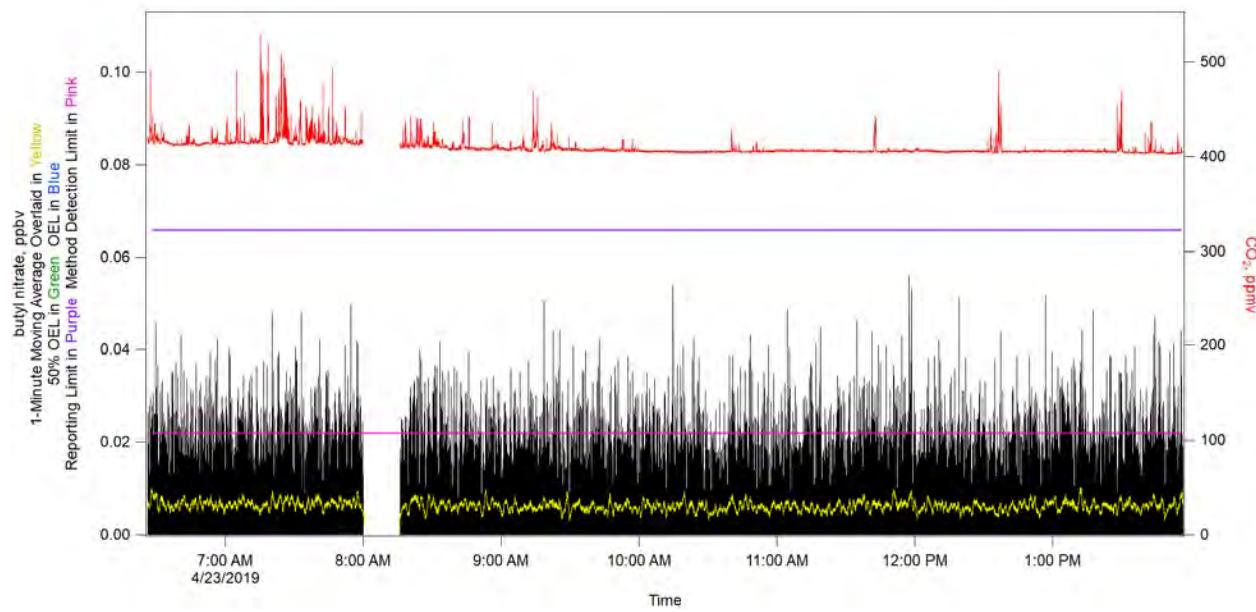
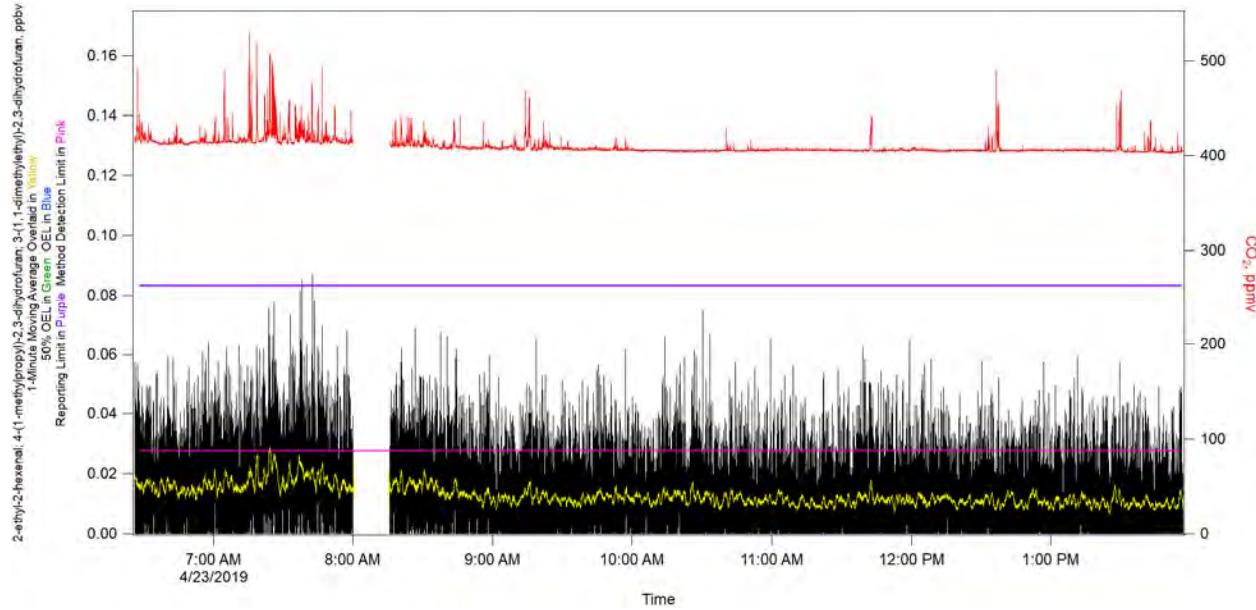


Figure 3-37. Butyl Nitrate.



**Figure 3-38. 2-ethyl-2-hexenal; 4-(1-methylpropyl)-2,3-dihydrofuran
3-(1,1-dimethylethyl)-2,3-dihydrofuran.**

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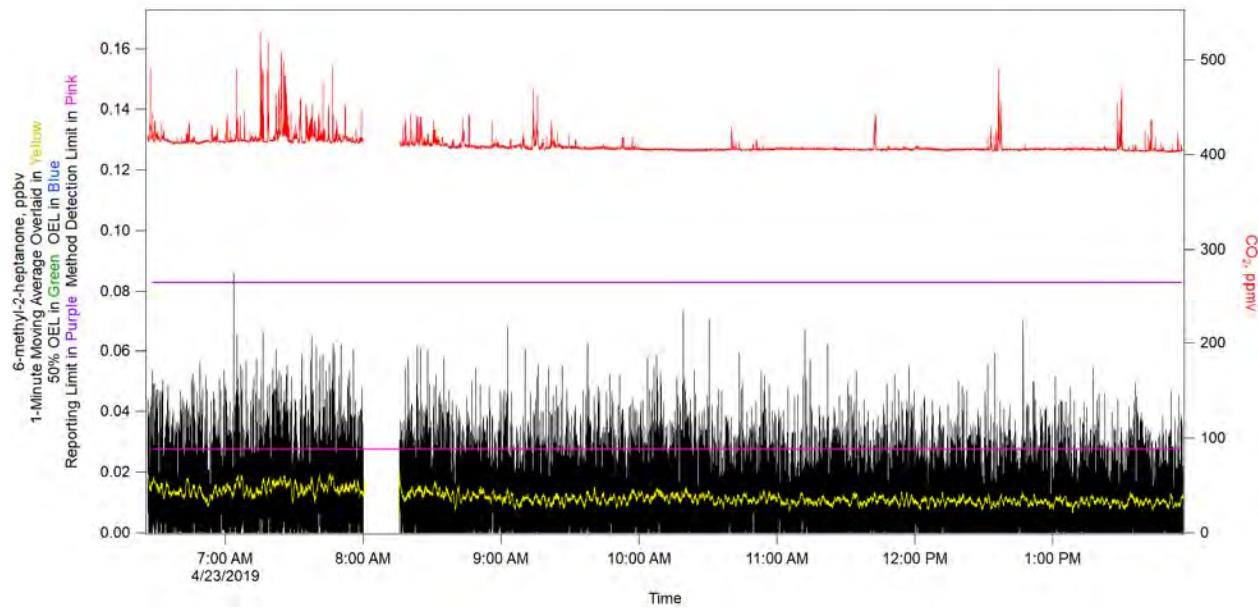


Figure 3-39. 6-methyl-2-heptanone.

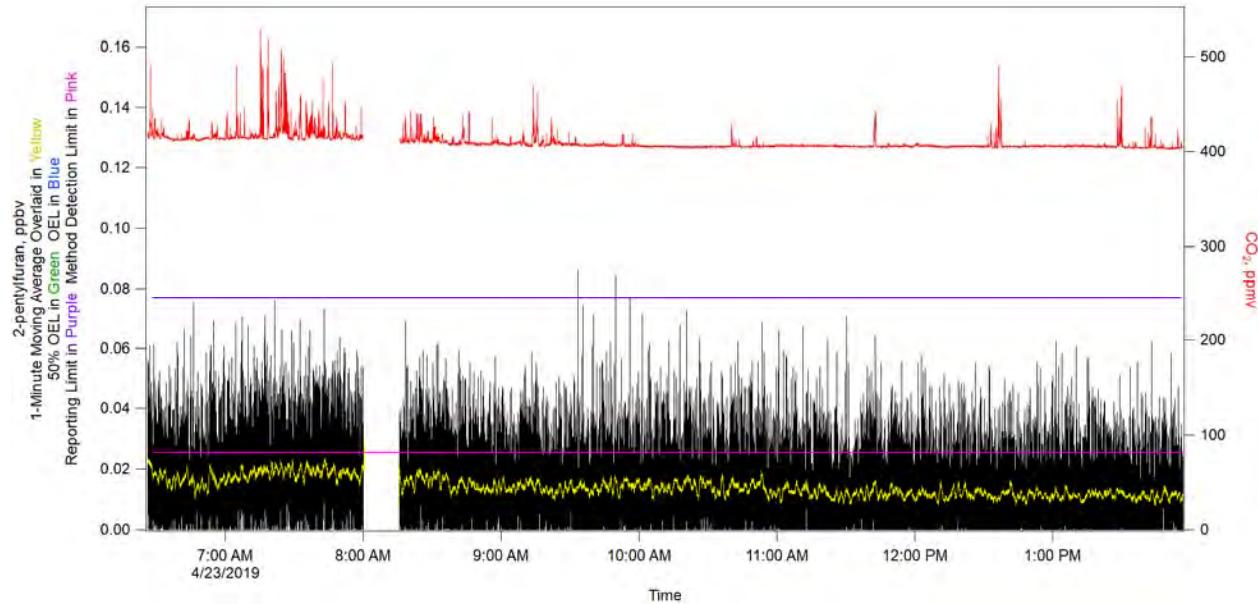


Figure 3-40. 2-pentylfuran.

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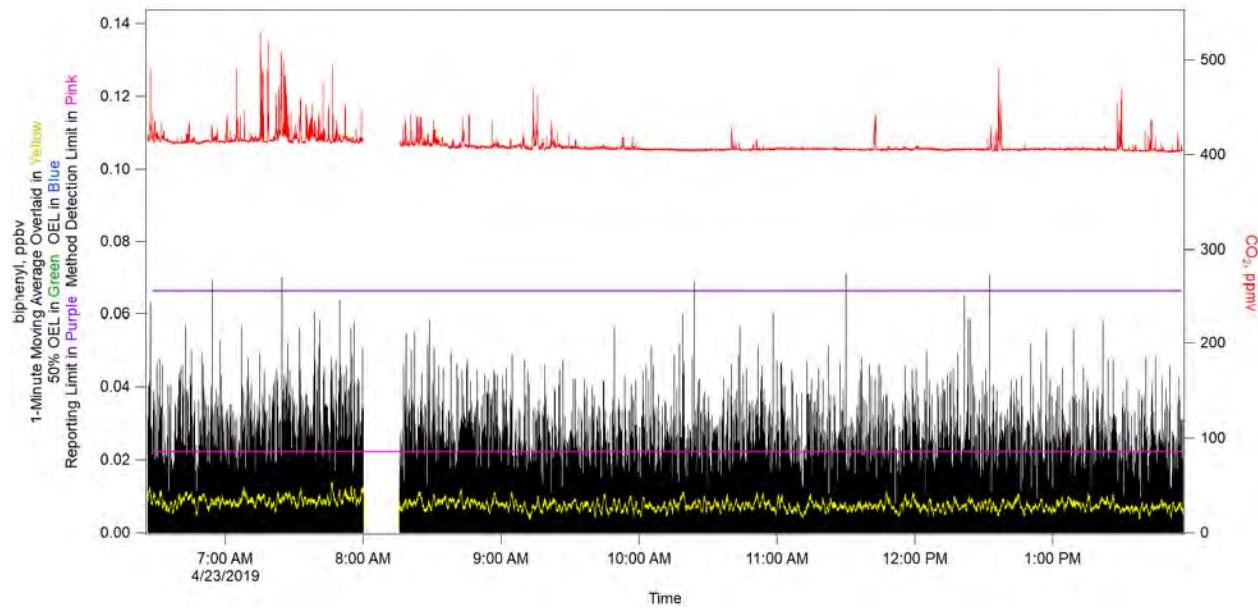


Figure 3-41. Biphenyl.

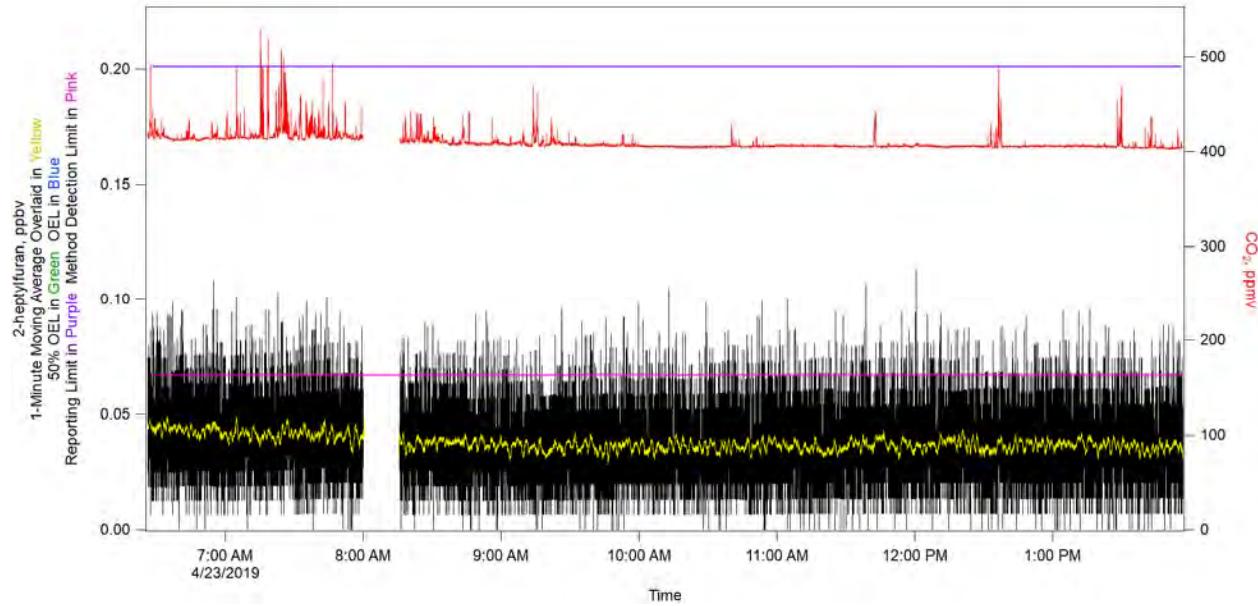


Figure 3-42. 2-heptylfuran.

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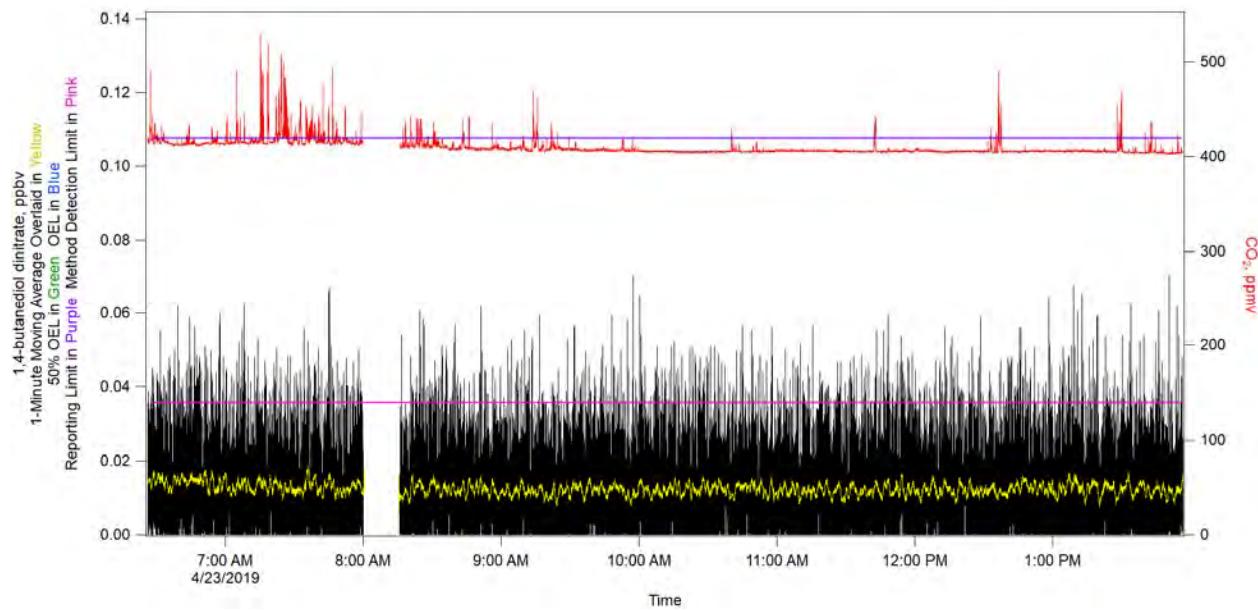


Figure 3-43. 1,4-butanediol Dinitrate.

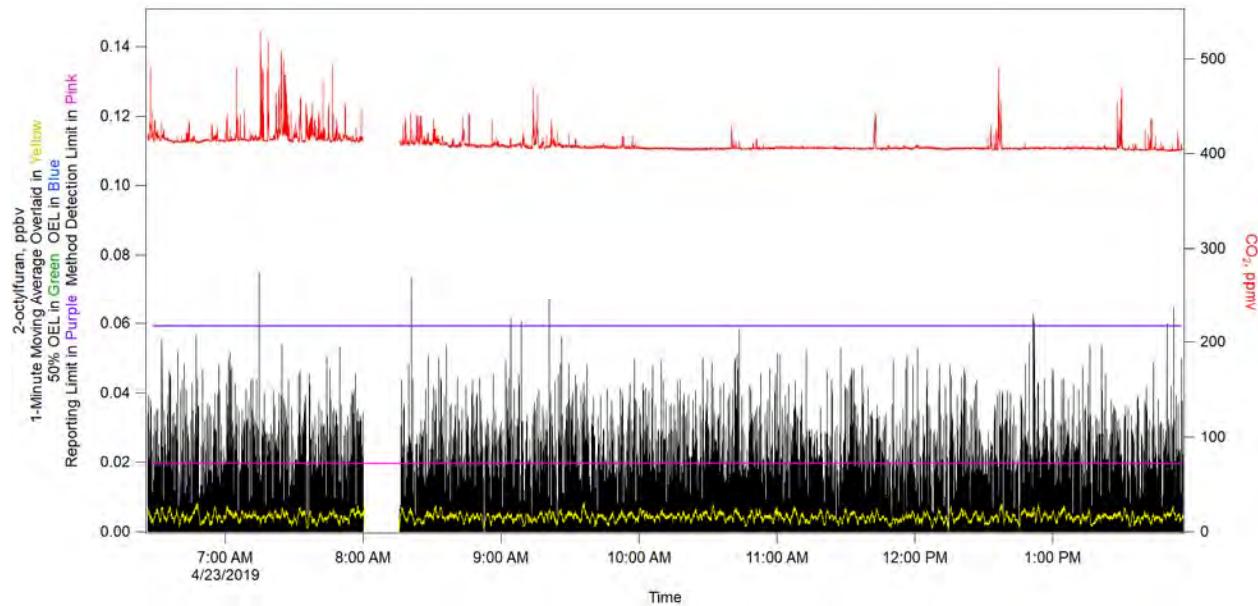


Figure 3-44. 2-octylfuran.

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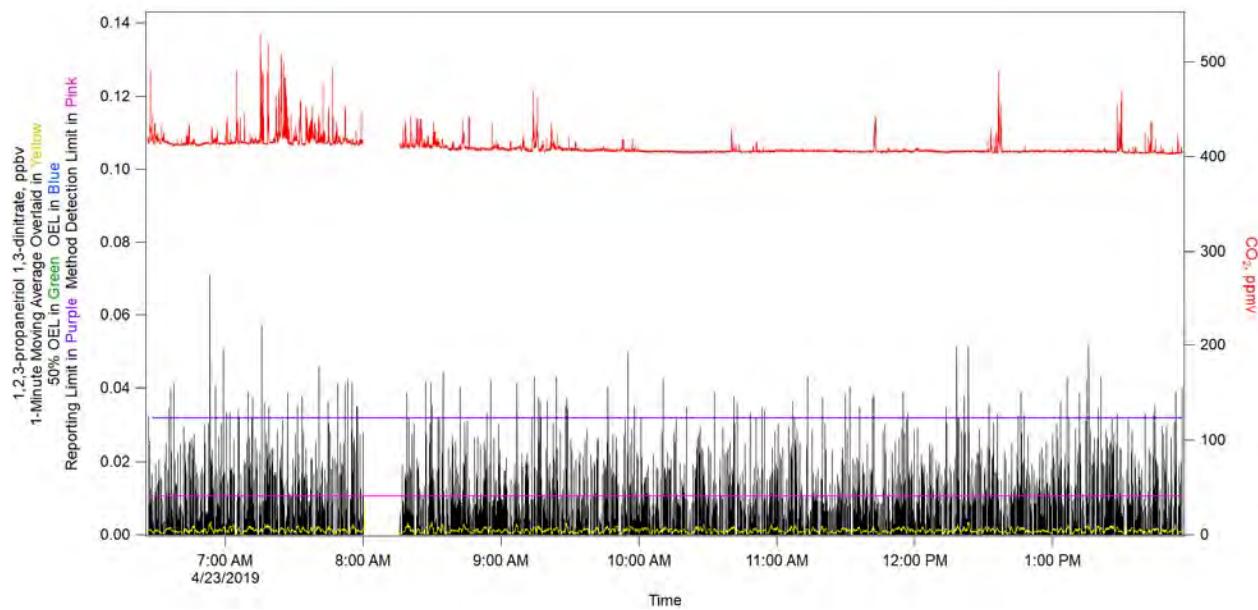


Figure 3-45. 1,2,3-propanetriol 1,3-dinitrate.

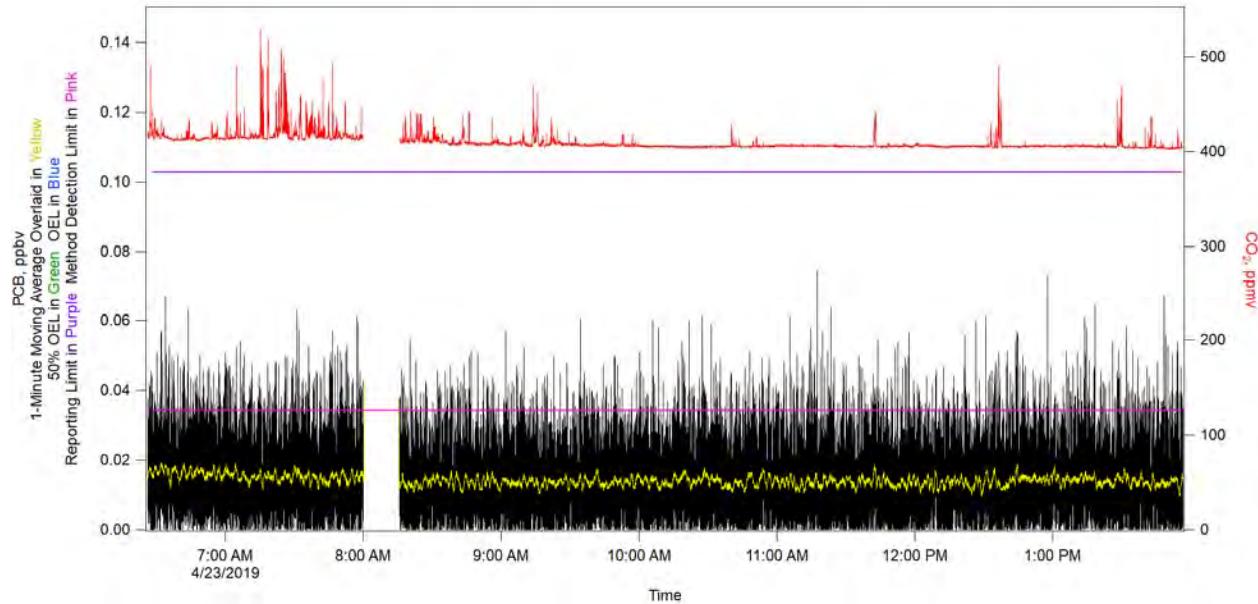


Figure 3-46. PCB.

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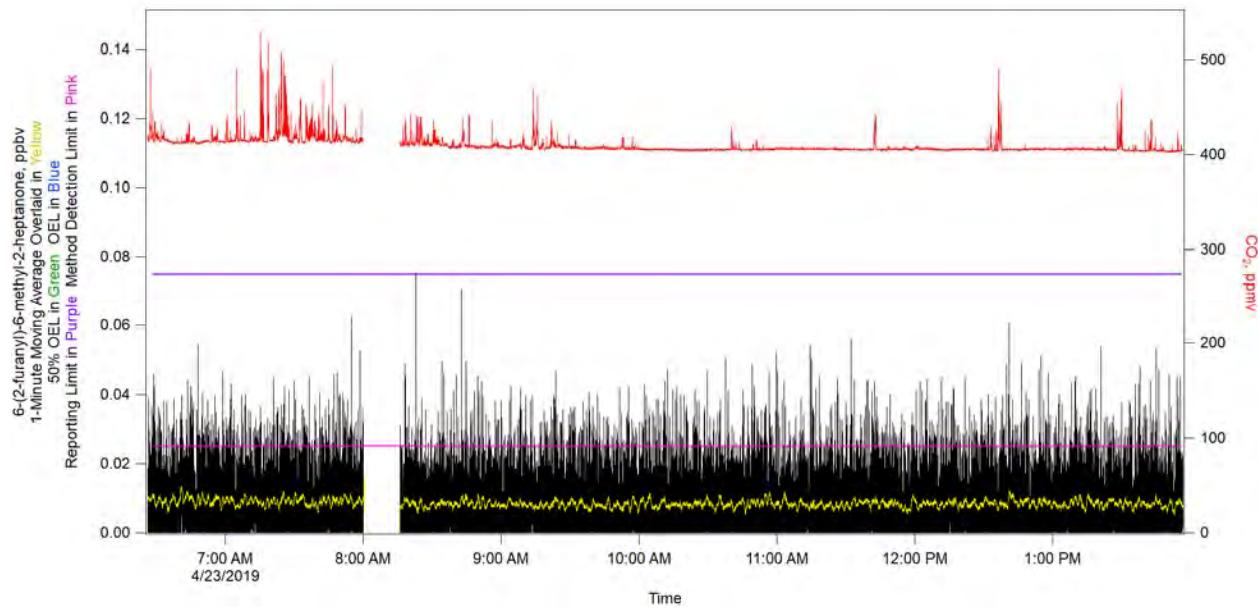


Figure 3-47. 6-(2-furanyl)-6-methyl-2-heptanone.

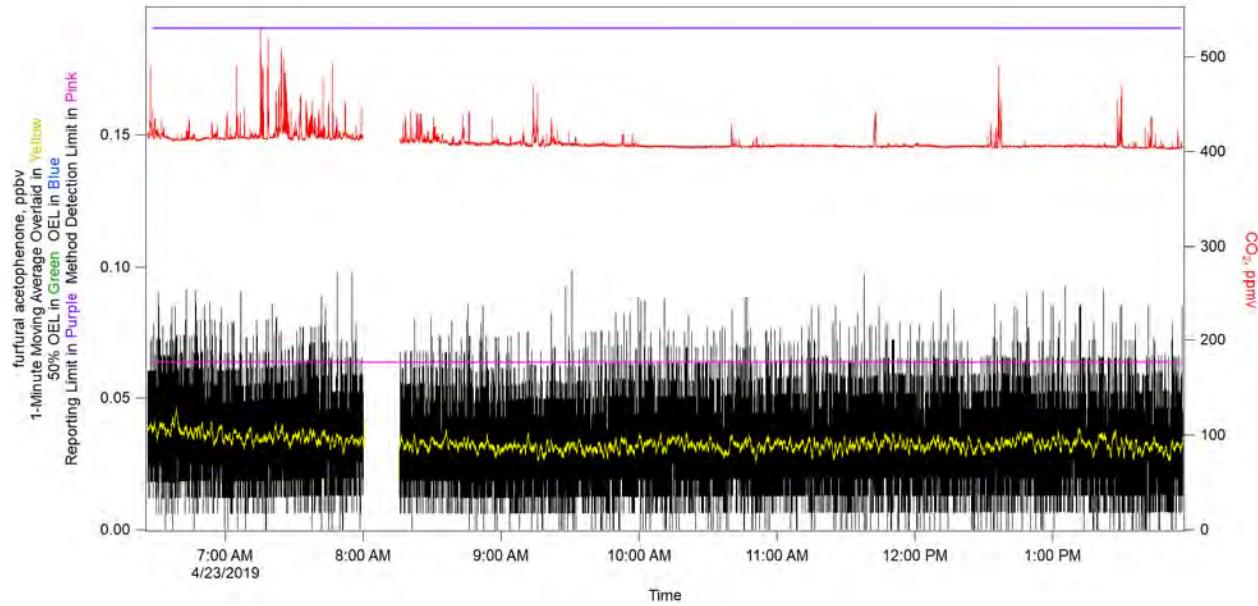


Figure 3-48. Furfural Acetophenone.

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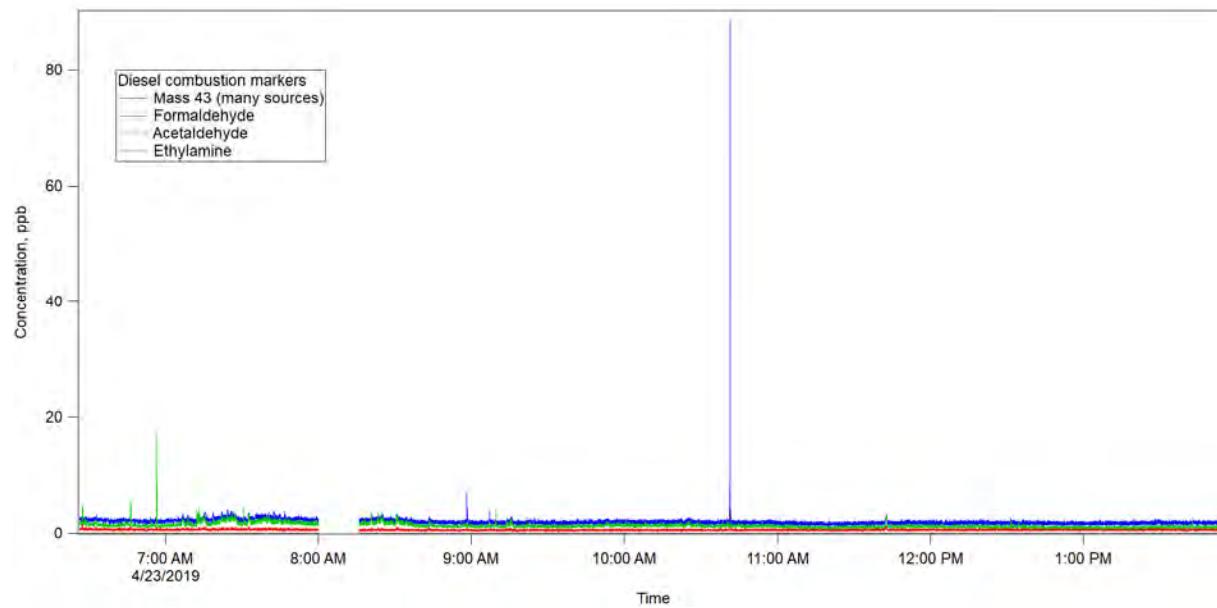


Figure 3-49. Diesel Combustion Markers.

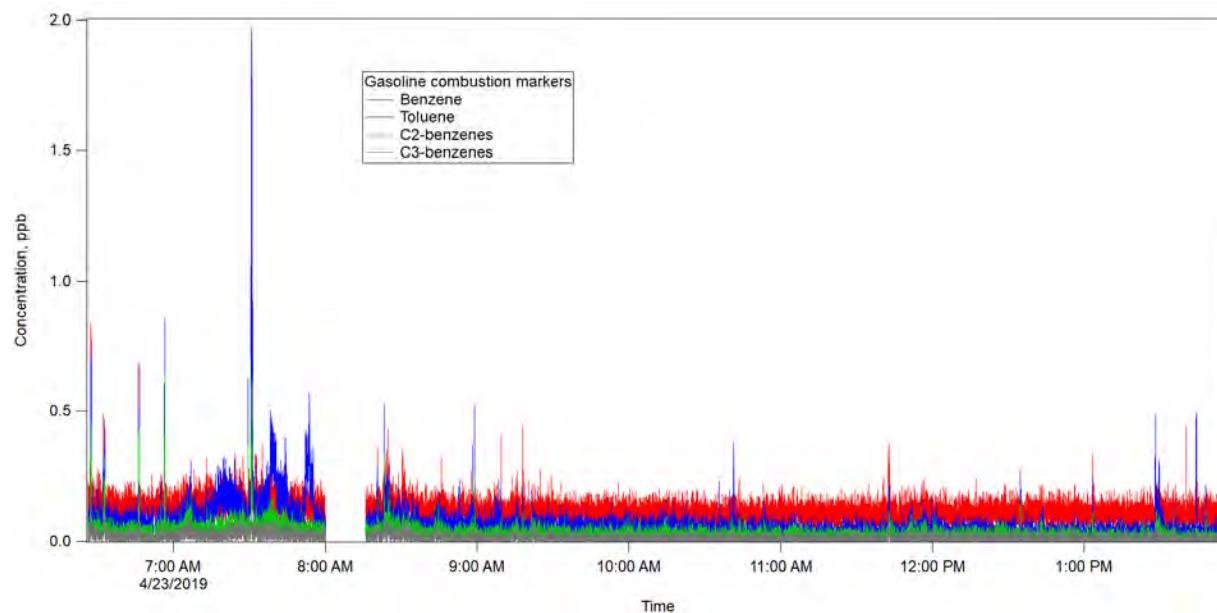


Figure 3-50. Gasoline Combustion Markers.

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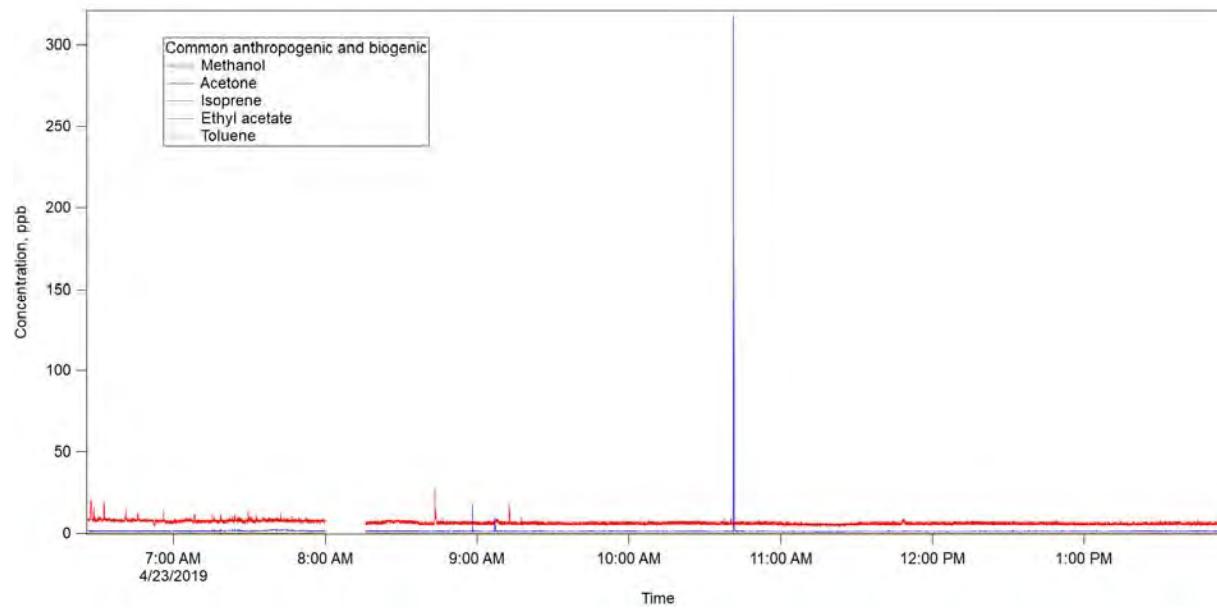


Figure 3-51. Plant and Human Markers

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4.0 APRIL 24, 2019 – AREA MONITORING

4.1 Quality Assessment

Data from April 24, 2019, were assessed using Procedure 17124-DOE-HS-102. A Data Exchange Checklist was completed. The data were accepted by TerraGraphics with the following comments. Report No. 66409-RPT-004 was adequately documented and all checks passed the acceptance limits.

4.2 Summary

On April 24, 2019, ML personnel arrived at the TerraGraphics warehouse at 04:58. The QA/QC zero-air/span checks were performed on the LI-COR CO₂ monitor, the Picarro NH₃ analyzer, and the PTR-MS beginning at 05:14. The ML arrived on the Hanford Site and checked in with the CSM at 06:20. The ML began a site survey loop of A Farms at 06:30. The ML parked on the east side of A Farms from 07:02 to 09:30. The ML initiated a site survey loop at 09:30 until parking on the southeast corner of A Farm. The ML was relocated to the southeast side of C Farm at 11:08, due to varying wind directions. At 12:09, the ML Operators started a site survey loop until stopping on the northeast side of AP Farm, downwind of workers working on a pump within the farm. The Operators checked out with the CSM at 13:55 and the ML departed the site. The ML arrived back at the TerraGraphics warehouse at 14:54.

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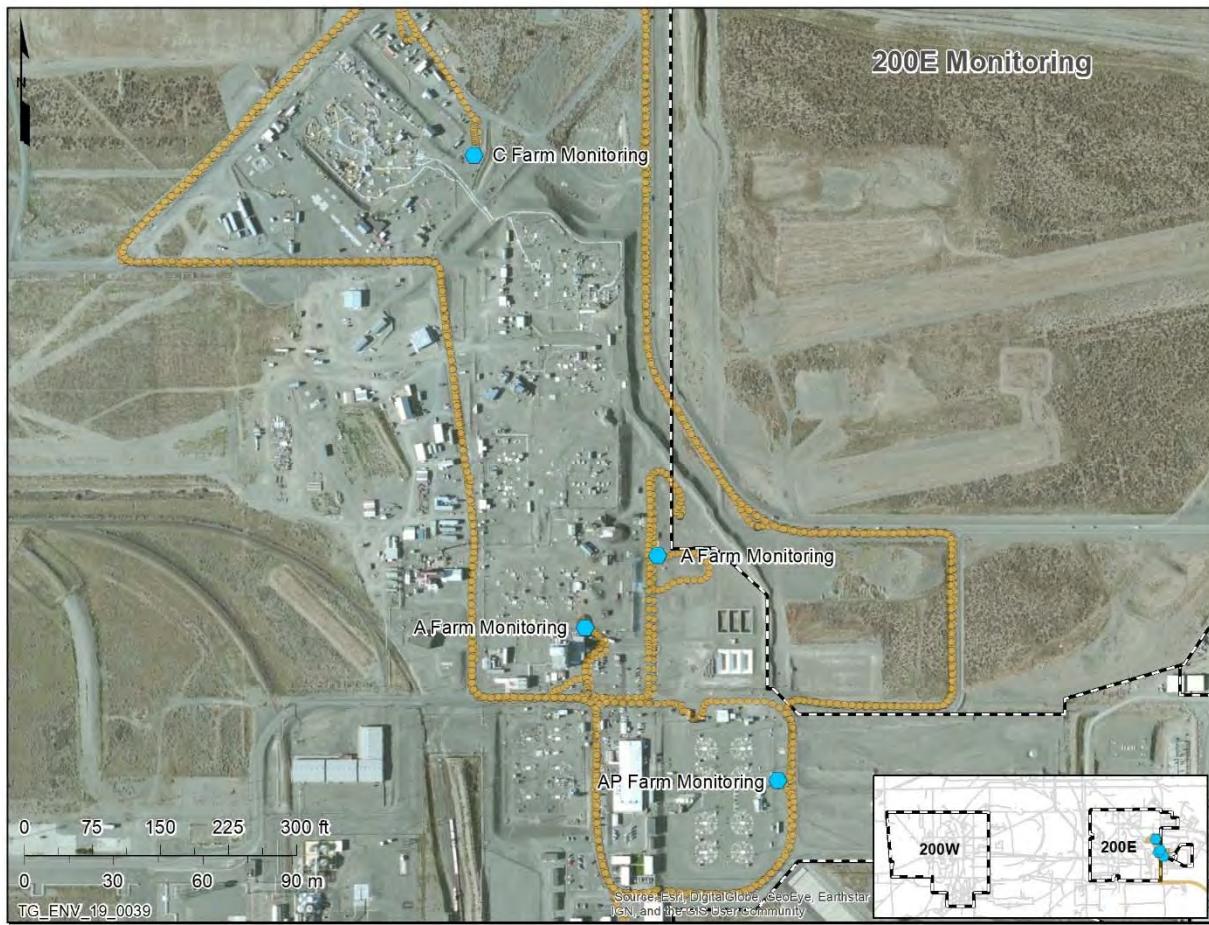


Figure 4-1. Location of the Mobile Laboratory for the Duration of the Monitoring Period.

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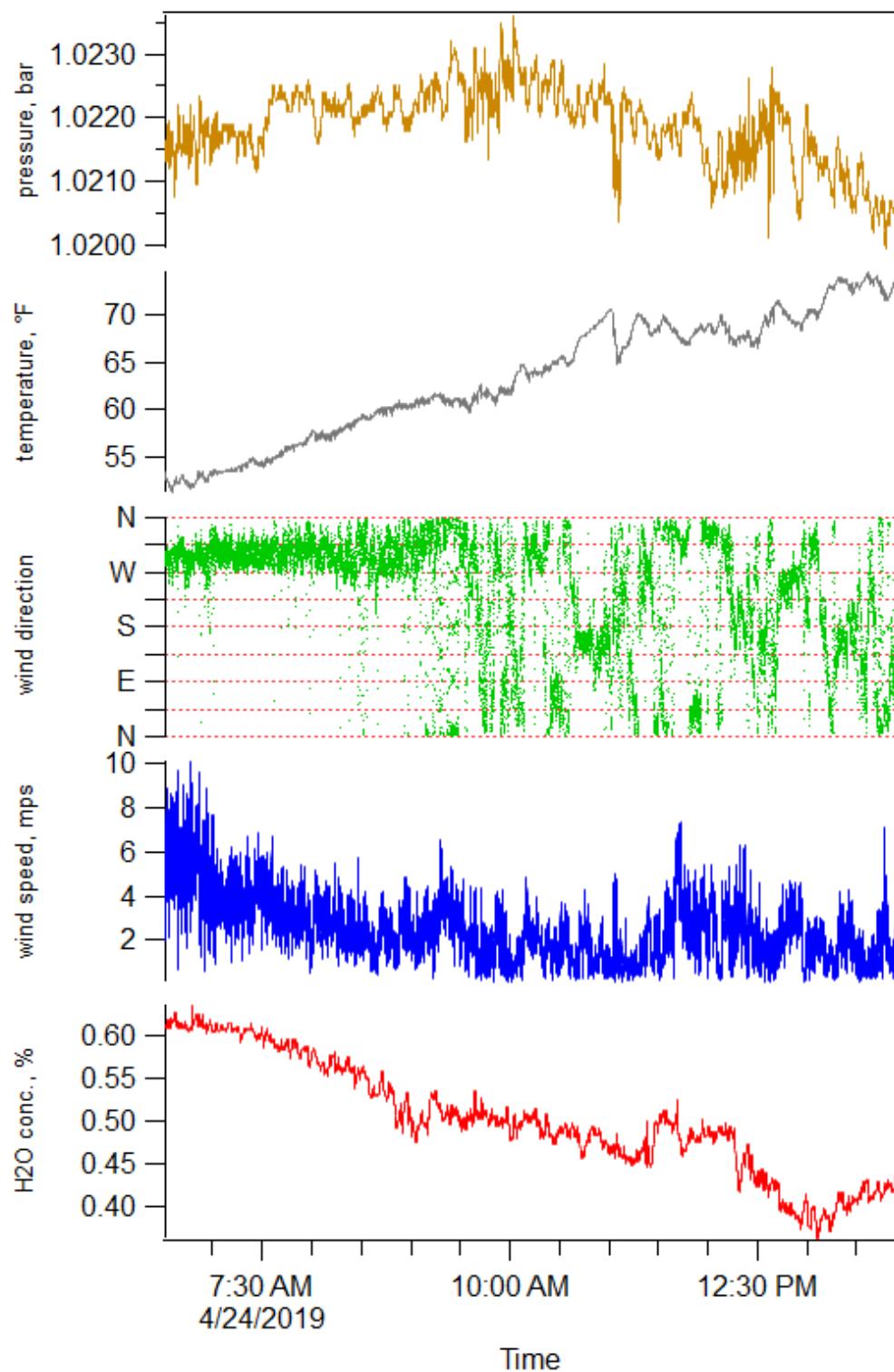


Figure 4-2. Weather Data for the Duration of the Monitoring Period.

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4.3 Samples Collected

Continuous air monitoring was performed using the following instrumentation:

- PTR-TOF 6000 X2,
- LI-COR CO₂ Monitor,
- Picarro Ammonia Monitor, and
- Airmar Weather Station.

Confirmatory air samples were not collected during this period.

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4.4 Area Monitoring

Table 4-1. Chemical of Potential Concern Statistical Information for the Monitoring Period of April 24, 2019. (2 Sheets)

COPC #	COPC Name	OEL (ppb)	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel St. Dev. (%)	Max. (ppb)	Median (ppb)
1	ammonia	25000	6.225	6.433†	1.619	25.173	12.918	5.807†
2	formaldehyde	300	1.721	<1.721	0.192	24.970	3.004	<1.721
3	methanol	200000	5.758	<5.758	0.537	9.661	11.743	<5.758
4	acetonitrile	20000	0.085	0.106†	0.034	32.614	0.749	0.103†
5	acetaldehyde	25000	1.027	1.089†	0.400	36.712	12.713	1.019†
6	ethylamine	5000	0.069	<0.069	0.021	63.894	0.149	<0.069
7	1,3-butadiene	1000	0.183	<0.183	0.070	53.208	0.860	<0.183
8	propanenitrile	6000	0.107	<0.107	0.029	59.594	0.648	<0.107
9	2-propenal	100	0.340	<0.340	0.056	71.789	0.830	<0.340
10	1-butanol + butenes	20000	0.214	<0.214	0.099	101.016	2.552	<0.214
11	methyl isocyanate	20	0.069	<0.069	0.024	59.737	0.172	<0.069
12	methyl nitrite	100	0.098	<0.098	0.040	54.107	0.689	<0.098
13	furan	1	0.062	<0.062	0.017	64.582	0.177	<0.062
14	Butanenitrile	8000	0.039	<0.039	0.017	79.130	0.278	<0.039
15	but-3-en-2-one + 2,3-dihydrofuran + 2,5-dihydrofuran	200, 1, 1	0.041	<0.041	0.020	72.524	N/A*	N/A*
16	butanal	25000	0.061	0.083†	0.028	33.325	0.257	0.081†
17	NDMA**	0.3	0.082	<0.082	0.020	162.759	0.174	<0.082
18	benzene	500	0.236	<0.236	0.245	156.516	10.653	<0.236
19	2,4-pentadienenitrile + pyridine	300, 1000	0.085	<0.085	0.024	58.787	0.816	<0.085
20	2-methylene butanenitrile	300	0.036	<0.036	0.011	80.193	0.096	<0.036
21	2-methylfuran	1	0.043	<0.043	0.020	75.306	0.282	<0.043
22	pentanenitrile	6000	0.036	<0.036	0.012	89.225	0.131	<0.036
23	3-methyl-3-butene-2-one + 2-methyl-2-butenal	20, 30	0.043	<0.043	0.018	76.310	0.199	<0.043
24	NEMA**	0.3	0.058	<0.058	0.015	139.362	0.108	<0.058
25	2,5-dimethylfuran	1	0.032	<0.032	0.015	89.864	0.133	<0.032
26	hexanenitrile	6000	0.031	<0.031	0.011	96.632	0.116	<0.031
27	2-hexanone (MBK)	5000	0.036	<0.036	0.012	95.091	0.097	<0.036
28	NDEA**	0.1	0.034	<0.034	0.010	148.635	0.097	<0.034

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Table 4-1. Chemical of Potential Concern Statistical Information for the Monitoring Period of April 24, 2019. (2 Sheets)

COPC #	COPC Name	OEL (ppb)	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel St. Dev. (%)	Max. (ppb)	Median (ppb)
29	butyl nitrite + 2-nitro-2-methylpropane	100, 300	0.058	<0.058	0.014	53.135	0.146	<0.058
30	2,4-dimethylpyridine	500	0.036	<0.036	0.018	151.138	0.513	<0.036
31	2-propylfuran + 2-ethyl-5-methylfuran	1	0.027	<0.027	0.011	110.342	0.111	<0.027
32	heptanenitrile	6000	0.027	<0.027	0.009	99.241	0.071	<0.027
33	4-methyl-2-hexanone	500	0.033	<0.033	0.010	96.276	0.070	<0.033
34	NMOR**	0.6	0.021	<0.021	0.009	176.297	0.098	<0.021
35	butyl nitrate	2500	0.022	<0.022	0.007	130.281	0.061	<0.022
36	2-ethyl-2-hexenal + 4-(1-methylpropyl)-2,3-dihydrofuran + 3-(1,1-dimethylethyl)-2,3-dihydrofuran	100, 1, 1	0.028	<0.028	0.010	95.048	0.079	<0.028
37	6-methyl-2-heptanone	8000	0.028	<0.028	0.009	94.952	0.069	<0.028
38	2-pentylfuran	1	0.026	<0.026	0.010	97.657	0.066	<0.026
39	biphenyl	200	0.022	<0.022	0.009	130.082	0.065	<0.022
40	2-heptylfuran	1	0.067	<0.067	0.015	43.444	0.121	<0.067
41	1,4-butanediol dinitrate	50	0.036	<0.036	0.010	89.723	0.073	<0.036
42	2-octylfuran	1	0.020	<0.02	0.008	209.628	0.068	<0.02
43	1,2,3-propanetriol 1,3-dinitrate	50	0.011	<0.011	0.004	373.112	0.053	<0.011
44	PCB	1000	0.034	<0.034	0.010	73.982	0.075	<0.034
45	6-(2-furanyl)-6-methyl-2-heptanone	1	0.025	<0.025	0.008	103.426	0.067	<0.025
46	furfural acetophenone	1	0.064	<0.064	0.015	44.099	0.104	<0.064
N/A*	The maximum peak value for but-3-en-2-one + 2,3 dihydrofuran + 2,5 dihydrofuran was 0.274 ppb and the median value was <0.041 ppb. The PTR-MS results for but-3-en-2-one + 2,3 dihydrofuran + 2,5 dihydrofuran are not compared to OEL concentrations because: 1) the result is suspect due to a known biogenic interferant (methacrolein) that is expected to be in concentrations that occasionally exceed the dihydrofuran OEL, and 2) this combination of COPCs have OEL concentrations that differ by a factor of 200, which provide widely variant bases for these numbers.							
**	Nitrosamine results are suspect due to isobaric interferants causing positive bias that have been encountered during previous background [53005-81-RPT-007, PTR-MS Mobile Laboratory Vapor Monitoring Background Study, (3/18/2018 – 4/20/2018), and Fiscal Year 2017 Mobile Laboratory Vapor Monitoring at the Hanford Site: Monitoring During Waste Disturbing Activities and Background Study, RJ Lee Group, Inc.].							
<	COPC Averages below the MDL.							
†	COPC Averages between the RL and the MDL.							
	COPC Averages >100% of the OEL.							
	COPC Averages 50-100% of the OEL.							
	COPC Averages 10-50% of the OEL.							

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Figure 4-3 through Figure 4-51 display 46 COPC signals, overlaid with the same signal smoothed using a 1-minute moving average (in cases where a moving average assists with data visualization), and CO₂, for the monitoring period April 24, 2019. If within range of the plot's left axis, a green horizontal line representing 50% of the COPC's OEL, a blue horizontal line representing the COPC's OEL, a horizontal purple line representing the RL, and a pink horizontal line representing the MDL are shown.

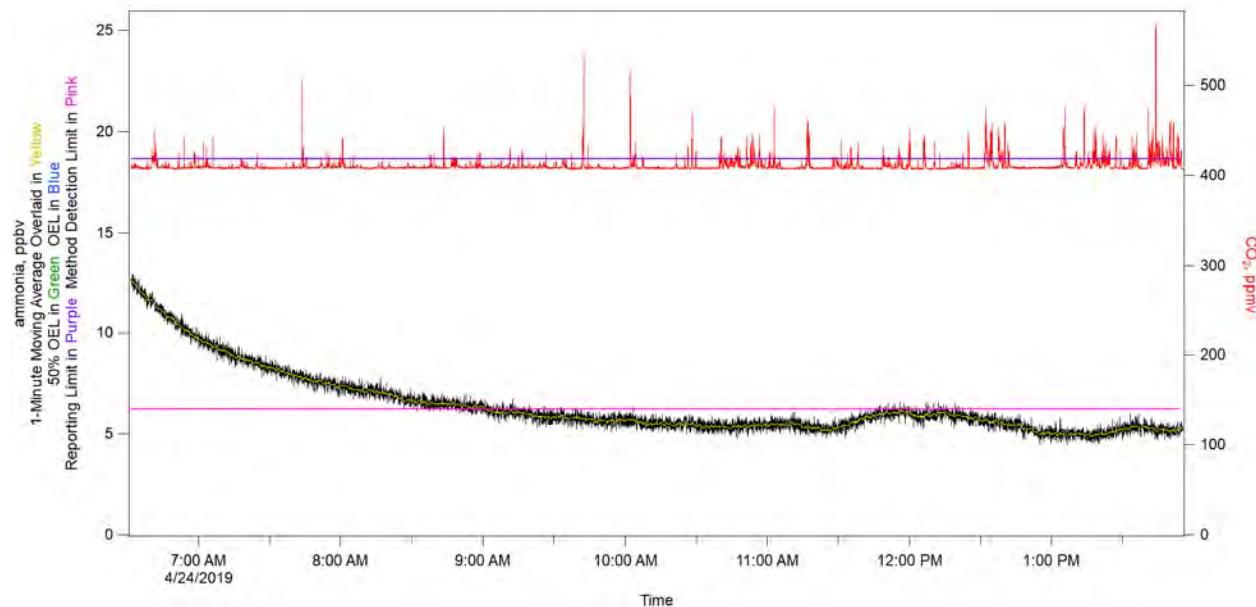


Figure 4-3. Ammonia.

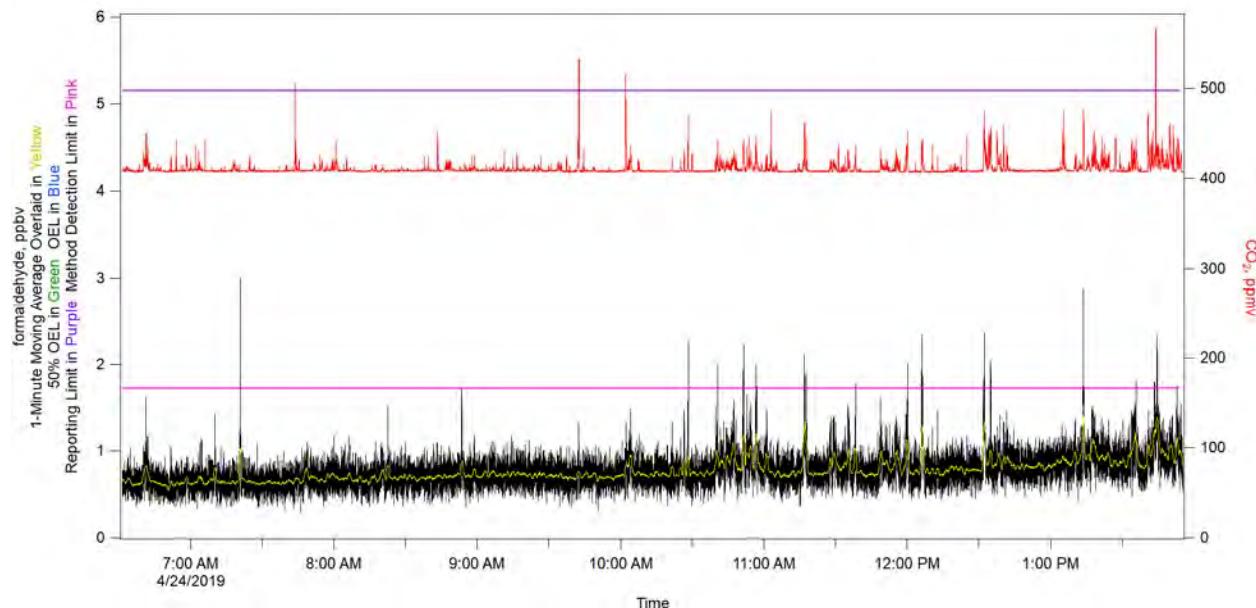


Figure 4-4. Formaldehyde.

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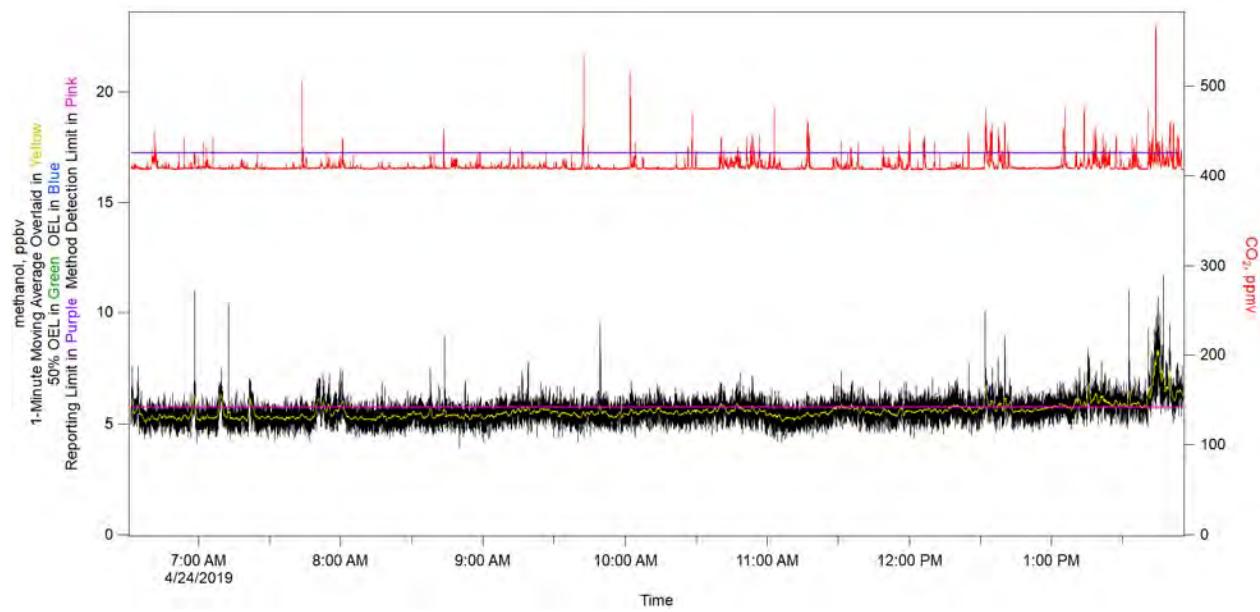


Figure 4-5. Methanol.

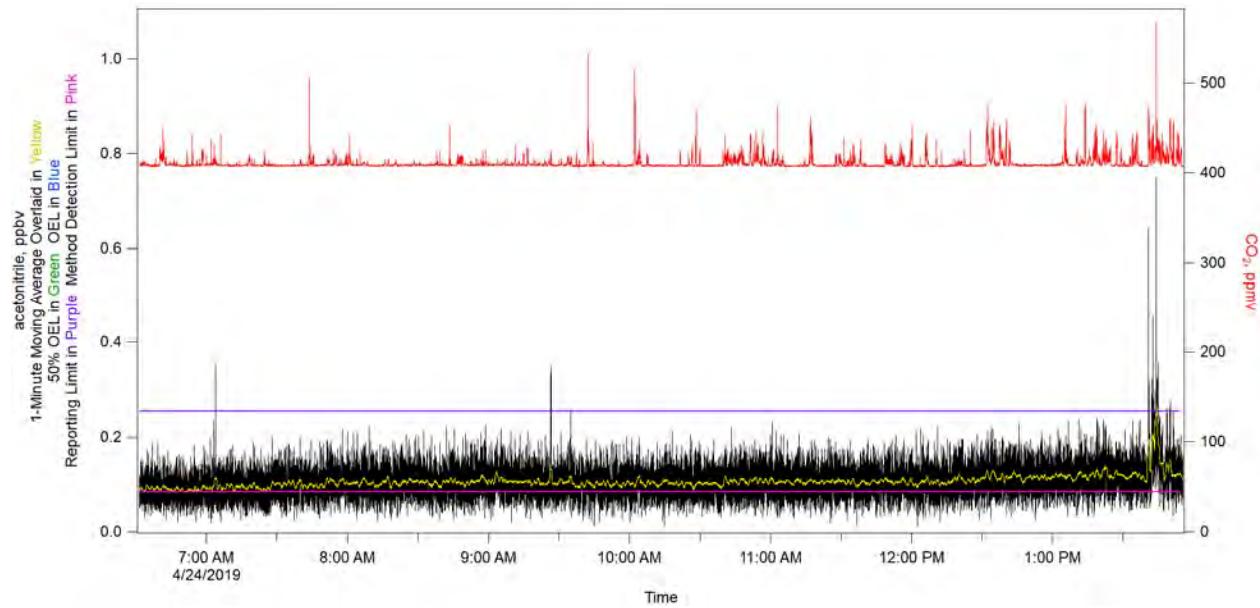


Figure 4-6. Acetonitrile.

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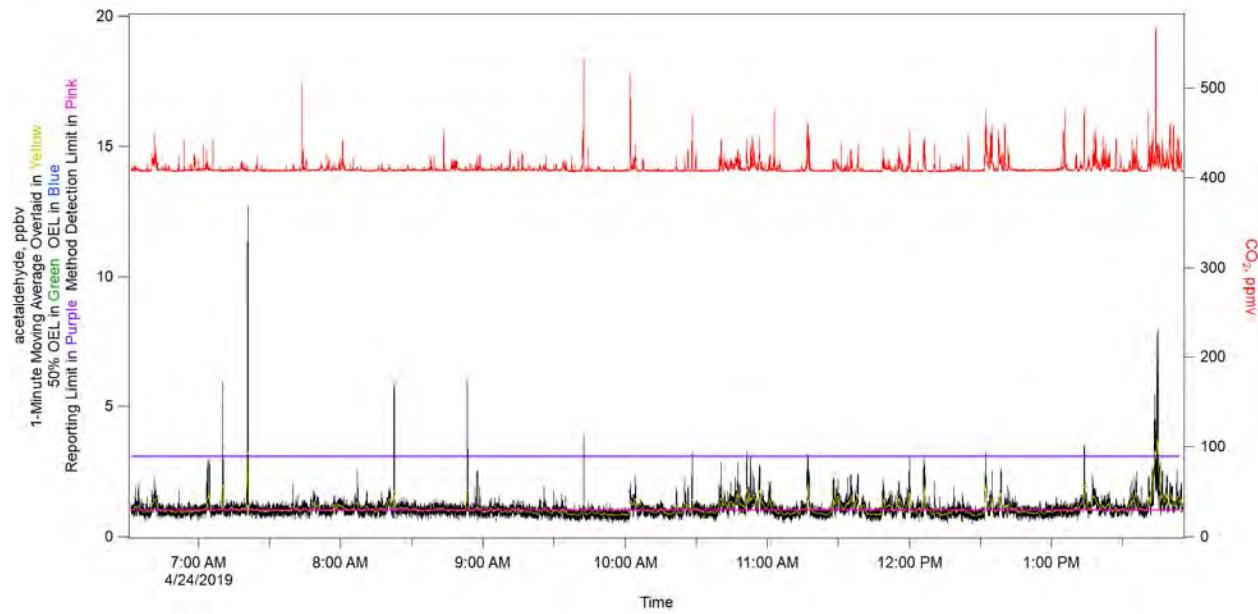


Figure 4-7. Acetaldehyde.

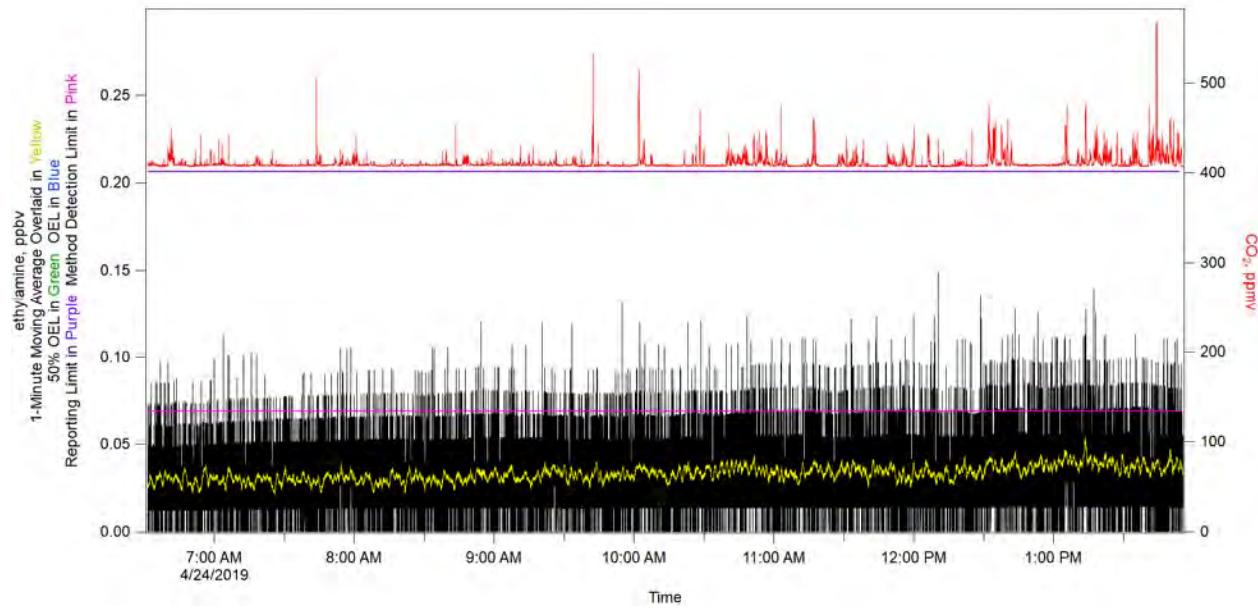


Figure 4-8. Ethylamine.

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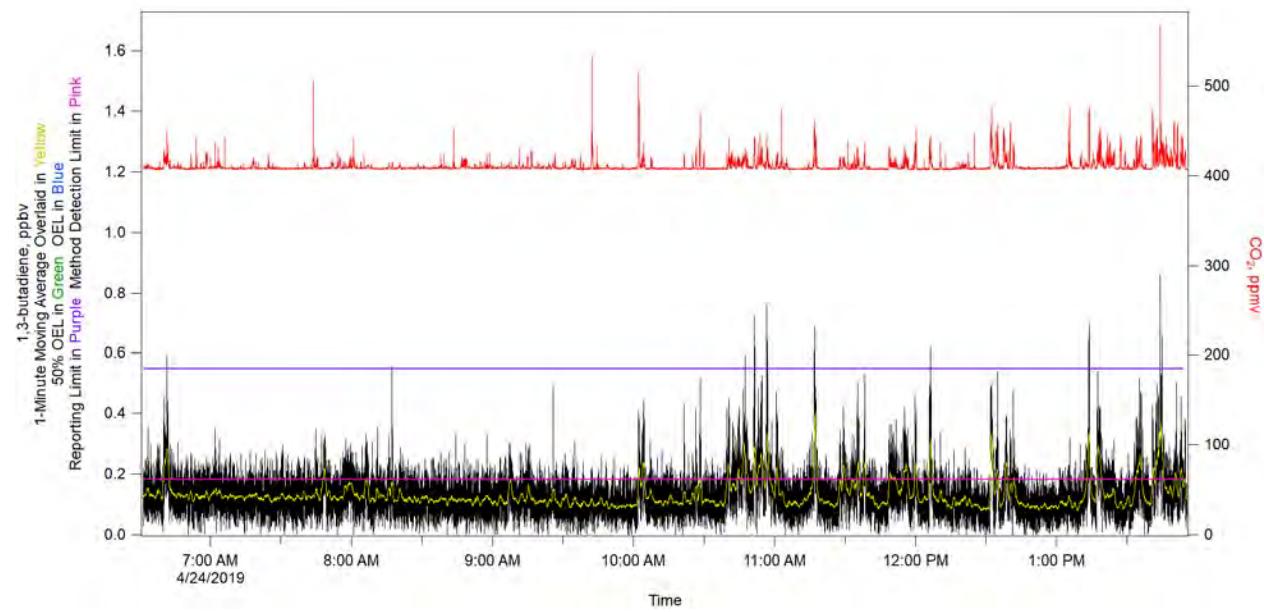


Figure 4-9. 1,3-butadiene.

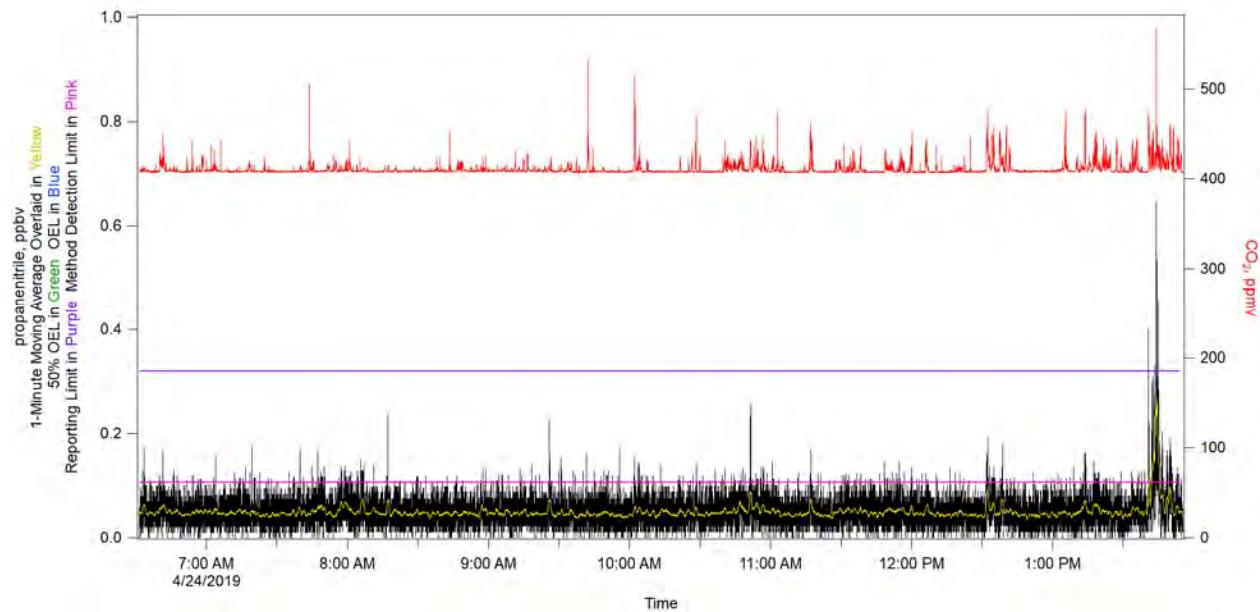


Figure 4-10. Propanenitrile.

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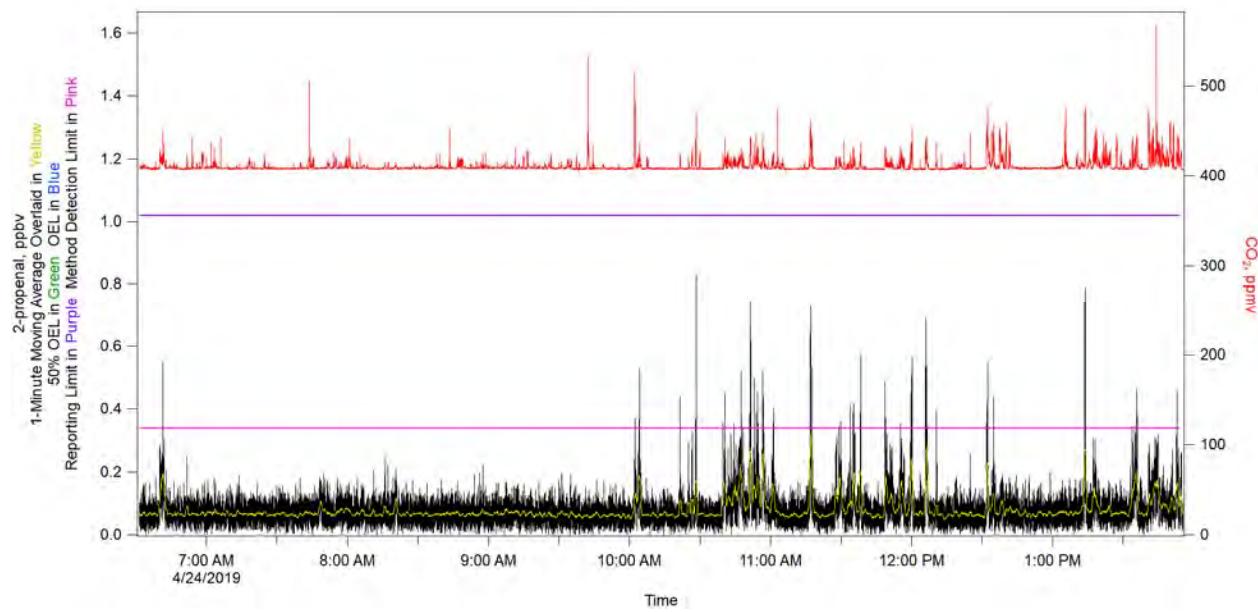


Figure 4-11. 2-propenal.

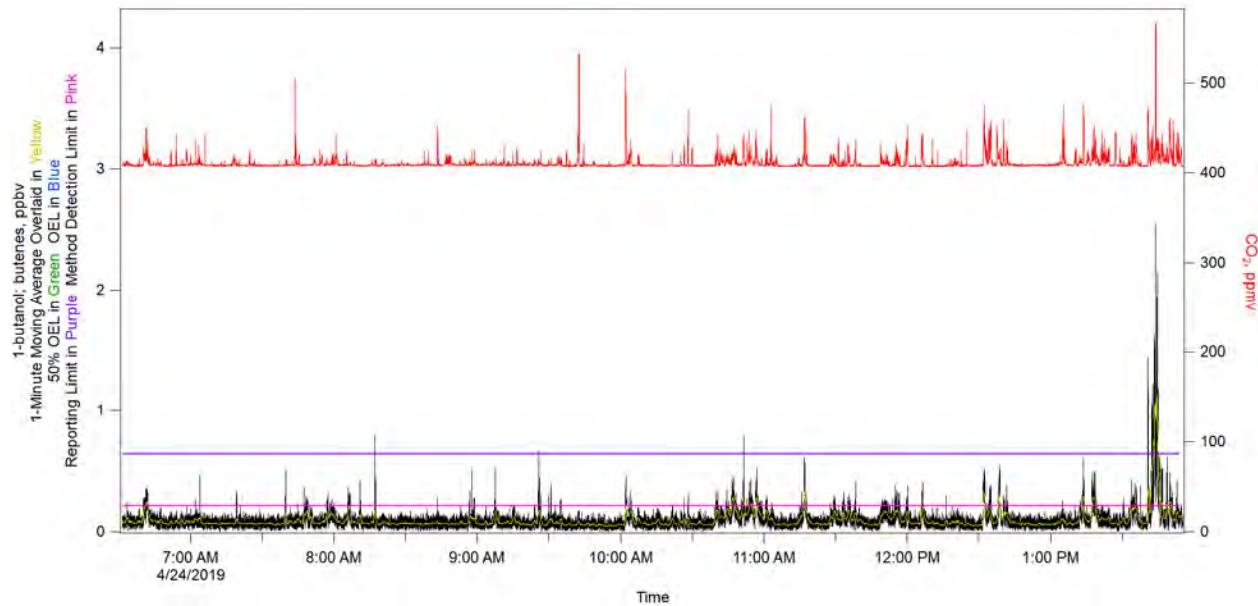


Figure 4-12. 1-butanol; Butenes.

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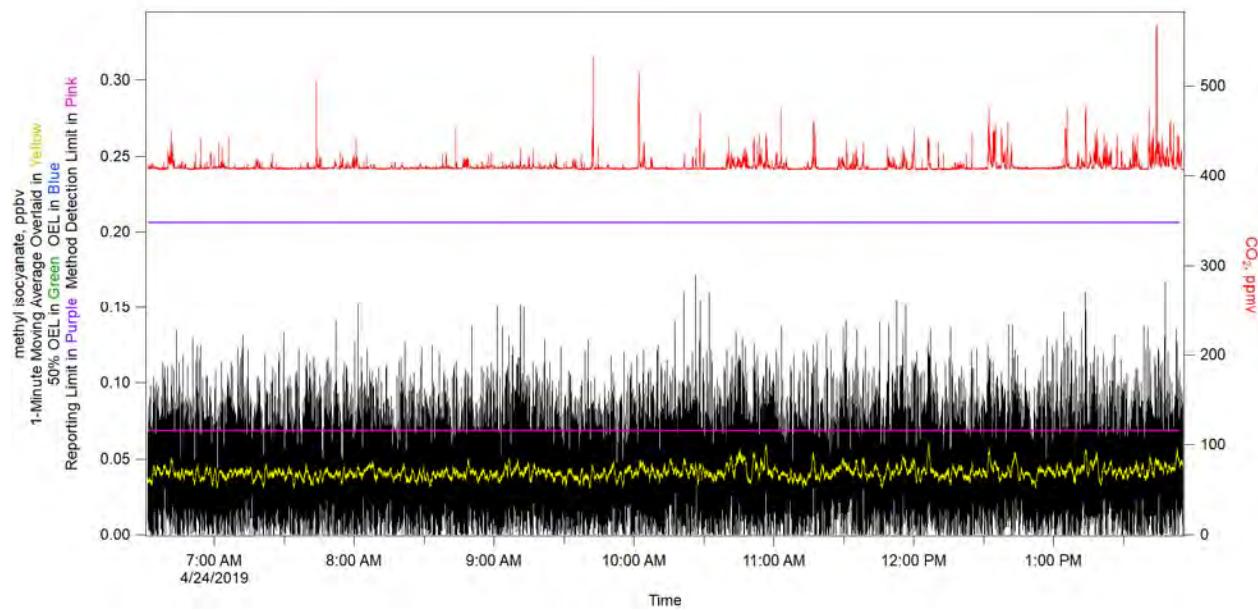


Figure 4-13. Methyl Isocyanate.

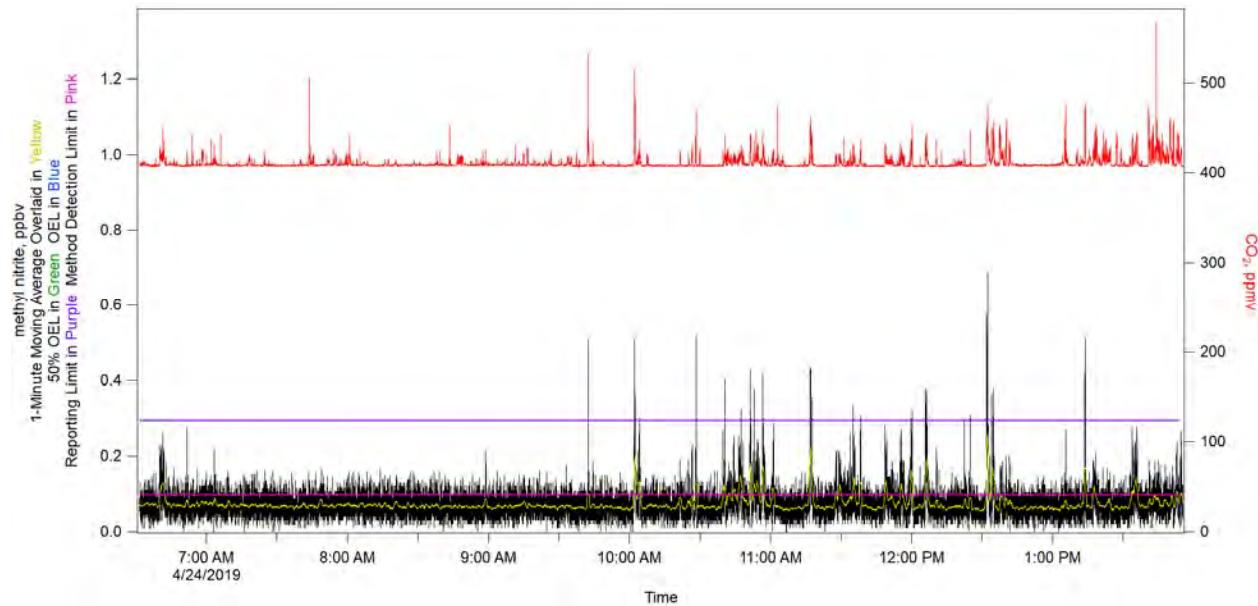


Figure 4-14. Methyl Nitrite.

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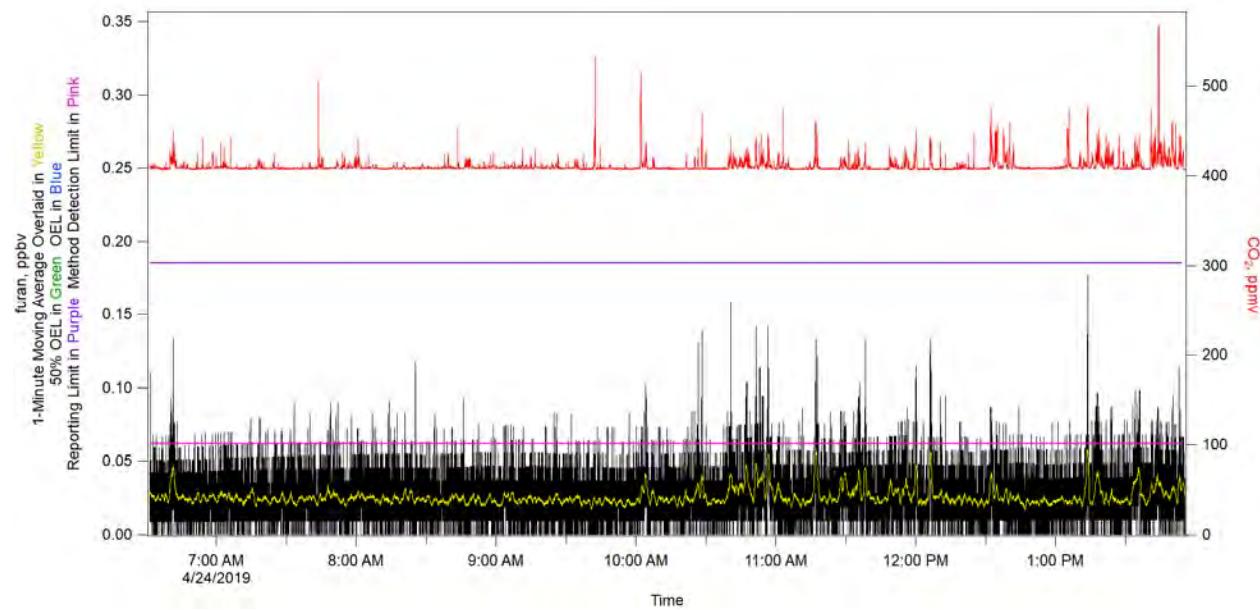


Figure 4-15. Furan.

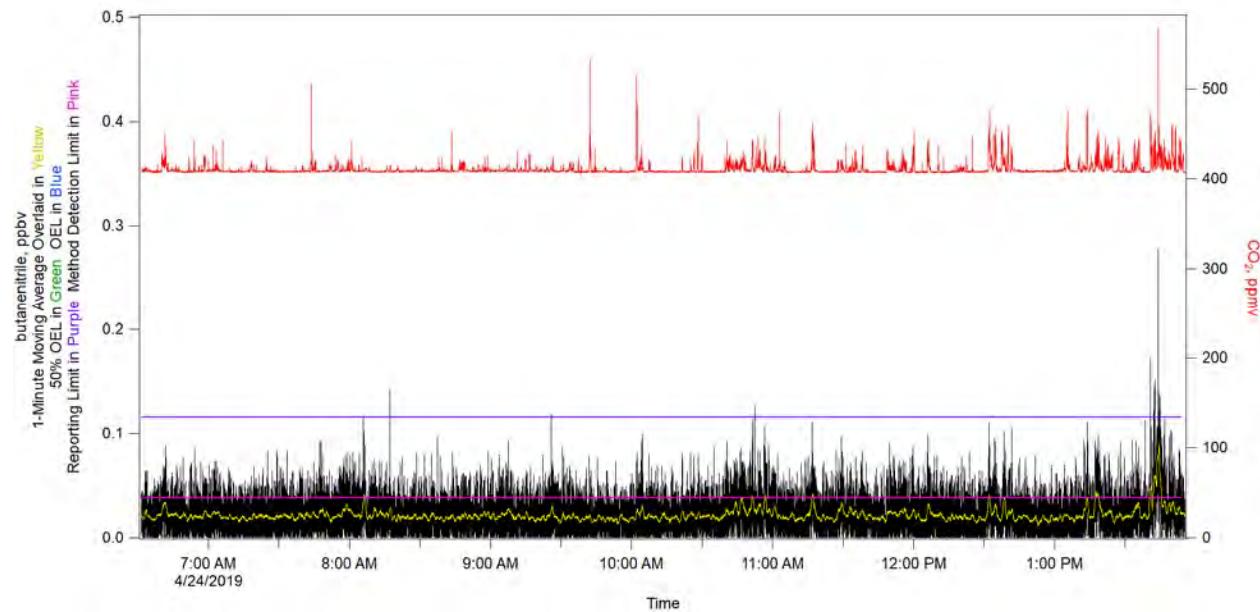


Figure 4-16. Butanenitrile.

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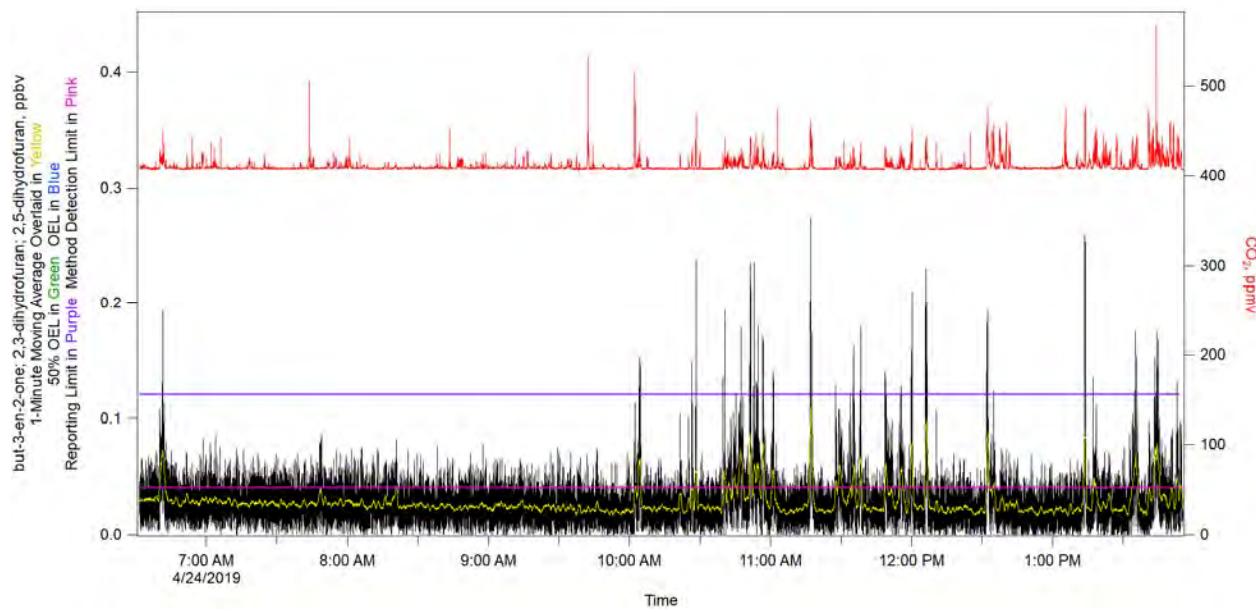


Figure 4-17. But-3-en-2-one; 2,3-dihydrofuran; 2,5-dihydrofuran.

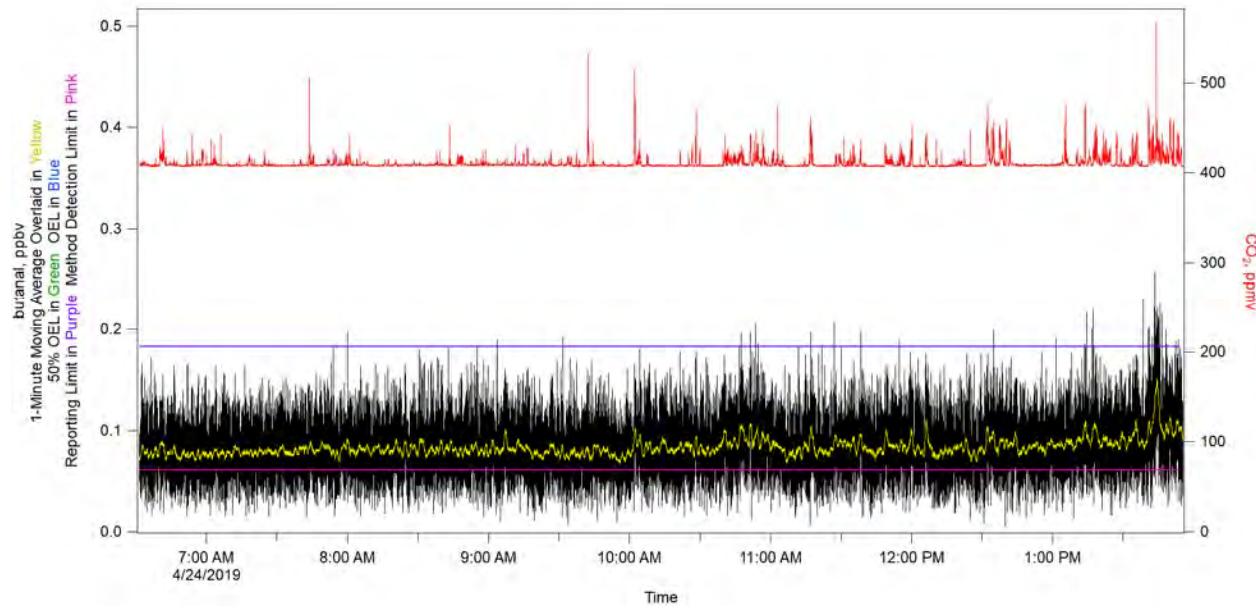


Figure 4-18. Butanal.

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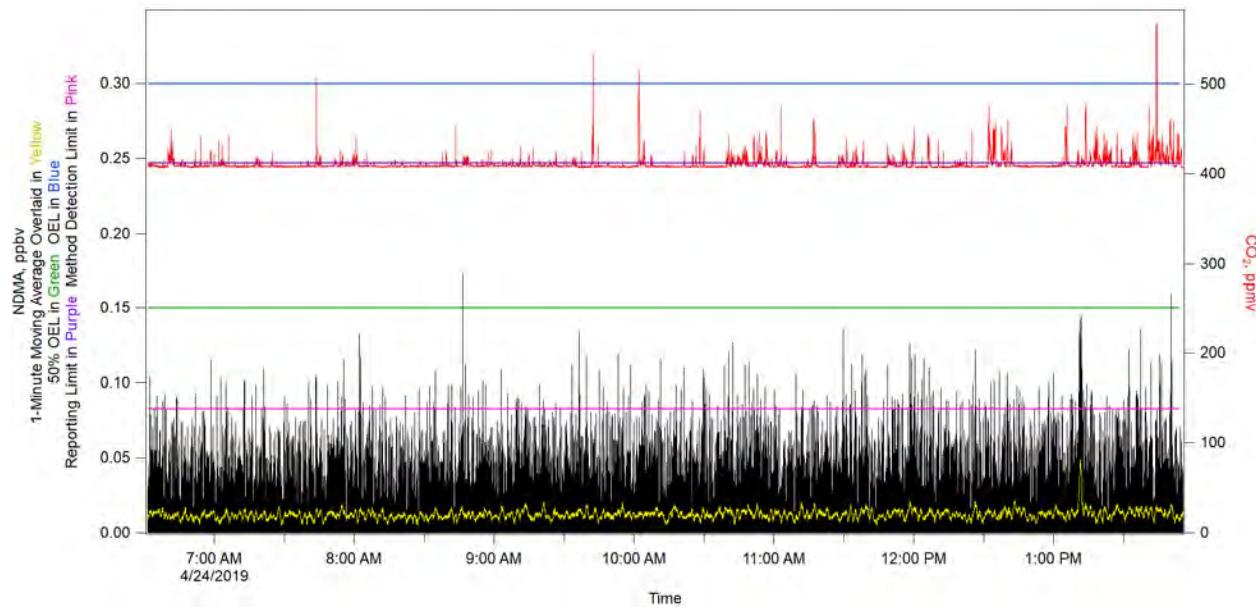


Figure 4-19. N-nitrosodimethylamine (NDMA).

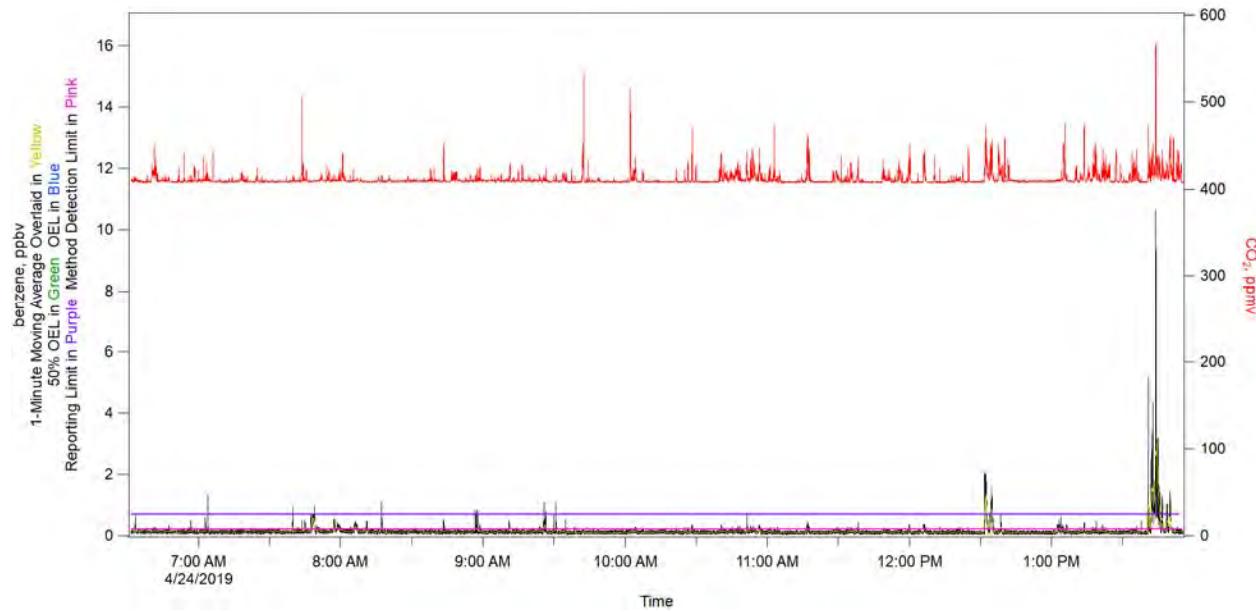


Figure 4-20. Benzene.

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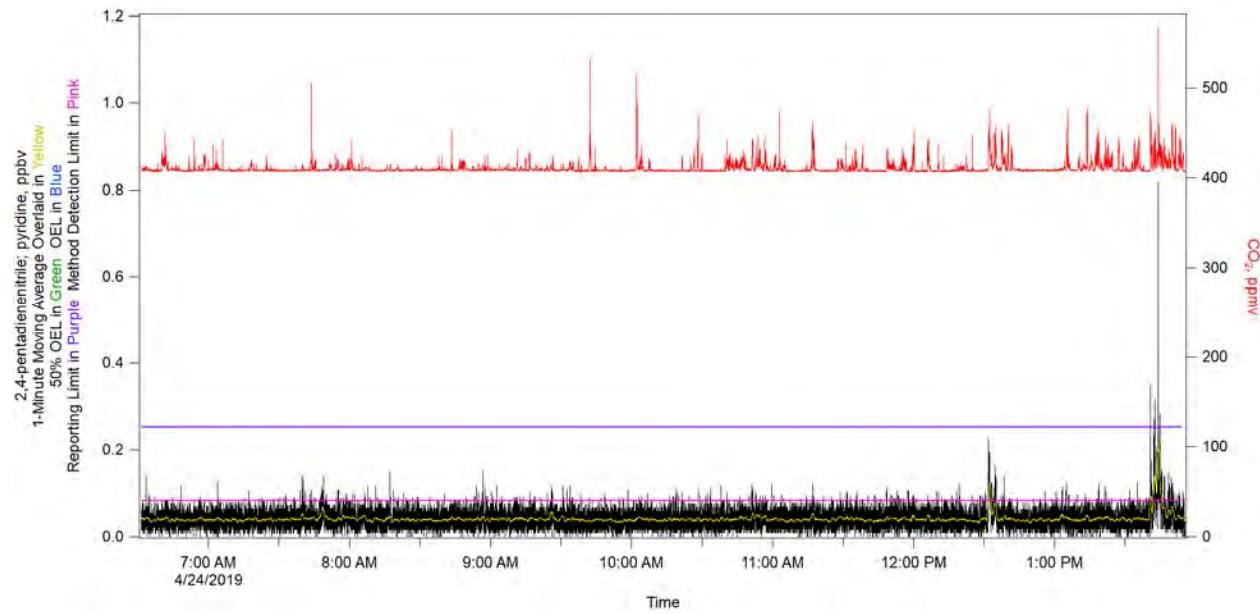


Figure 4-21. 2,4-pentadienenitrile; Pyridine.

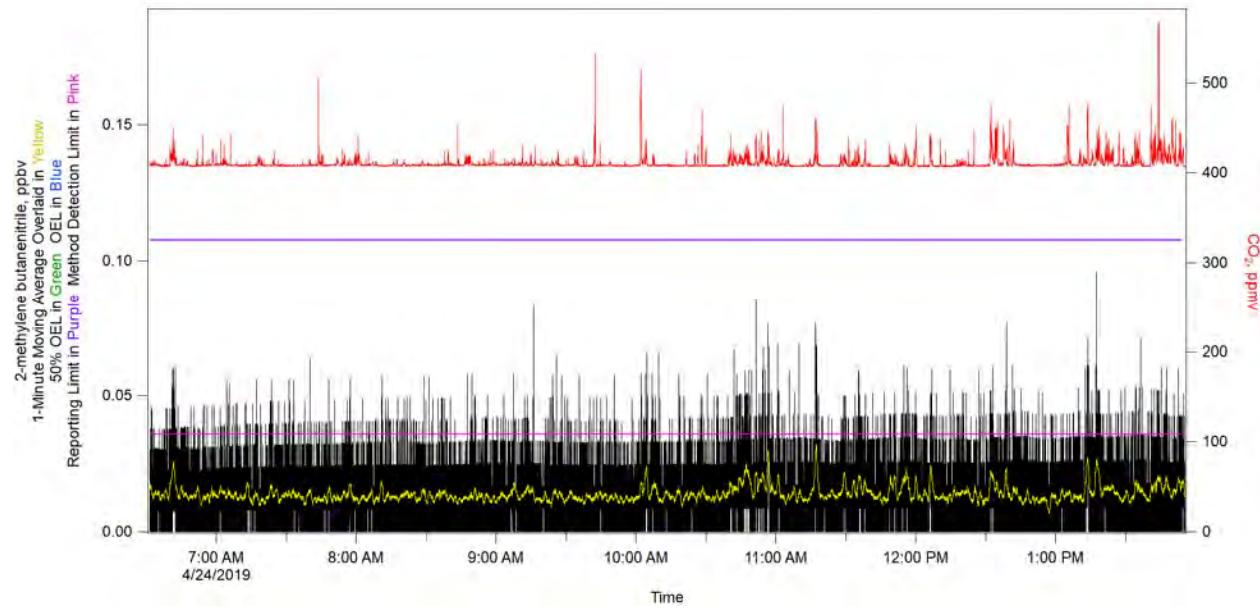


Figure 4-22. 2-methylene Butanenitrile.

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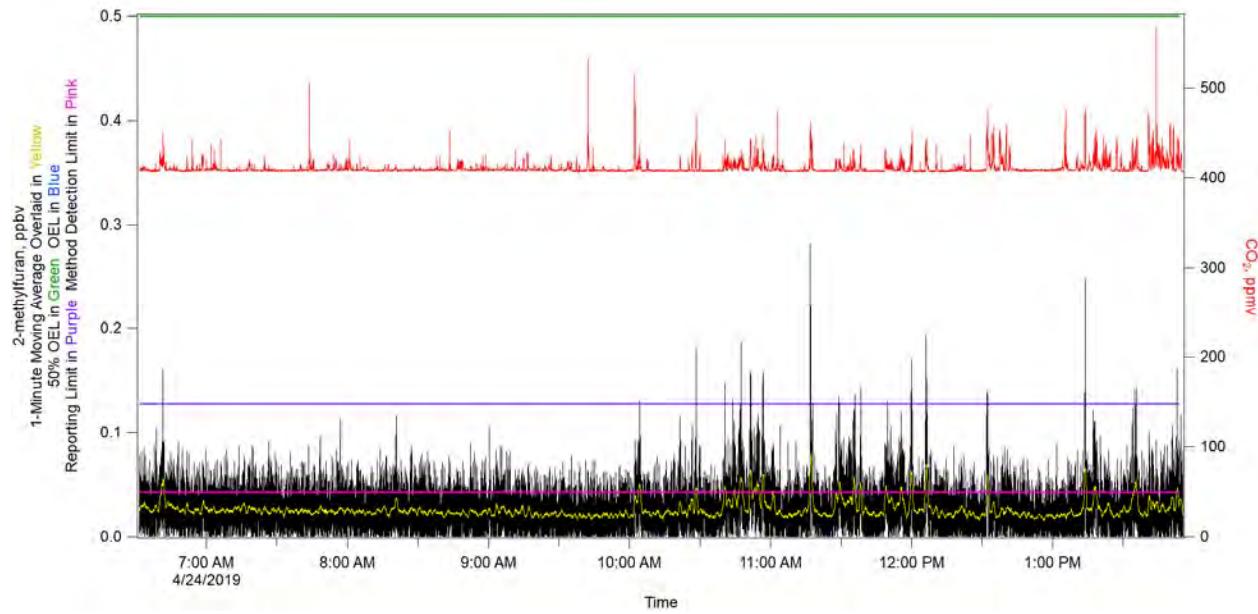


Figure 4-23. 2-methylfuran.

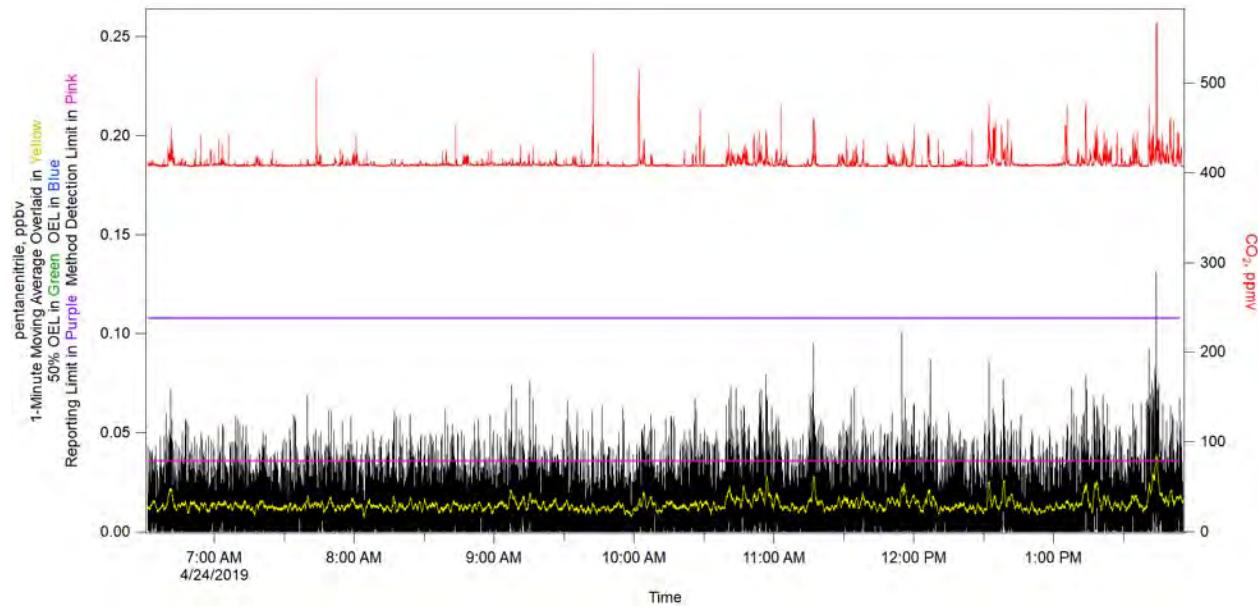


Figure 4-24. Pentanenitrile.

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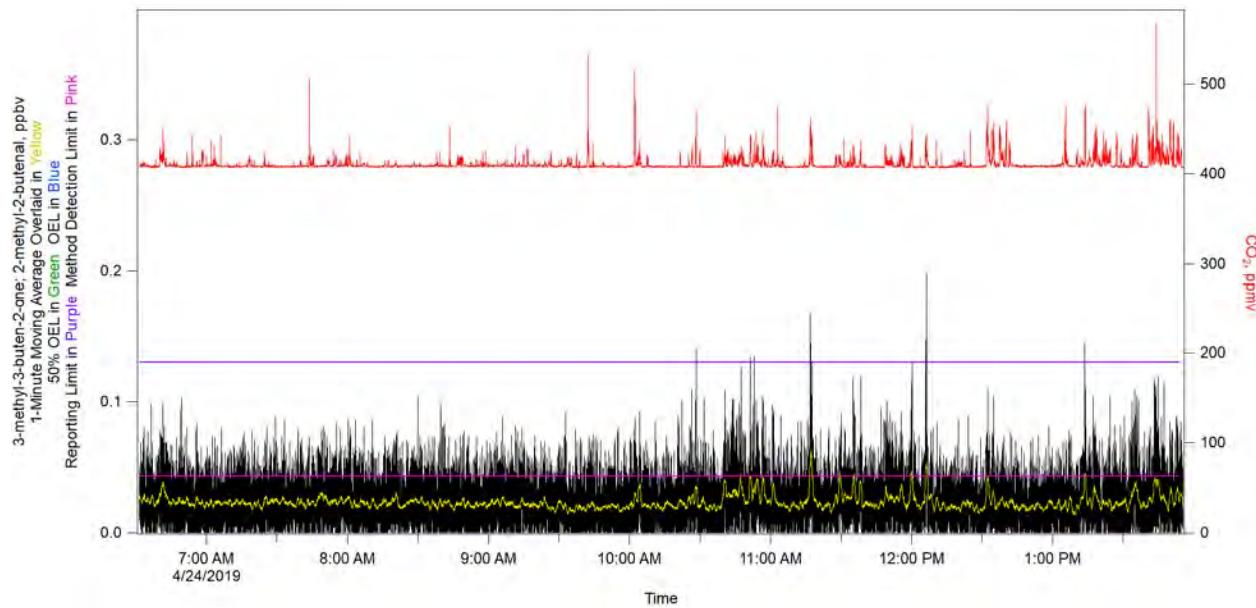


Figure 4-25. 3-methyl-3-buten-2-one; 2-methyl-2-butenal.

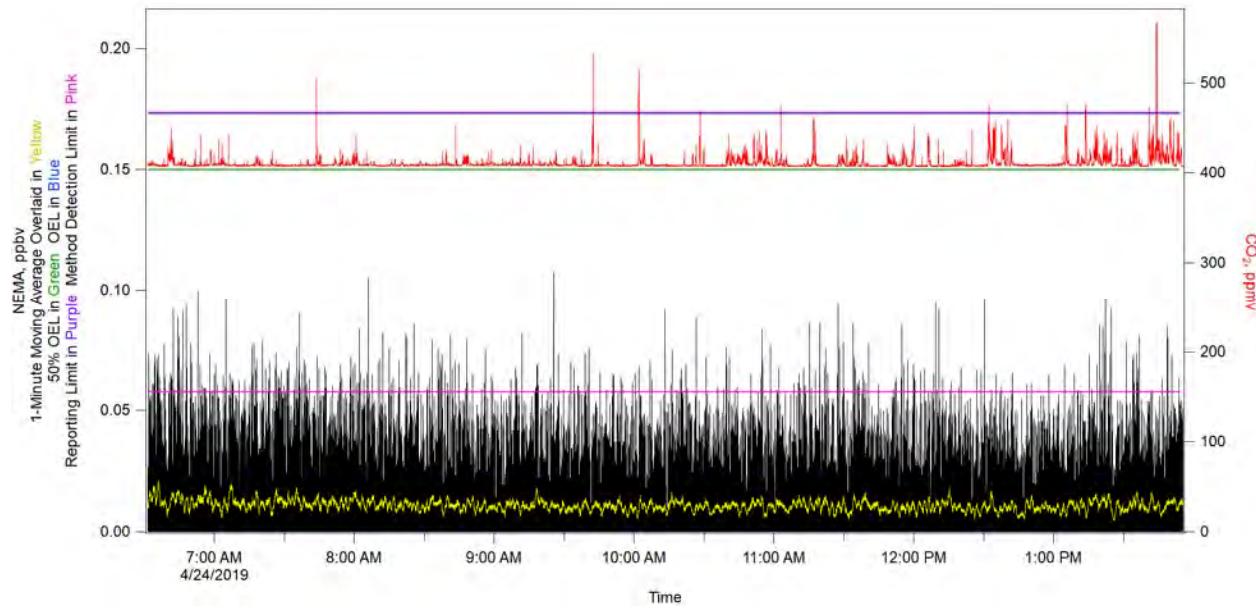


Figure 4-26. N-nitrosomethylethylamine (NEMA).

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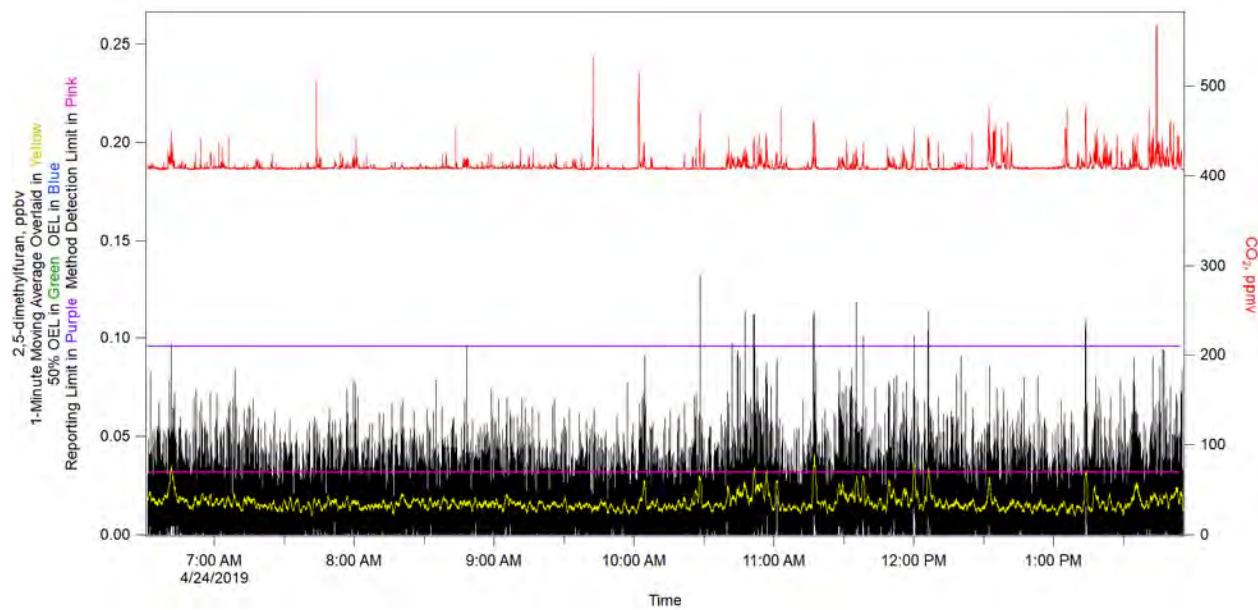


Figure 4-27. 2,5-dimethylfuran.

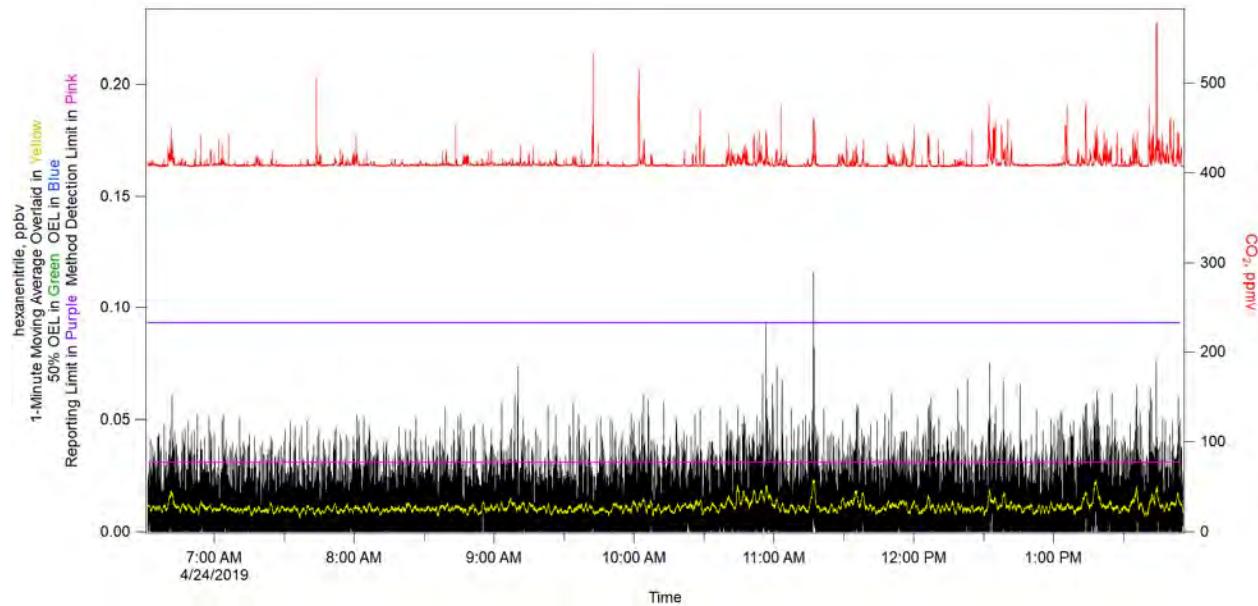


Figure 4-28. Hexanenitrile.

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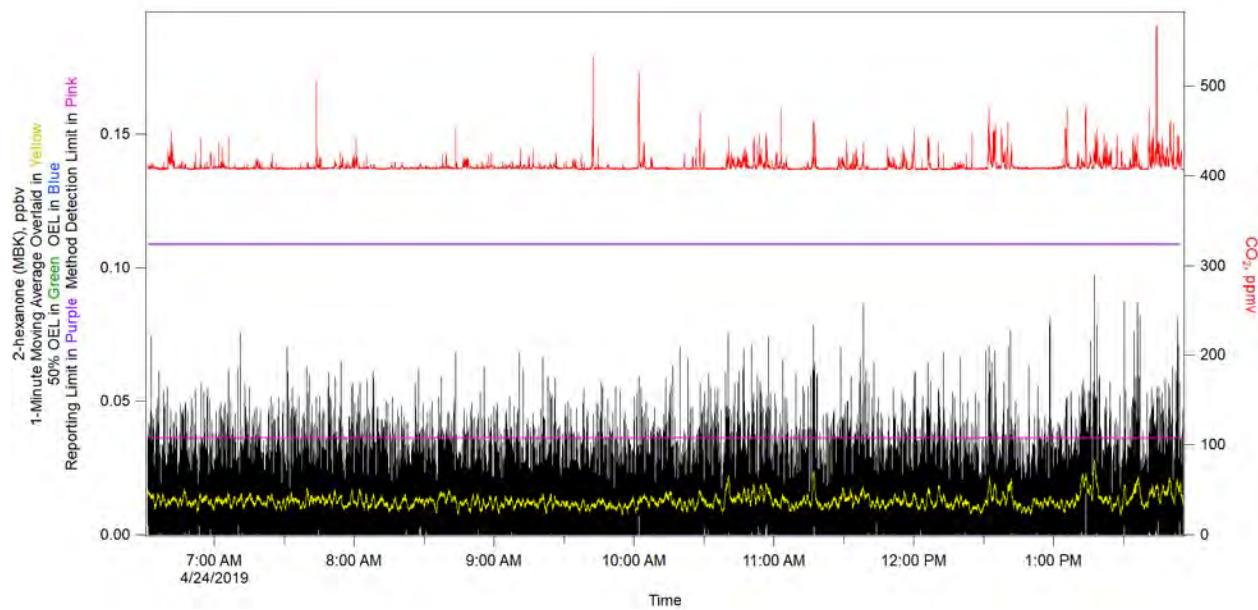


Figure 4-29. 2-hexanone (MBK).

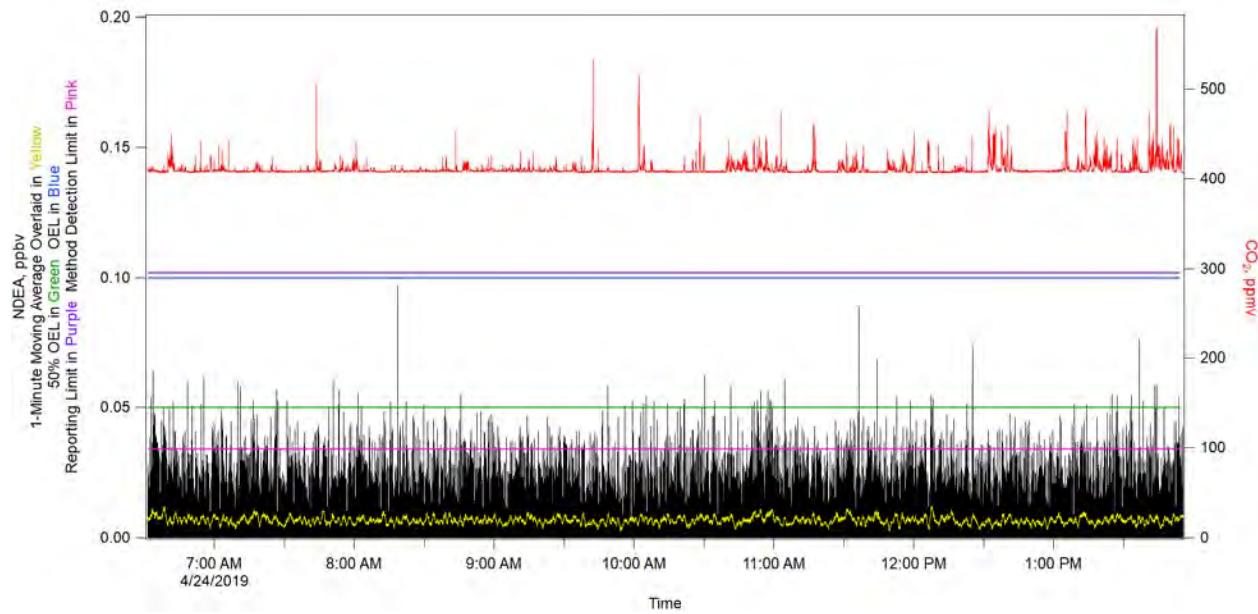


Figure 4-30. N-nitrosodiethylamine (NDEA).

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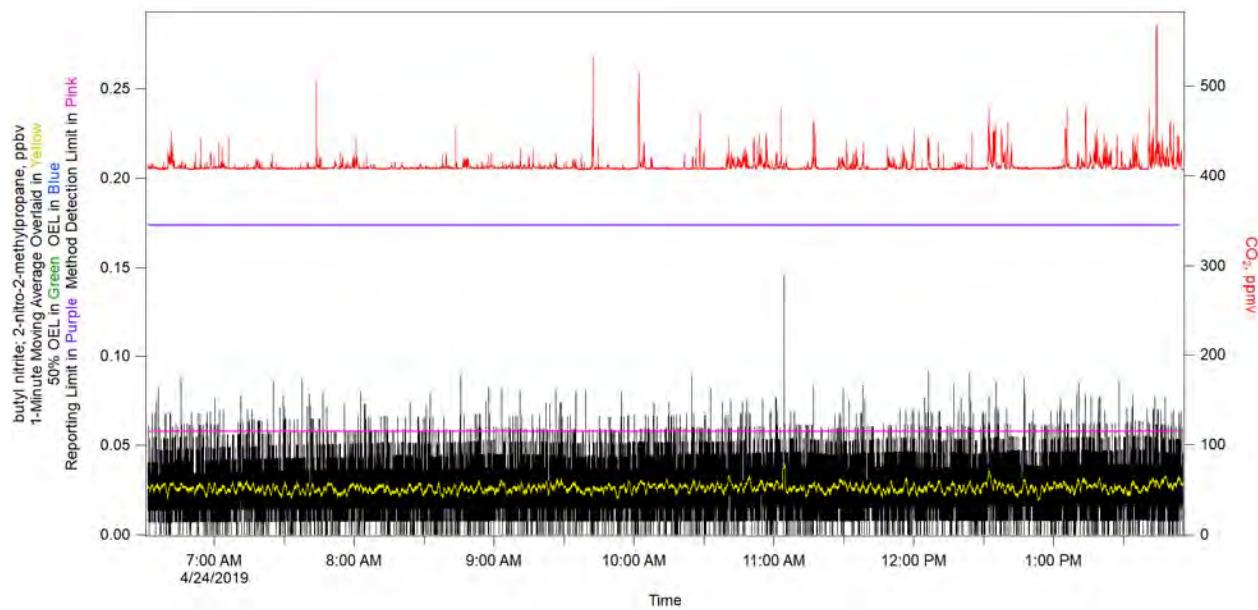


Figure 4-31. Butyl Nitrite; 2-nitro-2-methylpropane.

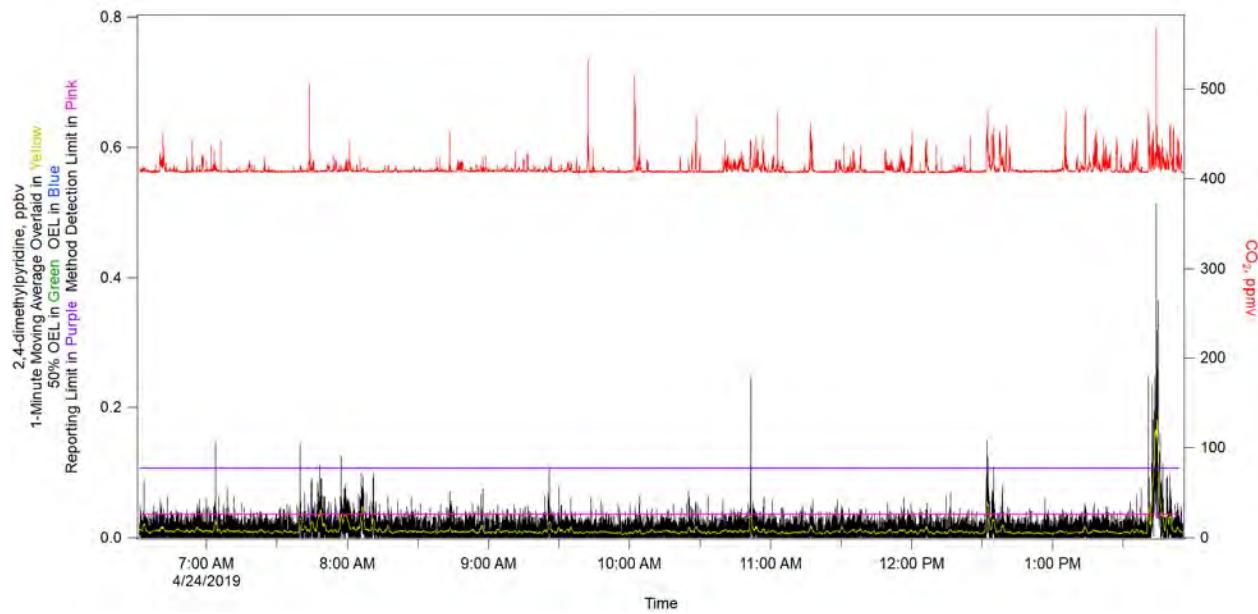


Figure 4-32. 2,4-dimethylpyridine.

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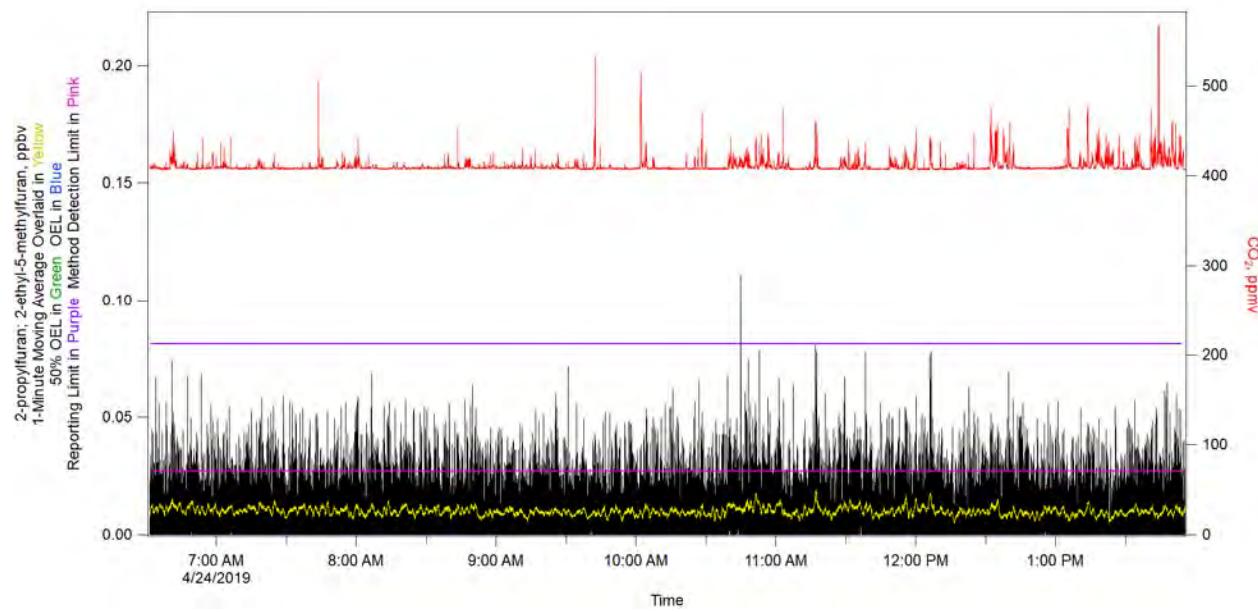


Figure 4-33. 2-propylfuran; 2-ethyl-5-methylfuran.

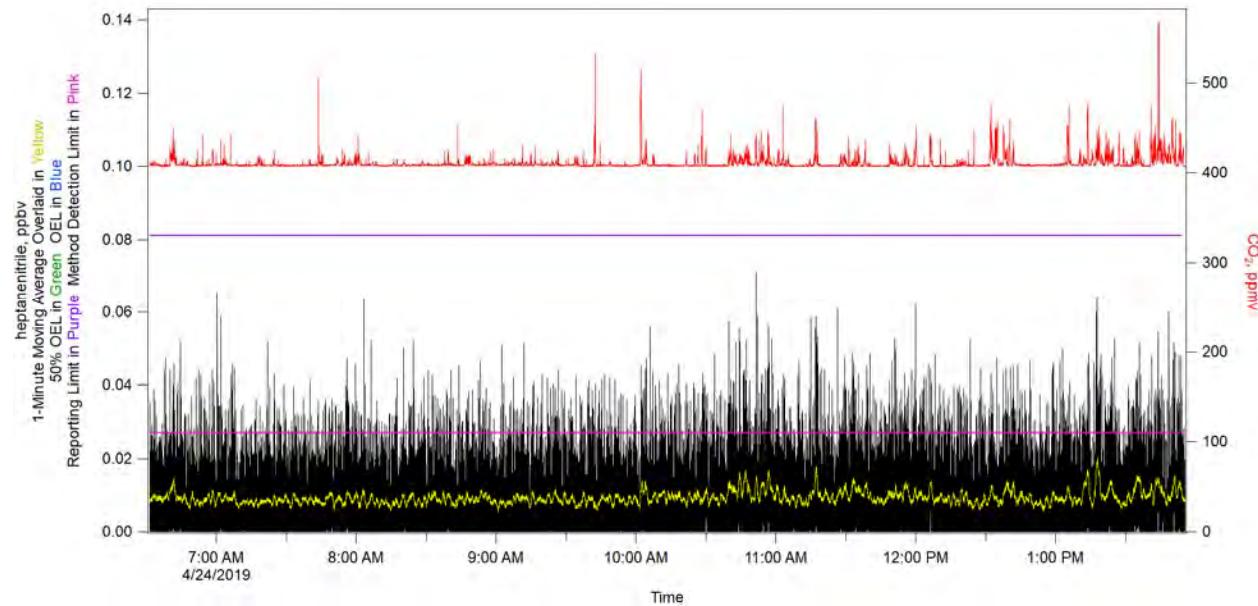


Figure 4-34. Heptanenitrile.

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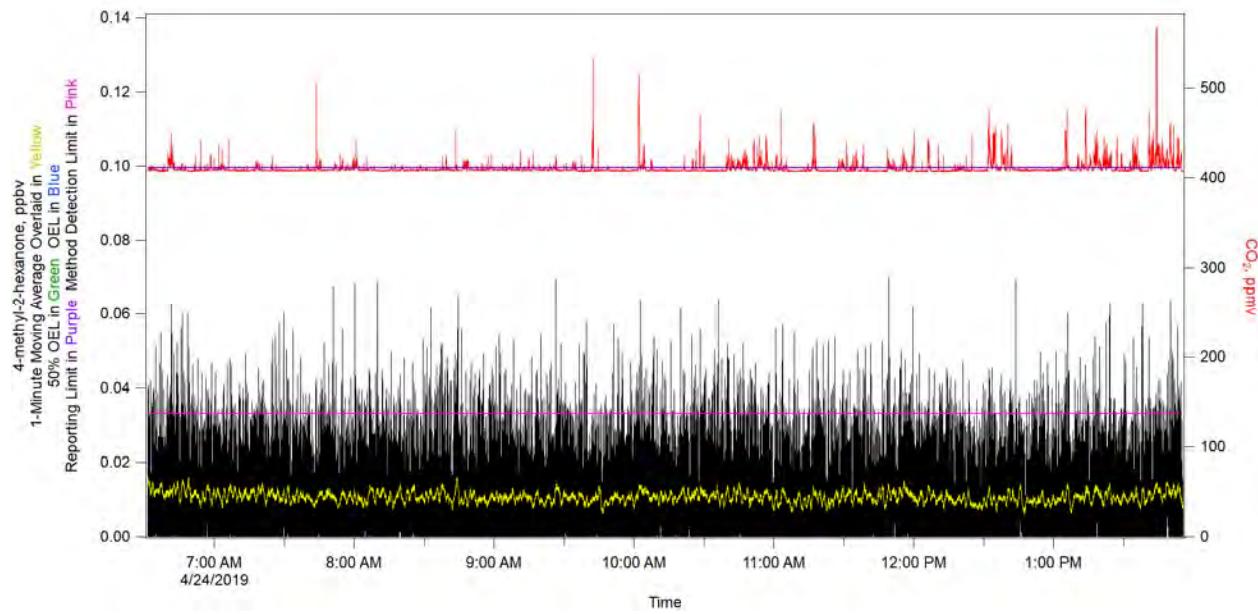


Figure 4-35. 4-methyl-2-hexanone.

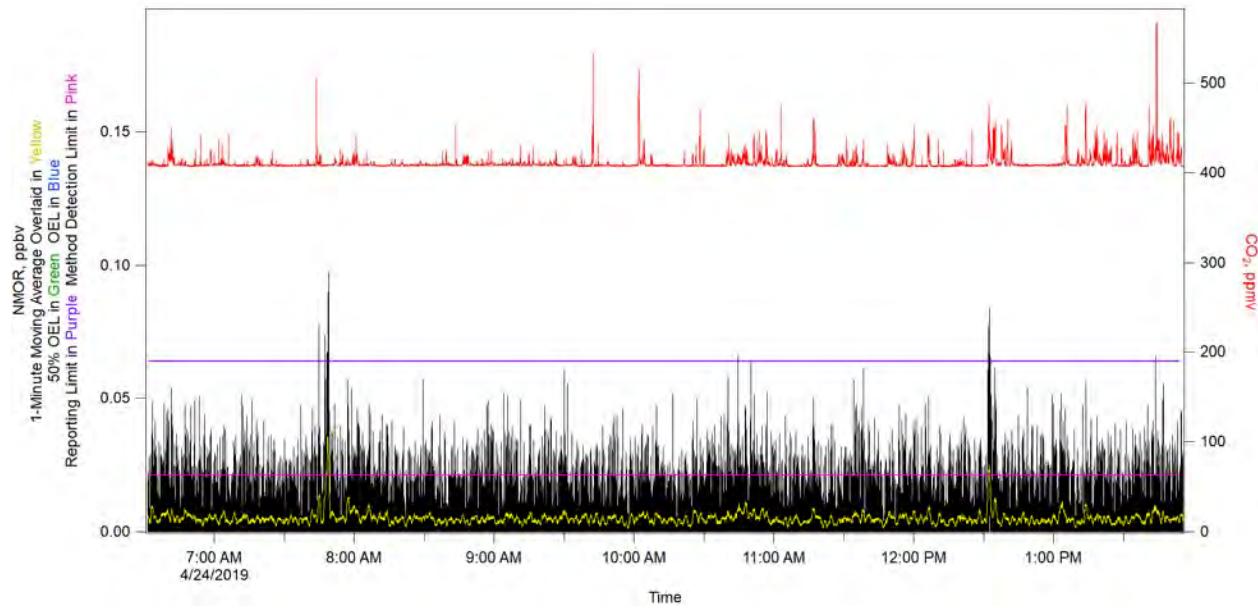


Figure 4-36. N-nitrosomorpholine (NMOR).

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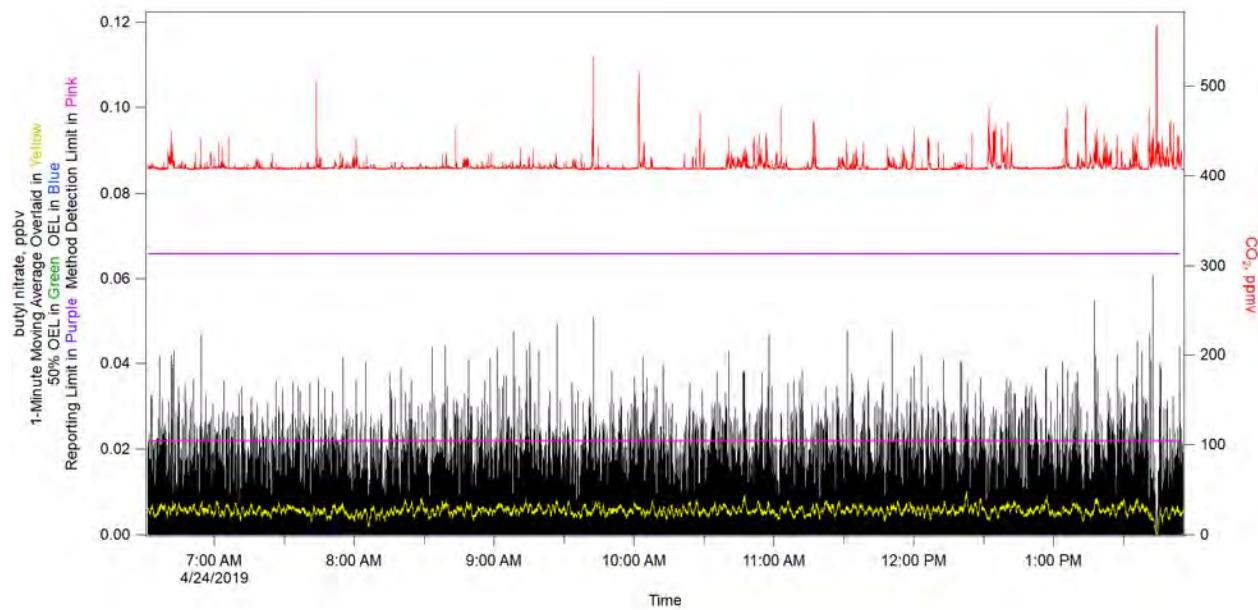
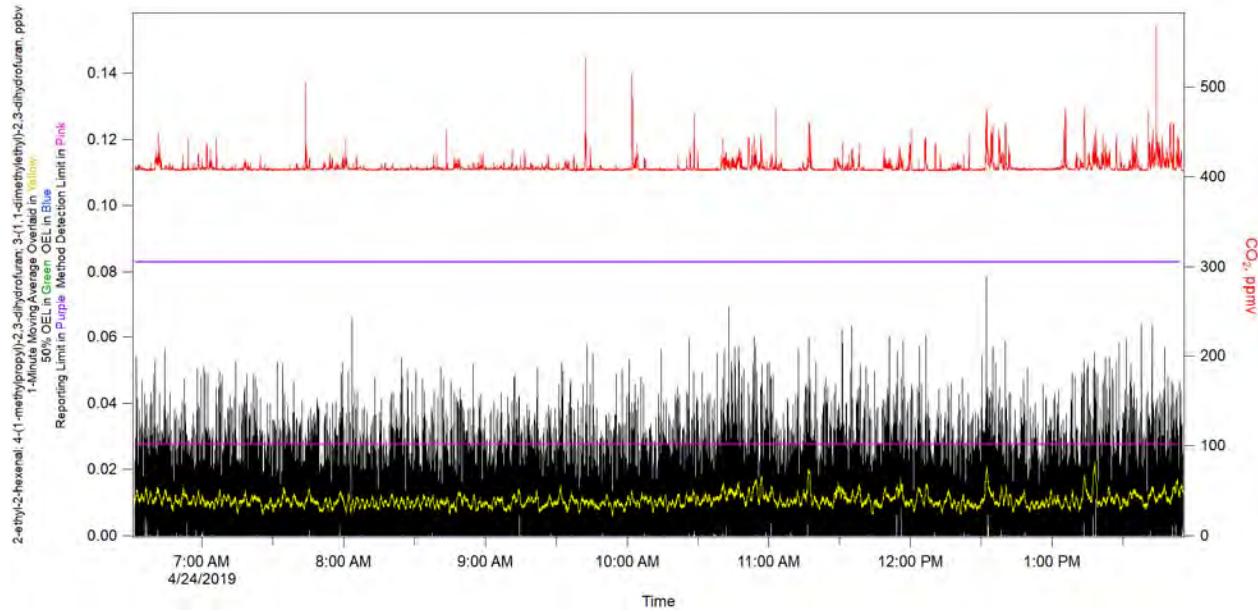


Figure 4-37. Butyl Nitrate.



**Figure 4-38. 2-ethyl-2-hexenal; 4-(1-methylpropyl)-2,3-dihydrofuran
3-(1,1-dimethylethyl)-2,3-dihydrofuran.**

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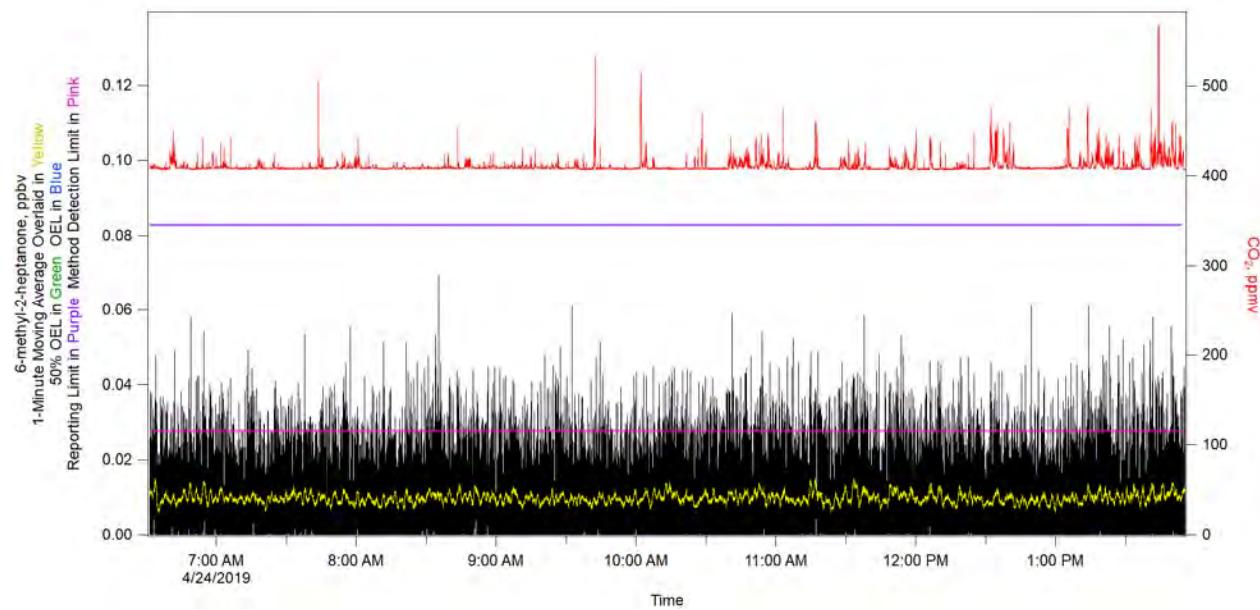


Figure 4-39. 6-methyl-2-heptanone.

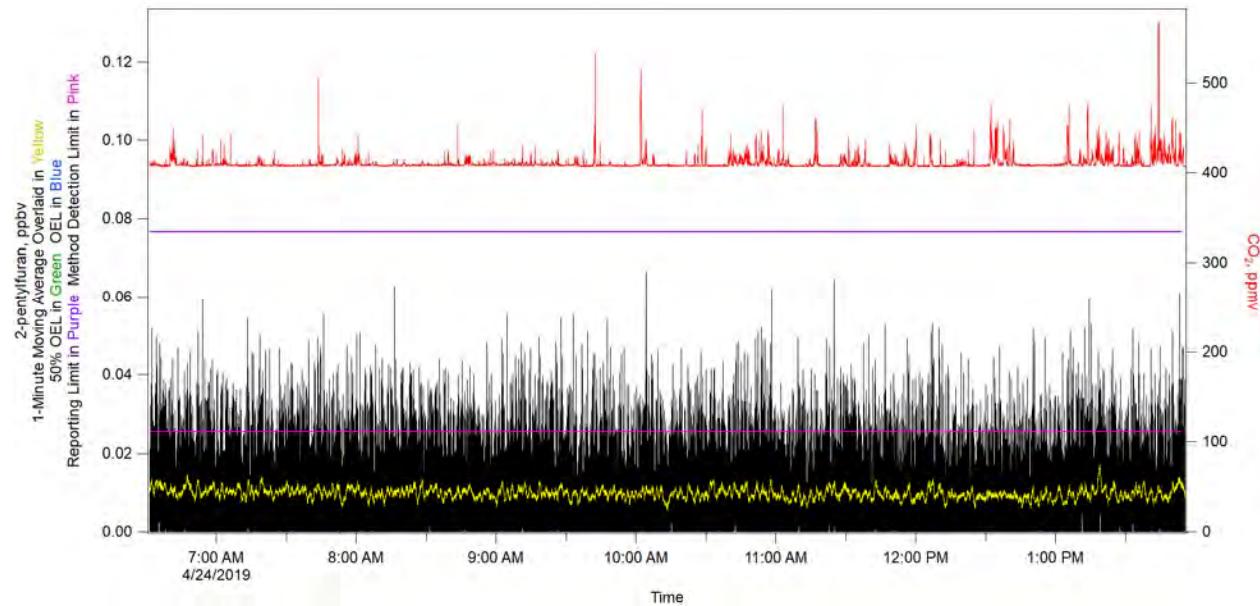


Figure 4-40. 2-pentylfuran.

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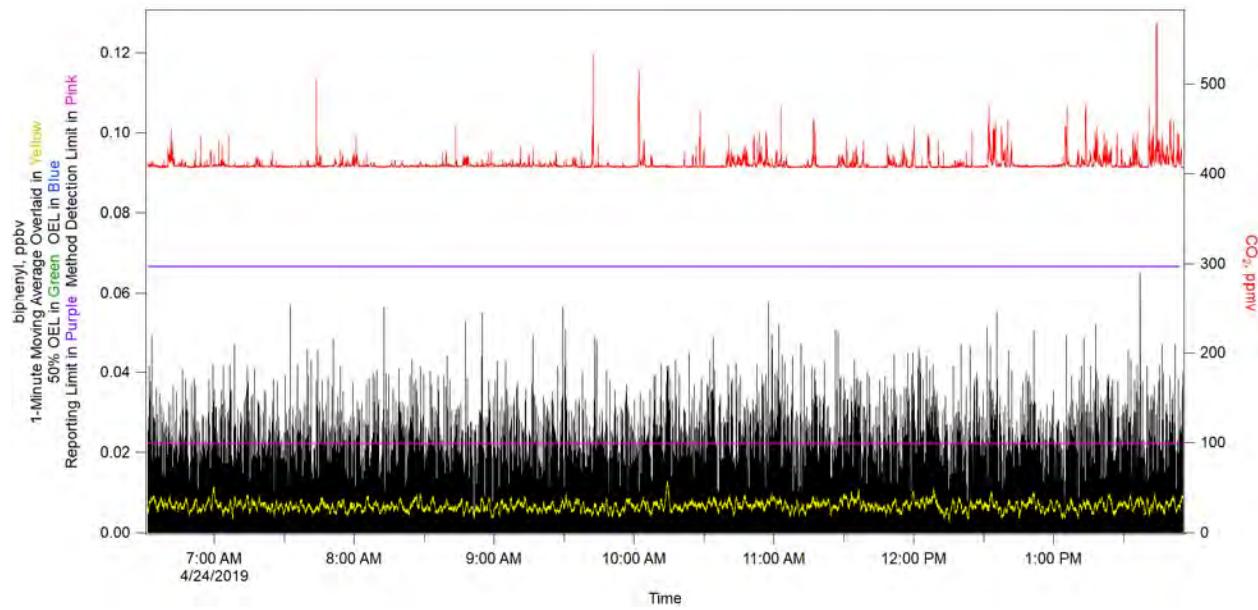


Figure 4-41. Biphenyl.

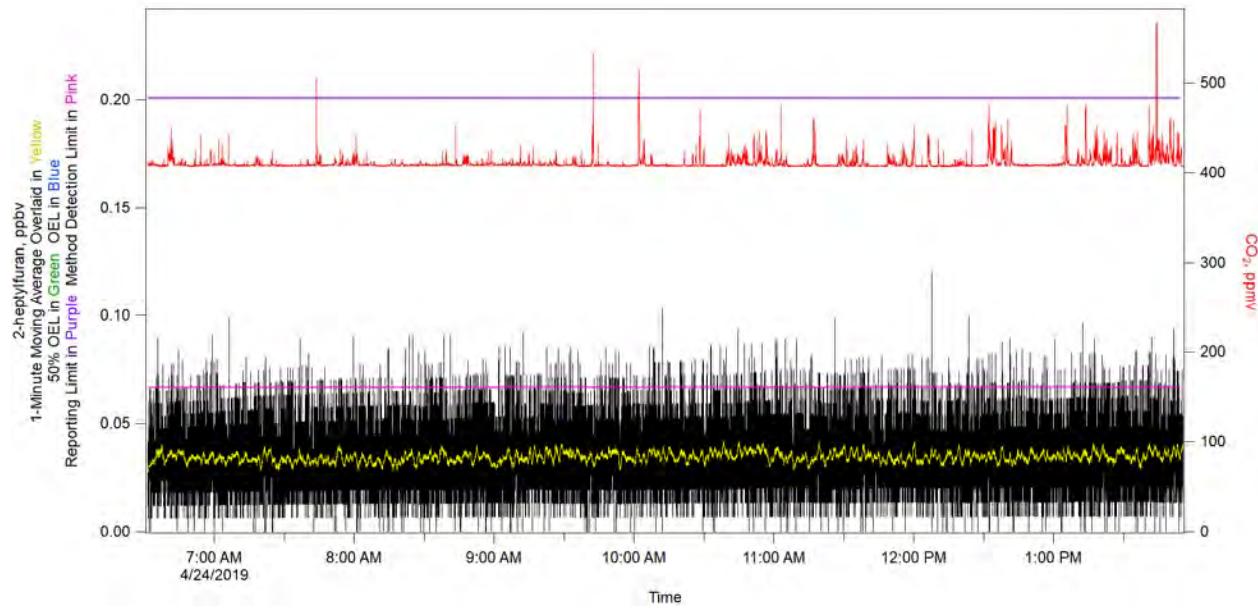


Figure 4-42. 2-heptylfuran.

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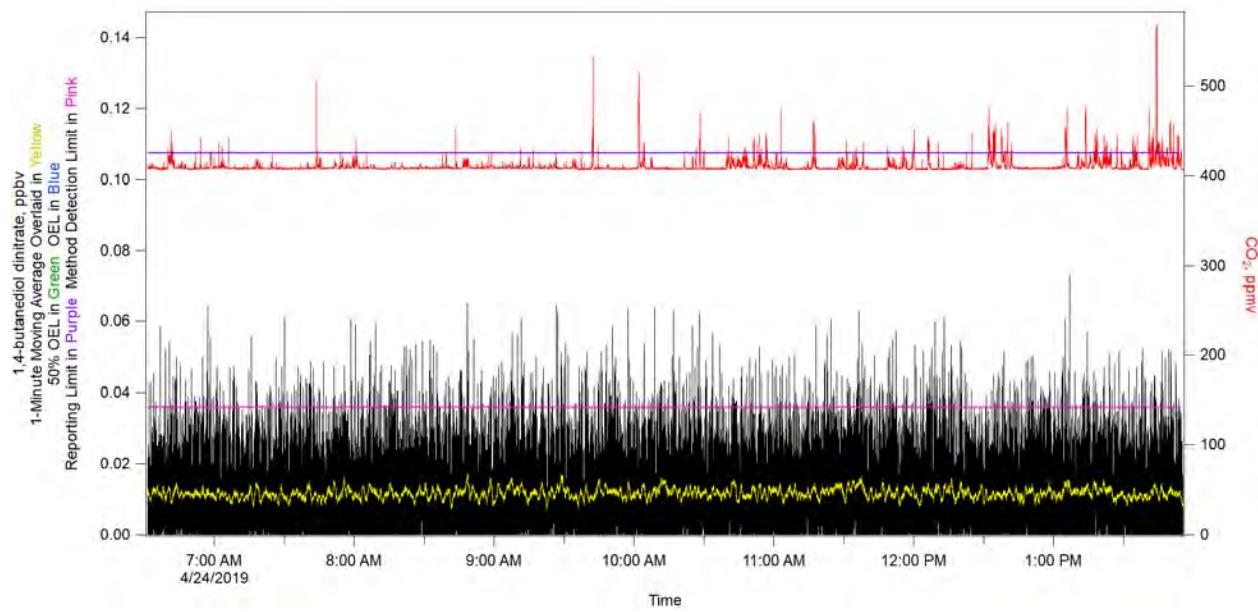


Figure 4-43. 1,4-butanediol Dinitrate.

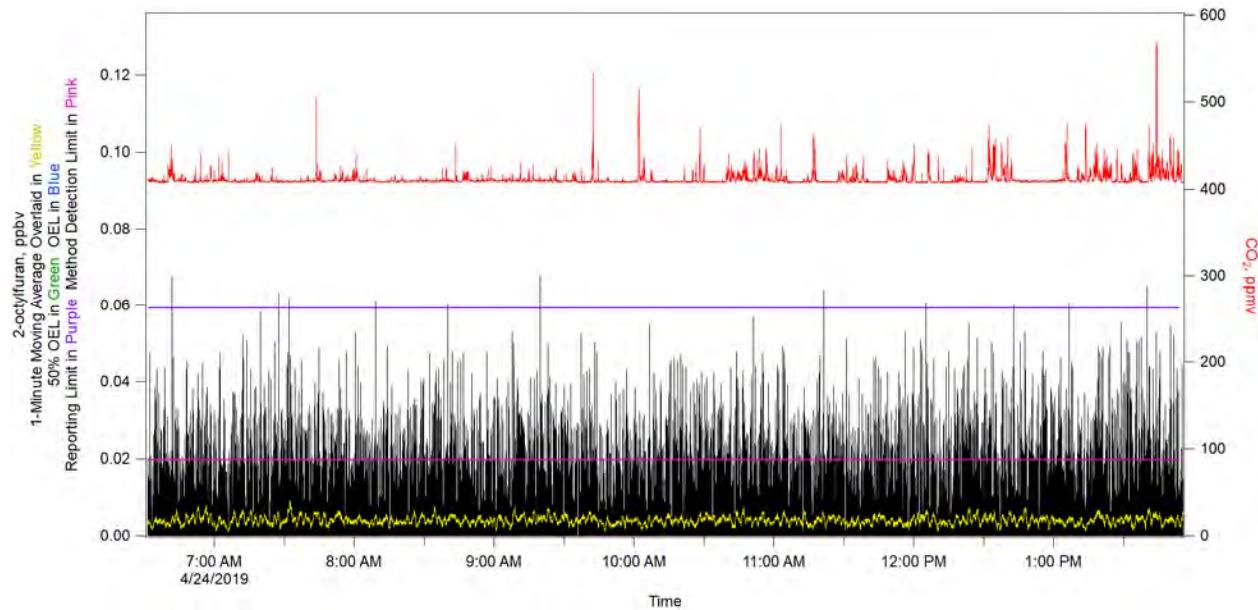


Figure 4-44. 2-octylfuran.

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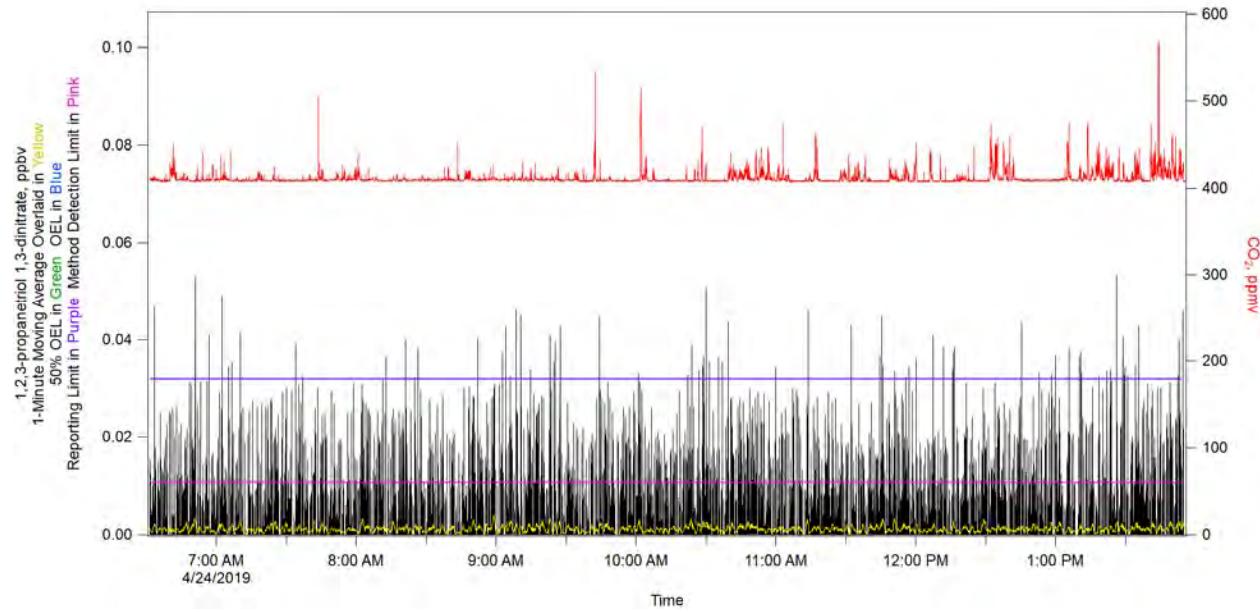


Figure 4-45. 1,2,3-propanetriol 1,3-dinitrate.

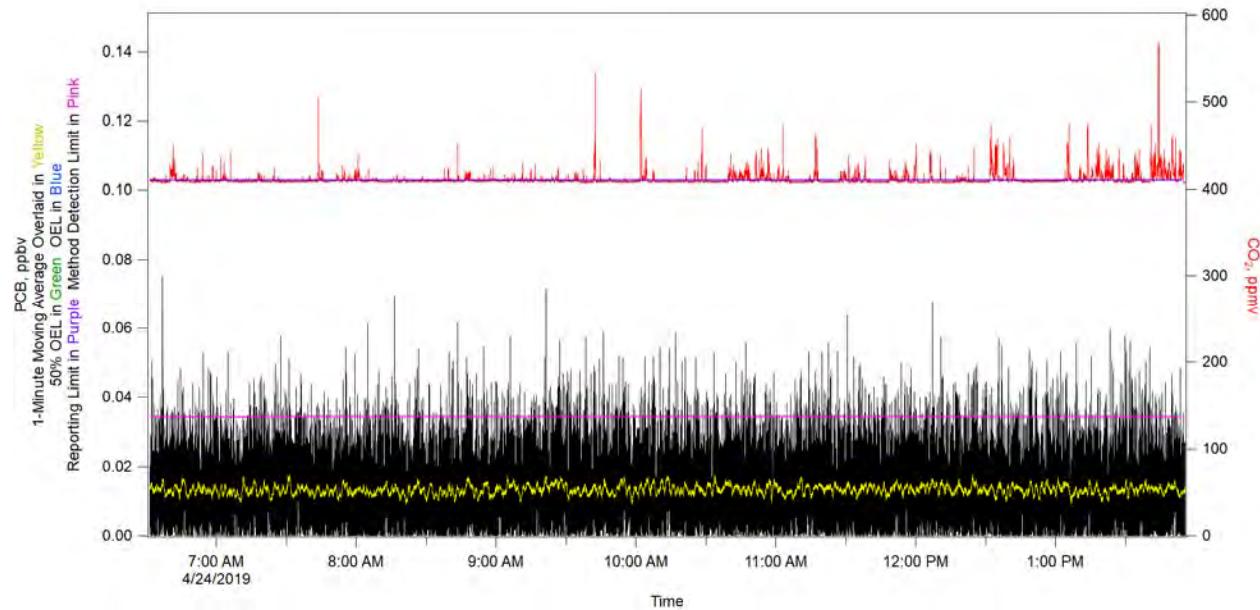


Figure 4-46. PCB.

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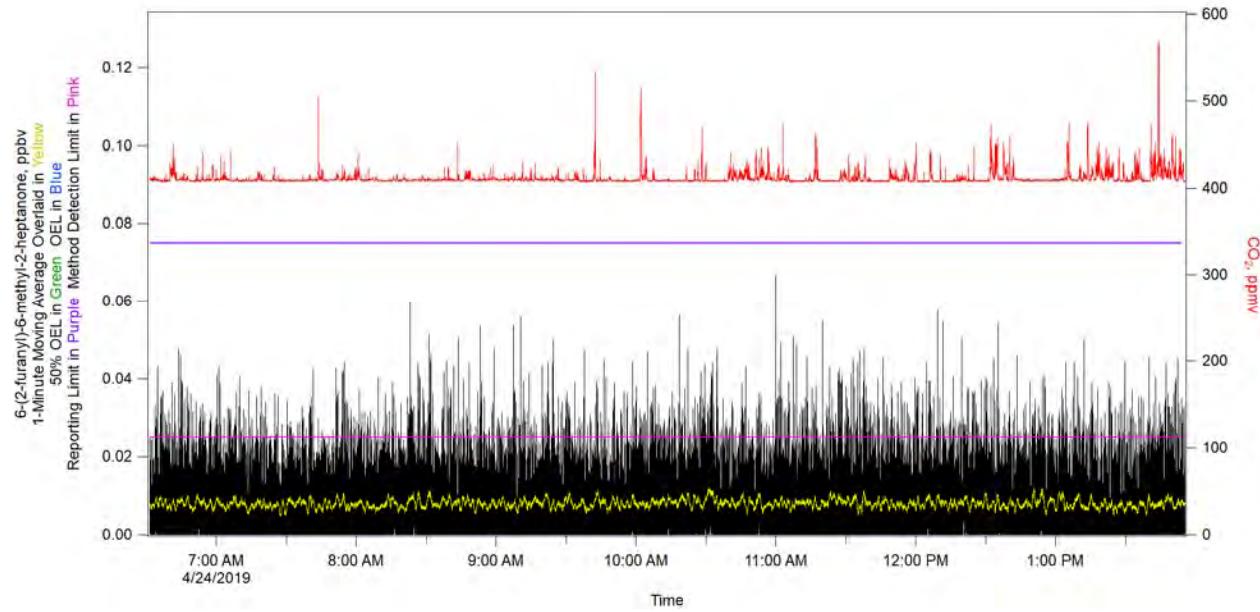


Figure 4-47. 6-(2-furanyl)-6-methyl-2-heptanone.

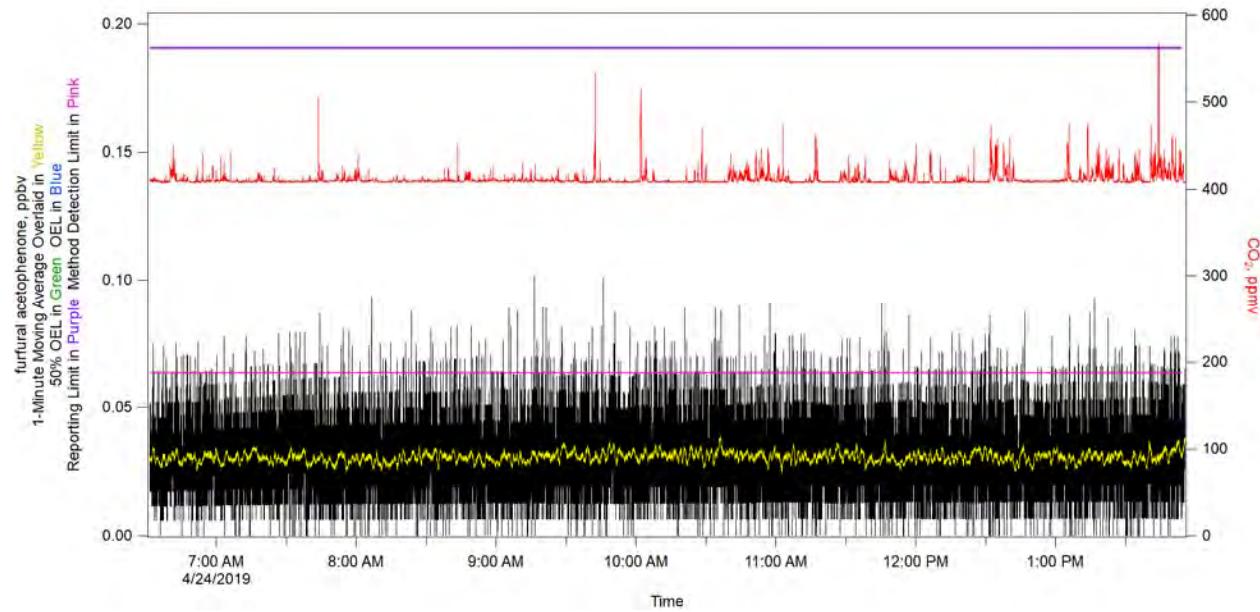


Figure 4-48. Furfural Acetophenone.

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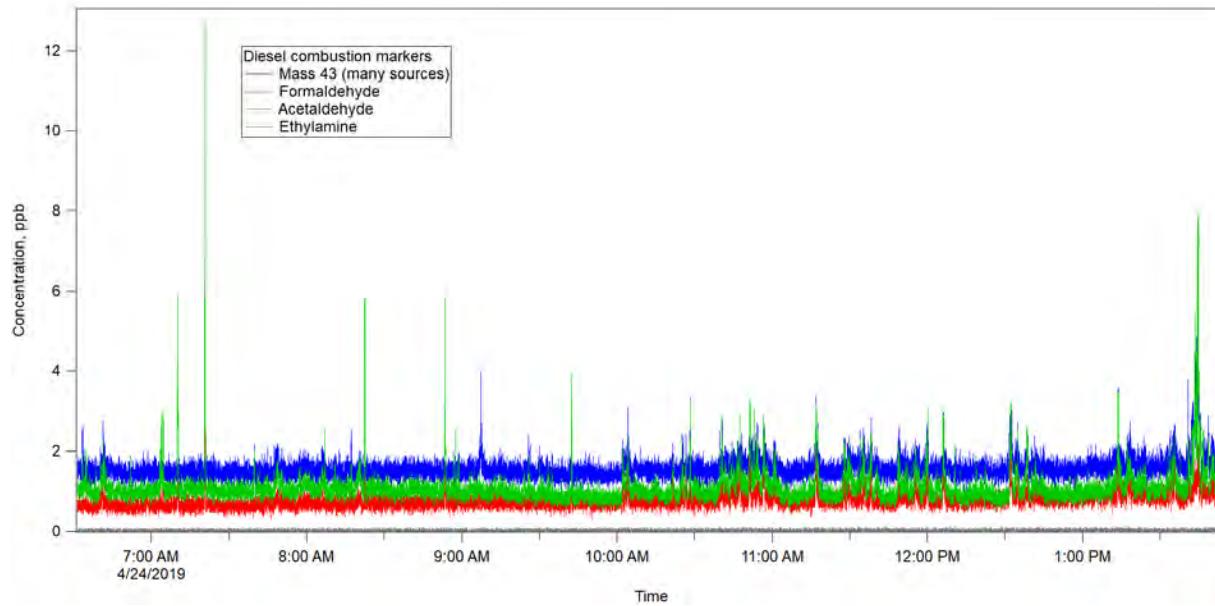


Figure 4-49. Diesel Combustion Markers.

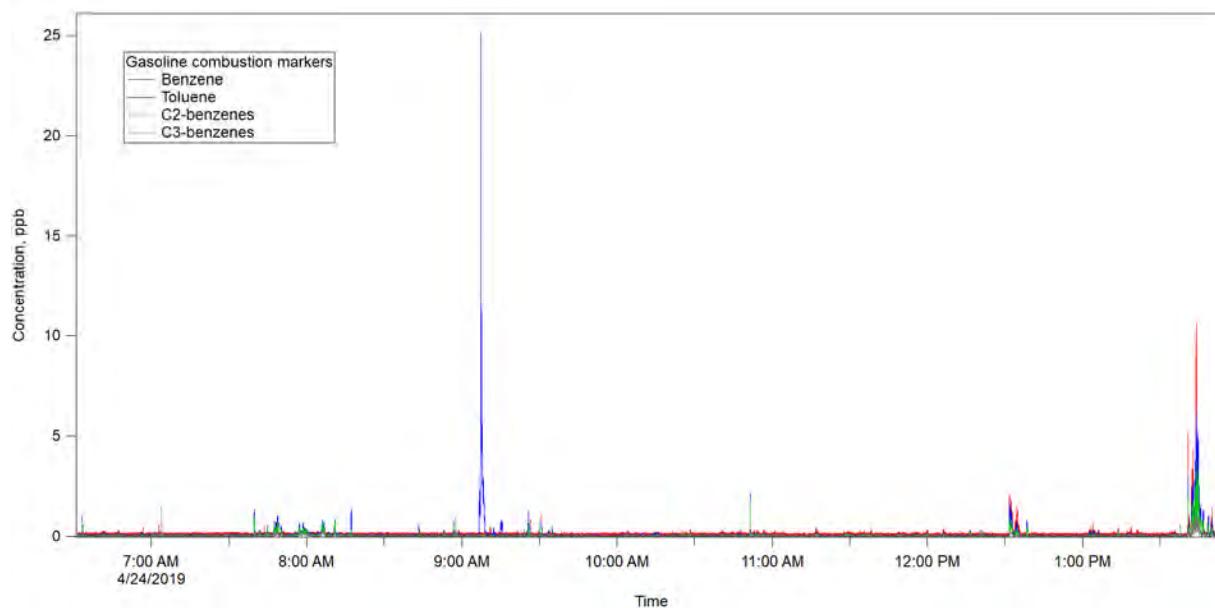


Figure 4-50. Gasoline Combustion Markers.

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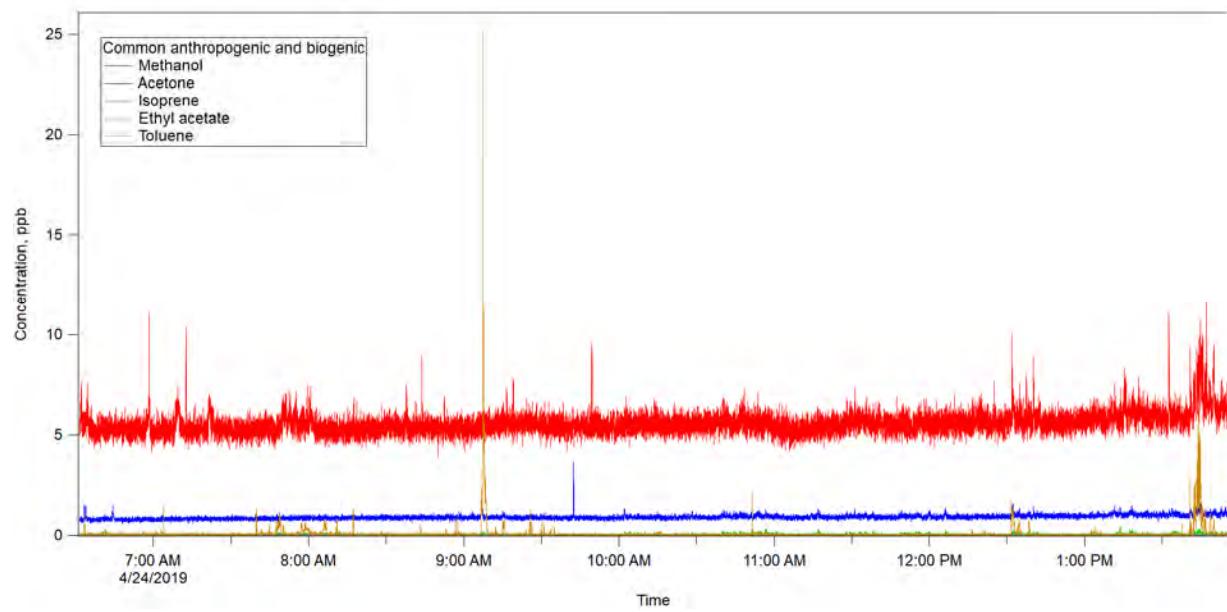


Figure 4-51. Plant and Human Markers.

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5.0 APRIL 25, 2019 – AREA MONITORING

5.1 Quality Assessment

Data from April 25, 2019, were assessed using Procedure 17124-DOE-HS-102. A Data Exchange Checklist was completed. The data were accepted by TerraGraphics with the following comments. Report No. 66409-RPT-004 was adequately documented and all checks passed the acceptance limits.

5.2 Summary

On April 25, 2019, ML personnel arrived at the TerraGraphics warehouse at 05:09. The QA/QC zero-air/span checks were performed on the LI-COR CO₂ monitor, the Picarro NH₃ analyzer, and the PTR-MS beginning at 05:26. The ML arrived on the Hanford Site and checked in with the CSM at 06:25. The ML began a site survey loop of A Farms at 06:40. The ML parked east of AP Farm at 06:40, monitoring AP-102 pump manual rotation activities. At 08:03, the ML arrived at U Farm and attempted to stay downwind of the change trailer, but wind conditions were varying. The ML departed U Farm at 08:29 and performed a site survey loop. The ML parked north of C Farm and began tracking an herbicide truck. At 11:19, the ML parked at the southeast corner of the TX/TY Farm area. At 12:08, the ML left the area and began performing site survey loops. The ML personnel checked out with the CSM at 14:00 and the ML departed the site. The ML arrived back at the TerraGraphics warehouse at 14:59.

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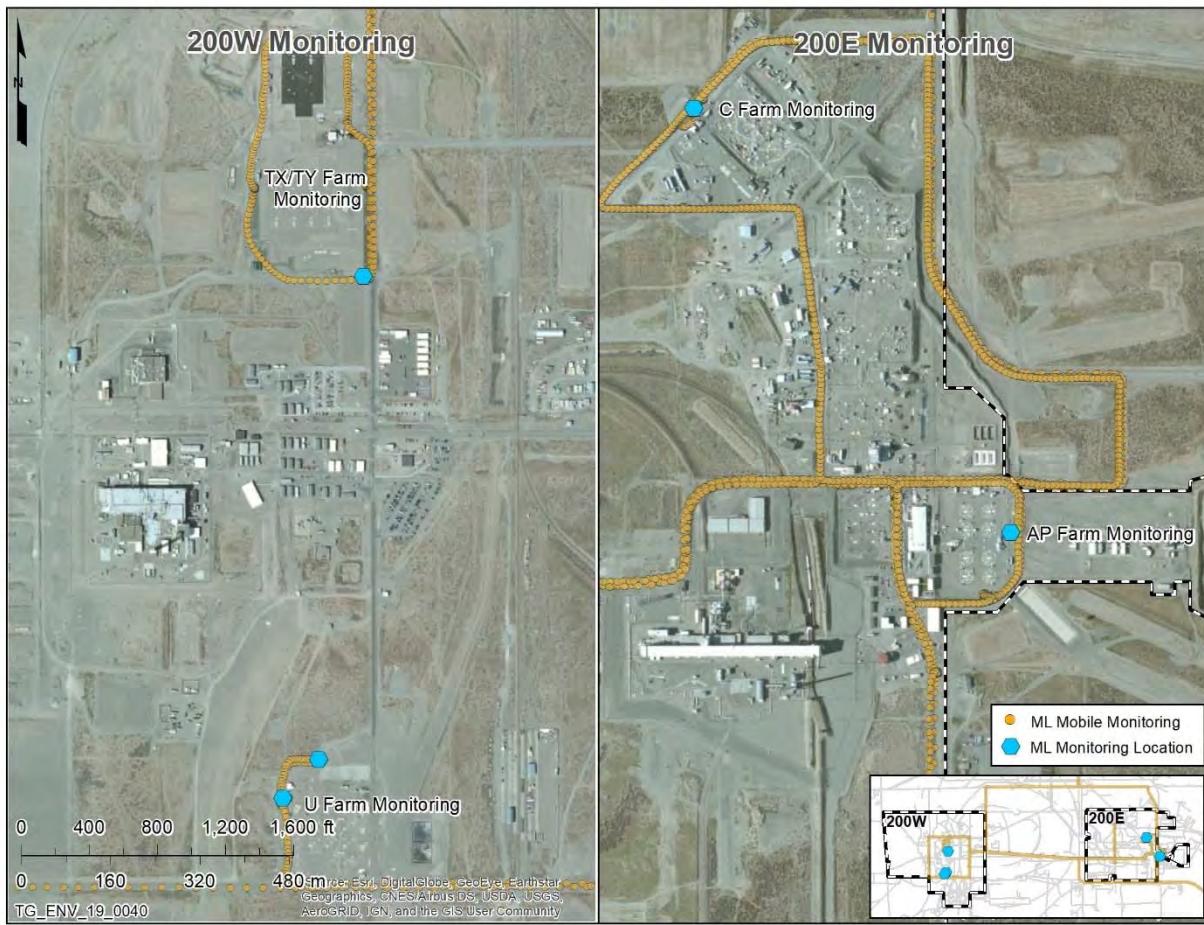


Figure 5-1. Location of the Mobile Laboratory for the Duration of the Monitoring Period.

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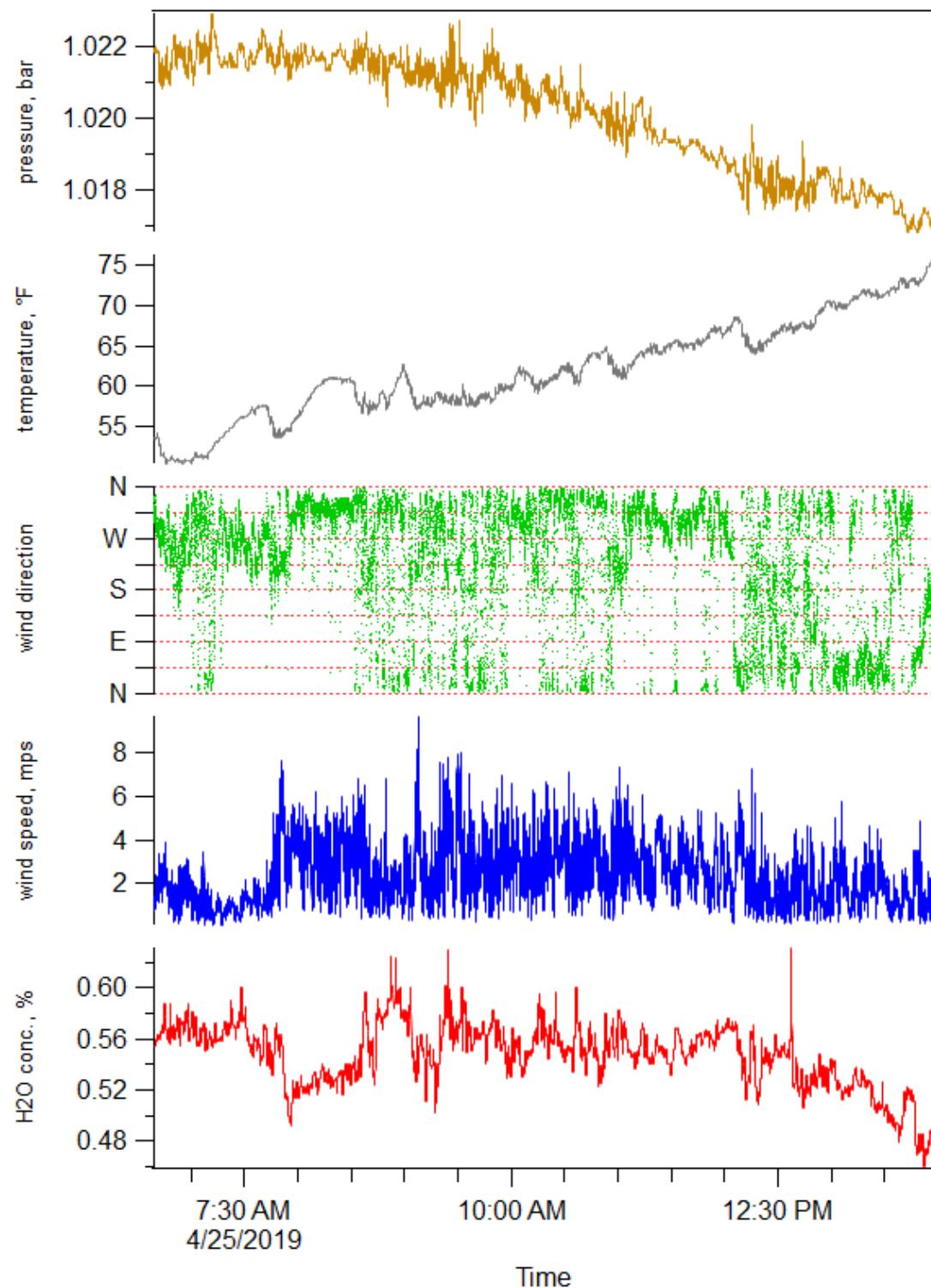


Figure 5-2. Weather Data for the Duration of the Monitoring Period.

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5.3 Samples Collected

Continuous air monitoring was performed using the following instrumentation:

- PTR-TOF 6000 X2,
- LI-COR CO₂ Monitor,
- Picarro Ammonia Monitor, and
- Airmar Weather Station.

Confirmatory air samples were not collected during this period.

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5.4 Area Monitoring

Table 5-1. Chemical of Potential Concern Statistical Information for the Monitoring Period of April 25, 2019. (2 Sheets)

COPC #	COPC Name	OEL (ppb)	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel St. Dev. (%)	Max. (ppb)	Median (ppb)
1	ammonia	25000	6.225	8.726†	1.328	15.221	14.661	8.312†
2	formaldehyde	300	1.721	<1.721	0.174	17.168	6.191	<1.721
3	methanol	200000	5.758	7.901†	1.419	17.955	116.233	7.832†
4	acetonitrile	20000	0.085	0.141†	0.035	24.837	0.549	0.139†
5	acetaldehyde	25000	1.027	1.447†	0.302	20.906	11.869	1.406†
6	ethylamine	5000	0.069	<0.069	0.021	63.319	0.149	<0.069
7	1,3-butadiene	1000	0.183	<0.183	0.055	49.557	0.714	<0.183
8	propanenitrile	6000	0.107	<0.107	0.024	46.163	0.218	<0.107
9	2-propenal	100	0.340	<0.34	0.058	53.396	1.078	<0.34
10	1-butanol + butenes	20000	0.214	<0.214	0.050	51.897	1.483	<0.214
11	methyl isocyanate	20	0.069	<0.069	0.025	57.418	0.174	<0.069
12	methyl nitrite	100	0.098	<0.098	0.037	39.216	0.572	<0.098
13	furan	1	0.062	<0.062	0.017	58.547	0.170	<0.062
14	butanenitrile	8000	0.039	<0.039	0.016	72.177	0.120	<0.039
15	but-3-en-2-one + 2,3-dihydrofuran + 2,5-dihydrofuran	200, 1, 1	0.041	<0.041	0.020	56.315	N/A*	N/A*
16	butanal	25000	0.061	0.143†	0.036	25.154	0.341	0.142†
17	NDMA**	0.3	0.082	<0.082	0.024	171.852	0.187	<0.082
18	benzene	500	0.236	<0.236	0.050	34.311	1.665	<0.236
19	2,4-pentadienenitrile + pyridine	300, 1000	0.085	<0.085	0.019	44.203	0.151	<0.085
20	2-methylene butanenitrile	300	0.036	<0.036	0.011	77.220	0.076	<0.036
21	2-methylfuran	1	0.043	<0.043	0.020	64.666	0.249	<0.043
22	pentanenitrile	6000	0.036	<0.036	0.012	81.538	0.094	<0.036
23	3-methyl-3-buten-2-one + 2-methyl-2-butenal	20, 30	0.043	<0.043	0.020	66.140	0.174	<0.043
24	NEMA**	0.3	0.058	<0.058	0.021	154.378	0.155	<0.058
25	2,5-dimethylfuran	1	0.032	<0.032	0.015	81.441	0.181	<0.032
26	hexanenitrile	6000	0.031	<0.031	0.010	93.082	0.073	<0.031
27	2-hexanone (MBK)	5000	0.036	<0.036	0.013	91.578	0.103	<0.036
28	NDEA**	0.1	0.034	<0.034	0.011	146.664	0.091	<0.034

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Table 5-1. Chemical of Potential Concern Statistical Information for the Monitoring Period of April 25, 2019. (2 Sheets)

COPC #	COPC Name	OEL (ppb)	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel St. Dev. (%)	Max. (ppb)	Median (ppb)
29	butyl nitrite + 2-nitro-2-methylpropane	100, 300	0.058	<0.058	0.015	49.992	0.103	<0.058
30	2,4-dimethylpyridine	500	0.036	<0.036	0.010	105.675	0.162	<0.036
31	2-propylfuran + 2-ethyl-5-methylfuran	1	0.027	<0.027	0.012	102.766	0.090	<0.027
32	heptanenitrile	6000	0.027	<0.027	0.009	97.437	0.061	<0.027
33	4-methyl-2-hexanone	500	0.033	<0.033	0.011	95.943	0.074	<0.033
34	NMOR**	0.6	0.021	<0.021	0.009	172.696	0.070	<0.021
35	butyl nitrate	2500	0.022	<0.022	0.007	125.048	0.053	<0.022
36	2-ethyl-2-hexenal + 4-(1-methylpropyl)-2,3-dihydrofuran + 3-(1,1-dimethylethyl)-2,3-dihydrofuran	100, 1, 1	0.028	<0.028	0.010	93.744	0.086	<0.028
37	6-methyl-2-heptanone	8000	0.028	<0.028	0.009	92.542	0.074	<0.028
38	2-pentylfuran	1	0.026	<0.026	0.010	91.077	0.082	<0.026
39	biphenyl	200	0.022	<0.022	0.009	127.802	0.060	<0.022
40	2-heptylfuran	1	0.067	<0.067	0.015	42.904	0.105	<0.067
41	1,4-butanediol dinitrate	50	0.036	<0.036	0.010	87.985	0.074	<0.036
42	2-octylfuran	1	0.020	<0.02	0.008	205.792	0.071	<0.02
43	1,2,3-propanetriol 1,3-dinitrate	50	0.011	<0.011	0.004	365.357	0.055	<0.011
44	PCB	1000	0.034	<0.034	0.010	72.343	0.069	<0.034
45	6-(2-furanyl)-6-methyl-2-heptanone	1	0.025	<0.025	0.008	100.891	0.059	<0.025
46	furfural acetophenone	1	0.064	<0.064	0.015	43.559	0.116	<0.064
N/A*	The maximum peak value for but-3-en-2-one + 2,3 dihydrofuran + 2,5 dihydrofuran was 0.343 ppb and the median value was <0.041 ppb. The PTR-MS results for but-3-en-2-one + 2,3 dihydrofuran + 2,5 dihydrofuran are not compared to OEL concentrations because: 1) the result is suspect due to a known biogenic interferant (methacrolein) that is expected to be in concentrations that occasionally exceed the dihydrofuran OEL, and 2) this combination of COPCs have OEL concentrations that differ by a factor of 200, which provide widely variant bases for these numbers.							
**	Nitrosamine results are suspect due to isobaric interferants causing positive bias that have been encountered during previous background [53005-81-RPT-007, <i>PTR-MS Mobile Laboratory Vapor Monitoring Background Study</i> , (3/18/2018 – 4/20/2018), and <i>Fiscal Year 2017 Mobile Laboratory Vapor Monitoring at the Hanford Site: Monitoring During Waste Disturbing Activities and Background Study</i> , RJ Lee Group, Inc.].							
<	COPC Averages below the MDL.							
†	COPC Averages between the RL and the MDL.							
	COPC Averages >100% of the OEL.							
	COPC Averages 50-100% of the OEL.							
	COPC Averages 10-50% of the OEL.							

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Figure 5-3 through Figure 5-51 display 46 COPC signals, overlaid with the same signal smoothed using a 1-minute moving average (in cases where a moving average assists with data visualization), and CO₂, for the monitoring period April 25, 2019. If within range of the plot's left axis, a green horizontal line representing 50% of the COPC's OEL, a blue horizontal line representing the COPC's OEL, a horizontal purple line representing the RL, and a pink horizontal line representing the MDL are shown.

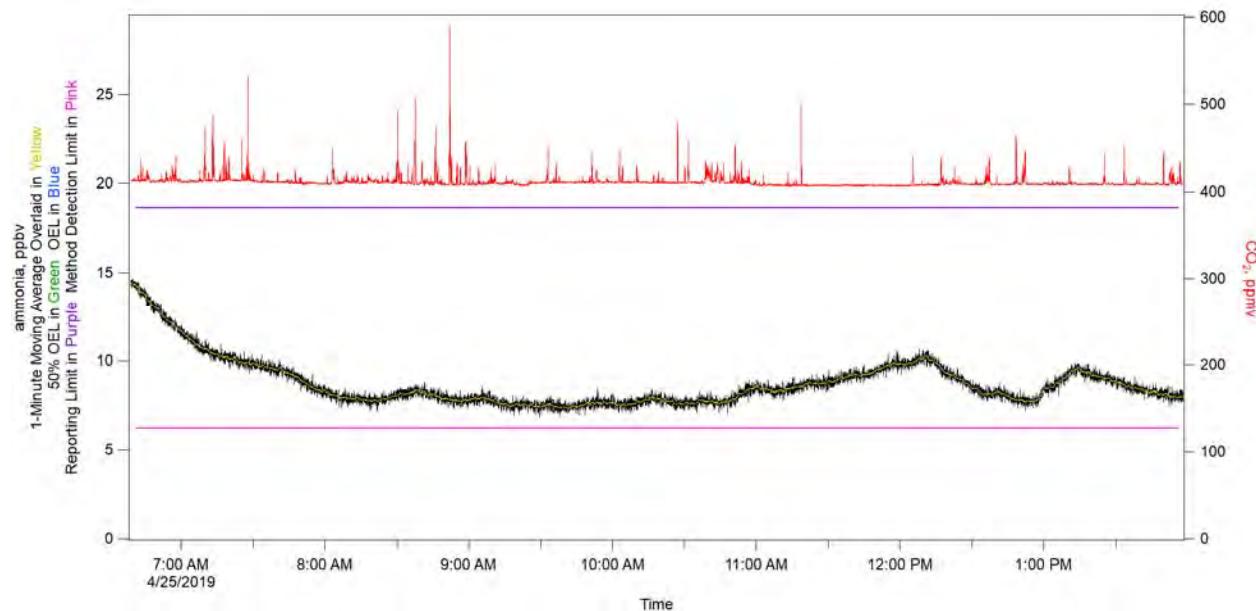


Figure 5-3. Ammonia.

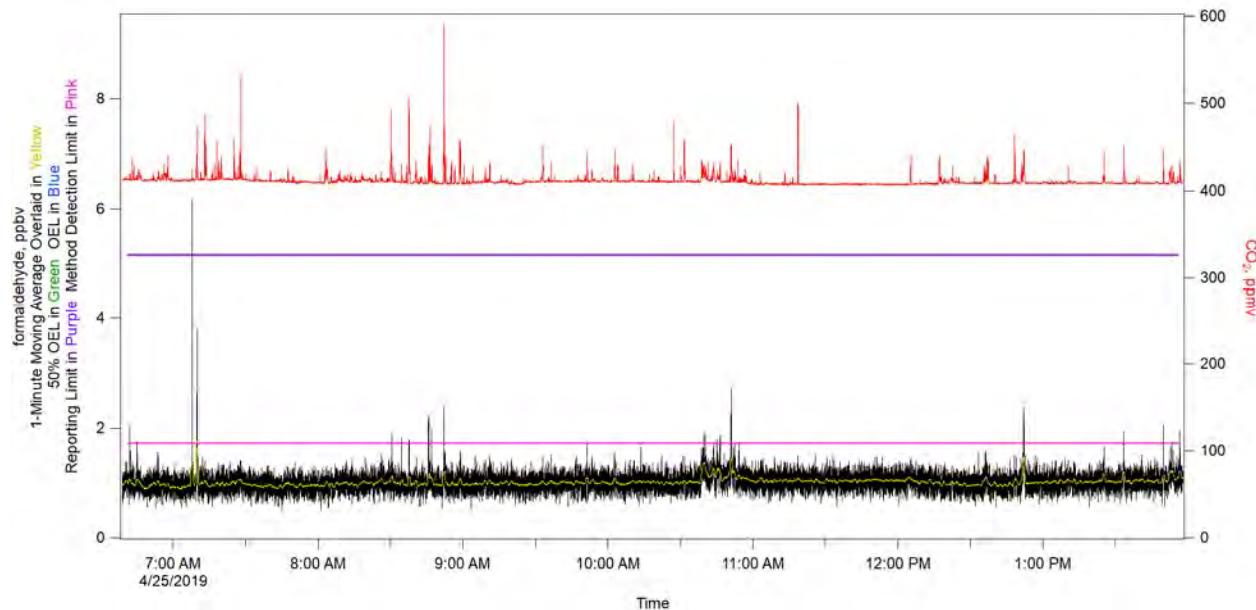


Figure 5-4. Formaldehyde.

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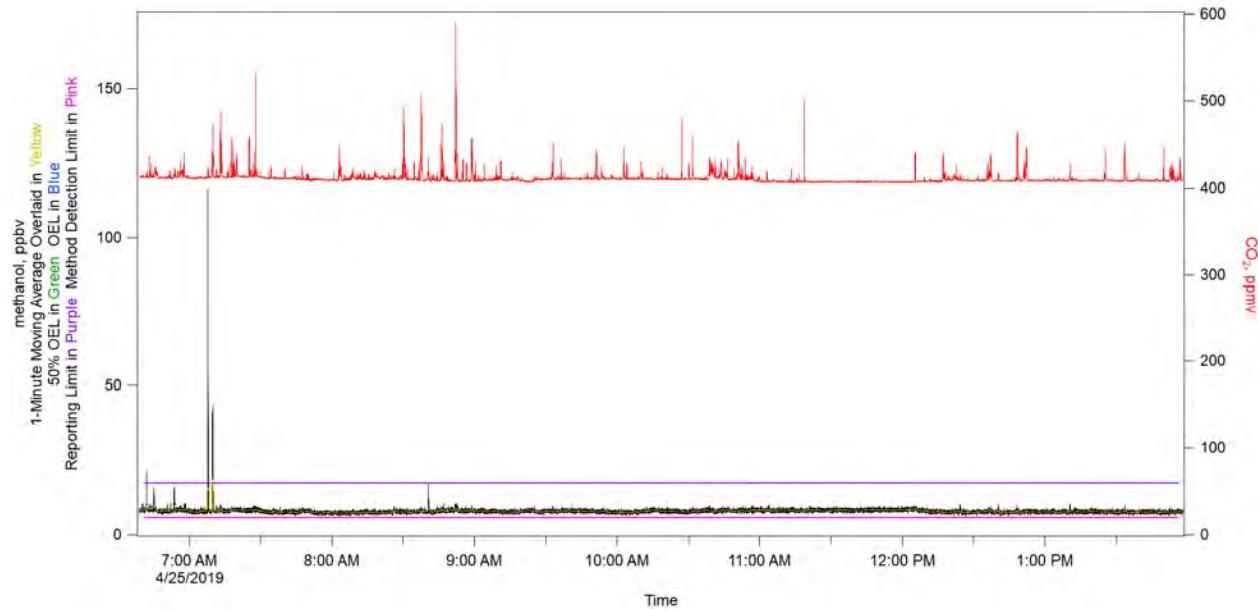


Figure 5-5. Methanol.

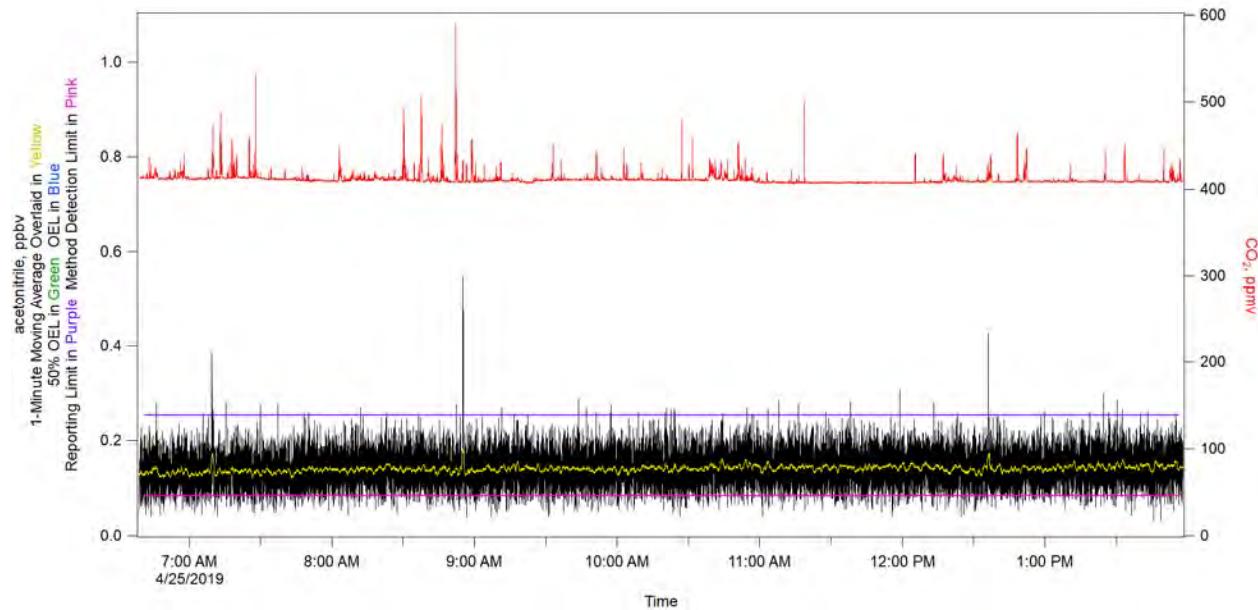


Figure 5-6. Acetonitrile.

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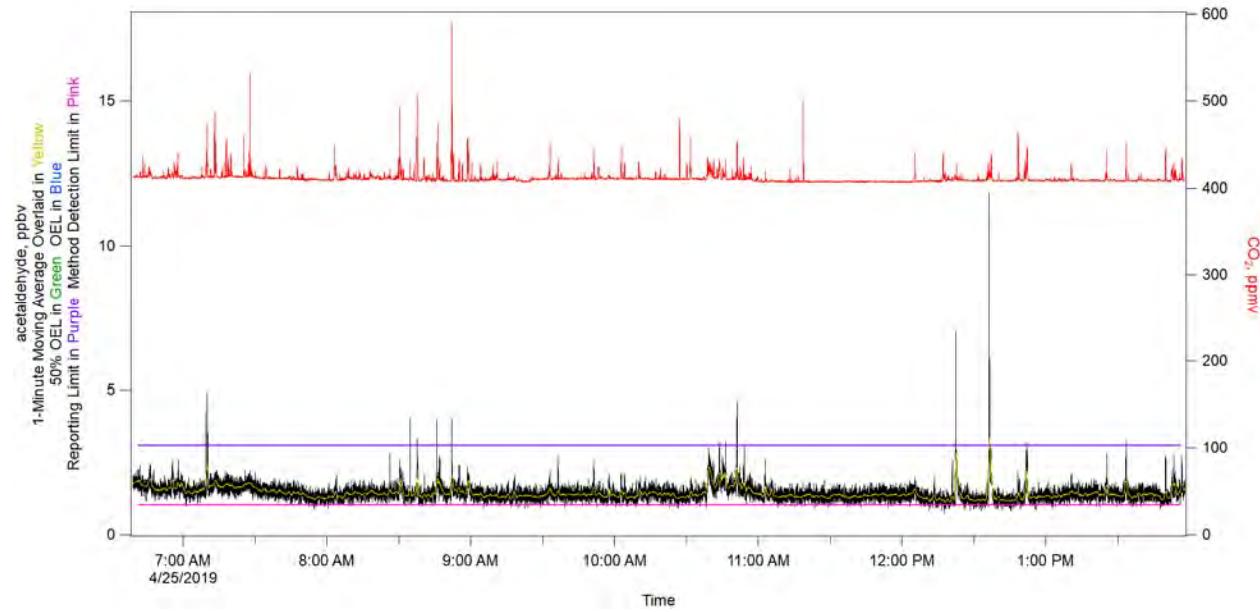


Figure 5-7. Acetaldehyde.

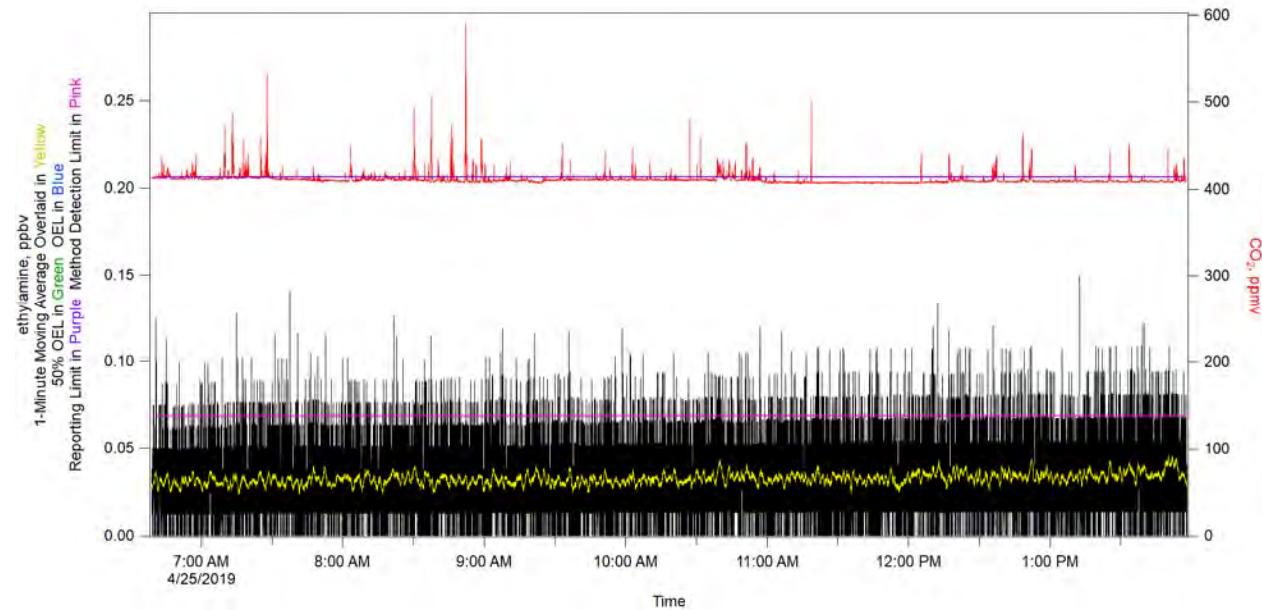


Figure 5-8. Ethylamine.

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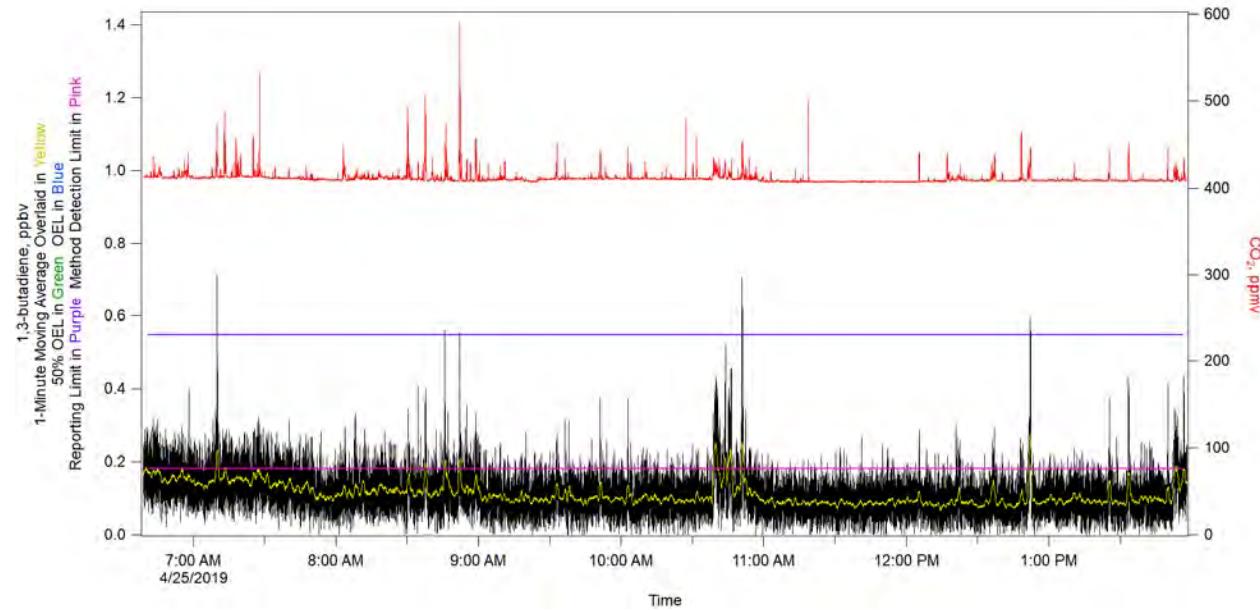


Figure 5-9. 1,3-butadiene.

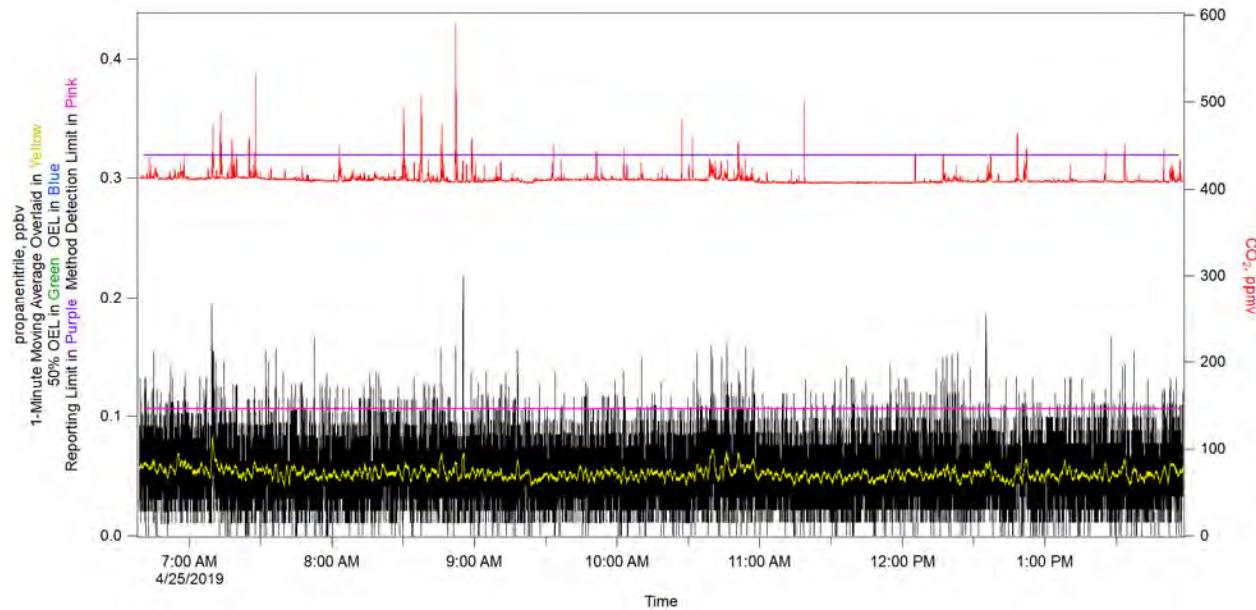


Figure 5-10. Propanenitrile.

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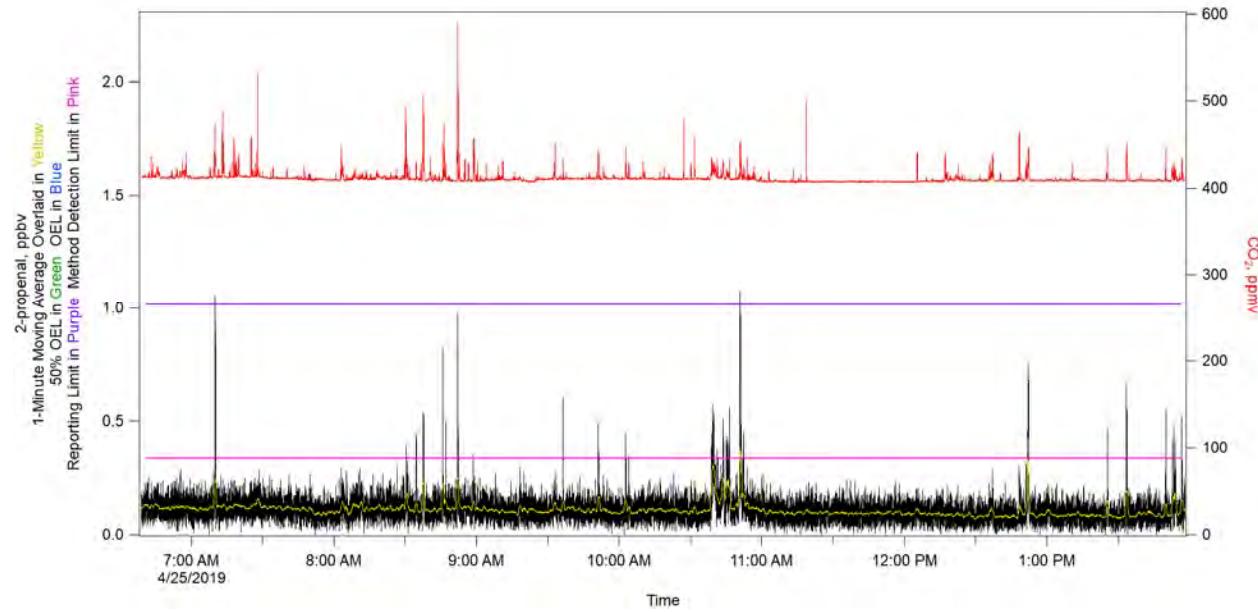


Figure 5-11. 2-propenal.

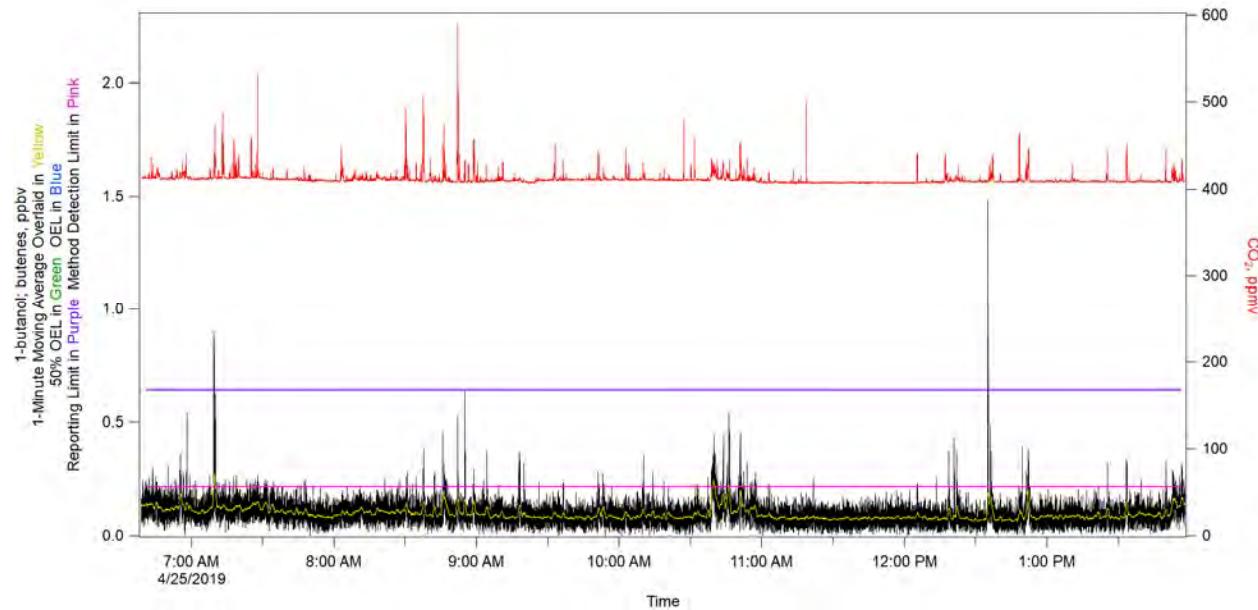


Figure 5-12. 1-butanol; Butenes.

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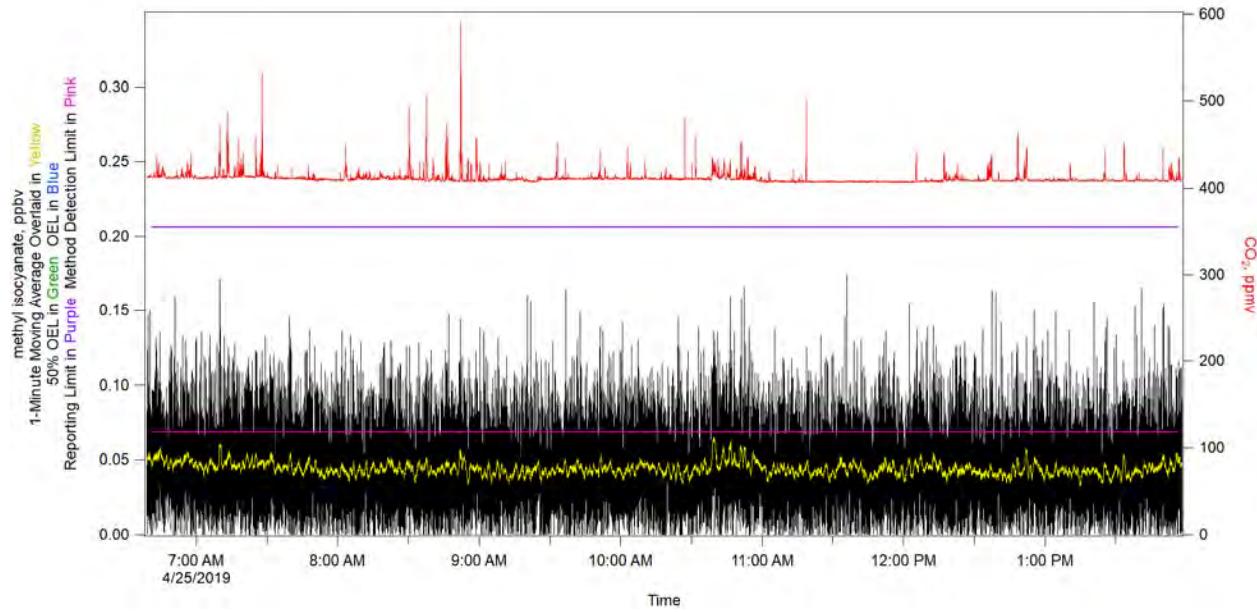


Figure 5-13. Methyl Isocyanate.

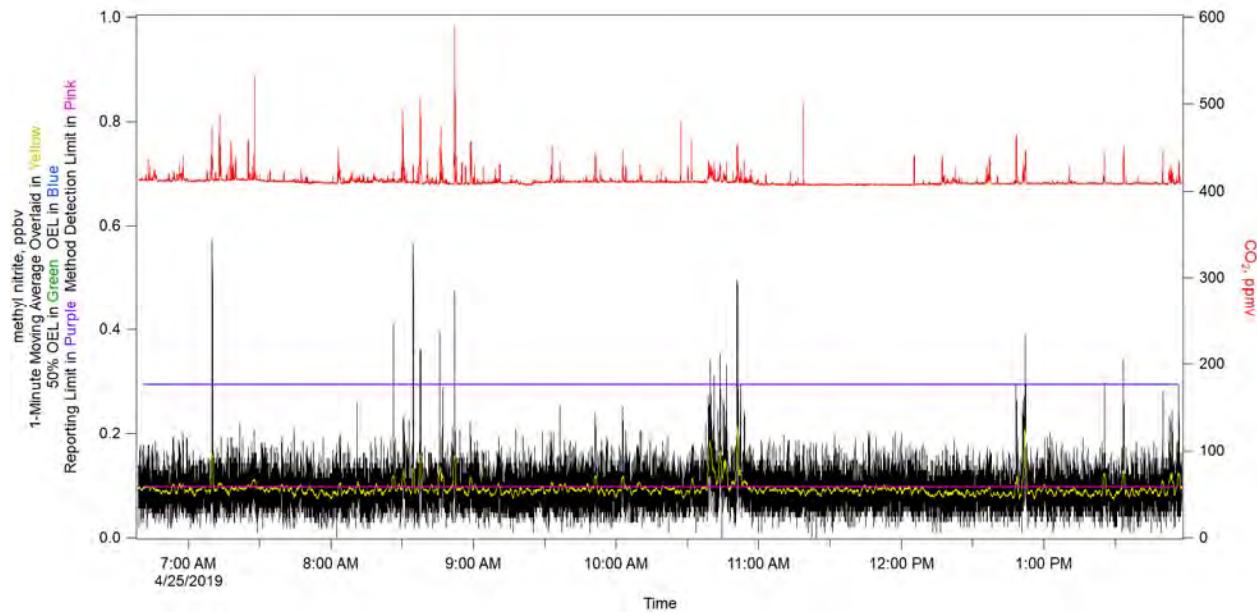


Figure 5-14. Methyl Nitrite.

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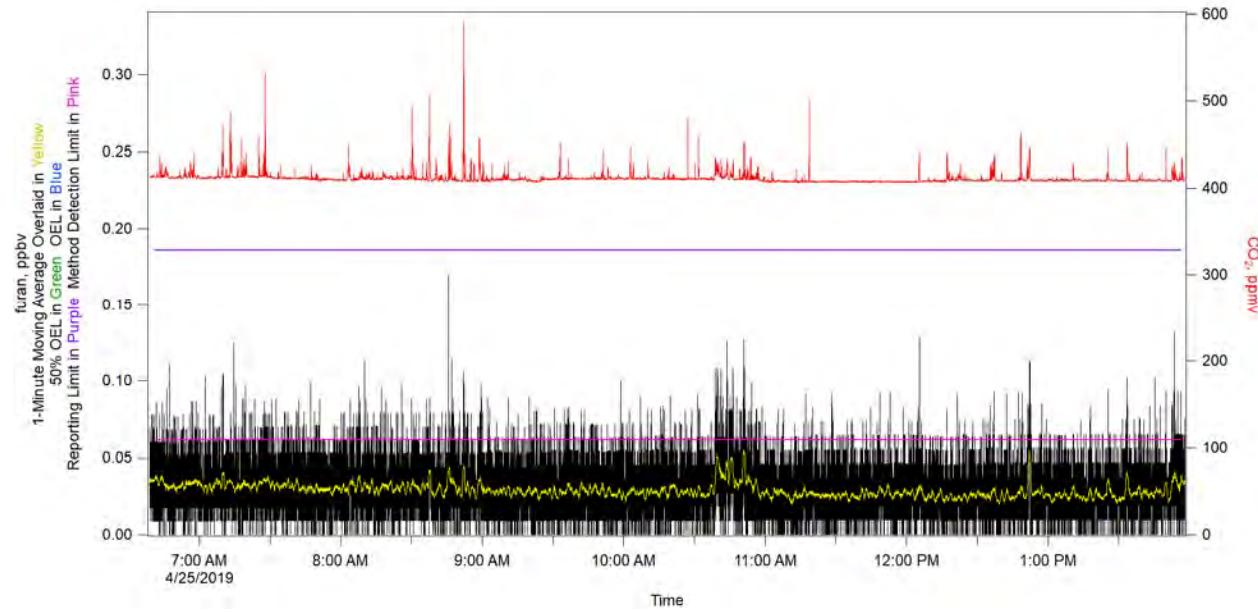


Figure 5-15. Furan.

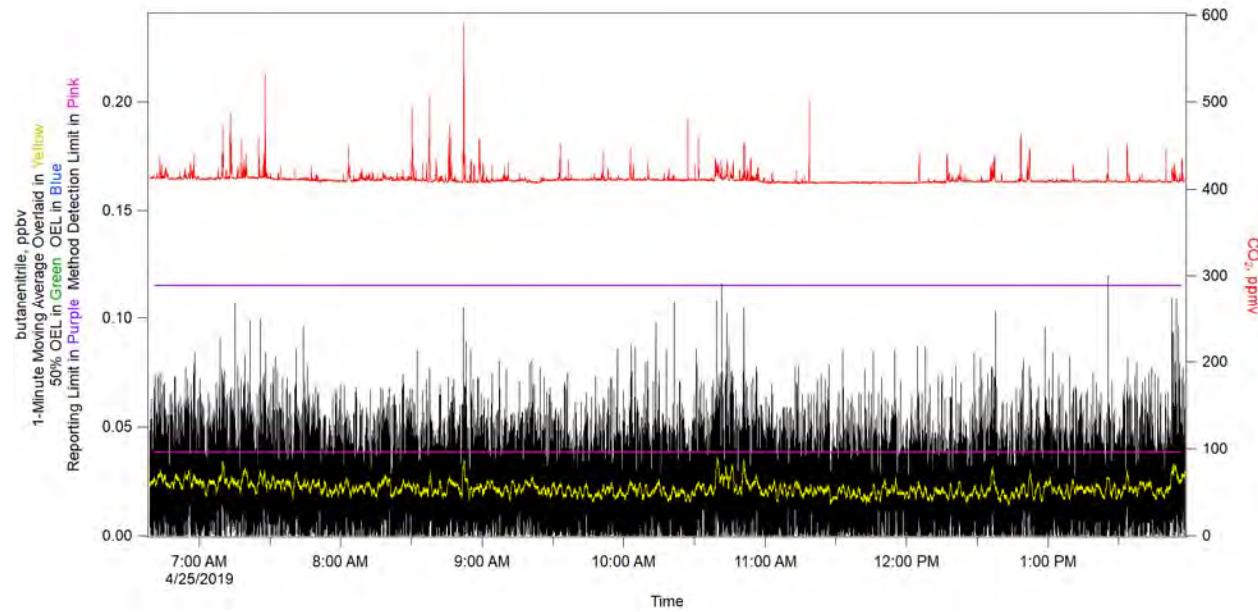


Figure 5-16. Butanenitrile.

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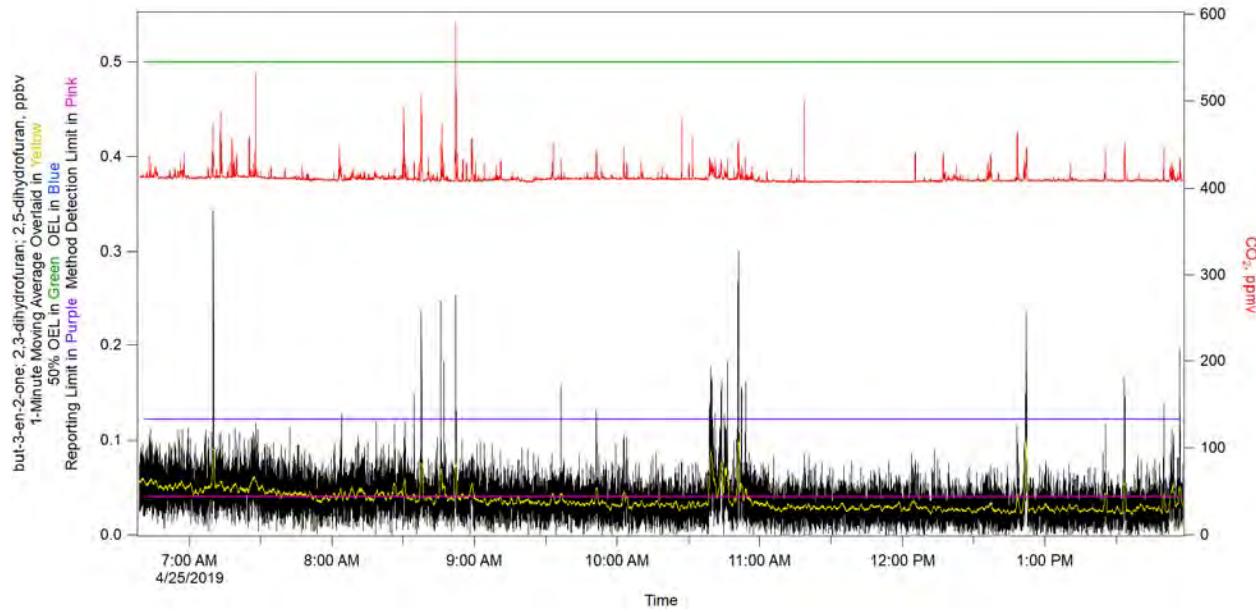


Figure 5-17. But-3-en-2-one; 2,3-dihydrofuran; 2,5-dihydrofuran.

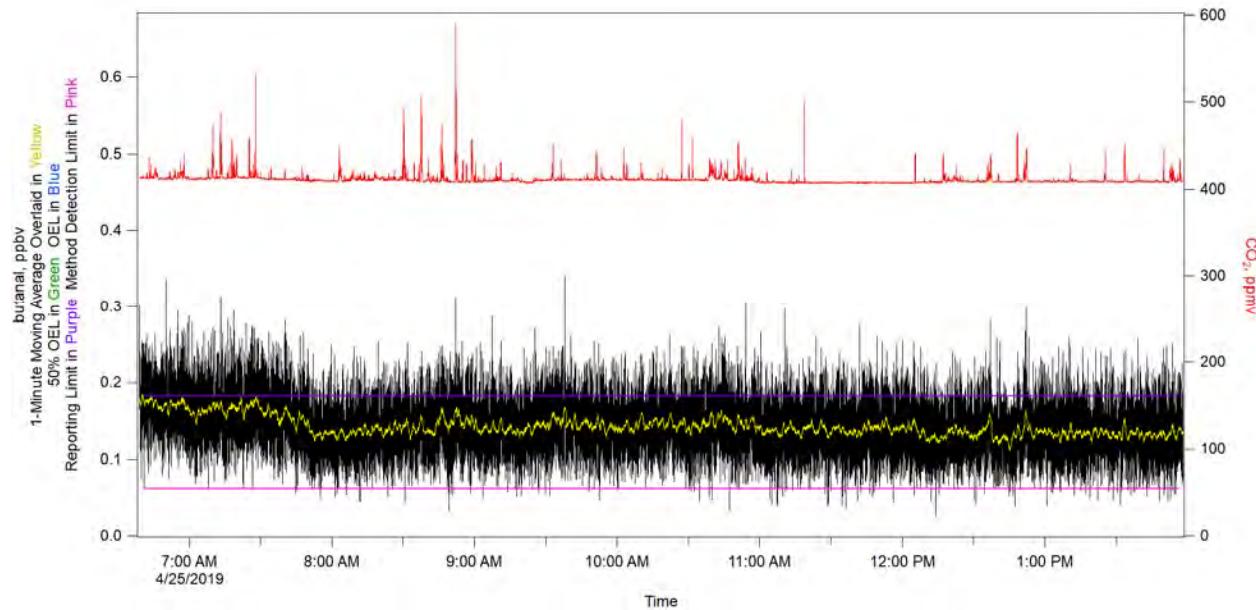


Figure 5-18. Butanal.

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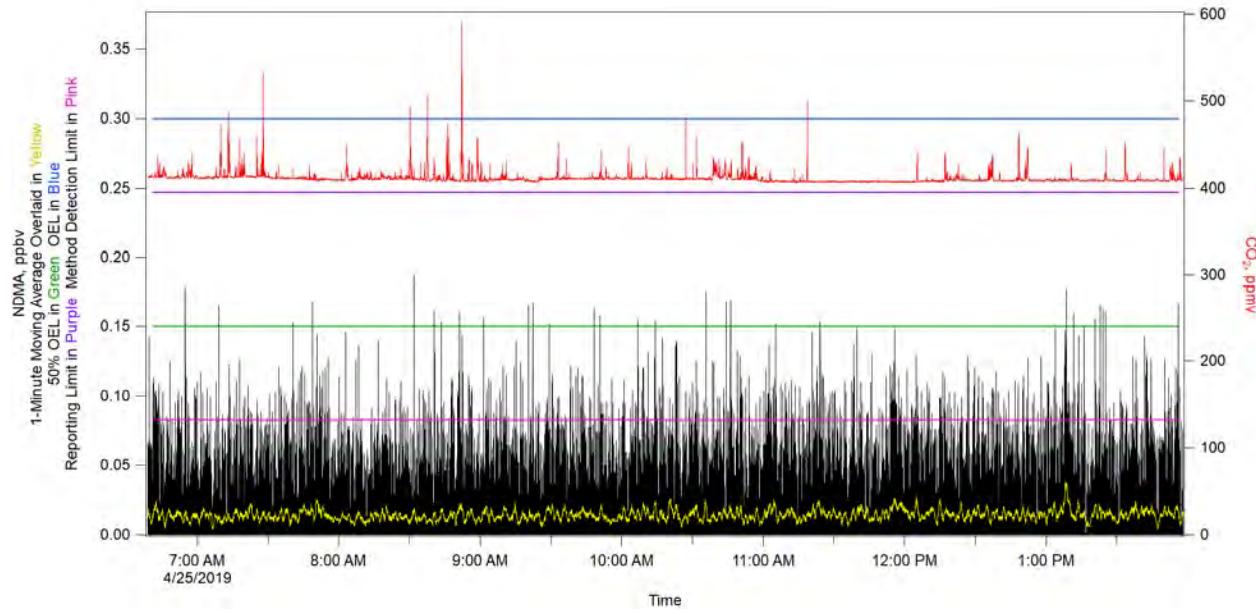


Figure 5-19. N-nitrosodimethylamine (NDMA).

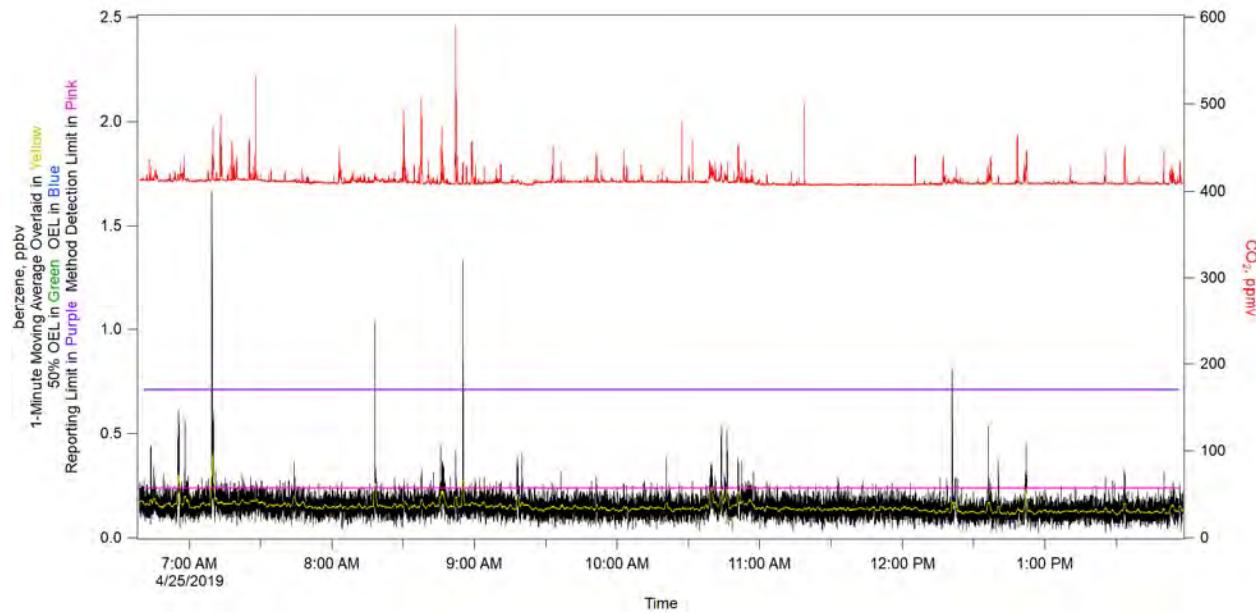


Figure 5-20. Benzene.

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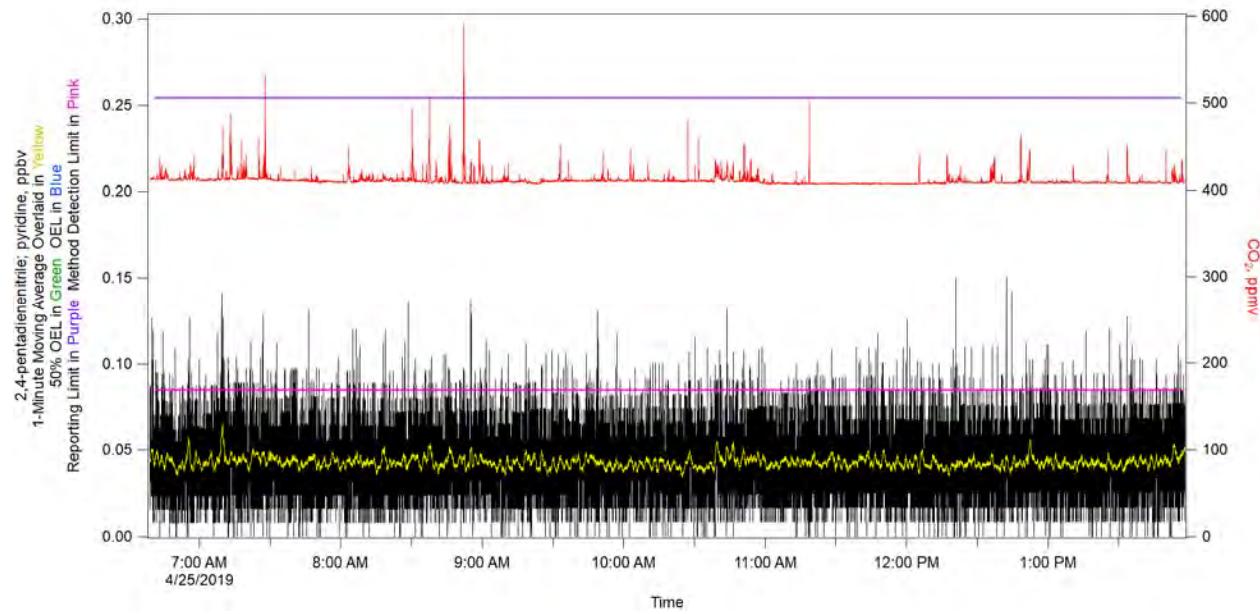


Figure 5-21. 2,4-pentadienenitrile; Pyridine.

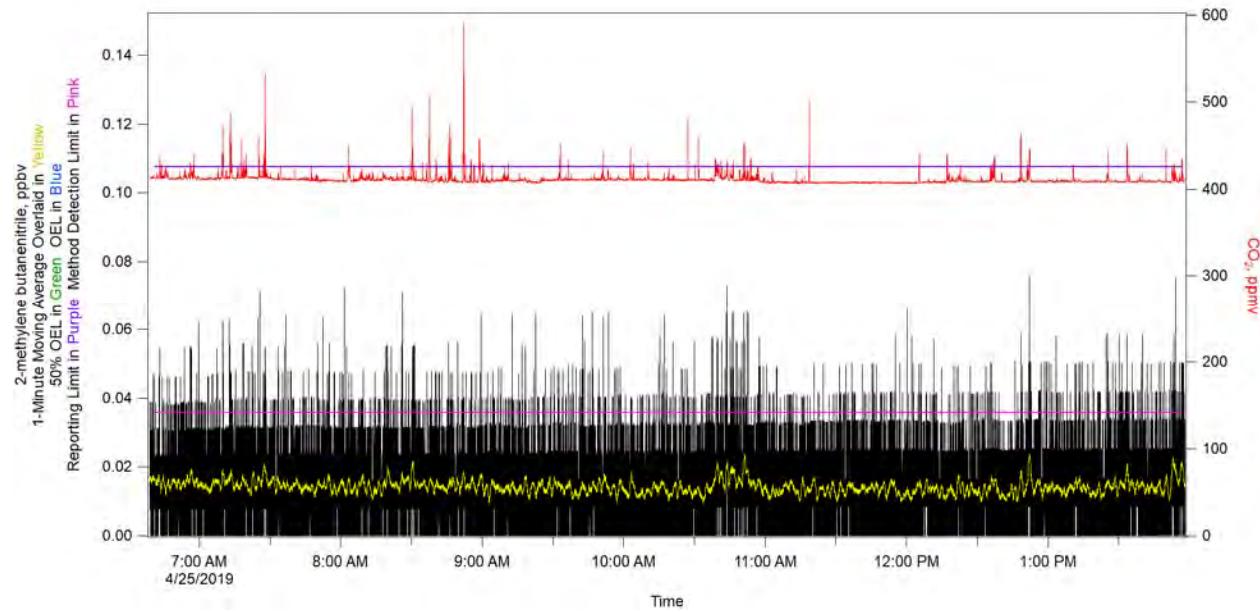


Figure 5-22. 2-methylene Butanenitrile.

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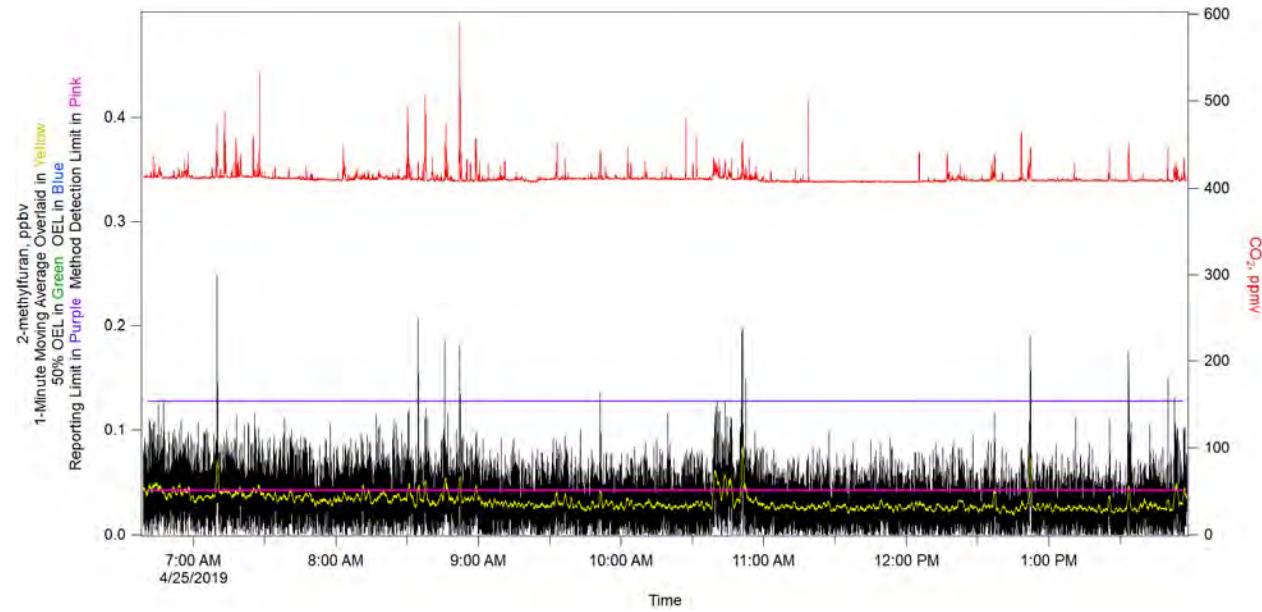


Figure 5-23. 2-methylfuran.

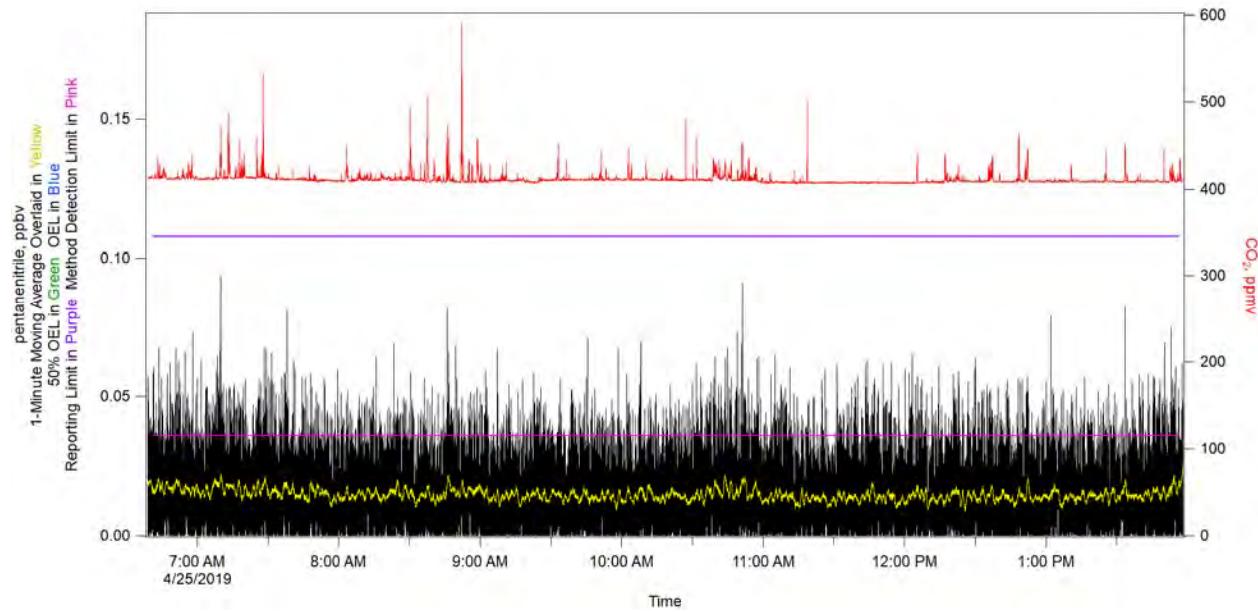


Figure 5-24. Pentanenitrile.

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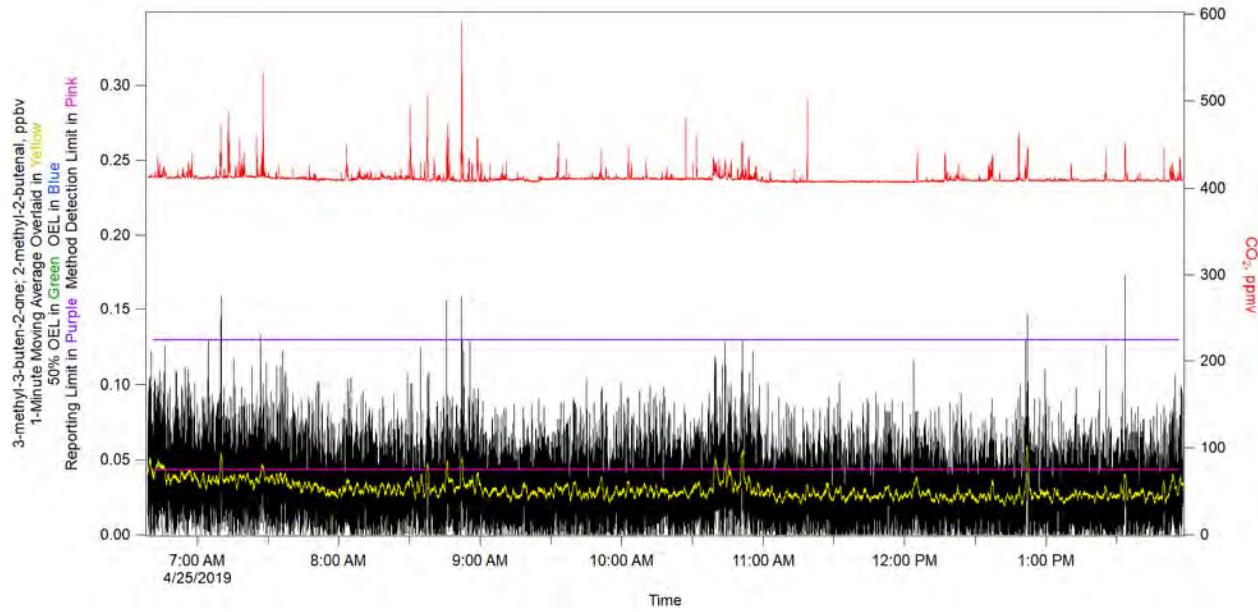


Figure 5-25. 3-methyl-3-buten-2-one; 2-methyl-2-butenal.

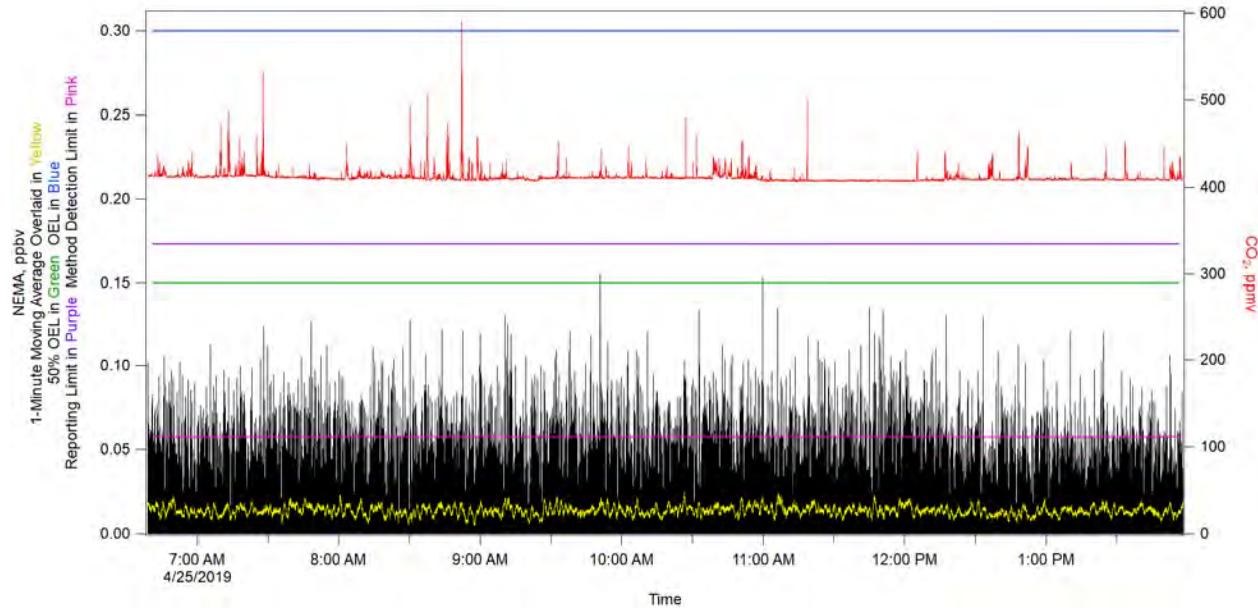


Figure 5-26. N-nitrosomethylethylamine (NEMA).

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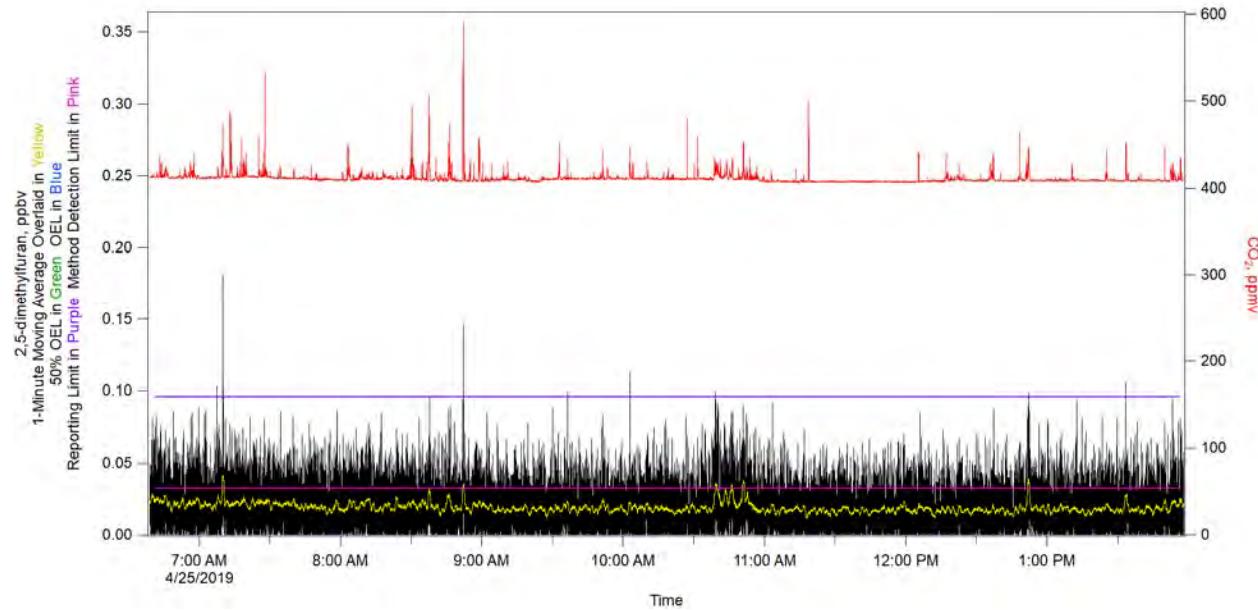


Figure 5-27. 2,5-dimethylfuran.

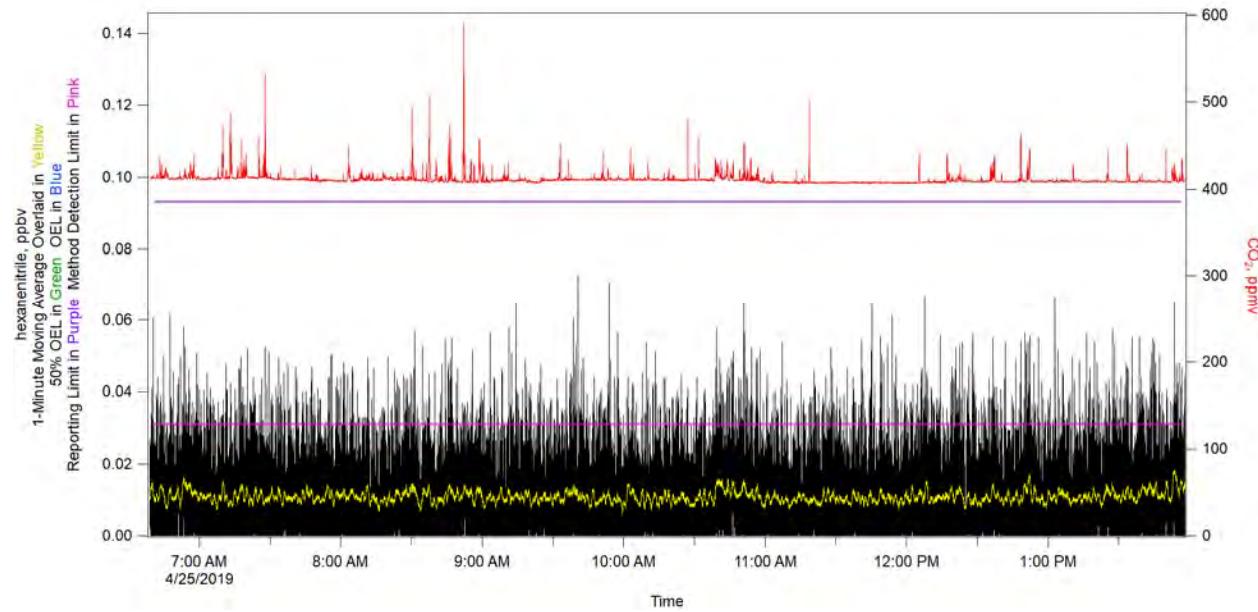


Figure 5-28. Hexanenitrile.

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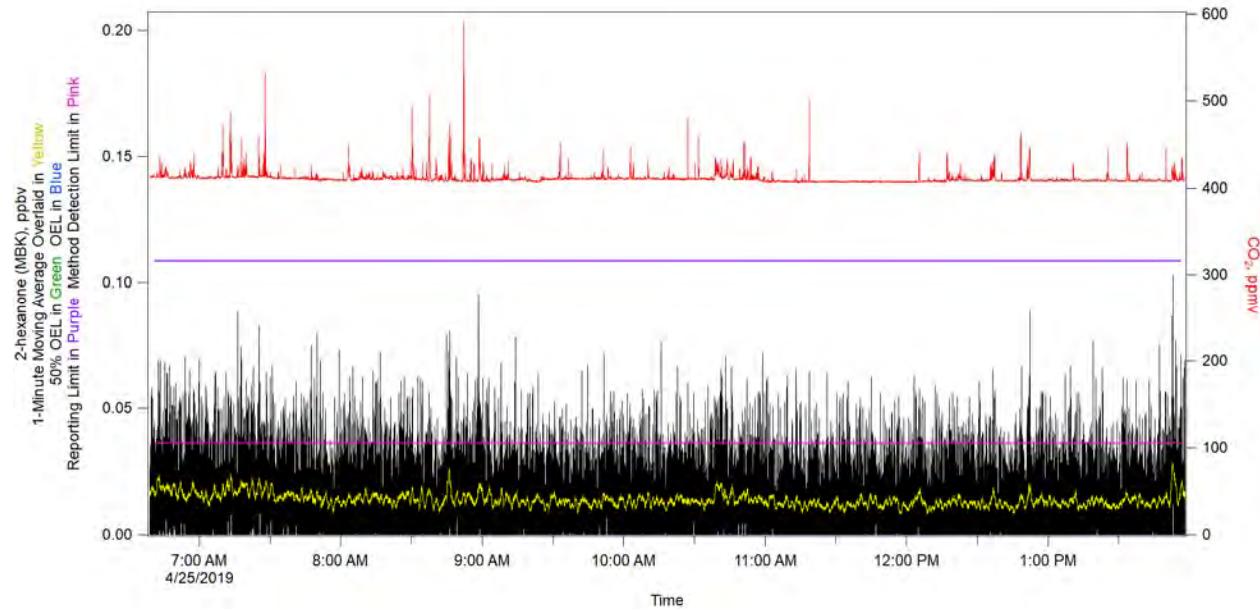


Figure 5-29. 2-hexanone (MBK).

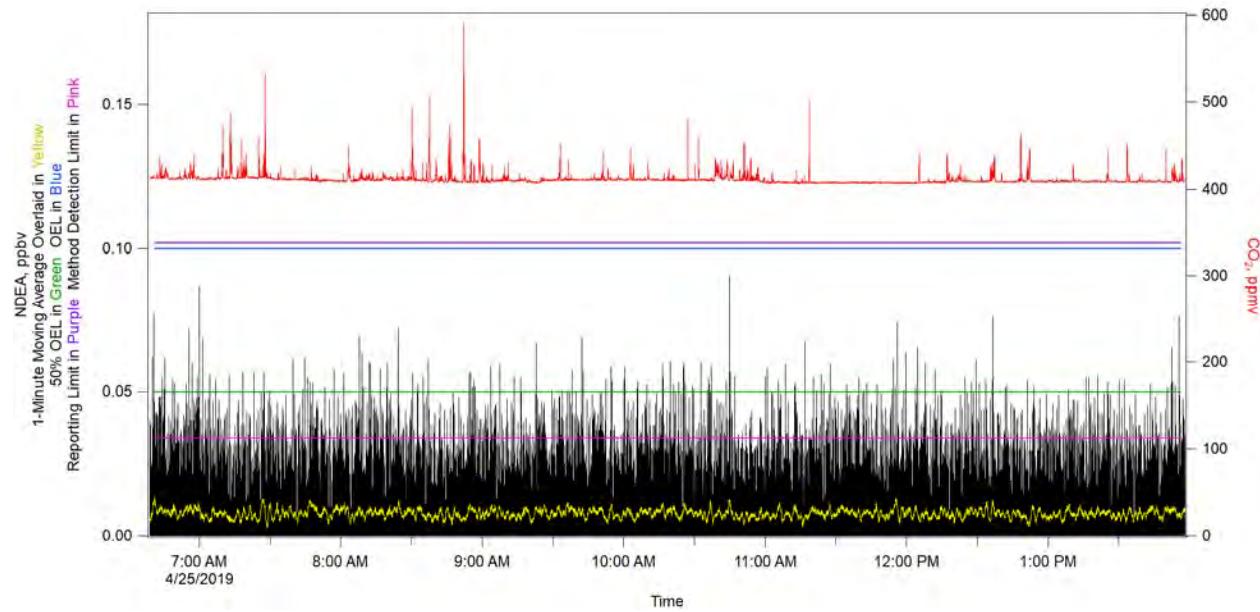


Figure 5-30. N-nitrosodiethylamine (NDEA).

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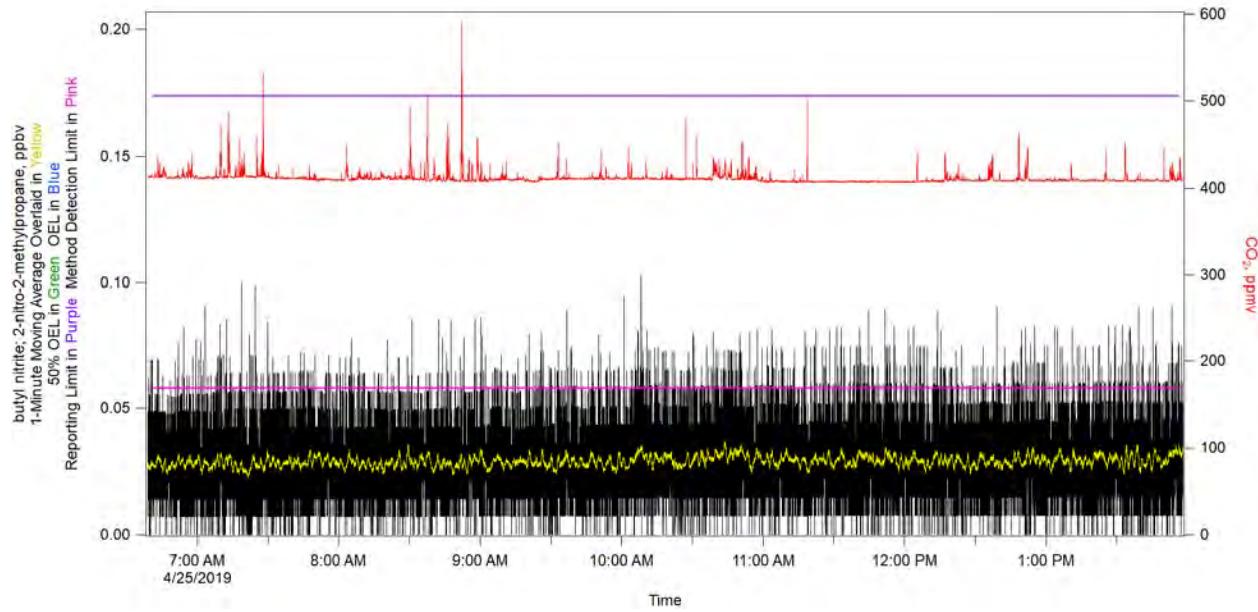


Figure 5-31. Butyl Nitrite; 2-nitro-2-methylpropane.

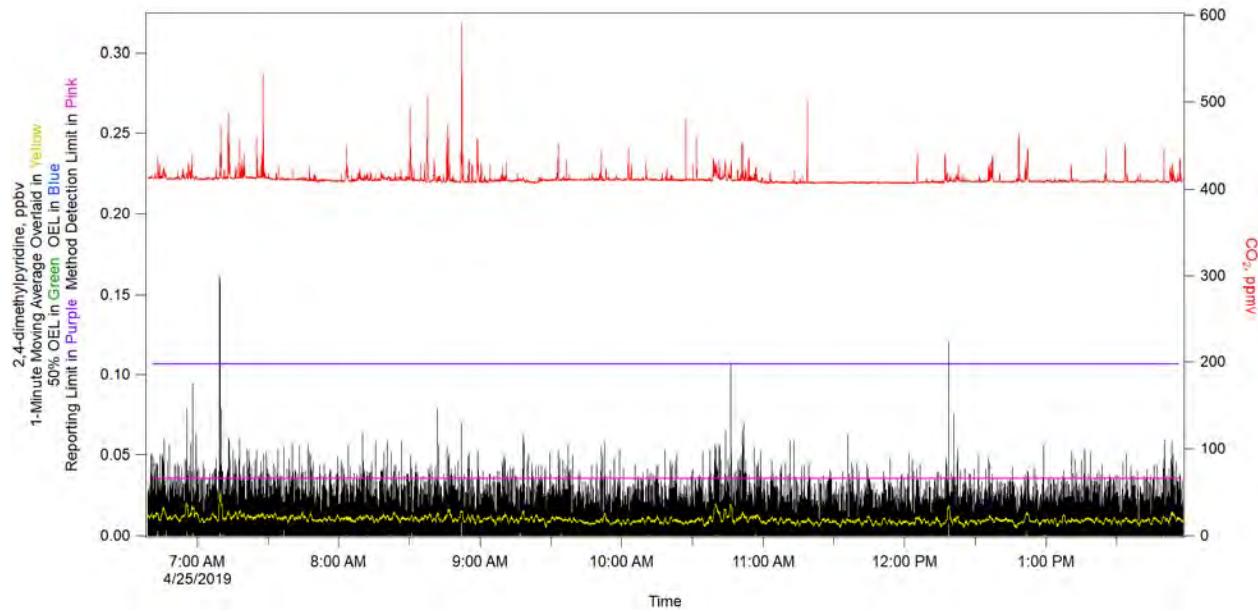


Figure 5-32. 2,4-dimethylpyridine.

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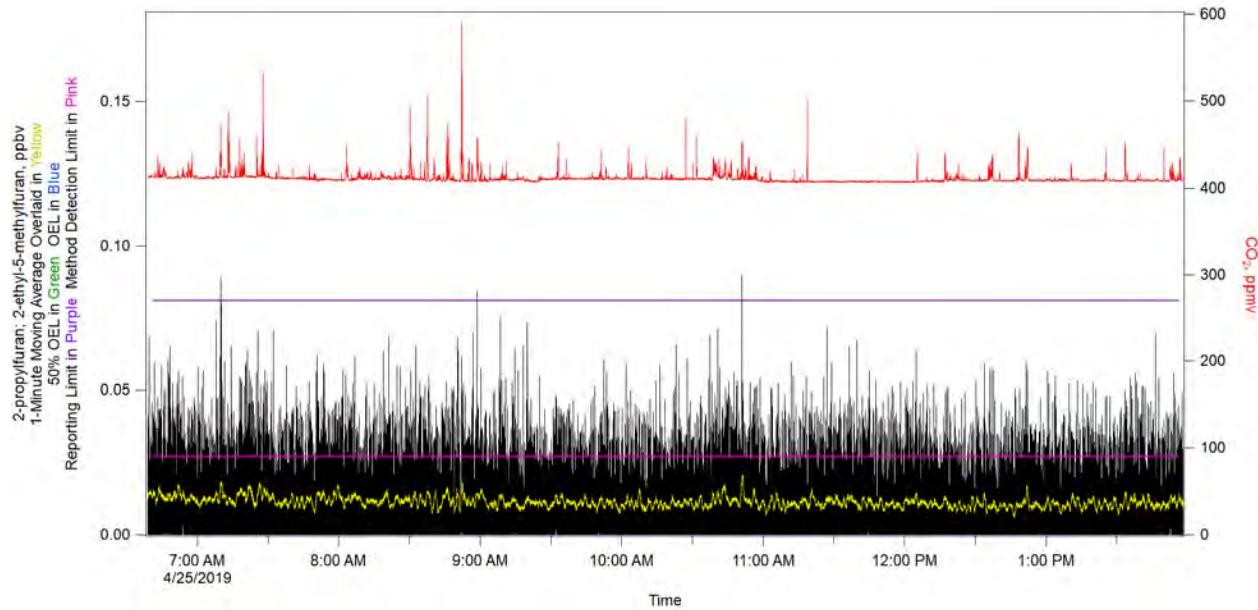


Figure 5-33. 2-propylfuran; 2-ethyl-5-methylfuran.

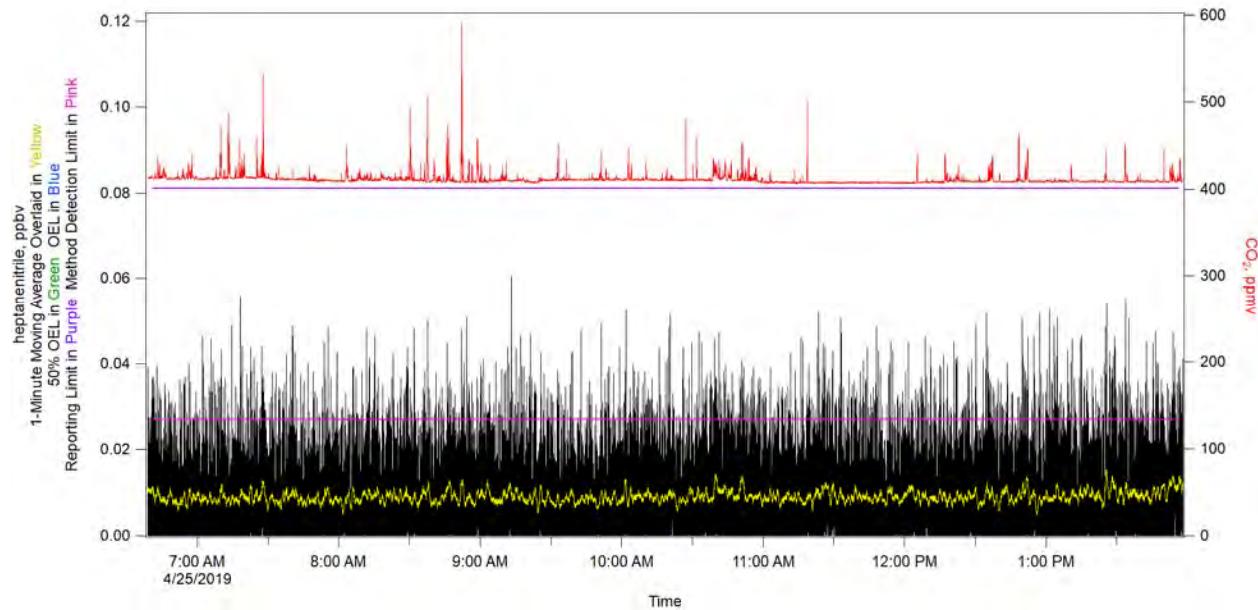


Figure 5-34. Heptanenitrile.

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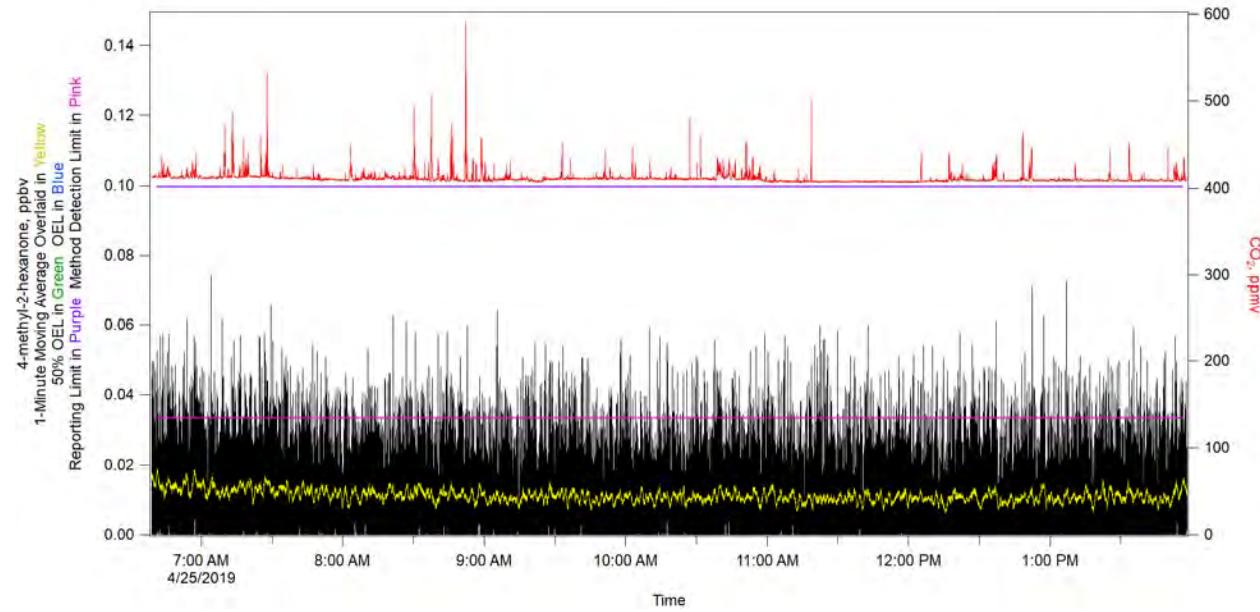


Figure 5-35. 4-methyl-2-hexanone.

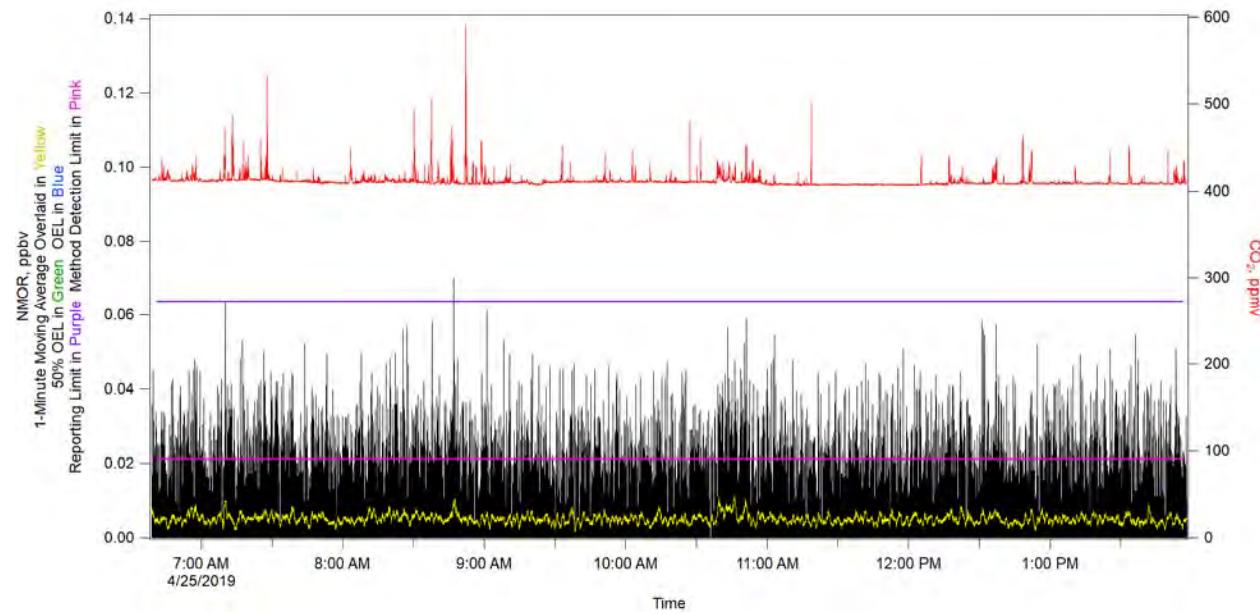


Figure 5-36. N-nitrosomorpholine (NMOR).

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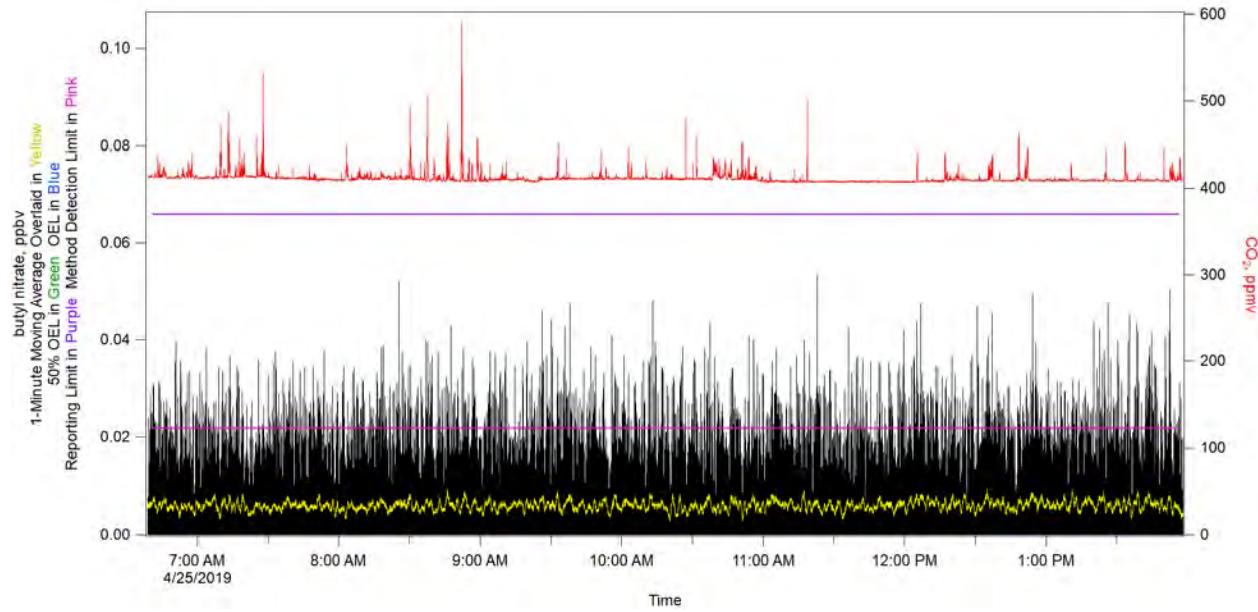
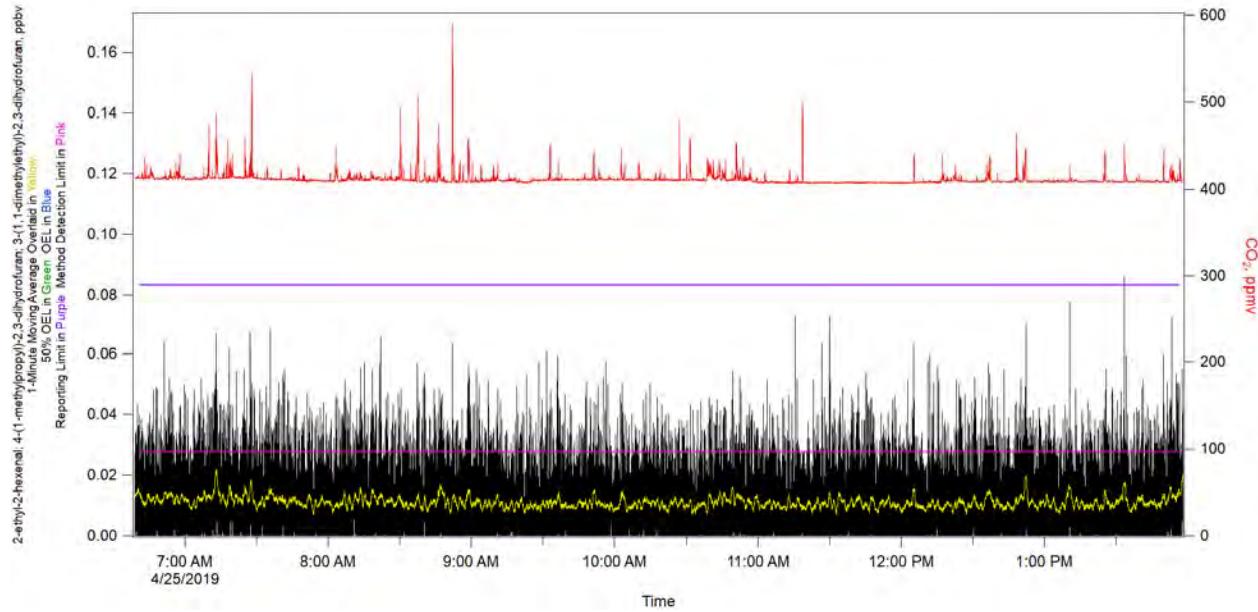


Figure 5-37. Butyl Nitrate.



**Figure 5-38. 2-ethyl-2-hexenal; 4-(1-methylpropyl)-2,3-dihydrofuran
3-(1,1-dimethylethyl)-2,3-dihydrofuran.**

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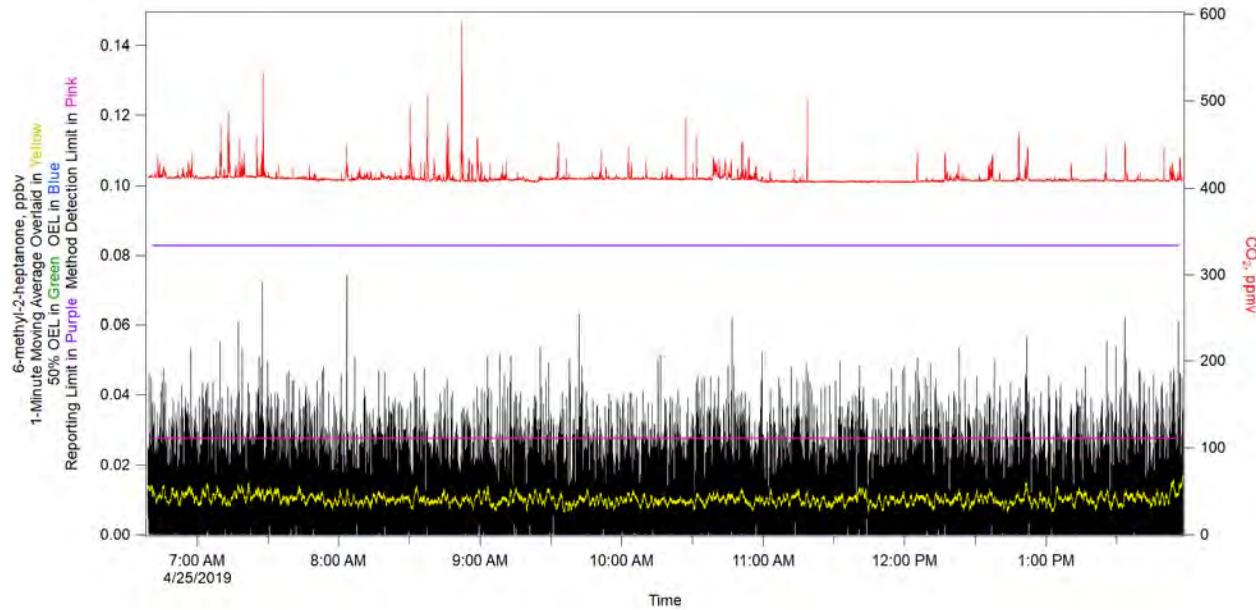


Figure 5-39. 6-methyl-2-heptanone.

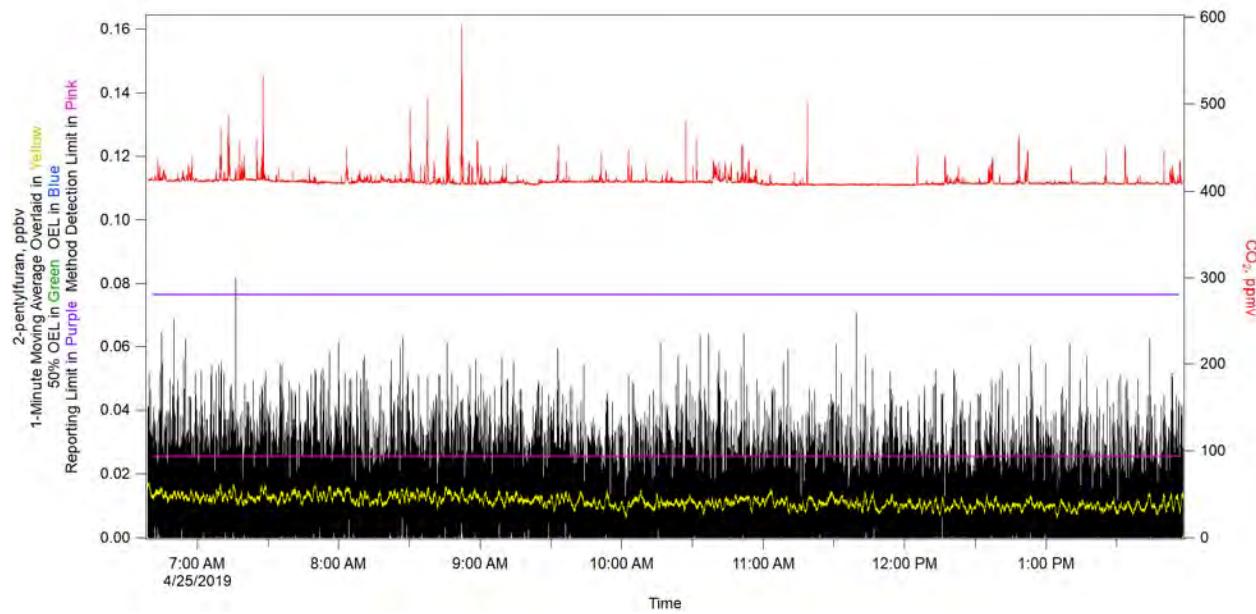


Figure 5-40. 2-pentylfuran.

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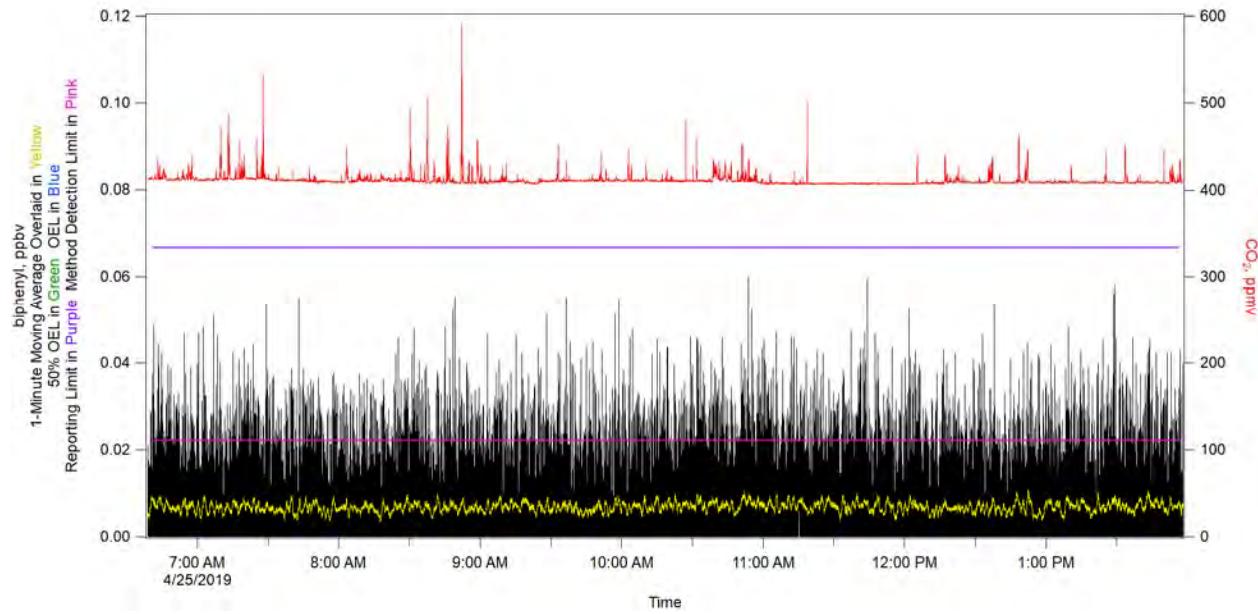


Figure 5-41. Biphenyl.

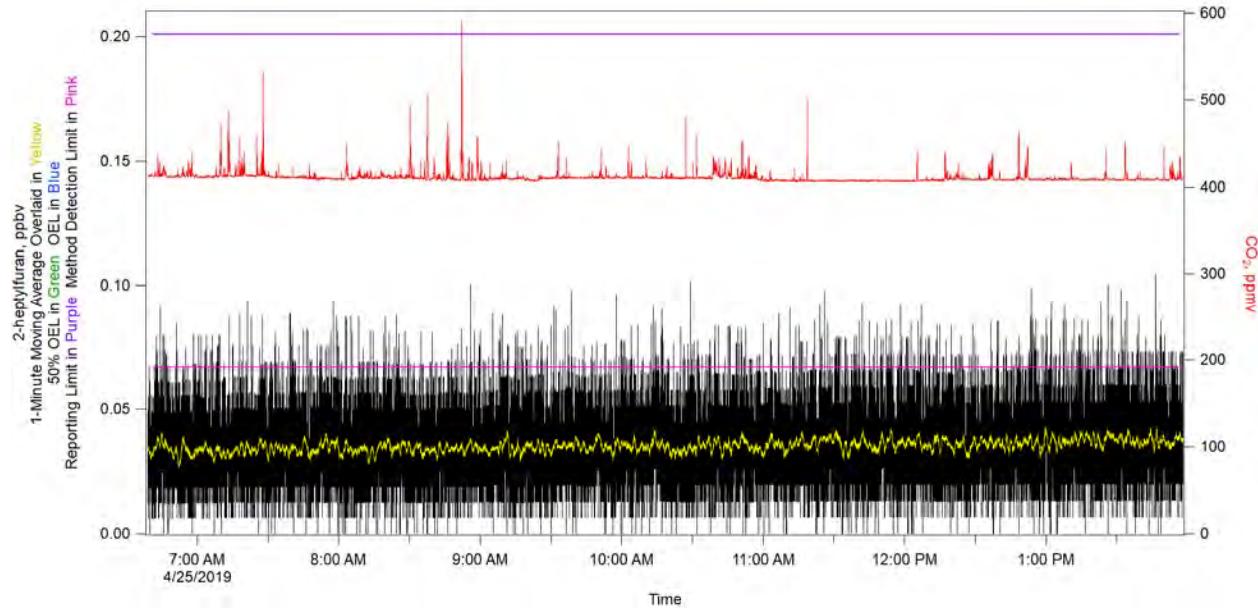


Figure 5-42. 2-heptylfuran.

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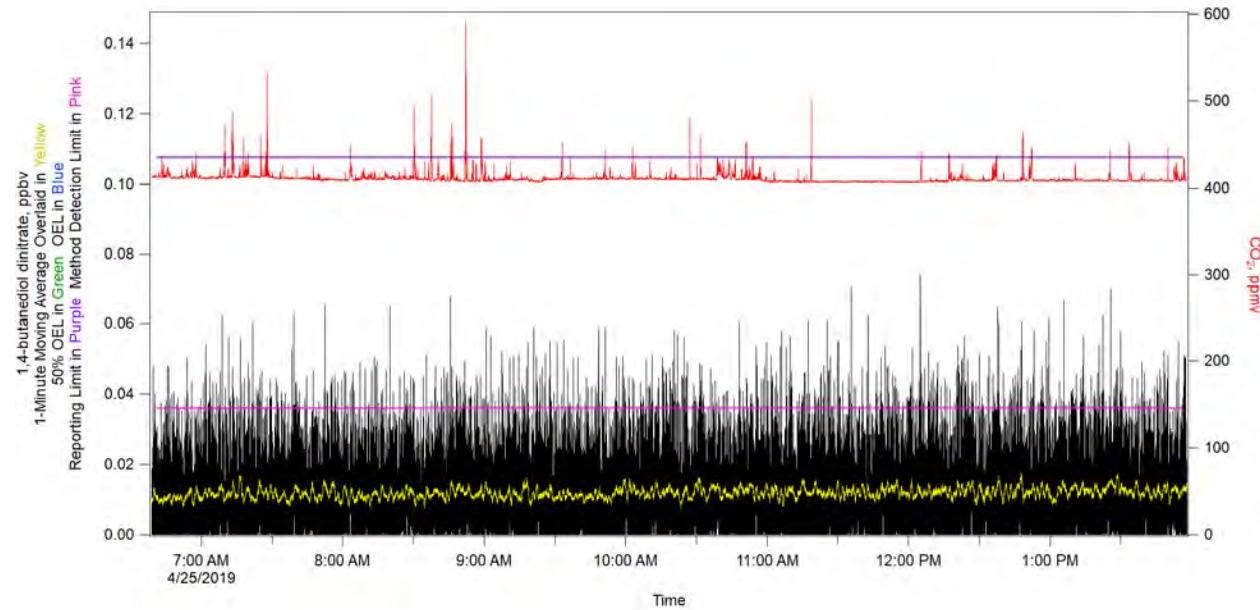


Figure 5-43. 1,4-butanediol Dinitrate.

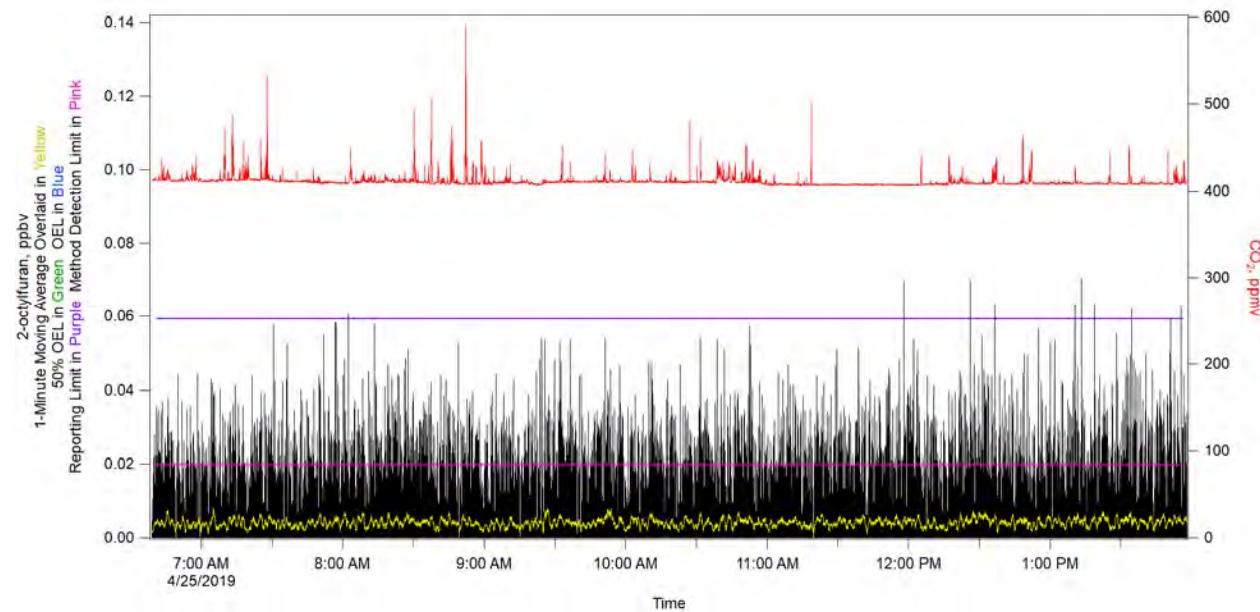


Figure 5-44. 2-octylfuran.

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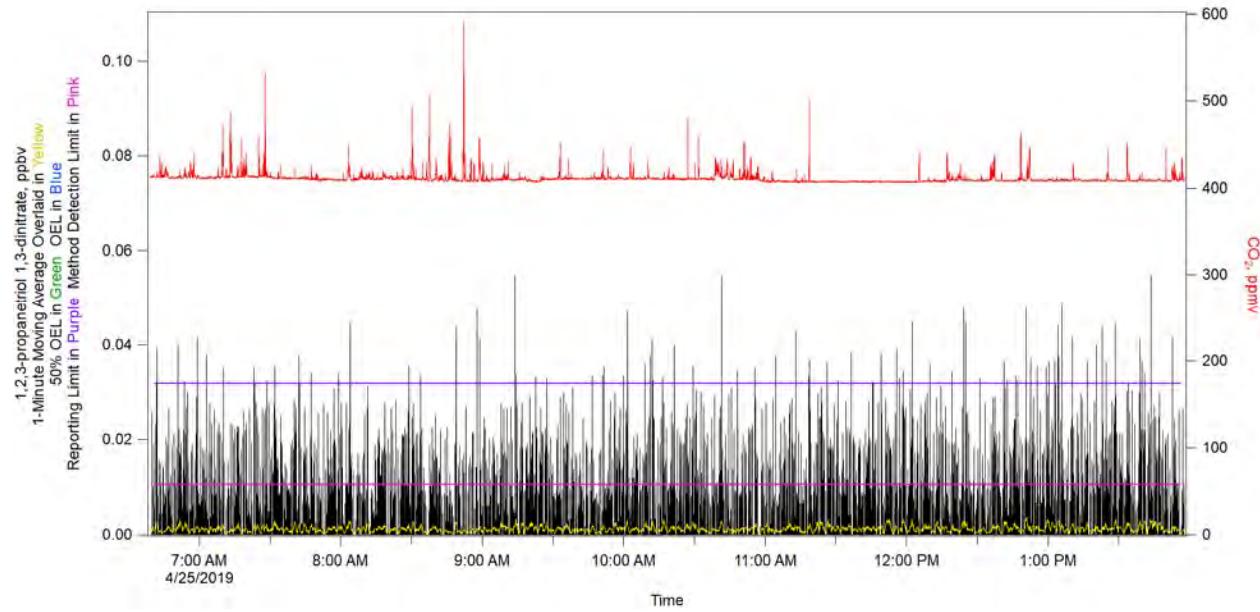


Figure 5-45. 1,2,3-propanetriol 1,3-dinitrate.

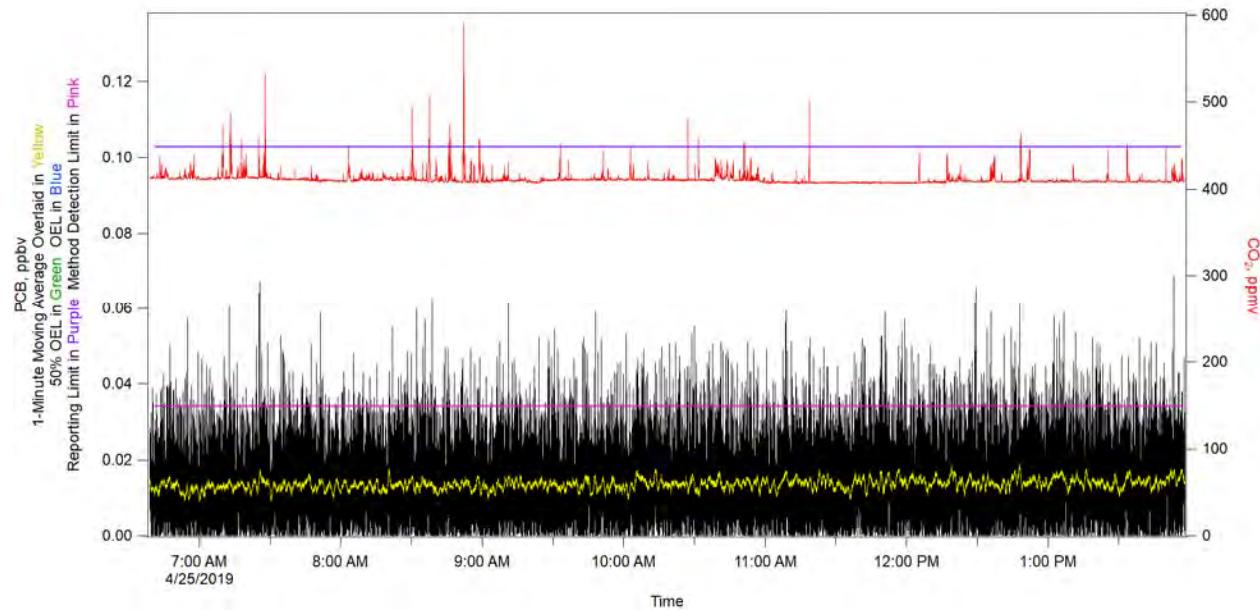


Figure 5-46. PCB.

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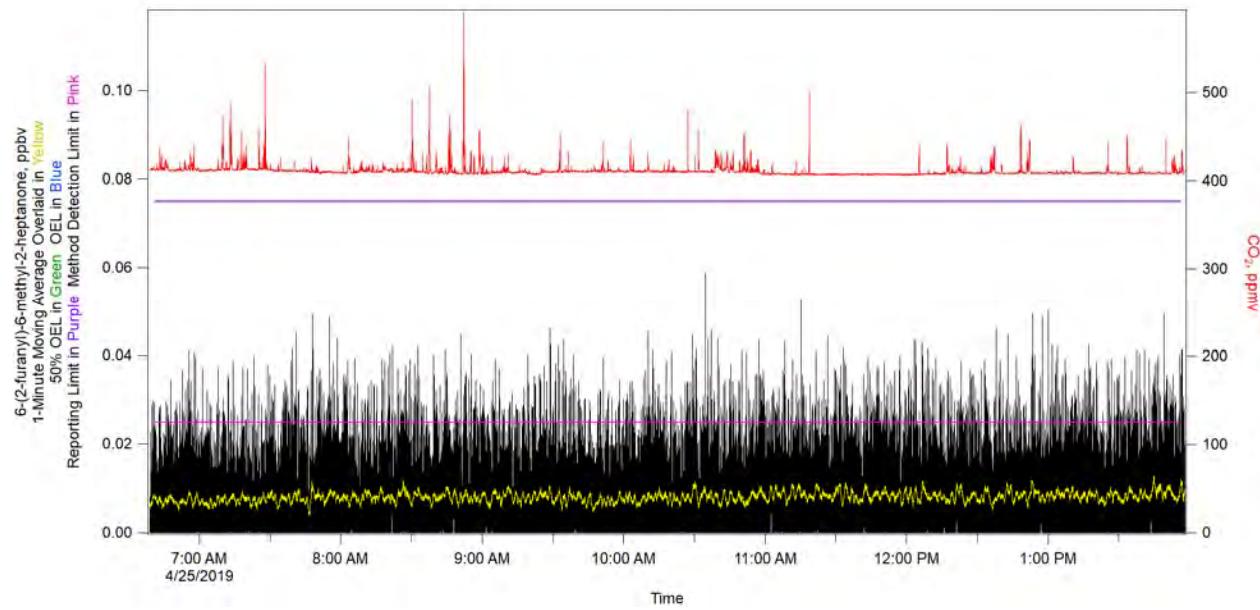


Figure 5-47. 6-(2-furanyl)-6-methyl-2-heptanone.

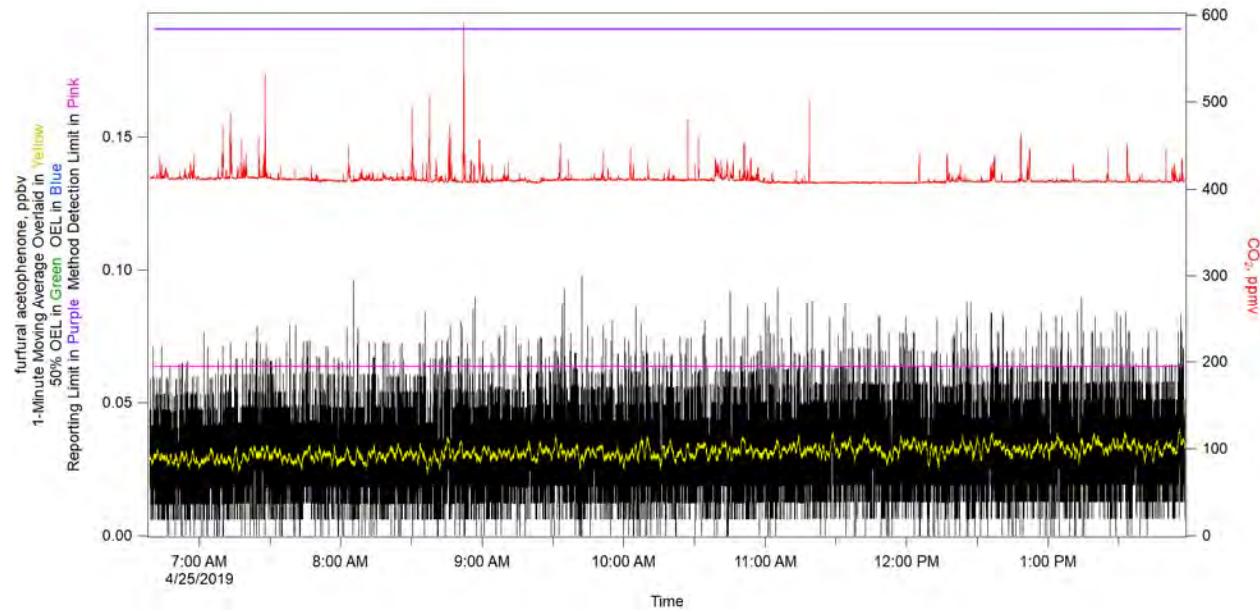


Figure 5-48. Furfural Acetophenone.

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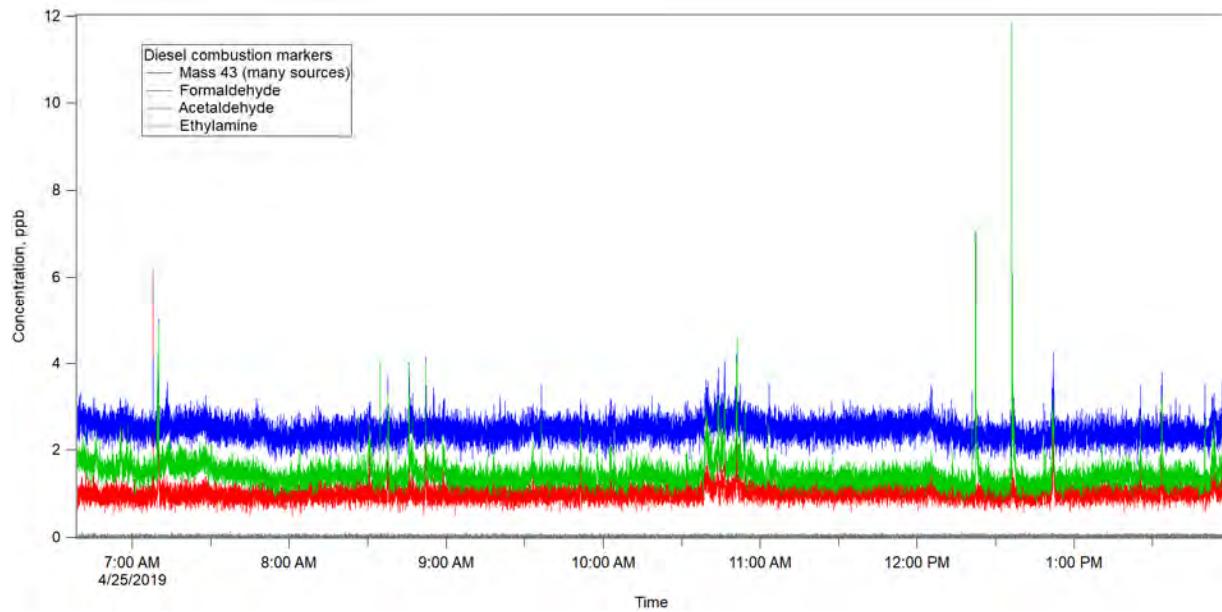


Figure 5-49. Diesel Combustion Markers.

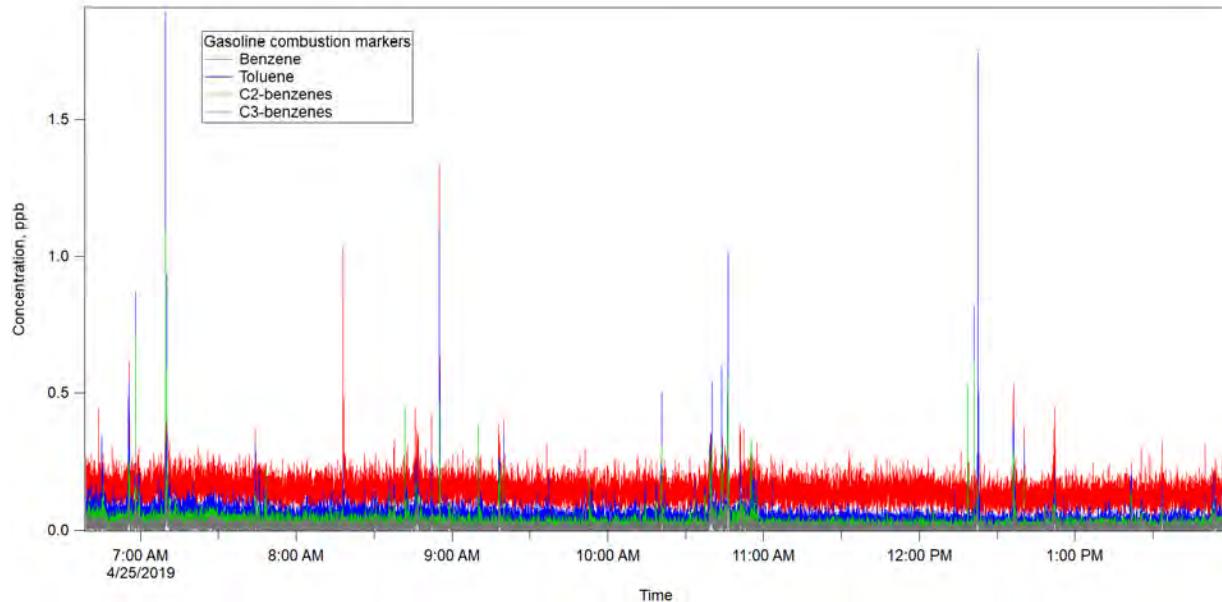


Figure 5-50. Gasoline Combustion Markers.

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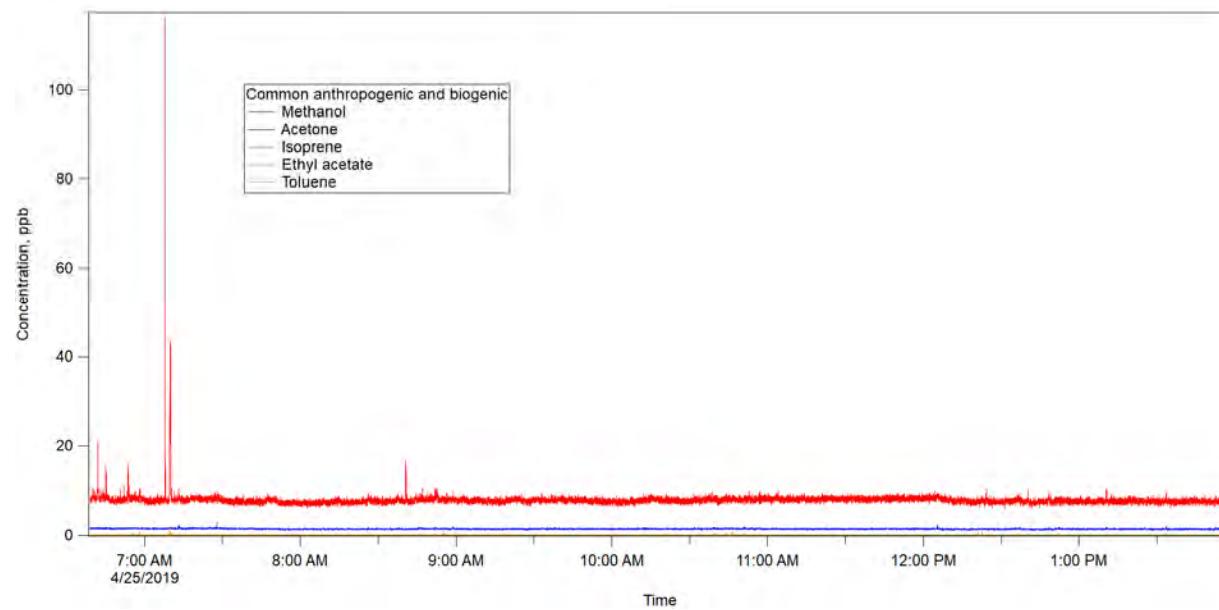


Figure 5-51. Plant and Human Markers.

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6.0 APRIL 26, 2019 – TESTING

6.1 Summary

On April 26, 2019, the ML personnel arrived at the TerraGraphics warehouse at 05:50. An ML Operator initiated a zero-air and span check on the PTR-MS using the volatile organic compound (VOC) gas standard (ID: 160-401380144-1) at 05:59. Figure 6-1 shows the response of toluene for the duration of the days testing. Once complete, the operator switched to the R&D gas standard and began a multipoint calibration at 06:36. The Operator connected calibration box (ID: CZ-MHE-001) line to the mast and began a 2-hour, zero-air check at 07:42. The purpose of the long zero is to provide a better data set for calculation of detection limits. After this, the R&D cylinder (ID: 160-401265983-1) was set to a flow of ~40 sccm and was allowed time to mix with the zero-air until all the species reached a stable concentration. The Operator connected two Carbotraps^{®4} to the sampling system and initiated sampling at 10:30. After an hour of sampling, the Carbotraps were disconnected from the sampling line and a Chain of Custody (COC) was completed. The purpose of this test was to collect a known amount of mass onto the Carbotraps and the results from the ML can be compared to this known mass and the PTR-MS values that were recorded simultaneously. The Operator replaced the almost empty zero-air gas standard with a new zero-air gas standard at 11:45. The ML was prepped for the April 29, 2019, deployment prior to ML personnel departing the TerraGraphics warehouse for the day.

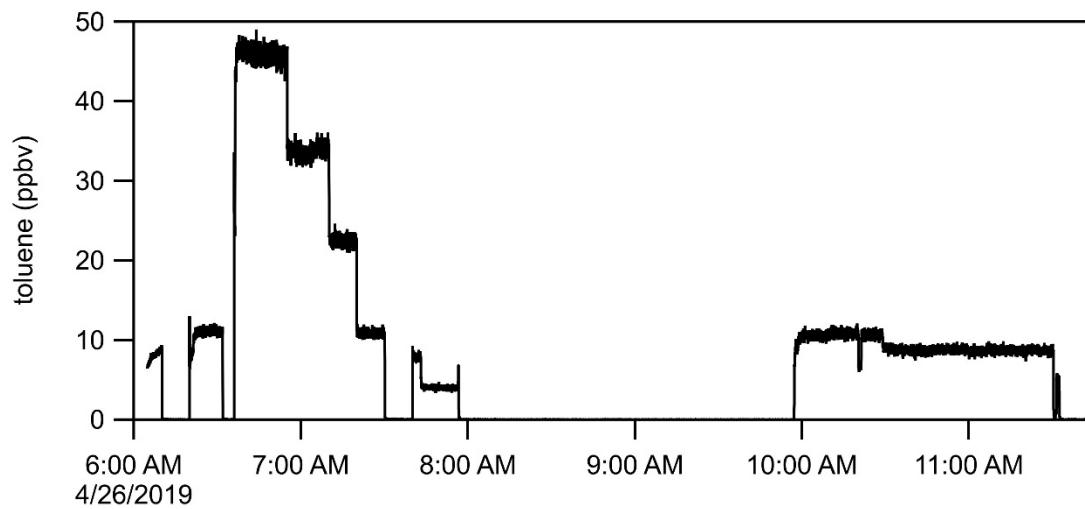


Figure 6-1. Time-series of Toluene During the Testing Performed on April 26, 2019.

⁴ Carbotrap is a registered trademark of Sigma-Aldrich Co., LLC, St. Louis, Missouri.

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7.0 ZERO-AIR AND SENSITIVITY VERIFICATION

Tables 7-1 through 7-6 display the zero-air and sensitivity checks.

Table 7-1. Zero-air Checks for the LI-COR CO₂ Monitor.

Date	Time	Instrument Check	Observed Result (ppm)	Expected Result (ppm)	% Difference	Acceptance Criteria (%)	Pass/Fail
04/22/19	05:43	Zero	0.009	<50	N/A	N/A	Pass
04/23/19	05:34	Zero	0.041	<50	N/A	N/A	Pass
04/24/19	05:46	Zero	0.198	<50	N/A	N/A	Pass
04/25/19	05:56	Zero	0.282	<50	N/A	N/A	Pass

Table 7-2. Span Checks for the LI-COR CO₂ Monitor.

Date	Time	Instrument Check	Observed Result (ppm)	Expected Result (ppm)	% Difference	Acceptance Criteria (%)	Pass/Fail
04/22/19	05:53	Span	361	385	-6.2	20	Pass
04/23/19	05:36	Span	365	384	-5.1	20	Pass
04/24/19	05:49	Span	365	383	-4.9	20	Pass
04/25/19	05:59	Span	365	384.7	-5.2	20	Pass

Table 7-3. Zero-air Checks for the Proton Transfer Reaction - Time-of-Flight.

Date	Time	Instrument Check	Observed Result (ppb)	Expected Result (ppb)	% Difference	Acceptance Criteria (%)	Pass/Fail
04/22/19	05:53	Zero	0.08	<0.5 ppb	N/A	N/A	Pass
04/23/19	05:46	Zero	0.075	<0.5 ppb	N/A	N/A	Pass
04/24/19	05:59	Zero	0.075	<0.5 ppb	N/A	N/A	Pass
04/25/19	06:09	Zero	0.07	<0.5 ppb	N/A	N/A	Pass

Table 7-4. Span Checks for the Proton Transfer Reaction - Time-of-Flight .

Date	Time	Instrument Check	Observed Result (ppb)	Expected Result (ppb)	% Difference	Acceptance Criteria (%)	Pass/Fail
04/22/19	06:03	Span	9.5	10.8	12	30	Pass
04/23/19	05:56	Span	9.75	10.8	9.7	30	Pass
04/24/19	06:10	Span	10	10.8	7.4	30	Pass
04/25/19	06:20	Span	9.8	10.8	9	30	Pass

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Table 7-5. Zero-air Checks for the Picarro.

Date	Time	Instrument Check	Observed Result (ppb)	Expected Result (ppb)	% Difference	Acceptance Criteria (%)	Pass/Fail
04/22/19	05:23	Zero	6.6	< 20 ppb	N/A	N/A	Pass
04/23/19	05:17	Zero	6.8	< 20 ppb	N/A	N/A	Pass
04/24/19	05:29	Zero	7.1	< 20 ppb	N/A	N/A	Pass
04/25/19	05:37	Zero	7.0	< 20 ppb	N/A	N/A	Pass

Table 7-6. Span Checks for the Picarro.

Date	Time	Instrument Check	Observed Result (ppb)	Expected Result (ppb)	% Difference	Acceptance Criteria (%)	Pass/Fail
04/22/19	05:36	Span	3650	3250	12.3	20	Pass
04/23/19	05:17	Span	3680	3250	13.0	20	Pass
04/24/19	05:42	Span	3690	3250	13.5	20	Pass
04/25/19	05:51	Span	3636	3250	11.9	20	Pass

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8.0 DATA PROCESSING AND REPORTING

During the Week of April 22, 2019 through April 26, 2019, the data processing team continued processing data from the previous week and current week. The reporting team worked towards the completion of reports for Week 34, Week 36, Week 37, Week 38, Month 5, Month 6, Month 7, and the report for the AOP-015 Event. The reporting team submitted the draft weekly summary report, 53005-81-RPT-046, *Weekly Report for Week 33 (March 18, 2019 – March 22, 2019)*.

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9.0 REFERENCES

17124-DOE-HS-102, 2018, "Mobile Laboratory Data Processing – Analysis," Revision 2, TerraGraphics Environmental Engineering, Inc., Pasco, Washington.

53005-81-COM-0519-005, *Subcontract 53005, Release 81 – Transmittal of Revision 1 of Special Communication Report for (AOP)-015 Event on April 22, 2019*, TerraGraphics Environmental Engineering, Inc., Pasco, Washington.

53005-81-RPT-007, 2018, *PTR-MS Mobile Laboratory Vapor Monitoring Background Study, (3/18/2018 – 4/20/2018)*, Revision 0, TerraGraphics Environmental Engineering, Inc., Pasco, Washington.

66409-RPT-004, 2019, *Mobile Laboratory Operational Procedure*, Revision 11, TerraGraphics Environmental Engineering, Inc., Pasco, Washington.

Fiscal Year 2017 Mobile Laboratory Vapor Monitoring at the Hanford Site: Monitoring During Waste Disturbing Activities and Background Study, 2017, RJ Lee Group, Inc., Pasco, Washington.

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APPENDIX A

**53005-81-COM-0519-005, REVISION 1 OF SPECIAL COMMUNICATION REPORT
FOR (AOP)-015 EVENT ON APRIL 22, 2019**

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www.terrographics.com

VIA E-MAIL – TOCVND@rl.gov / Kyle_R_Dickman@rl.gov

53005-81-COM-0519-005

Revision 1

May 9, 2019

Kyle Dickman
 Buyer's Technical Representative
 Washington River Protection Solutions, LLC
 Post Office Box 850
 Richland, Washington 99352

Dear Mr. Dickman:

SUBJECT: SUBCONTRACT 53005, RELEASE 81 – TRANSMITTAL OF REVISION 1 OF THE SPECIAL COMMUNICATION REPORT FOR AOP-015 EVENT ON APRIL 22, 2019

Reference: Letter, From Mr. Rich Westberg to Mr. Kyle Dickman, *Subcontract 53005, Release 81 – Transmittal of Special Communication Report for AOP-015 Event on April 22, 2019*, 53005-81-COM-0519-005, dated May 6, 2019.

TerraGraphics is pleased to transmit Revision 1 of the attached letter report for the Abnormal Operating Procedure (AOP)-015 event that took place on April 22, 2019. There was a mass calibration issue discovered in the referenced report above that has now been corrected in the attached letter report. Please accept our apologies for this oversight.

Please note that TOCVND@rl.gov is included as a recipient of this submittal for reference only. This submittal is not listed on the Master Submittal Register (MSR) for this task.

Thank you for the opportunity to support this task. If you have any questions, please feel free to contact me at (509) 547-3883.

Sincerely,

Rich Westberg
 Business Unit Manager, Health, Safety, & Analytical Services
 Pasco Office

/kls

cc:

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Attachment 1

LETTER REPORT FOR AOP-015 EVENT ON APRIL 22, 2019

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1.0 INTRODUCTION

On April 22, 2019, the TerraGraphics Mobile Laboratory (ML) was deployed to the Hanford Site for area monitoring around the 200 East location. The ML's objective was to perform vapor monitoring for a range of compounds, primarily focused on chemicals of potential concern (COPCs) and odor-causing compounds. At approximately 9:18 Pacific Standard Time (PST), an Abnormal Operating Procedure (AOP)-015 event was reported near the U-Farm change trailer. There were no significant responses of COPCs or odor compounds, with levels remaining below occupational exposure limits (OELs) or at typical background levels at the initial time of the AOP-015 event.

The ML operators were notified of the AOP-015 event by Central Shift Office (CSO) announcement at 9:18 PST, while monitoring the northwest side of 241-A Farm, and immediately headed to the CSO for briefing. By 9:50 PST, following further instruction from the CSO, the ML operators traveled to 200 West and positioned approximately 500 feet to the south of the entrance of the U-Farm change trailer. At this time, the ML operators set up the 208-foot heated inlet line for vapor monitoring. The ML was instructed by the CSO to maintain its position. After the arrival of the Industrial Hygiene Technicians (IHTs) at approximately 10:17 PST, the heated inlet line was extended north by the IHTs toward the U-Farm change trailer with the end of the line falling less than halfway between the ML position and the source of the AOP-015 event. The ML operators continued vapor monitoring in this position, sampling from the 208-foot heated inlet line, until 14:02 PST, at which time the CSO announced an exit from the AOP-015. During this monitoring period, the wind direction was primarily based from the south. Consequently, the ML inlet was positioned upwind and 300 feet away from the U-Farm change trailer for the majority of the monitoring.

2.0 April 22, 2019 – AOP-015 EVENT

2.1 Quality Assessment

Data from April 16, 2019, were assessed using Procedure 17124-DOE-HS-102, "Mobile Laboratory Data Processing-Analysis." A Data Acceptance Checklist was completed. The data were accepted by TerraGraphics with the following comments. Report No. 66409-RPT-004, *Mobile Laboratory Operational Procedure*, was adequately documented and all checks passed the acceptance limits.

2.2 Summary

Figure 2-1 below shows the position of the ML and its 208-foot heated line inlet for the duration of the monitoring period.

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Figure 2-1. Position of the Mobile Laboratory for the Duration of Monitoring (10:17 PST – 14:02 PST).

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Figure 2-2 below shows meteorological data from the ML mast for the duration of the monitoring period.

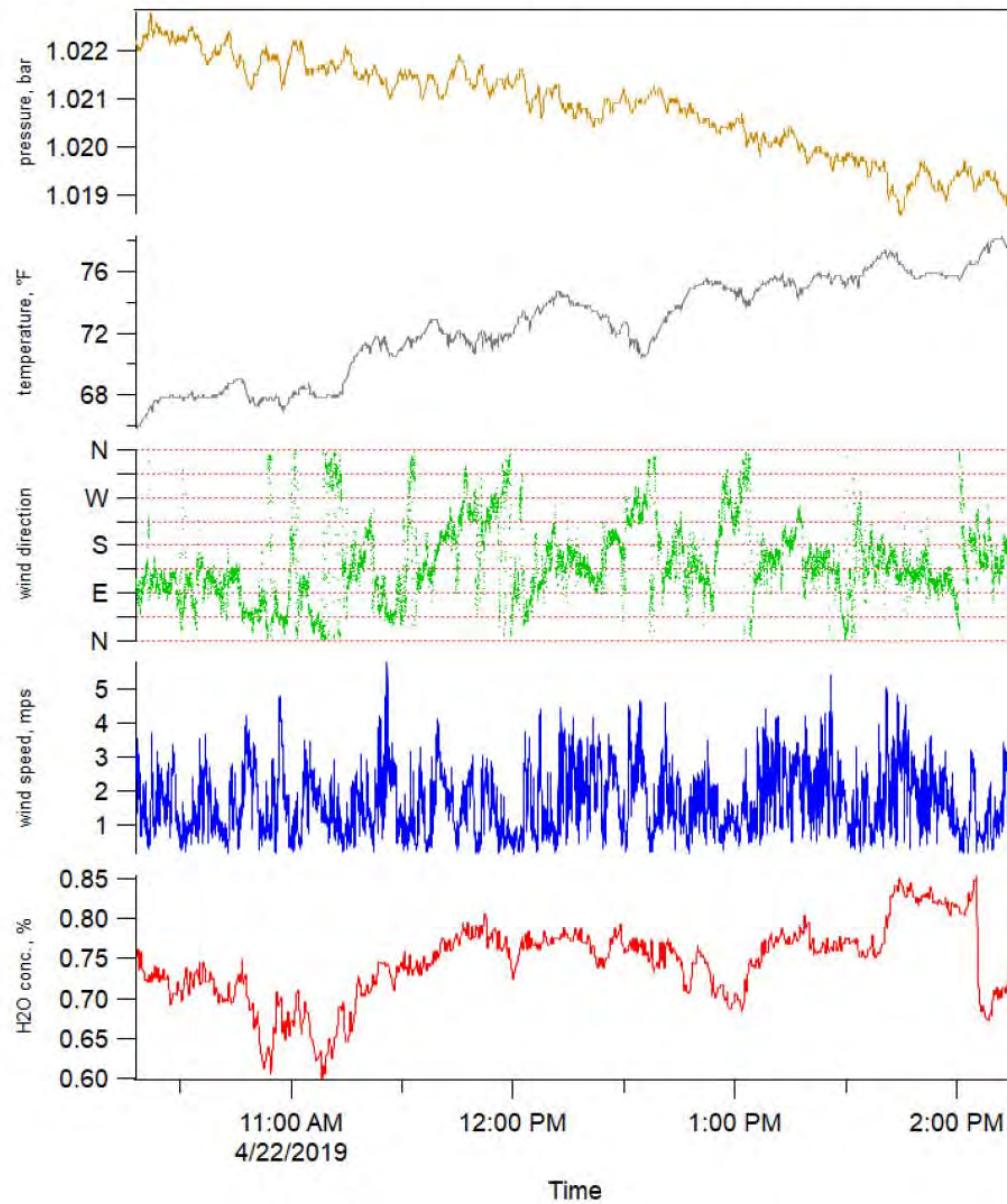


Figure 2-2. Meteorological Data from Mobile Laboratory Mast for the Duration of Monitoring.

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2.3 Chemicals of Potential Concern

Table 2-1. Chemical of Potential Concern Statistical Information for the Monitoring Period of April 22, 2019. (2 Sheets)

COPC #	COPC Name	OEL (ppb)	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel. St. Dev. %	Max (ppb)	Median (ppb)
1	Ammonia	25000	6.225	8.602†	0.401	4.658	10.190	8.579†
2	Formaldehyde	300	1.721	2.225†	0.291	13.097	3.324	2.22†
3	Methanol	200000	5.758	20.292	2.157	10.629	26.112	20.378
4	Acetonitrile	20000	0.085	0.121†	0.033	27.005	0.260	<0.119†
5	Acetaldehyde	25000	1.027	2.438†	0.302	12.387	3.625	2.425†
6	Ethylamine	5000	0.069	<0.069	0.024	58.427	0.166	<0.069
7	1,3-butadiene	1000	0.183	0.285†	0.078	27.319	0.714	0.282†
8	Propanenitrile	6000	0.107	<0.107	0.030	38.903	0.226	<0.107
9	2-propenal	100	0.340	<0.34	0.049	30.420	0.398	<0.34
10	1-butanol + butenes	20000	0.214	<0.214	0.050	31.860	0.410	<0.214
11	methyl isocyanate	20	0.069	0.083†	0.035	42.416	0.266	0.08†
12	methyl nitrite	100	0.098	0.217†	0.050	22.983	0.433	0.217†
13	Furan	1	0.062	<0.062	0.022	44.090	0.155	<0.062
14	Butanenitrile	8000	0.039	<0.039	0.021	58.294	0.147	<0.039
15	but-3-en-2-one + 2,3-dihydrofuran + 2,5-dihydrofuran	100, 1, 1	0.041	0.051†	0.019	38.469	N/A*	N/A*
16	Butanal	25000	0.061	0.264	0.054	20.465	0.518	0.262
17	NDMA**	0.3	0.082	<0.082	0.030	198.986	0.269	<0.082
18	Benzene	500	0.236	0.413†	0.075	18.184	0.699	0.412†
19	2,4-pentadienenitrile + pyridine	300, 1000	0.085	<0.085	0.024	36.674	0.192	<0.085
20	2-methylene butanenitrile	30	0.036	<0.036	0.014	59.873	0.103	<0.036
21	2-methylfuran	1	0.043	0.046†	0.024	52.210	0.178	0.044†
22	Pantanenitrile	6000	0.036	<0.036	0.016	67.023	0.119	<0.036
23	3-methyl-3-buten-2-one + 2-methyl-2-butenal	20, 30	0.043	<0.043	0.024	58.595	0.161	<0.043
24	NEMA**	0.3	0.058	<0.058	0.029	154.500	0.181	<0.058
25	2,5-dimethylfuran	1	0.032	<0.032	0.020	72.220	0.125	<0.032
26	Hexanenitrile	6000	0.031	<0.031	0.014	77.287	0.098	<0.031
27	2-hexanone (MBK)	5000	0.036	<0.036	0.021	66.718	0.149	<0.036
28	NDEA**	0.1	0.034	<0.034	0.016	134.378	0.107	<0.034

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Table 2-1. Chemical of Potential Concern Statistical Information for the Monitoring Period of April 22, 2019. (2 Sheets)

COPC #	COPC Name	OEL (ppb)	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel. St. Dev. %	Max (ppb)	Median (ppb)
29	butyl nitrite + 2-nitro-2-methylpropane	100, 30	0.058	<0.058	0.020	39.035	0.132	<0.058
30	2,4-dimethylpyridine	500	0.036	<0.036	0.016	76.457	0.118	<0.036
31	2-propylfuran + 2-ethyl-5-methylfuran	1	0.027	<0.027	0.014	113.359	0.102	<0.027
32	Heptanenitrile	6000	0.027	<0.027	0.012	81.009	0.084	<0.027
33	4-methyl-2-hexanone	500	0.033	<0.033	0.015	81.918	0.092	<0.033
34	NMOR**	0.6	0.021	<0.021	0.012	200.988	0.093	<0.021
35	butyl nitrate	2500	0.022	<0.022	0.010	110.338	0.067	<0.022
36	2-ethyl-2-hexenal + 4-(1-methylpropyl)-2,3-dihydrofuran + 3-(1,1-dimethylethyl)-2,3-dihydrofuran	100,1,1	0.028	<0.028	0.014	80.121	0.100	<0.028
37	6-methyl-2-heptanone	8000	0.028	<0.028	0.013	81.765	0.089	<0.028
38	2-pentylfuran	1	0.026	<0.026	0.013	81.122	0.078	<0.026
39	Biphenyl	200	0.022	<0.022	0.014	106.903	0.079	<0.022
40	2-heptylfuran	1	0.067	<0.067	0.020	33.915	0.150	<0.067
41	1,4-butanediol dinitrate	50	0.036	<0.036	0.016	64.364	0.101	<0.036
42	2-octylfuran	1	0.020	<0.02	0.010	204.239	0.093	<0.02
43	1,2,3-propanetriol 1,3-dinitrate	50	0.011	<0.011	0.010	264.422	0.080	<0.011
44	PCB	1000	0.034	<0.034	0.016	53.239	0.109	<0.034
45	6-(2-furanyl)-6-methyl-2-heptanone	1	0.025	<0.025	0.010	93.233	0.080	<0.025
46	furfural acetophenone	1	0.064	<0.064	0.019	36.634	0.135	<0.064
N/A*	The maximum peak value for but-3-en-2-one + 2,3 dihydrofuran + 2,5 dihydrofuran was 0.143 ppb and the median value was 0.049† ppb. The PTR-MS results for but-3-en-2-one + 2,3 dihydrofuran + 2,5 dihydrofuran are not compared to OEL concentrations because: 1) the result is suspect due to a known biogenic interferant (methacrolein) that is expected to be in concentrations that occasionally exceed the dihydrofuran OEL, and 2) this combination of COPCs have OEL concentrations that differ by a factor of 200, which provide widely variant bases for these numbers.							
**	Nitrosamine results are suspect due to isobaric interferants causing positive bias that have been encountered during previous background studies the Spring 2018 background study [53005-81-RPT-007, PTR-MS Mobile Laboratory Vapor Monitoring Background Study, (3/18/2018 – 4/20/2018), and Fiscal Year 2017 Mobile Laboratory Vapor Monitoring at the Hanford Site: Monitoring During Waste Disturbing Activities and Background Study, RJ Lee Group, Inc., 2017].							
<	COPC Average/Median Below the MDL.							
†	COPC Average/Median Between the MDL and the RL (i.e., 3*MDL).							

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2.4 Odor Compounds

Table 2-2. Odor Statistical Information for the Monitoring Period of April 22, 2019.

Odor #	Odor Name	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel. St. Dev. (ppb)	Max (ppb)	Median (ppb)
1	hydrogen sulfide	2.801	<2.801	0.199	8.924	3.108	<2.801
2	methyl mercaptan	0.147	<0.147	0.044	30.407	0.366	<0.147
3	dimethyl sulfide; ethanethiol	0.088	0.193†	0.045	23.122	0.402	0.193†
4	allyl mercaptan	0.033	<0.033	0.013	201.445	0.126	<0.033
5	1-propanethiol; isopropyl mercaptan	0.041	<0.041	0.023	139.480	0.182	<0.041
6	2-butene-1-thiol	0.062	<0.062	0.012	263.193	0.147	<0.062
7	diethyl sulfide; 2-methylpropane-2-thiol	0.281	0.373†	0.089	23.980	0.824	0.365†
8	thiopropanal sulfuroxide	0.024	0.126	0.051	40.564	0.353	0.123
9	dimethyl disulfide	0.027	0.053†	0.030	57.077	0.203	0.051†
10	1-pentanethiol; 2,2-dimethylpropane-1-thiol	0.046	<0.046	0.019	121.401	0.121	<0.046
11	benzenethiol	0.028	0.107	0.057	52.758	0.320	0.105
12	diallyl sulfide	0.022	<0.022	0.013	193.852	0.130	<0.022
13	methyl propyl disulfide	0.016	0.026†	0.032	124.046	0.187	0.011†
14	methylbenzenethiol	0.027	<0.027	0.017	114.929	0.111	<0.027
15	dimethyl trisulfide	0.020	<0.02	0.010	157.053	0.069	<0.02
16	(1-oxoethyl) thiophene	0.040	<0.04	0.024	170.062	0.164	<0.04
17	(1-oxopropyl) thiophene	0.029	0.031†	0.024	75.714	0.138	0.028†
18	dipropyl disulfide	0.024	<0.024	0.013	102.344	0.087	<0.024
19	methyl propyl trisulfide	0.019	<0.019	0.013	160.881	0.089	<0.019
20	dimethyl tetrasulfide	0.021	<0.021	0.009	92.339	0.057	<0.021
21	dipropyl trisulfide	0.022	<0.022	0.013	177.118	0.096	<0.022
22	diphenyl sulfide	0.026	<0.026	0.014	137.747	0.099	<0.026
<	COPC Average/Median Below the MDL.						
†	COPC Average/Median Between the MDL and the RL (i.e., 3*MDL).						

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2.5 All Other Mass Signals Detected by Proton Transfer Reaction – Mass Spectrometer

Table 2-3. Other Mass Signal Statistical Information for the Monitoring Period of April 22, 2019. (12 Sheets)

#	Species Name	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	ReL St. Dev. (ppb)	Max (ppb)	Median (ppb)
1	nominal m/z 26	0.189	<0.189	0.056	41.944	0.423	<0.189
2	nominal m/z 27	0.435	<0.435	0.092	24.673	0.792	<0.435
3	nominal m/z 36	1.631	1.911†	0.184	9.616	2.648	1.909†
4	nominal m/z 38	0.678	4.531	0.383	8.460	5.972	4.54
5	nominal m/z 39	1.530	6.899	0.434	6.295	8.448	6.91
6	nominal m/z 40	0.272	0.332†	0.073	22.025	0.629	0.332†
7	nominal m/z 41	0.551	0.747†	0.129	17.318	1.270	0.743†
8	unknown m/z 42	0.298	0.304†	0.081	26.649	0.667	0.3†
9	nominal m/z 43	1.411	7.81	0.803	10.281	10.266	7.872
10	nominal m/z 44	0.708	0.715†	0.104	14.588	1.126	0.713†
11	nominal m/z 45	5.841	6.62†	1.204	18.194	12.005	6.378†
12	nominal m/z 46	2.018	3.023†	0.429	14.181	8.074	2.963†
13	formamide	0.240	0.264†	0.063	23.961	0.523	0.264†
14	formic acid	0.908	7.271	0.629	8.649	9.198	7.331
15	ethanol	0.050	<0.05	0.021	164.305	0.170	<0.05
16	nominal m/z 48	0.267	0.331†	0.067	20.086	0.639	0.331†
17	nominal m/z 50	0.473	0.499†	0.081	16.301	0.850	0.497†
18	nominal m/z 51	0.536	0.615†	0.104	16.938	1.125	0.612†
19	nominal m/z 52	0.588	<0.588	0.060	20.750	0.541	<0.588
20	nominal m/z 53	0.280	0.283†	0.064	22.441	0.623	0.278†
21	nominal m/z 54	0.308	<0.308	0.043	27.589	0.357	<0.308
22	unknown m/z 55	0.091	<0.091	0.040	57.514	0.255	<0.091
23	unknown m/z 58	0.270	<0.27	0.026	49.651	0.191	<0.27
24	C ₃ H ₇ N	0.089	0.093†	0.037	40.220	0.271	0.09†
25	acetone	0.274	6.635	0.775	11.686	9.543	6.659
26	nominal m/z 60	0.292	0.768†	0.110	14.296	1.245	0.766†
27	acetic acid + acetate fragment	0.592	4.247	0.472	11.102	5.907	4.287
28	nominal m/z 64	0.160	0.241†	0.052	21.707	0.484	0.242†
29	nominal m/z 65	0.154	0.291†	0.057	19.611	0.542	0.29†

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Table 2-3. Other Mass Signal Statistical Information for the Monitoring Period of April 22, 2019. (12 Sheets)

#	Species Name	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel. St. Dev. (ppb)	Max (ppb)	Median (ppb)
30	nominal m/z 66	0.126	<0.126	0.032	31.989	0.259	<0.126
31	nominal m/z 67	0.154	0.159†	0.041	25.770	0.361	0.157†
32	nominal m/z 68	0.132	<0.132	0.031	31.179	0.252	<0.132
33	isoprene	0.108	0.115†	0.038	32.727	0.301	0.115†
34	unknown m/z 70	0.032	<0.032	0.016	74.956	0.102	<0.032
35	C ₅ H ₁₀	0.074	<0.074	0.030	52.058	0.198	<0.074
36	unknown m/z 71	0.037	0.071†	0.031	44.067	0.223	0.069†
37	nominal m/z 72	0.211	<0.211	0.035	26.983	0.290	<0.211
38	nominal m/z 73	0.113	<0.113	0.016	57.765	0.111	<0.113
39	C ₃ H ₄ O ₂	0.099	0.182†	0.051	28.271	0.401	0.18†
40	nominal m/z 74	0.226	<0.226	0.043	23.023	0.361	<0.226
41	methyl acetate	0.056	0.365	0.081	22.070	0.679	0.363
42	nominal m/z 76	0.245	<0.245	0.035	26.105	0.280	<0.245
43	unknown m/z 77(a)	0.115	0.222†	0.057	25.738	0.475	0.22†
44	unknown m/z 77(b)	0.023	<0.023	0.014	116.799	0.089	<0.023
45	nominal m/z 78	0.236	<0.236	0.040	22.759	0.363	<0.236
46	nominal m/z 81	0.164	0.196†	0.045	22.730	0.409	0.193†
47	C ₆ H ₁₀	0.069	0.104†	0.040	38.336	0.284	0.101†
48	unknown m/z 84	0.039	<0.039	0.016	67.372	0.107	<0.039
49	nominal m/z 85	0.063	0.091†	0.028	30.539	0.211	0.092†
50	C ₄ H ₄ O ₂	0.049	0.131†	0.041	30.951	0.310	0.129†
51	C ₆ H ₁₂	0.049	<0.049	0.020	64.297	0.135	<0.049
52	nominal m/z 86	0.156	<0.156	0.032	26.840	0.250	<0.156
53	nominal m/z 87	0.155	0.357†	0.060	16.685	0.602	0.355†
54	nominal m/z 88	0.224	<0.224	0.031	26.453	0.275	<0.224
55	nominal m/z 89	0.068	<0.068	0.014	58.045	0.091	<0.068
56	ethyl acetate	0.026	0.11	0.052	47.030	0.348	0.109
57	unknown m/z 89	0.065	0.088†	0.034	38.627	0.264	0.086†
58	nominal m/z 90	0.395	<0.395	0.030	27.442	0.231	<0.395
59	unknown m/z 91(a)	0.472	<0.472	0.031	36.891	0.234	<0.472
60	unknown m/z 91(b)	0.037	0.06†	0.039	65.785	0.249	0.056†

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Table 2-3. Other Mass Signal Statistical Information for the Monitoring Period of April 22, 2019. (12 Sheets)

#	Species Name	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel. St. Dev. (ppb)	Max (ppb)	Median (ppb)
61	nominal m/z 92	0.143	0.561	0.109	19.474	1.029	0.551
62	unknown m/z 93	0.047	<0.047	0.040	132.426	0.272	<0.047
63	toluene	0.055	2.88	0.537	18.635	5.244	2.796
64	nominal m/z 94	0.223	0.399†	0.068	17.015	0.726	0.396†
65	unknown m/z 95(a)	0.079	0.176†	0.054	30.663	0.433	0.172†
66	unknown m/z 95(b)	0.063	<0.063	0.030	52.435	0.183	<0.063
67	unknown m/z 95(c)	0.083	0.136†	0.055	40.161	0.381	0.133†
68	nominal m/z 96	0.149	<0.149	0.034	24.573	0.273	<0.149
69	unknown m/z 97	0.036	<0.036	0.020	58.602	0.133	<0.036
70	C ₅ H ₄ O ₂	0.089	<0.089	0.027	48.368	0.186	<0.089
71	C ₇ H ₁₂	0.054	<0.054	0.025	51.552	0.183	<0.054
72	unknown m/z 98(a)	0.040	<0.04	0.013	71.873	0.089	<0.04
73	unknown m/z 98(b)	0.035	<0.035	0.017	69.936	0.117	<0.035
74	unknown m/z 98(c)	0.031	<0.031	0.015	89.979	0.106	<0.031
75	nominal m/z 99	0.149	0.21†	0.042	20.116	0.388	0.211†
76	nominal m/z 100	0.036	0.692	0.099	14.345	1.109	0.691
77	nominal m/z 101	0.089	0.056†	0.021	37.373	0.150	0.055†
78	unknown m/z 101	0.054	0.079†	0.030	37.438	0.223	0.078†
79	C ₅ H ₈ O ₂	0.049	0.081†	0.033	41.112	0.244	0.078†
80	nominal m/z 102	0.172	<0.172	0.029	27.092	0.235	<0.172
81	nominal m/z 103	0.041	0.042†	0.018	43.283	0.148	0.039†
82	ethyl propionate	0.025	<0.025	0.022	94.551	0.125	<0.025
83	unknown m/z 103	0.037	0.067†	0.027	40.100	0.188	0.066†
84	nominal m/z 104	0.142	<0.142	0.022	35.938	0.184	<0.142
85	unknown m/z 105	0.063	0.173†	0.046	26.654	0.390	0.171†
86	styrene	0.030	<0.03	0.010	374.427	0.145	<0.03
87	nominal m/z 106	1.046	<1.046	0.037	21.034	0.353	<1.046
88	nominal m/z 107	0.089	<0.089	0.019	41.078	0.146	<0.089
89	unknown m/z 107	0.079	<0.079	0.031	44.605	0.228	<0.079
90	C ₂ benzenes	0.053	0.116†	0.039	34.088	0.322	0.112†
91	nominal m/z 108	0.206	<0.206	0.022	35.282	0.161	<0.206

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#	Species Name	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel. St. Dev. (ppb)	Max (ppb)	Median (ppb)
92	unknown m/z 108	0.041	<0.041	0.020	53.336	0.126	<0.041
93	nominal m/z 109	0.184	0.559	0.079	14.191	0.898	0.558
94	nominal m/z 110	0.099	<0.099	0.027	29.081	0.235	<0.099
95	unknown m/z 111	0.025	0.076	0.045	59.513	0.280	0.073
96	C ₆ H ₆ O ₂	0.029	<0.029	0.024	130.718	0.170	<0.029
97	C ₈ H ₁₄	0.043	<0.043	0.020	63.358	0.135	<0.043
98	unknown m/z 112	0.032	<0.032	0.014	72.778	0.093	<0.032
99	C ₆ H ₉ NO	0.028	<0.028	0.012	101.557	0.075	<0.028
100	nominal m/z 113	0.180	0.319†	0.054	16.946	0.532	0.317†
101	nominal m/z 114	0.133	<0.133	0.030	26.907	0.239	<0.133
102	unknown m/z 115(a)	0.095	<0.095	0.018	66.306	0.117	<0.095
103	unknown m/z 115(b)	0.062	<0.062	0.022	72.684	0.129	<0.062
104	unknown m/z 115(c)	0.027	0.034†	0.023	66.898	0.156	0.032†
105	nominal m/z 116	0.078	<0.078	0.021	35.944	0.163	<0.078
106	C ₆ H ₁₂ O ₂	0.035	0.036†	0.021	58.416	0.141	0.034†
107	unknown m/z 117	0.020	<0.02	0.016	84.175	0.118	<0.02
108	nominal m/z 118	0.149	<0.149	0.025	29.722	0.208	<0.149
109	nominal m/z 119	0.128	0.17†	0.036	21.239	0.334	0.169†
110	nominal m/z 120	0.077	<0.077	0.017	43.792	0.110	<0.077
111	unknown m/z 120	0.027	<0.027	0.013	85.842	0.102	<0.027
112	unknown m/z 121(a)	0.031	0.039†	0.026	67.558	0.154	0.037†
113	unknown m/z 121(b)	0.036	0.274	0.065	23.765	0.558	0.272
114	C ₃ benzenes	0.033	<0.033	0.018	106.124	0.145	<0.033
115	nominal m/z 122	0.088	0.153†	0.034	22.366	0.291	0.153†
116	unknown m/z 123(a)	0.054	<0.054	0.031	68.653	0.189	<0.054
117	unknown m/z 123(b)	0.054	<0.054	0.031	68.653	0.189	<0.054
118	unknown m/z 123(c)	0.033	<0.033	0.016	62.838	0.117	<0.033
119	nominal m/z 124	0.202	0.274†	0.045	16.454	0.446	0.271†
120	unknown m/z 125(a)	0.025	<0.025	0.013	84.950	0.080	<0.025
121	unknown m/z 125(b)	0.025	<0.025	0.013	84.950	0.080	<0.025
122	unknown m/z 125(c)	0.027	<0.027	0.015	76.451	0.095	<0.027

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#	Species Name	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel. St. Dev. (ppb)	Max (ppb)	Median (ppb)
123	nominal m/z 126	0.101	0.128†	0.031	23.986	0.246	0.125†
124	unknown m/z 127(a)	0.019	0.103	0.043	41.604	0.281	0.104
125	unknown m/z 127(b)	0.028	<0.028	0.015	89.940	0.096	<0.028
126	nominal m/z 128	0.086	<0.086	0.023	30.583	0.173	<0.086
127	unknown m/z 129(a)	0.025	<0.025	0.015	111.770	0.111	<0.025
128	unknown m/z 129(b)	0.035	0.175	0.045	25.913	0.349	0.173
129	naphthalene	0.025	<0.025	0.018	101.843	0.108	<0.025
130	nominal m/z 130	0.081	<0.081	0.023	32.362	0.180	<0.081
131	nominal m/z 131	0.120	0.294†	0.053	18.014	0.529	0.293†
132	nominal m/z 132	0.074	<0.074	0.022	33.778	0.171	<0.074
133	nominal m/z 133	0.072	<0.072	0.022	33.135	0.173	<0.072
134	nominal m/z 134	0.078	<0.078	0.020	34.733	0.164	<0.078
135	C ₄ benzenes	0.052	<0.052	0.019	39.977	0.135	<0.052
136	nominal m/z 136	0.096	0.104†	0.036	34.638	0.256	0.1†
137	nominal m/z 137	0.079	0.098†	0.027	27.966	0.205	0.099†
138	nominal m/z 138	0.068	<0.068	0.020	34.760	0.157	<0.068
139	unknown m/z 139(a)	0.036	0.059†	0.023	38.876	0.185	0.058†
140	unknown m/z 139(b)	0.024	<0.024	0.011	94.582	0.072	<0.024
141	nominal m/z 140	0.071	0.071†	0.022	31.618	0.162	0.07†
142	unknown m/z 141(a)	0.026	<0.026	0.014	123.920	0.101	<0.026
143	unknown m/z 141(b)	0.026	<0.026	0.014	123.920	0.101	<0.026
144	unknown m/z 141(c)	0.029	<0.029	0.014	82.368	0.086	<0.029
145	nominal m/z 142	0.083	<0.083	0.022	31.427	0.169	<0.083
146	nominal m/z 143	0.075	0.092†	0.026	28.169	0.198	0.091†
147	nominal m/z 144	0.086	<0.086	0.022	31.494	0.174	<0.086
148	nominal m/z 145	0.093	0.094†	0.026	27.283	0.217	0.091†
149	nominal m/z 146	0.064	<0.064	0.019	36.453	0.141	<0.064
150	nominal m/z 147	0.079	0.126†	0.031	24.353	0.258	0.126†
151	nominal m/z 148	0.069	<0.069	0.021	33.503	0.153	<0.069
152	nominal m/z 149	0.078	0.089†	0.026	28.919	0.200	0.09†
153	nominal m/z 150	0.072	<0.072	0.022	31.604	0.173	<0.072

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#	Species Name	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel. St. Dev. (ppb)	Max (ppb)	Median (ppb)
154	unknown m/z 151(a)	0.021	<0.021	0.012	120.573	0.084	<0.021
155	unknown m/z 151(b)	0.021	<0.021	0.012	120.573	0.084	<0.021
156	unknown m/z 151(c)	0.026	<0.026	0.012	97.023	0.103	<0.026
157	nominal m/z 152	0.070	<0.07	0.020	34.568	0.153	<0.07
158	nominal m/z 153	0.074	0.087†	0.025	28.901	0.203	0.084†
159	nominal m/z 154	0.065	<0.065	0.020	36.039	0.153	<0.065
160	unknown m/z 155(a)	0.019	<0.019	0.012	163.973	0.097	<0.019
161	unknown m/z 155(b)	0.019	<0.019	0.012	163.973	0.097	<0.019
162	unknown m/z 155(c)	0.026	<0.026	0.012	91.749	0.094	<0.026
163	nominal m/z 156	0.068	<0.068	0.019	36.554	0.153	<0.068
164	nominal m/z 157	0.078	0.087†	0.025	28.524	0.194	0.084†
165	nominal m/z 158	0.072	<0.072	0.020	34.231	0.144	<0.072
166	unknown m/z 159(a)	0.037	<0.037	0.017	49.893	0.129	<0.037
167	unknown m/z 159(b)	0.029	<0.029	0.014	60.253	0.108	<0.029
168	nominal m/z 160	0.063	<0.063	0.018	37.734	0.130	<0.063
169	nominal m/z 161	0.077	0.085†	0.024	28.699	0.198	0.083†
170	nominal m/z 162	0.064	<0.064	0.018	37.231	0.151	<0.064
171	nominal m/z 163	0.081	<0.081	0.023	30.189	0.191	<0.081
172	nominal m/z 164	0.080	<0.08	0.021	33.408	0.151	<0.08
173	nominal m/z 165	0.099	0.099†	0.027	26.768	0.218	0.096†
174	nominal m/z 166	0.075	<0.075	0.020	34.001	0.157	<0.075
175	nominal m/z 168	0.067	<0.067	0.019	35.223	0.143	<0.067
176	nominal m/z 169	0.074	0.074†	0.023	30.296	0.185	0.075†
177	nominal m/z 170	0.069	<0.069	0.019	35.334	0.144	<0.069
178	nominal m/z 171	0.072	<0.072	0.022	31.337	0.170	<0.072
179	nominal m/z 172	0.070	<0.07	0.019	34.769	0.143	<0.07
180	nominal m/z 173	0.045	<0.045	0.017	39.621	0.116	<0.045
181	nominal m/z 174	0.066	<0.066	0.018	37.418	0.149	<0.066
182	nominal m/z 175	0.073	<0.073	0.022	32.162	0.168	<0.073
183	nominal m/z 176	0.063	<0.063	0.018	37.439	0.142	<0.063
184	nominal m/z 177	0.082	0.083†	0.024	28.702	0.202	0.082†

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#	Species Name	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel. St. Dev. (ppb)	Max (ppb)	Median (ppb)
185	nominal m/z 178	0.072	<0.072	0.020	34.260	0.150	<0.072
186	nominal m/z 179	0.082	<0.082	0.023	29.994	0.177	<0.082
187	nominal m/z 180	0.086	<0.086	0.022	31.341	0.156	<0.086
188	unknown m/z 181(a)	0.026	<0.026	0.013	122.401	0.099	<0.026
189	unknown m/z 181(b)	0.026	<0.026	0.013	122.401	0.099	<0.026
190	unknown m/z 181(c)	0.030	<0.03	0.013	93.774	0.084	<0.03
191	nominal m/z 182	0.078	<0.078	0.021	32.580	0.156	<0.078
192	unknown m/z 183(a)	0.028	<0.028	0.013	97.497	0.088	<0.028
193	unknown m/z 183(b)	0.032	<0.032	0.013	71.171	0.083	<0.032
194	unknown m/z 183(c)	0.025	<0.025	0.016	105.609	0.097	<0.025
195	nominal m/z 184	0.091	<0.091	0.022	30.343	0.175	<0.091
196	nominal m/z 185	0.089	<0.089	0.023	29.560	0.180	<0.089
197	nominal m/z 186	0.084	<0.084	0.021	32.072	0.155	<0.084
198	unknown m/z 187(a)	0.030	<0.03	0.016	80.055	0.099	<0.03
199	unknown m/z 187(b)	0.034	<0.034	0.013	70.090	0.093	<0.034
200	nominal m/z 188	0.095	<0.095	0.023	29.507	0.189	<0.095
201	unknown m/z 189	0.034	<0.034	0.013	63.604	0.089	<0.034
202	nominal m/z 190	0.081	<0.081	0.021	32.287	0.155	<0.081
203	nominal m/z 191	0.086	<0.086	0.024	27.998	0.190	<0.086
204	nominal m/z 192	0.081	<0.081	0.021	32.026	0.169	<0.081
205	nominal m/z 193	0.087	<0.087	0.022	30.117	0.180	<0.087
206	nominal m/z 194	0.082	<0.082	0.021	32.034	0.160	<0.082
207	unknown m/z 195(a)	0.032	0.034†	0.017	50.589	0.115	0.033†
208	unknown m/z 195(b)	0.026	<0.026	0.010	88.190	0.068	<0.026
209	nominal m/z 196	0.075	<0.075	0.020	33.716	0.174	<0.075
210	nominal m/z 197	0.079	<0.079	0.021	32.208	0.168	<0.079
211	nominal m/z 198	0.082	<0.082	0.021	32.168	0.159	<0.082
212	nominal m/z 200	0.081	<0.081	0.021	31.839	0.159	<0.081
213	nominal m/z 201	0.071	<0.071	0.020	34.409	0.167	<0.071
214	nominal m/z 202	0.064	<0.064	0.018	37.915	0.127	<0.064
215	nominal m/z 203	0.045	<0.045	0.015	45.230	0.100	<0.045

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#	Species Name	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel. St. Dev. (ppb)	Max (ppb)	Median (ppb)
216	nominal m/z 205	0.078	<0.078	0.022	30.736	0.168	<0.078
217	nominal m/z 206	0.054	<0.054	0.016	41.373	0.112	<0.054
218	nominal m/z 207	0.040	<0.04	0.014	47.237	0.120	<0.04
219	nominal m/z 208	0.039	<0.039	0.013	50.580	0.087	<0.039
220	nominal m/z 209	0.039	<0.039	0.014	47.364	0.113	<0.039
221	nominal m/z 210	0.038	<0.038	0.013	51.565	0.087	<0.038
222	nominal m/z 211	0.076	<0.076	0.020	32.397	0.160	<0.076
223	nominal m/z 212	0.037	<0.037	0.012	53.991	0.086	<0.037
224	nominal m/z 213	0.043	<0.043	0.016	40.208	0.114	<0.043
225	nominal m/z 214	0.083	<0.083	0.021	31.499	0.165	<0.083
226	nominal m/z 215	0.095	<0.095	0.023	28.719	0.193	<0.095
227	nominal m/z 216	0.078	<0.078	0.020	32.541	0.165	<0.078
228	nominal m/z 217	0.087	<0.087	0.022	30.716	0.166	<0.087
229	nominal m/z 218	0.042	<0.042	0.014	46.371	0.100	<0.042
230	nominal m/z 219	0.050	<0.05	0.016	42.688	0.131	<0.05
231	nominal m/z 220	0.049	<0.049	0.015	43.906	0.099	<0.049
232	nominal m/z 221	0.087	<0.087	0.022	29.644	0.172	<0.087
233	nominal m/z 224	0.040	<0.061	0.017	37.505	0.125	<0.061
234	nominal m/z 225	0.042	<0.04	0.013	50.942	0.093	<0.04
235	nominal m/z 226	0.041	<0.042	0.014	47.620	0.100	<0.042
236	nominal m/z 227	0.038	<0.041	0.013	48.945	0.099	<0.041
237	nominal m/z 228	0.037	<0.038	0.013	49.325	0.087	<0.038
238	nominal m/z 229	0.034	<0.037	0.012	53.325	0.073	<0.037
239	nominal m/z 230	0.034	<0.034	0.012	53.793	0.085	<0.034
240	nominal m/z 231	0.041	<0.034	0.012	56.394	0.073	<0.034
241	nominal m/z 232	0.036	<0.041	0.014	48.680	0.091	<0.041
242	nominal m/z 233	0.036	<0.036	0.012	52.600	0.086	<0.036
243	nominal m/z 234	0.034	<0.036	0.012	52.945	0.091	<0.036
244	nominal m/z 235	0.046	<0.034	0.012	56.086	0.085	<0.034
245	nominal m/z 236	0.043	<0.046	0.014	45.116	0.104	<0.046
246	nominal m/z 237	0.034	<0.043	0.014	48.160	0.092	<0.043

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Table 2-3. Other Mass Signal Statistical Information for the Monitoring Period of April 22, 2019. (12 Sheets)

#	Species Name	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel. St. Dev. (ppb)	Max (ppb)	Median (ppb)
247	nominal m/z 238	0.038	<0.034	0.012	56.296	0.085	<0.034
248	nominal m/z 239	0.047	<0.038	0.012	52.234	0.085	<0.038
249	nominal m/z 240	0.050	<0.047	0.015	44.415	0.117	<0.047
250	nominal m/z 241	0.046	<0.05	0.015	43.688	0.111	<0.05
251	nominal m/z 242	0.042	<0.046	0.014	44.218	0.097	<0.046
252	nominal m/z 243	0.037	<0.042	0.013	48.475	0.085	<0.042
253	nominal m/z 244	0.038	<0.037	0.013	50.302	0.085	<0.037
254	nominal m/z 245	0.042	<0.038	0.013	52.361	0.085	<0.038
255	nominal m/z 246	0.041	<0.042	0.014	47.974	0.098	<0.042
256	nominal m/z 247	0.038	<0.041	0.013	49.239	0.097	<0.041
257	nominal m/z 248	0.043	<0.038	0.013	50.431	0.079	<0.038
258	nominal m/z 249	0.036	<0.043	0.014	47.795	0.092	<0.043
259	nominal m/z 250	0.035	<0.036	0.012	53.962	0.084	<0.036
260	nominal m/z 251	0.036	<0.035	0.012	53.941	0.084	<0.035
261	nominal m/z 252	0.034	<0.036	0.012	53.114	0.079	<0.036
262	nominal m/z 253	0.041	<0.034	0.011	55.941	0.084	<0.034
263	nominal m/z 254	0.037	<0.041	0.013	49.604	0.103	<0.041
264	nominal m/z 255	0.040	<0.037	0.012	53.559	0.085	<0.037
265	nominal m/z 256	0.035	<0.04	0.013	49.329	0.085	<0.04
266	nominal m/z 257	0.043	<0.035	0.012	54.848	0.072	<0.035
267	nominal m/z 258	0.033	<0.043	0.014	47.417	0.098	<0.043
268	nominal m/z 259	0.046	<0.033	0.011	56.427	0.084	<0.033
269	nominal m/z 260	0.036	<0.046	0.014	45.401	0.124	<0.046
270	nominal m/z 261	0.042	<0.036	0.012	53.321	0.084	<0.036
271	nominal m/z 262	0.039	<0.042	0.013	48.120	0.091	<0.042
272	nominal m/z 263	0.041	<0.039	0.013	51.042	0.084	<0.039
273	nominal m/z 264	0.037	<0.041	0.013	48.022	0.091	<0.041
274	nominal m/z 265	0.038	<0.037	0.012	52.020	0.077	<0.037
275	nominal m/z 266	0.038	<0.038	0.013	51.789	0.085	<0.038
276	nominal m/z 267	0.037	<0.038	0.012	50.955	0.090	<0.038
277	nominal m/z 268	0.037	<0.037	0.012	53.095	0.090	<0.037

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Table 2-3. Other Mass Signal Statistical Information for the Monitoring Period of April 22, 2019. (12 Sheets)

#	Species Name	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel. St. Dev. (ppb)	Max (ppb)	Median (ppb)
278	nominal m/z 270	0.039	<0.037	0.012	51.984	0.091	<0.037
279	nominal m/z 271	0.043	<0.039	0.013	50.909	0.097	<0.039
280	nominal m/z 272	0.039	<0.043	0.014	46.535	0.096	<0.043
281	nominal m/z 273	0.041	<0.039	0.013	50.371	0.090	<0.039
282	nominal m/z 274	0.046	<0.041	0.013	48.871	0.103	<0.041
283	nominal m/z 275	0.046	<0.046	0.014	44.914	0.134	<0.046
284	nominal m/z 276	0.031	<0.046	0.014	45.684	0.090	<0.046
285	nominal m/z 277	0.038	<0.031	0.011	60.092	0.071	<0.031
286	nominal m/z 278	0.040	<0.038	0.013	51.868	0.084	<0.038
287	nominal m/z 279	0.033	<0.04	0.013	49.916	0.096	<0.04
288	nominal m/z 280	0.040	<0.033	0.012	56.432	0.083	<0.033
289	nominal m/z 281	0.042	<0.04	0.013	49.719	0.090	<0.04
290	nominal m/z 282	0.031	<0.042	0.013	49.076	0.078	<0.042
291	nominal m/z 283	0.031	<0.031	0.011	59.097	0.064	<0.031
292	nominal m/z 284	0.037	<0.031	0.011	58.821	0.071	<0.031
293	nominal m/z 285	0.048	<0.037	0.012	52.348	0.083	<0.037
294	nominal m/z 286	0.037	<0.048	0.015	43.502	0.097	<0.048
295	nominal m/z 287	0.038	<0.037	0.012	52.596	0.083	<0.037
296	nominal m/z 288	0.037	<0.038	0.012	51.355	0.084	<0.038
297	nominal m/z 289	0.038	<0.037	0.012	52.142	0.095	<0.037
298	nominal m/z 290	0.031	<0.038	0.013	50.316	0.096	<0.038
299	nominal m/z 291	0.044	<0.031	0.011	58.435	0.077	<0.031
300	nominal m/z 292	0.044	<0.044	0.014	45.893	0.089	<0.044
301	nominal m/z 293	0.042	<0.044	0.014	46.947	0.095	<0.044
302	nominal m/z 294	0.036	<0.042	0.013	47.570	0.114	<0.042
303	nominal m/z 295	0.049	<0.036	0.012	53.895	0.089	<0.036
304	nominal m/z 296	0.040	<0.049	0.015	43.398	0.103	<0.049
305	nominal m/z 297	0.073	<0.04	0.013	49.579	0.096	<0.04
306	nominal m/z 298	0.035	<0.073	0.020	32.459	0.158	<0.073
307	nominal m/z 299	0.032	<0.035	0.012	54.028	0.083	<0.035
308	nominal m/z 300	0.031	<0.032	0.011	58.386	0.077	<0.032

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Table 2-3. Other Mass Signal Statistical Information for the Monitoring Period of April 22, 2019. (12 Sheets)

#	Species Name	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel. St. Dev. (ppb)	Max (ppb)	Median (ppb)
309	nominal m/z 301	0.045	<0.031	0.011	58.400	0.076	<0.031
310	nominal m/z 302	0.035	<0.045	0.014	45.644	0.096	<0.045
311	nominal m/z 303	0.044	<0.035	0.012	54.842	0.083	<0.035
312	nominal m/z 304	0.039	<0.044	0.014	46.838	0.090	<0.044
313	nominal m/z 305	0.039	<0.039	0.013	50.691	0.090	<0.039
314	nominal m/z 306	0.039	<0.039	0.013	50.629	0.095	<0.039
315	nominal m/z 307	0.037	<0.039	0.013	50.564	0.090	<0.039
316	nominal m/z 308	0.037	<0.037	0.012	51.132	0.083	<0.037
317	nominal m/z 309	0.044	<0.037	0.012	51.824	0.083	<0.037
318	nominal m/z 310	0.041	<0.044	0.014	46.122	0.090	<0.044
319	nominal m/z 311	0.057	<0.041	0.013	49.083	0.090	<0.041
320	nominal m/z 312	0.048	<0.057	0.016	39.580	0.121	<0.057
321	nominal m/z 313	0.068	<0.048	0.015	44.167	0.102	<0.048
322	nominal m/z 314	0.048	<0.068	0.018	35.107	0.140	<0.068
323	nominal m/z 315	0.050	<0.048	0.015	43.475	0.096	<0.048
324	nominal m/z 316	0.047	<0.05	0.015	42.425	0.107	<0.05
325	nominal m/z 317	0.047	<0.047	0.014	44.753	0.096	<0.047
326	nominal m/z 318	0.048	<0.047	0.014	45.381	0.109	<0.047
327	nominal m/z 319	0.041	<0.048	0.015	44.262	0.103	<0.048
328	nominal m/z 320	0.047	<0.041	0.013	48.309	0.090	<0.041
329	nominal m/z 321	0.035	<0.047	0.014	44.119	0.096	<0.047
330	nominal m/z 322	0.038	<0.035	0.012	54.346	0.083	<0.035
331	nominal m/z 323	0.034	<0.038	0.012	51.646	0.084	<0.038
332	nominal m/z 324	0.038	<0.034	0.012	55.202	0.076	<0.034
333	nominal m/z 325	0.032	<0.038	0.012	52.065	0.089	<0.038
334	nominal m/z 326	0.032	<0.032	0.011	57.514	0.083	<0.032
335	nominal m/z 327	0.032	<0.032	0.011	57.480	0.070	<0.032
336	nominal m/z 328	0.032	<0.032	0.011	56.928	0.077	<0.032
337	nominal m/z 329	0.033	<0.032	0.011	57.293	0.076	<0.032
338	nominal m/z 330	0.039	<0.033	0.011	57.197	0.096	<0.033
339	nominal m/z 332	0.055	<0.039	0.014	44.449	0.108	<0.039

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#	Species Name	MDL (ppb)	Ave. (ppb)	St. Dev. (ppb)	Rel. St. Dev. (ppb)	Max (ppb)	Median (ppb)
340	nominal m/z 333	0.032	<0.055	0.016	39.208	0.132	<0.055
341	nominal m/z 334	0.032	<0.032	0.011	56.983	0.070	<0.032
342	nominal m/z 335	0.032	<0.032	0.011	57.294	0.083	<0.032
343	nominal m/z 336	0.033	<0.032	0.011	57.539	0.077	<0.032
344	nominal m/z 337	0.033	<0.033	0.011	56.544	0.076	<0.033
345	nominal m/z 338	0.032	<0.033	0.011	57.067	0.077	<0.033
346	nominal m/z 339	0.032	<0.032	0.011	57.053	0.076	<0.032
347	nominal m/z 340	0.033	<0.032	0.011	56.704	0.076	<0.032
348	nominal m/z 341	0.032	<0.033	0.011	57.538	0.076	<0.033
349	nominal m/z 342	0.033	<0.032	0.011	57.365	0.071	<0.032
350	nominal m/z 343	0.033	<0.033	0.011	56.469	0.070	<0.033
351	nominal m/z 344	0.033	<0.033	0.011	56.363	0.076	<0.033
352	nominal m/z 345	0.032	<0.033	0.011	56.683	0.077	<0.033
353	nominal m/z 346	0.033	<0.032	0.011	56.901	0.083	<0.032
354	nominal m/z 347	0.033	<0.033	0.011	56.822	0.076	<0.033
355	nominal m/z 348	0.033	<0.033	0.011	56.821	0.077	<0.033
356	nominal m/z 349	0.033	<0.033	0.011	56.525	0.071	<0.033
357	nominal m/z 350	0.033	<0.033	0.011	56.402	0.076	<0.033
<	Signal Average/Median Below the MDL.						
†	Signal Average/Median Between the MDL and the RL (i.e., 3*MDL).						

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3.0 PRELIMINARY ANALYSIS

The Subject Matter Expert (SME) reviewed the maximum values to identify species of interest which were then further investigated for trends and plumes. No COPCs or odor species responded during this time that warranted further investigation. The only response of interest was at nominal m/z 46. Figure 3-1 shows the response at nominal m/z 46 during the monitoring period. There were multiple short-lived spikes increasing well above the background level. Identification of the species responsible for the m/z 46 response needs more investigation, but the spikes were coupled with spikes in CO₂ which suggests a combustion source. The lack of response from aromatics and aldehydes typical of diesel and gasoline vehicle combustion engines suggests it is not vehicle related. Current work has observed a large response at nominal m/z 46 in generator exhaust, which is a plausible explanation for the short-lived spikes correlated with CO₂.

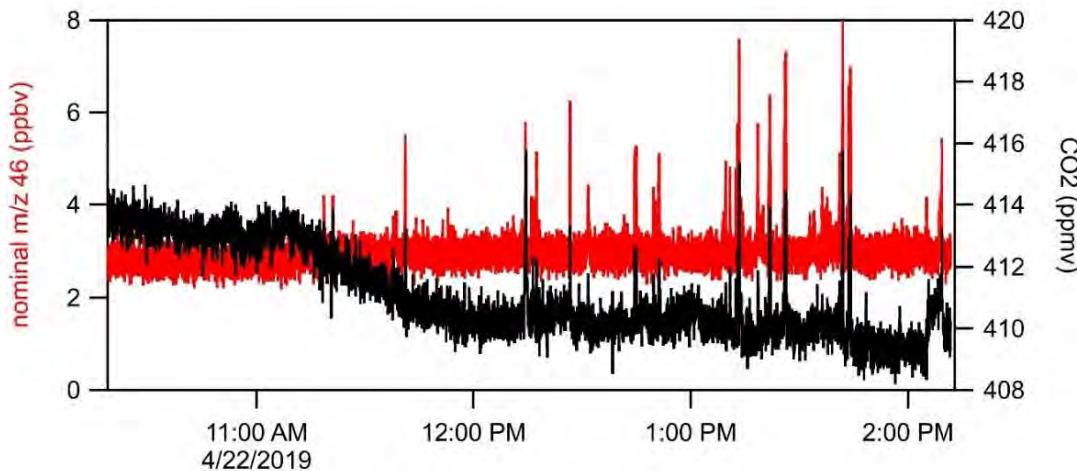


Figure 3-1. Response of CO₂ and Nominal m/z 46 During Monitoring after the AOP-015 Event on 04/22/2019.

4.0 REFERENCES

17124-DOE-HS-102, 2018, "Mobile Laboratory Data Processing-Analysis," Revision 0, TerraGraphics Environmental Engineering, Inc., Pasco, Washington.

66409-RPT-004, 2019, *Mobile Laboratory Operational Procedure*, Revision 11, TerraGraphics Environmental Engineering, Inc., Pasco, Washington.