LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049717.D
Acq On : 7 Aug 2021 23:39
Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Title : SVOA CALIBRATION

Signal : TIC: BG049717.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.079	3	7	16	rBV2	61414	138316	3.78%	0.718%
2	3.161	16	21	43	rVB	268413	531036	14.50%	2.758%
3	3.972	153	159	162	rBV3	3368	5838	0.16%	0.030%
4	4.189	192	196	200	rVB4	3110	5348	0.15%	0.028%
5	4.412	227	234	241	rBV2	39066	68084	1.86%	0.354%
6	4.559	256	259	264	rVB3	6407	10227	0.28%	0.053%
7	4.829	299	305	311	rBV2	28443	47844	1.31%	0.248%
8	5.117	351	354	359	rVB3	2823	4189	0.11%	0.022%
9	5.217	367	371	377	rBV3	3144	6424	0.18%	0.033%
10	6.028	499	509	511	rBV5	6156	11652	0.32%	0.061%
11	6.069	514	516	520	rVB3	4177	5213	0.14%	0.027%
12	7.203	699	709	717	rBV2	74927	148917	4.07%	0.773%
13	7.361	730	736	746	rBV	55961	103423	2.82%	0.537%
14	7.573	766	772	782	rVB	80971	147218	4.02%	0.765%
15	7.720	795	797	800	rBV3	4034	4652	0.13%	0.024%
16	8.043	843	852	859	rBV	127788	221161	6.04%	1.149%
17	8.178	870	875	879	rVB3	6218	9580	0.26%	0.050%
18	8.754	967	973	984	rVB2	85928	160034	4.37%	0.831%
19	8.871	988	993	997	rBV4	3958	5772	0.16%	0.030%
20	9.212	1041	1051	1059	rBV2	91285	164954	4.50%	0.857%
21	9.364	1072	1077	1085	rVB3	11807	21177	0.58%	0.110%
22	9.940	1167	1175	1183	rBV2	80657	145460	3.97%	0.755%
23	10.252	1225	1228	1230	rBV3	3227	4022	0.11%	0.021%
24	10.486	1260	1268	1277	rBV2	108766	208822	5.70%	1.084%
25	10.627	1287	1292	1293	rBV3	5041	5187	0.14%	0.027%
26	10.862	1325	1332	1339	rVV	164820	306387	8.37%	1.591%
27	10.945	1342	1346	1350	rVB2	3631	4545	0.12%	0.024%
28	11.438	1425	1430	1432	rBV4	2646	4270	0.12%	0.022%
29	11.661	1463	1468	1475	rVB4	5032	10575	0.29%	0.055%
30	11.802	1488	1492	1498	rVB4	6491	10895	0.30%	0.057%
31	11.890	1505	1507	1513	rVB5	4163	5399	0.15%	0.028%
32	12.308	1573	1578	1583	rVB4	2133	3926	0.11%	0.020%
33	12.402	1590	1594	1599	rBV4	7731	10601	0.29%	0.055%
34	12.584	1621	1625	1629	rBV4	2129	4226	0.12%	0.022%
35	13.054	1701	1705	1710	rVB5	4915	7797	0.21%	0.040%
36	13.136	1715	1719	1725	rBV5	8115	11976	0.33%	0.062%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049717.D
Acq On : 7 Aug 2021 23:39
Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Title : SVOA CALIBRATION

37	13.300	1739	1747	1750	rBV4	3221	7327	0.20%	0.038%
38	13.559	1785	1791	1795	rBV5	5209	11471	0.31%	0.060%
39	13.612	1797	1800	1805	rVB3	8283	13312	0.36%	0.069%
40	13.970	1855	1861	1865	rBV5	2468	4551	0.12%	0.024%
41	14.093	1875	1882	1889	rBV	204037	303392	8.28%	1.576%
42	14.164	1892	1894	1897	rVB2	4023	4341	0.12%	0.023%
43	14.322	1913	1921	1927	rBV4	12010	26203	0.72%	0.136%
44	14.387	1927	1932	1941	rVB	224377	362854	9.91%	1.884%
45	14.546	1956	1959	1962	rVV3	5776	5951	0.16%	0.031%
46	14.581	1962	1965	1968	rVB4	3797	5043	0.14%	0.026%
47	14.693	1977	1984	1992	rBV	258211	413233	11.28%	2.146%
48	14.904	2014	2020	2031	rBV4	24208	55041	1.50%	0.286%
49	15.098	2049	2053	2061	rBV6	11089	24215	0.66%	0.126%
50	15.345	2090	2095	2097	rBV6	3803	5482	0.15%	0.028%
51	15.427	2104	2109	2114	rVB	41403	54472	1.49%	0.283%
52	15.497	2114	2121	2128	rBV3	35401	73132	2.00%	0.380%
53	15.685	2147	2153	2160	rVB	286917	450398	12.30%	2.339%
54	15.809	2167	2174	2179	rBV2	62341	98976	2.70%	0.514%
55	15.973	2200	2202	2206	rBV3	2924	3930	0.11%	0.020%
56	16.243	2242	2248	2255	rVB6	17965	30443	0.83%	0.158%
57	16.390	2270	2273	2277	rBV6	5916	9283	0.25%	0.048%
58	16.572	2294	2304	2310	rBV6	20090	44397	1.21%	0.231%
59	16.725	2326	2330	2331	rBV4	4626	5114	0.14%	0.027%
60	16.831	2343	2348	2354	rBV4	27640	42739	1.17%	0.222%
61	17.195	2405	2410	2415	rBV4	28727	48278	1.32%	0.251%
62	17.330	2429	2433	2434	rBV4	22521	28182	0.77%	0.146%
63	17.448	2447	2453	2458	rVV	277000	431010	11.77%	2.238%
64	17.542	2462	2469	2476	rVB2	238741	396140	10.82%	2.057%
65	17.647	2484	2487	2491	rBV5	18187	21756	0.59%	0.113%
66	18.100	2560	2564	2569	rVB4	21872	29272	0.80%	0.152%
67	18.205	2580	2582	2587	rVB5	10287	14516	0.40%	0.075%
68	18.276	2590	2594	2599	rVB5	13327	19245	0.53%	0.100%
69	18.329	2599	2603	2612	rBV2	120885	183970	5.02%	0.955%
70	18.617	2649	2652	2654	rBV3	15530	19065	0.52%	0.099%
71	18.646	2654	2657	2666	rVB7	50954	79657	2.18%	0.414%
72	18.793	2679	2682	2684	rBV3	15173	18294	0.50%	0.095%
73	18.981	2711	2714	2719	rBV6	40922	56914	1.55%	0.296%
74	19.181	2746	2748	2752	rVB4	20043	21696	0.59%	0.113%
75	19.257	2756	2761	2765	rVV4	52592	93509	2.55%	0.486%
76	19.298	2765	2768	2771	rVB3	30049	33455	0.91%	0.174%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049717.D
Acq On : 7 Aug 2021 23:39
Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Title : SVOA CALIBRATION

77	19.480	2796	2799	2806	rVB5	42224	63961	1.75%	0.332%
78	19.598	2816	2819	2822	rBV4	29025	35746	0.98%	0.186%
79	19.833	2854	2859	2865	rVB2	308093	456105	12.45%	2.369%
80	20.320	2938	2942	2944	rBV2	113565	148571	4.06%	0.772%
81	20.350	2944	2947	2956	rVB	118938	165049	4.51%	0.857%
82	20.679	2999	3003	3010	rVB7	64679	137608	3.76%	0.715%
83	20.790	3019	3022	3026	rBV	88940	114764	3.13%	0.596%
84	21.025	3057	3062	3070	rBV4	116883	198501	5.42%	1.031%
85	21.354	3113	3118	3130	rBV	945704	1434656	39.18%	7.450%
86	21.730	3177	3182	3190	rVB	295053	491282	13.42%	2.551%
87	22.159	3250	3255	3261	rVB	203338	309928	8.46%	1.610%
88	22.553	3313	3322	3337	rVB2	473054	1426726	38.96%	7.409%
89	23.592	3492	3499	3509	rVB2	678349	1415765	38.66%	7.352%
90	24.056	3572	3578	3583	rBV2	147291	317262	8.66%	1.648%
91	24.121	3583	3589	3602	rVB	374705	948763	25.91%	4.927%
92	25.014	3733	3741	3749	rBV2	155211	429158	11.72%	2.229%
93	25.560	3825	3834	3846	rVB2	496013	1384539	37.81%	7.190%
94	26.330	3953	3965	3982	rBV	1063917	3662163	100.00%	19.018%

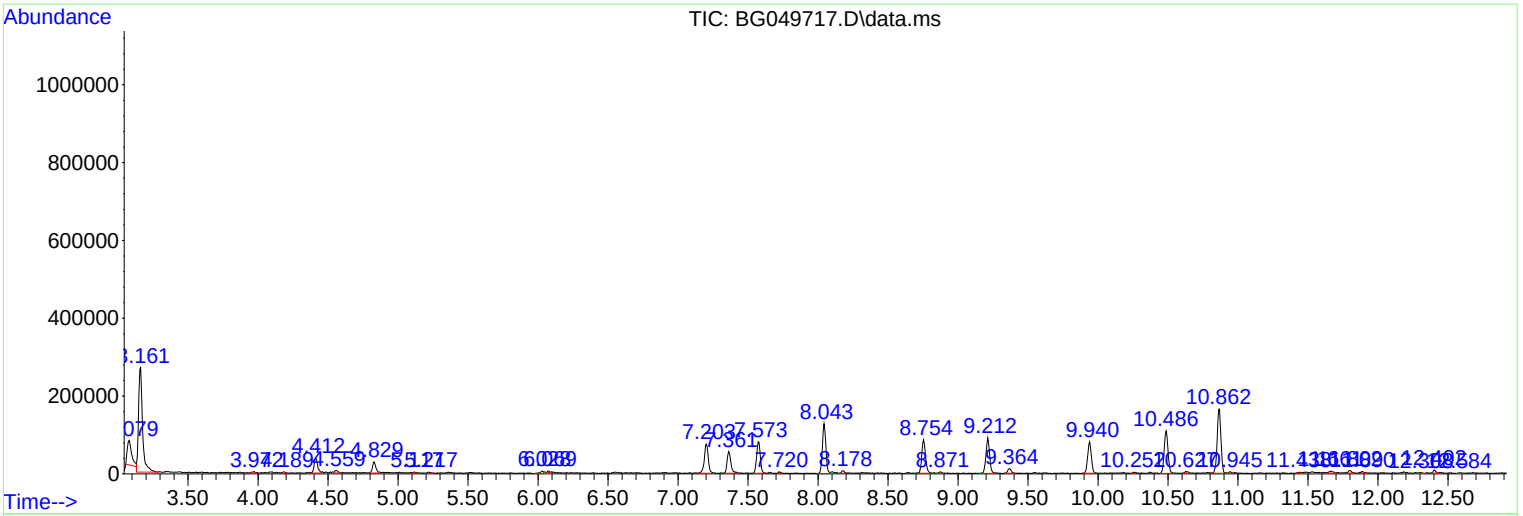
Sum of corrected areas: 19255953

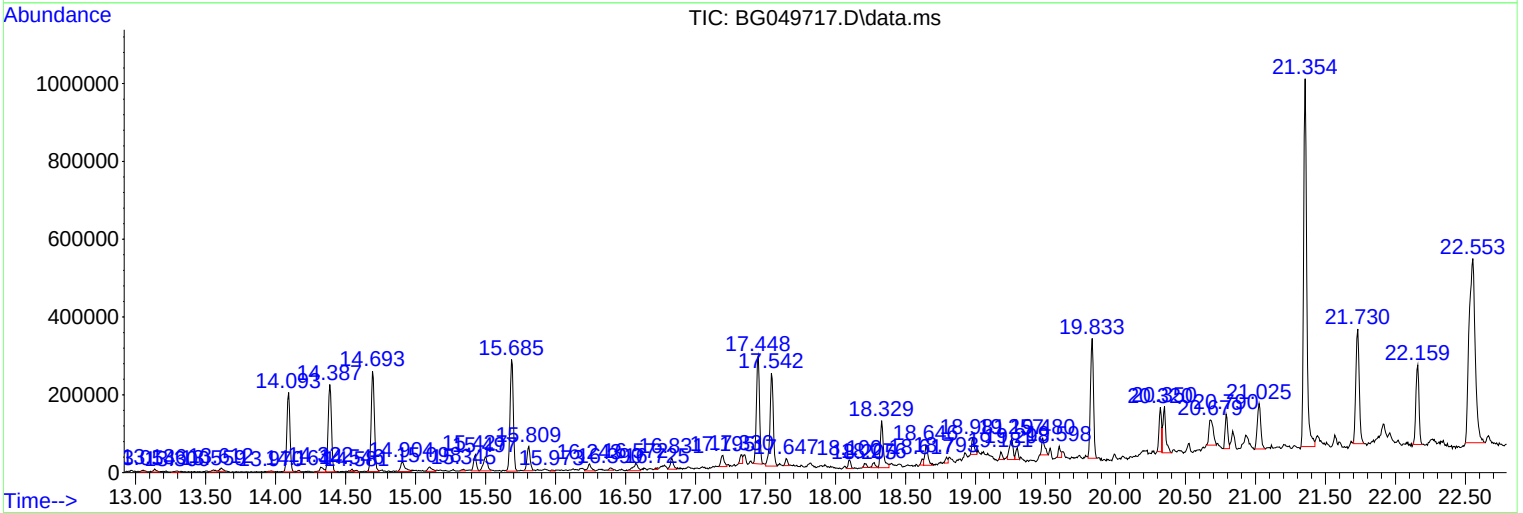
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049717.D
Acq On : 7 Aug 2021 23:39
Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

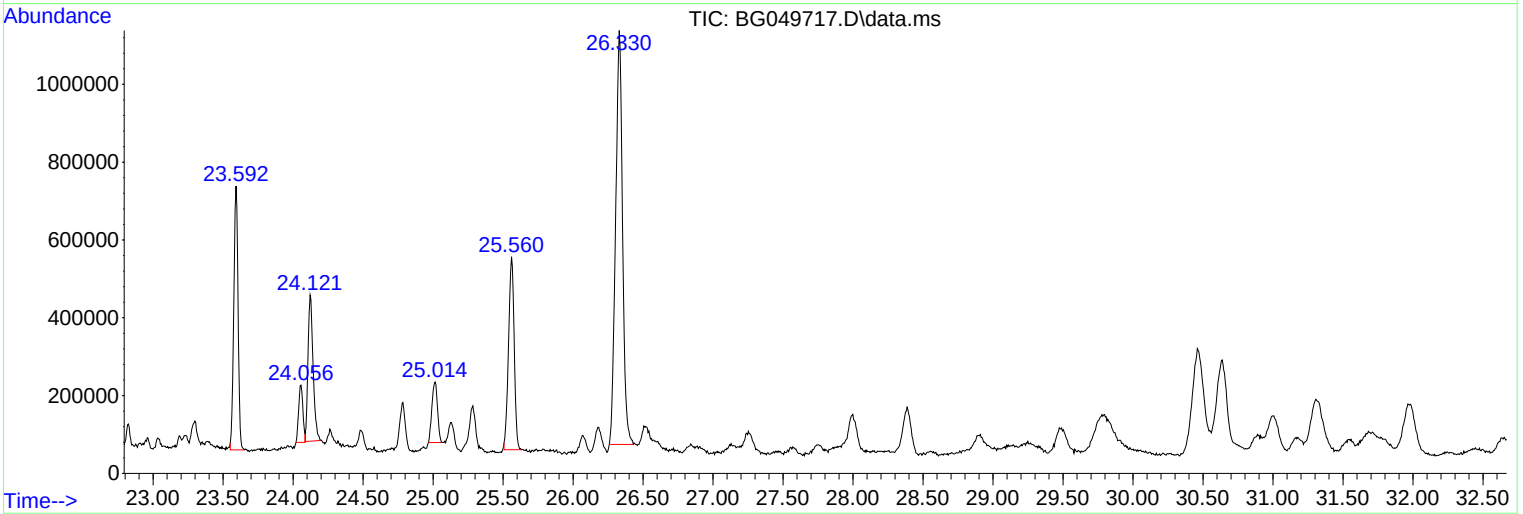
Instrument :
BNA_G
ClientSampleId :
JCE01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P







Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049717.D
Acq On : 7 Aug 2021 23:39
Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

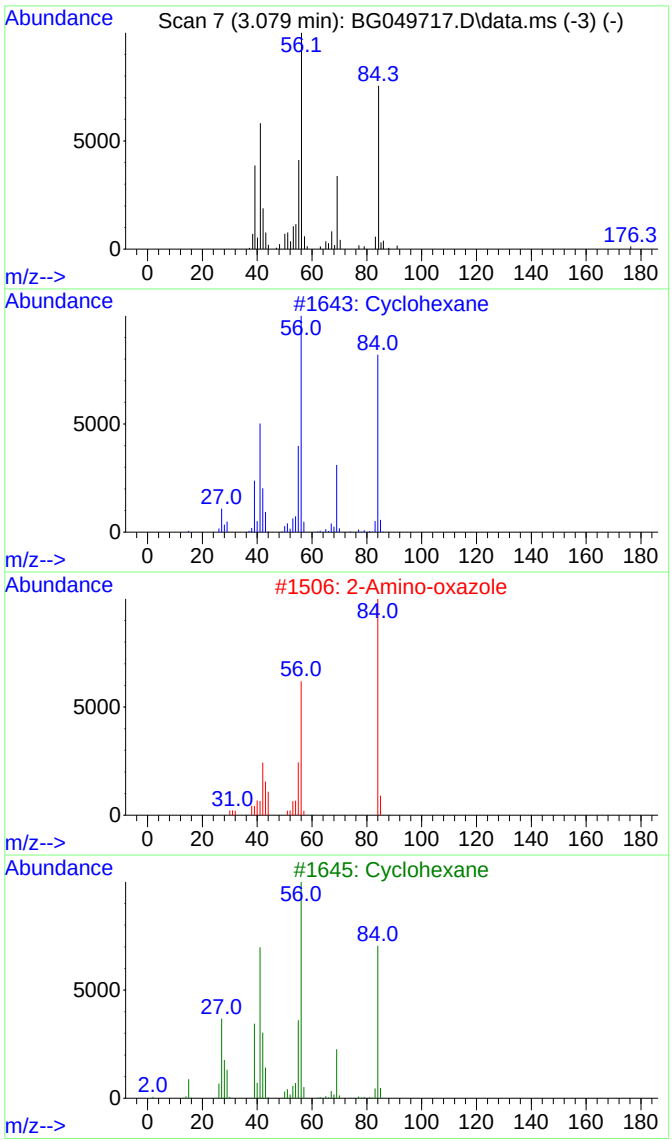
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

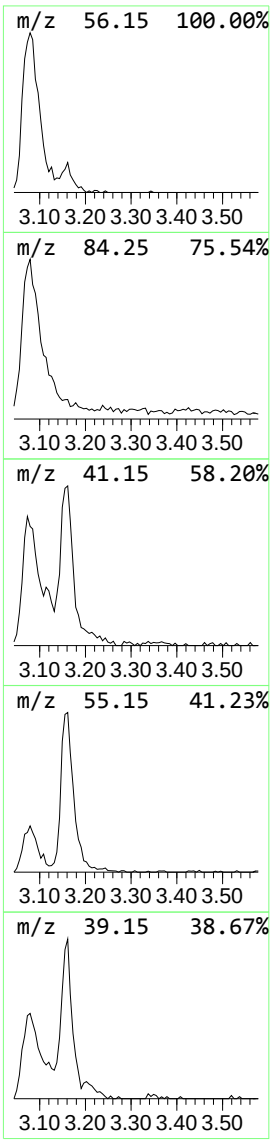
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Cyclic3.079 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.079	12.51 ng/ul	138316	1,4-Dichlorobenzene-d4	8.043

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane	84	C6H12	000110-82-7	81
2	2-Amino-oxazole	84	C3H4N2O	004570-45-0	64
3	Cyclohexane	84	C6H12	000110-82-7	59
4	Cyclohexane	84	C6H12	000110-82-7	59
5	Cyclohexane	84	C6H12	000110-82-7	58





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049717.D
Acq On : 7 Aug 2021 23:39
Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

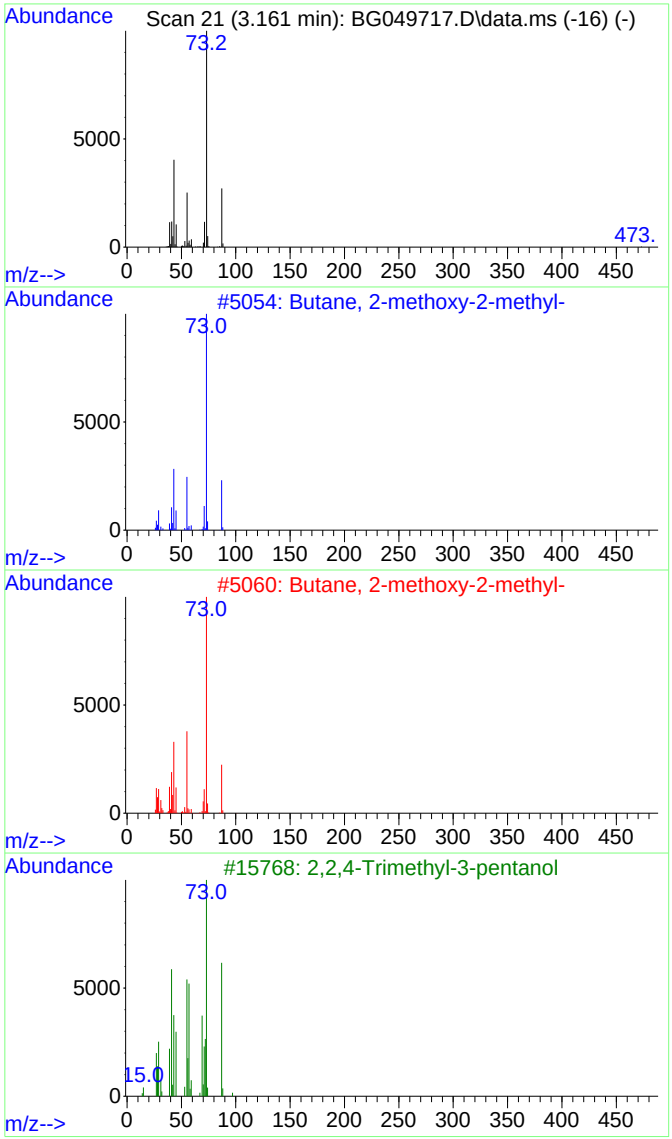
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

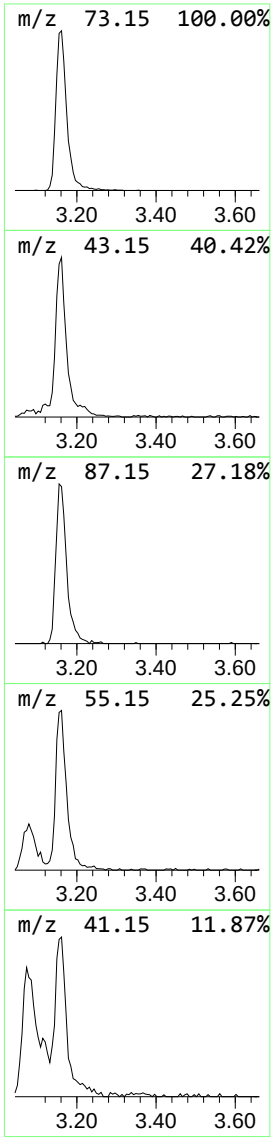
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.161	48.02 ng/ul	531036	1,4-Dichlorobenzene-d4	8.043

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	56
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	35
5	2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	32





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049717.D
Acq On : 7 Aug 2021 23:39
Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
JCE01

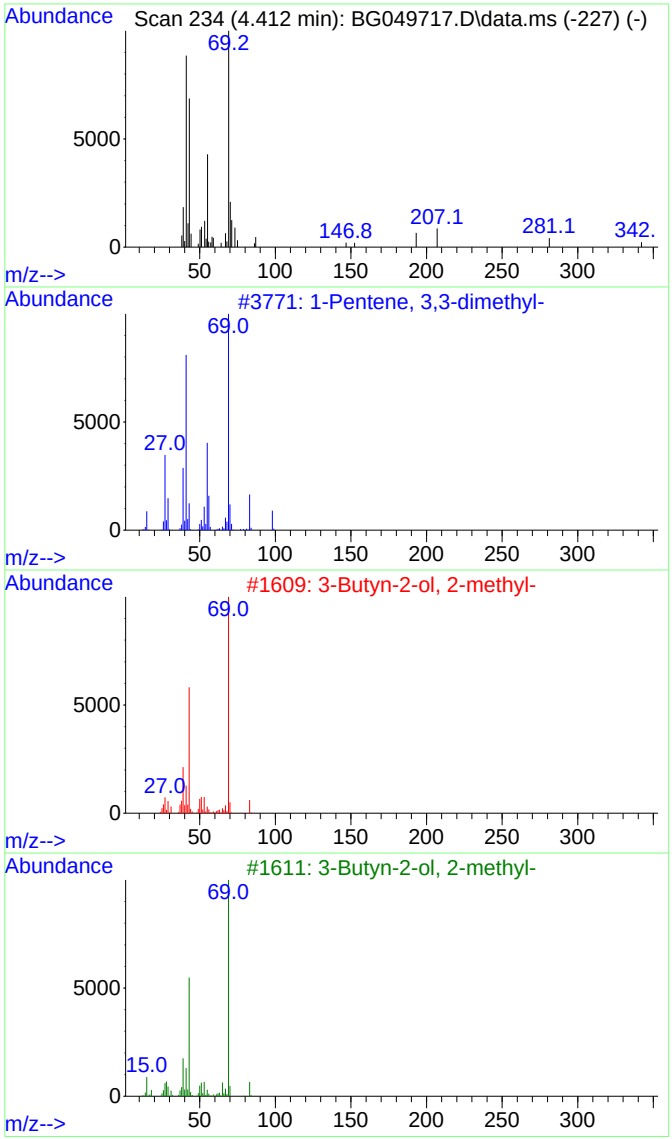
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

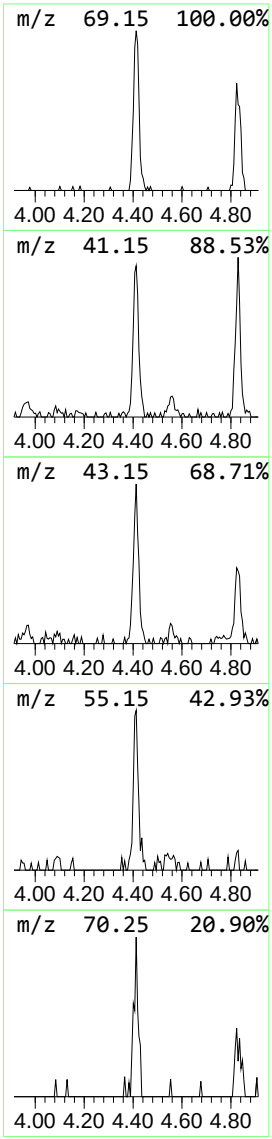
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.412	6.16 ng/ul	68084	1,4-Dichlorobenzene-d4	8.043

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentene, 3,3-dimethyl-	98	C7H14	003404-73-7	47
2	3-Butyn-2-ol, 2-methyl-	84	C5H8O	000115-19-5	43
3	3-Butyn-2-ol, 2-methyl-	84	C5H8O	000115-19-5	43
4	3-Methyl-2-hexene	98	C7H14	017618-77-8	42
5	1-Pentene, 3-ethyl-	98	C7H14	004038-04-4	40





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049717.D
Acq On : 7 Aug 2021 23:39
Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

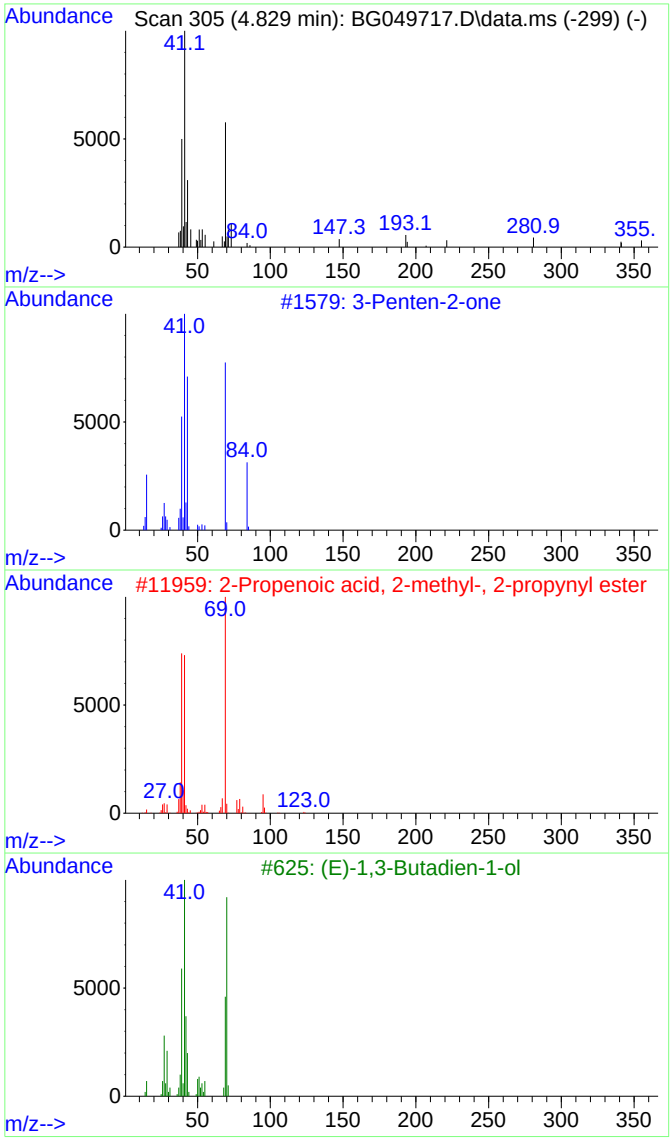
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

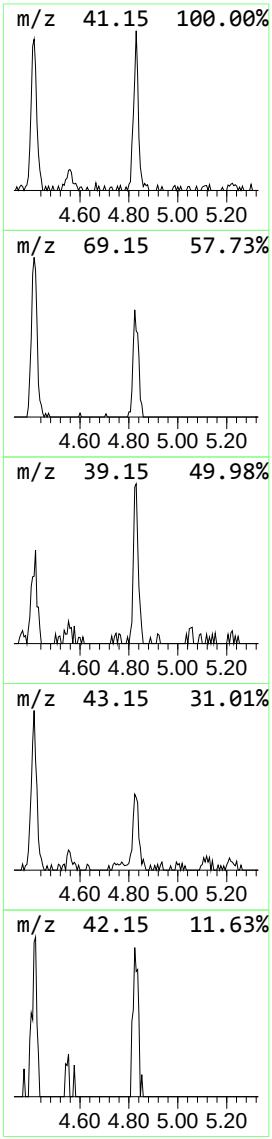
****

Peak Number 4 unknown-01 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.829	4.33 ng/ul	47844	1,4-Dichlorobenzene-d4	8.043

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Penten-2-one	84	C5H8O	000625-33-2	43
2	2-Propenoic acid, 2-methyl-, 2-p...	124	C7H8O2	013861-22-8	37
3	(E)-1,3-Butadien-1-ol	70	C4H6O	070411-98-2	32
4	2-Propenoic acid, 2-methyl-	86	C4H6O2	000079-41-4	25
5	3-Penten-2-one, (E)-	84	C5H8O	003102-33-8	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049717.D
Acq On : 7 Aug 2021 23:39
Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

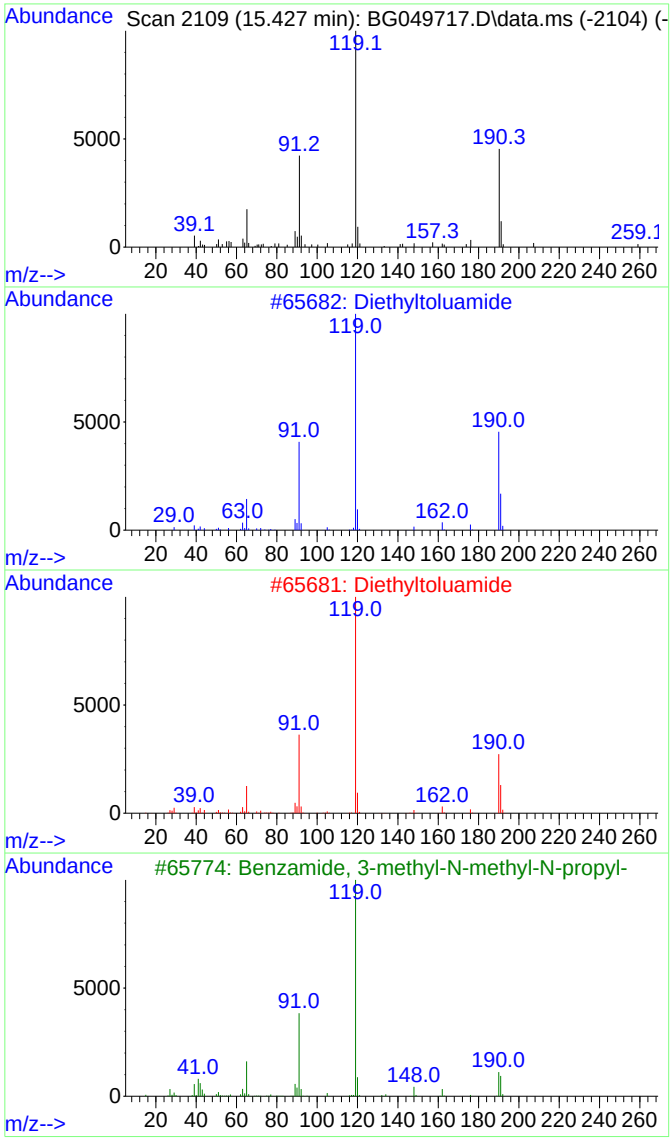
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

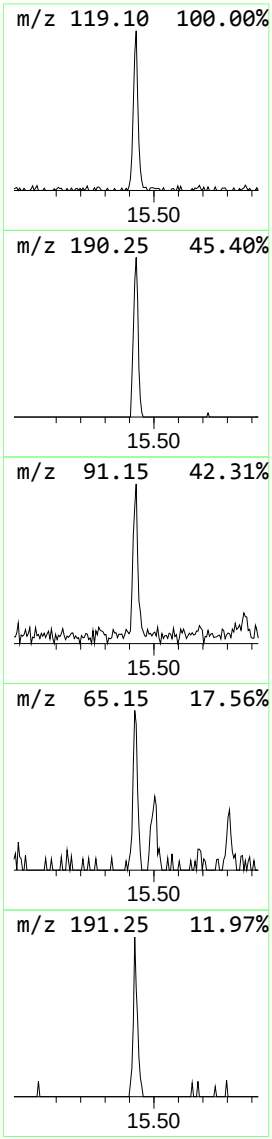
****

Peak Number 5 Diethyltoluamide Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.427	2.64 ng/ul	54472	Acenaphthene-d10	14.693

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethyltoluamide	191	C12H17NO	000134-62-3	91
2	Diethyltoluamide	191	C12H17NO	000134-62-3	87
3	Benzamide, 3-methyl-N-methyl-N-p...	191	C12H17NO	1000421-46-2	81
4	Isoxazole-3-carboximidamide, N'-...	259	C13H13N3O3	263869-49-4	70
5	Benzamide, N,N-diethyl-4-methyl-	191	C12H17NO	002728-05-4	68





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Acq On : 7 Aug 2021 23:39
Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

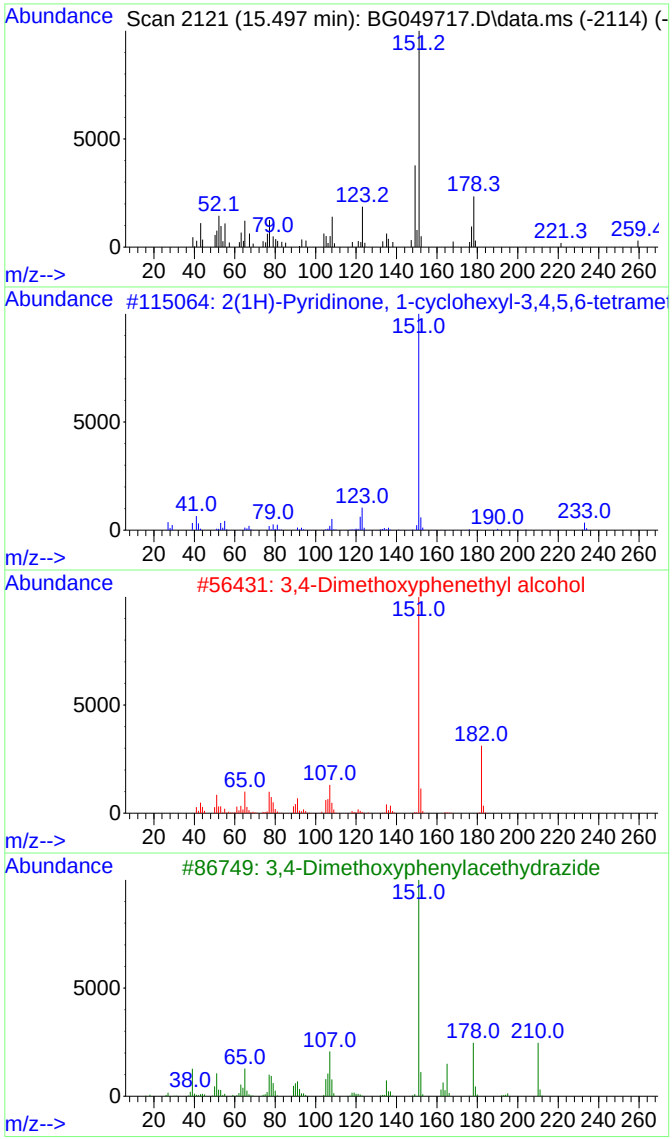
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

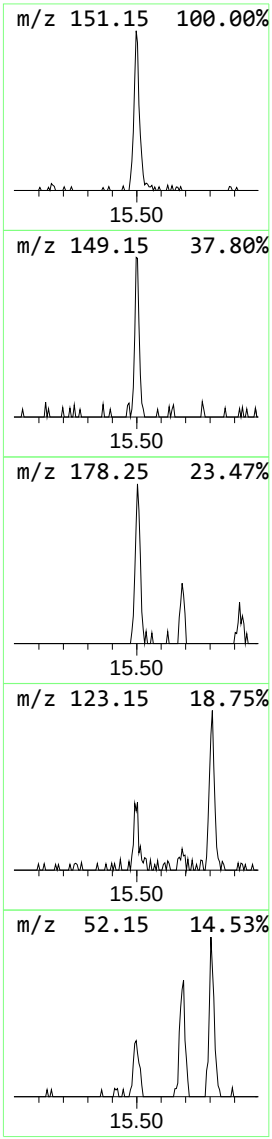
****

Peak Number 6 unknown-02 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.497	3.54 ng/ul	73132	Acenaphthene-d10	14.693

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(1H)-Pyridinone, 1-cyclohexyl-3...	233	C15H23NO	082754-44-7	47
2	3,4-Dimethoxyphenethyl alcohol	182	C10H14O3	007417-21-2	43
3	3,4-Dimethoxyphenylacethydrazide	210	C10H14N2O3	060075-23-2	40
4	1,4-Dimethoxy-2,3-dimethylbenzene	166	C10H14O2	039021-83-5	38
5	4-(Methylsulfanyl)benzohydrazide	182	C8H10N2OS	1000484-57-9	38





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Acq On : 7 Aug 2021 23:39
Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

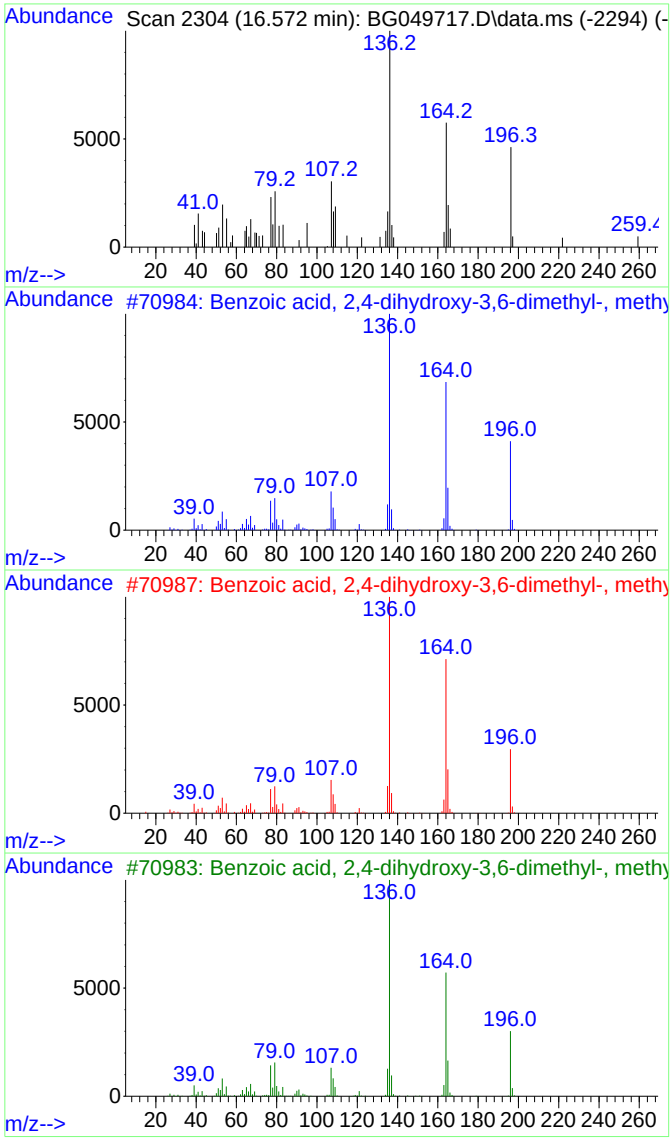
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

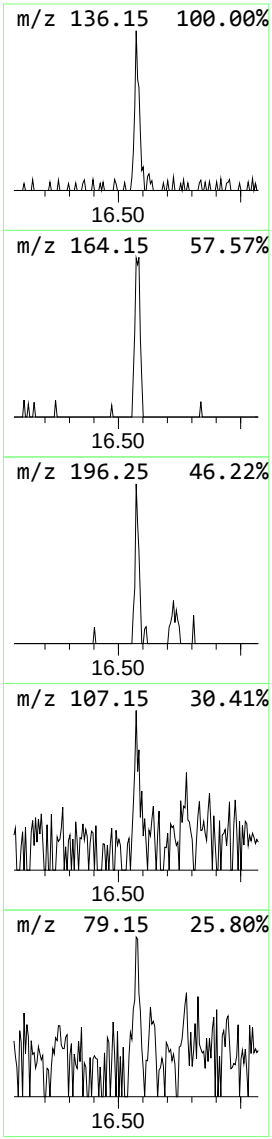
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzoic acid, 2,4-dihydroxy... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.572	2.06 ng/ul	44397	Phenanthrene-d10	17.448

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 2,4-dihydroxy-3,6-...	196	C10H12O4	004707-47-5	94
2	Benzoic acid, 2,4-dihydroxy-3,6-...	196	C10H12O4	004707-47-5	91
3	Benzoic acid, 2,4-dihydroxy-3,6-...	196	C10H12O4	004707-47-5	90
4	Benzoic acid, 2,4-dihydroxy-3,6-...	196	C10H12O4	004707-47-5	87
5	Benzoic acid, 2,4-dihydroxy-3,6-...	196	C10H12O4	004707-47-5	72





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Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
JCE01

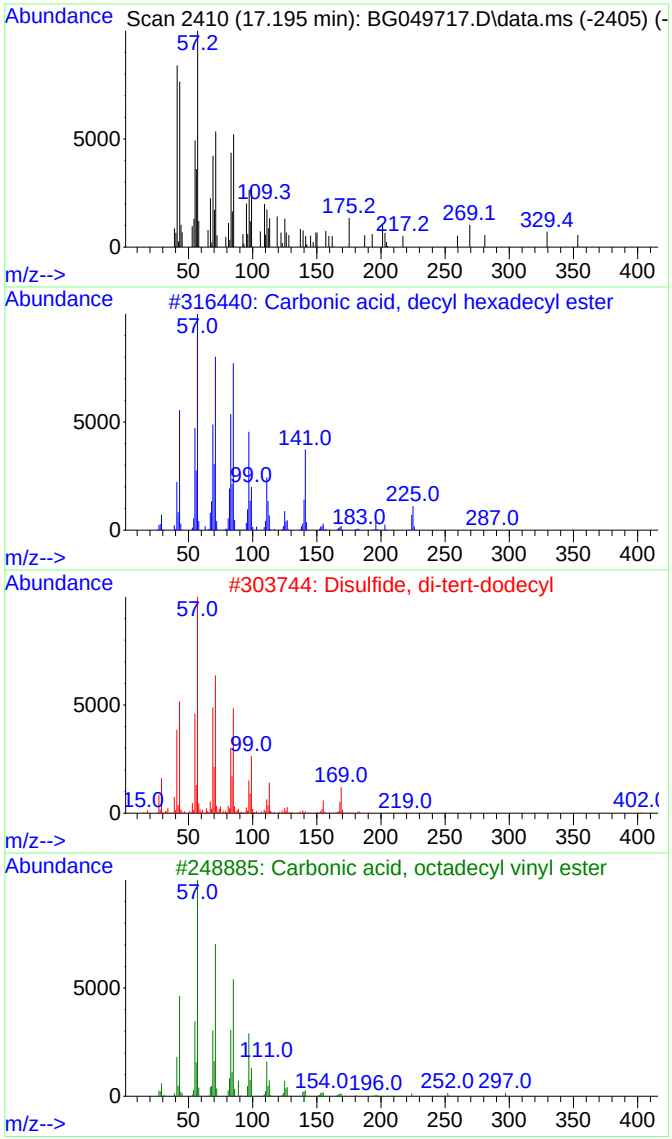
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

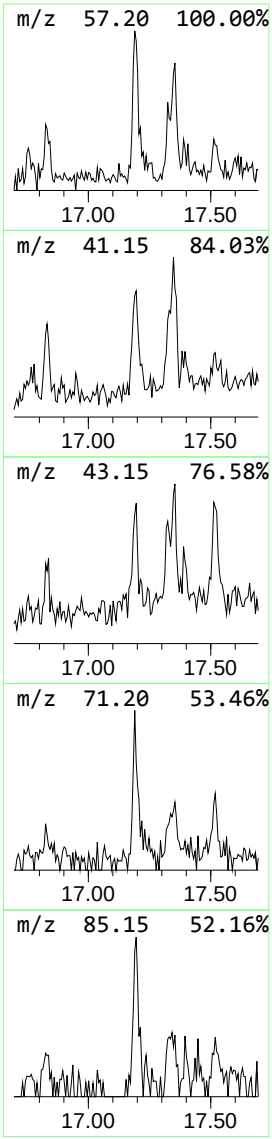
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Carbonic acid, decyl hexade... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.195	2.24 ng/ul	48278	Phenanthrene-d10	17.448

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Carbonic acid, decyl hexadecyl e...	426	C27H54O3	1000383-16-5	58
2	Disulfide, di-tert-dodecyl	402	C24H50S2	027458-90-8	49
3	Carbonic acid, octadecyl vinyl e...	340	C21H40O3	1000382-54-4	49
4	Carbonic acid, decyl undecyl ester	356	C22H44O3	1000383-16-0	47
5	Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	47





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Acq On : 7 Aug 2021 23:39
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Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
JCE01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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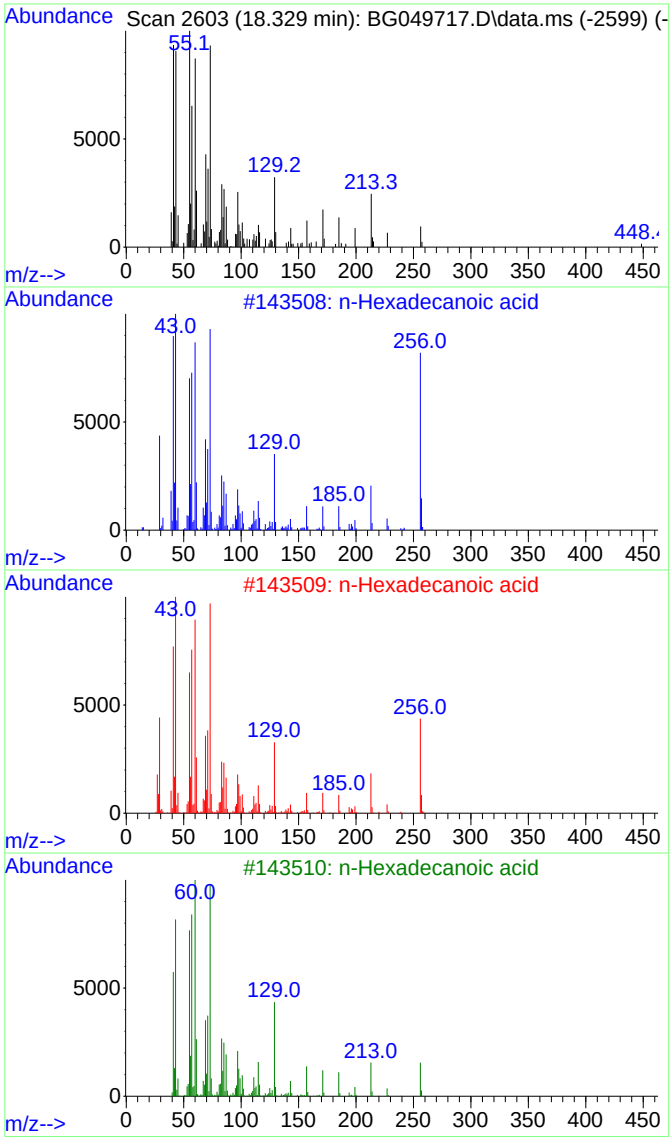
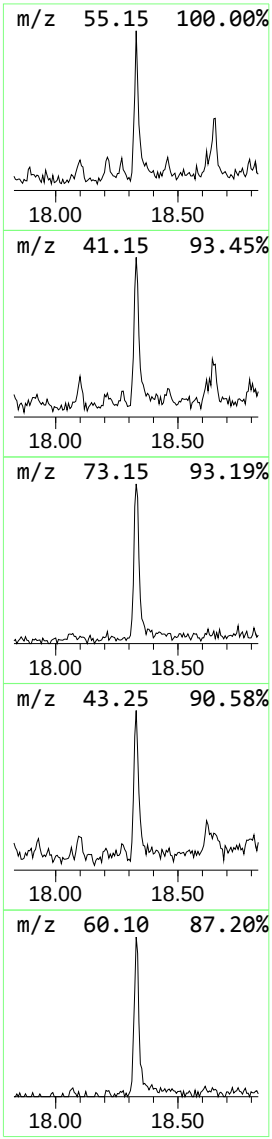
Peak Number 9 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.329	8.54 ng/ul	183970	Phenanthrene-d10	17.448

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	76
4	Tridecanoic acid	214	C13H26O2	000638-53-9	76
5	Tridecanoic acid	214	C13H26O2	000638-53-9	64





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Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

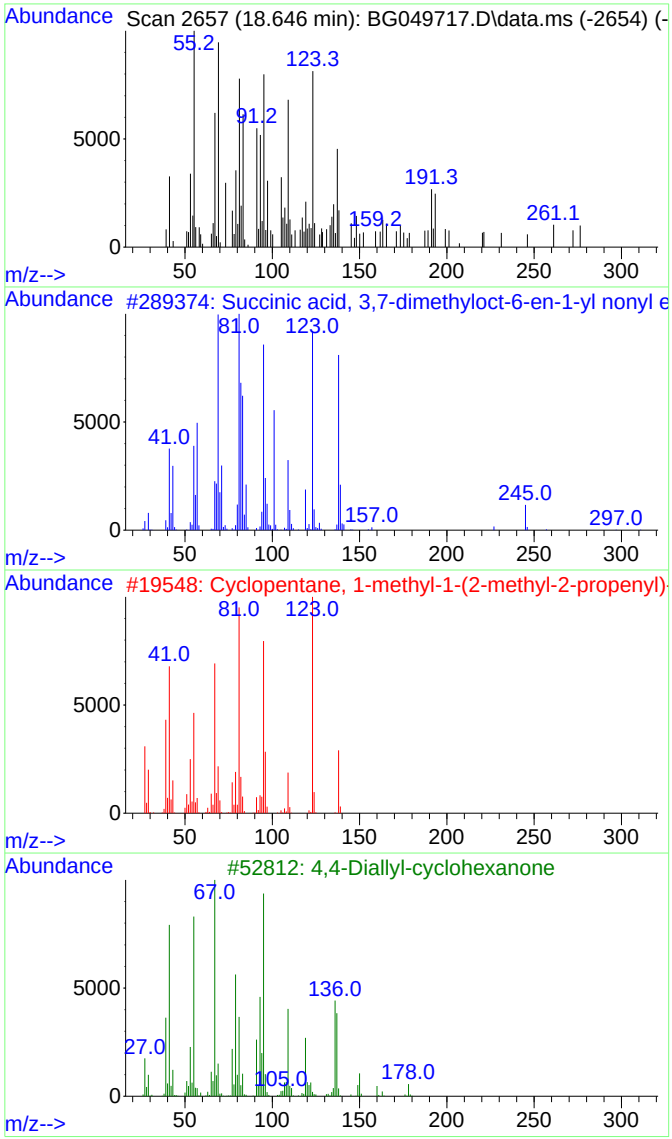
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

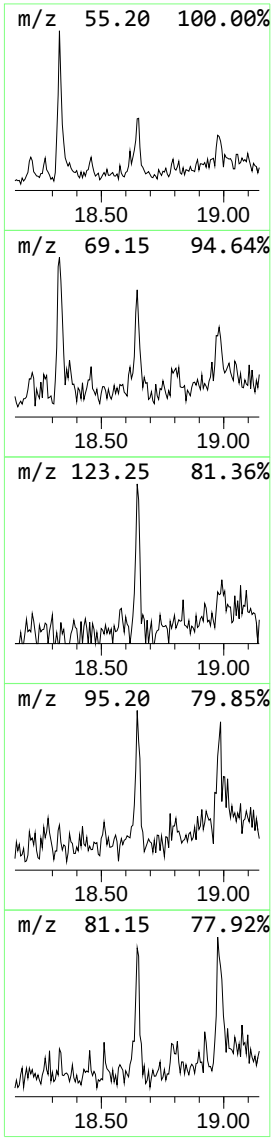
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Succinic acid, 3,7-dimethyl... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.646	3.70 ng/ul	79657	Phenanthrene-d10	17.448

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Succinic acid, 3,7-dimethyloct-6...	382	C23H42O4	1010353-34-3	58
2	Cyclopentane, 1-methyl-1-(2-meth...	138	C10H18	074764-47-9	46
3	4,4-Diallyl-cyclohexanone	178	C12H18O	1000186-50-2	44
4	.alpha.-Damascone	192	C13H20O	031089-90-4	30
5	3,7-Decadiyne, 2,2,5,5,6,6,9,9-o...	246	C18H30	017553-35-4	30





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Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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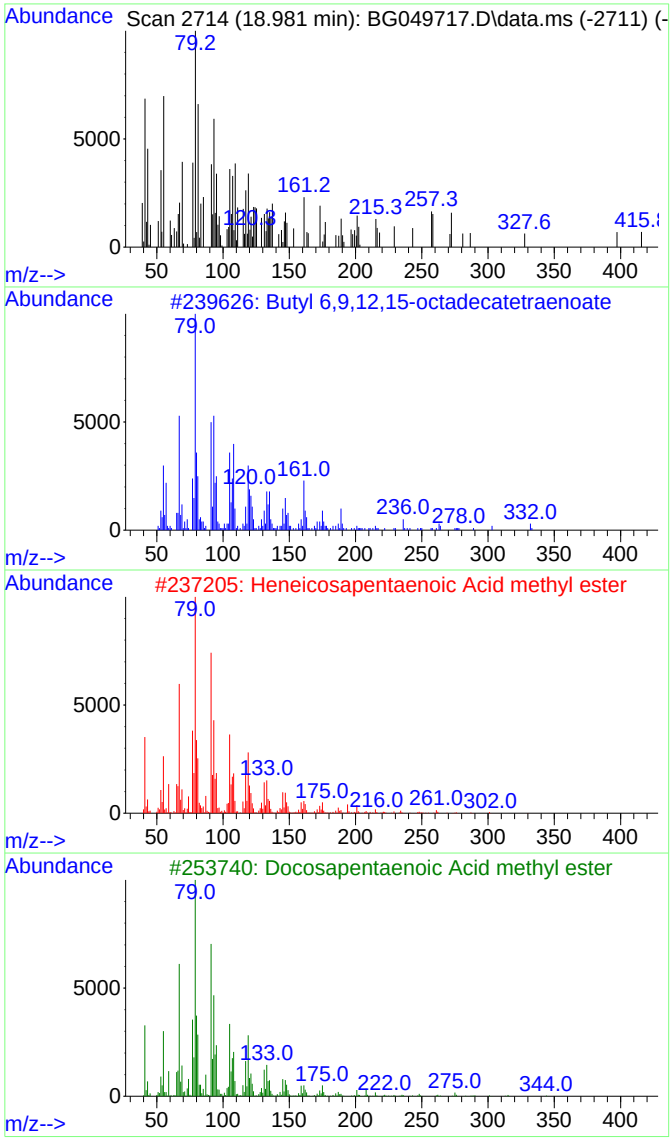
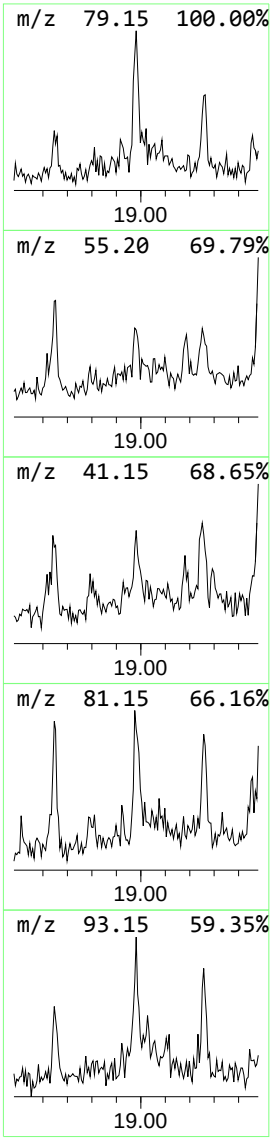
Peak Number 11 unknown-03 Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.981	2.64 ng/ul	56914	Phenanthrene-d10	17.448

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butyl 6,9,12,15-octadecatetraenoate	332	C22H36O2	1000336-71-3	38
2	Heneicosapentaenoic Acid methyl ...	330	C22H34O2	065919-53-1	35
3	Docosapentaenoic Acid methyl ester	344	C23H36O2	108698-02-8	35
4	1,5,9,11-Tridecatetraene, 12-met...	190	C14H22	062338-27-6	30
5	Octadecahydro-cyclopenta[e]pyrene	258	C19H30	1000371-48-4	30





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Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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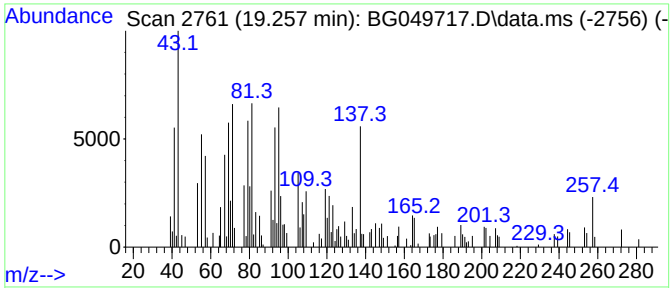
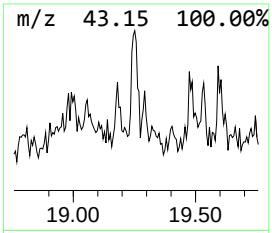
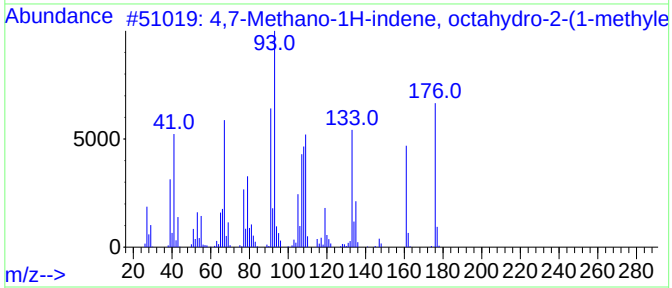
Peak Number 12 unknown-04 Concentration Rank 18

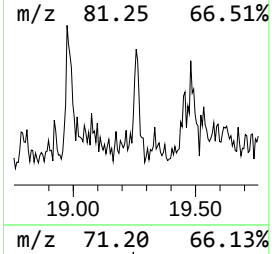
R.T.	EstConc	Area	Relative to ISTD	R.T.
19.257	4.34 ng/ul	93509	Phenanthrene-d10	17.448

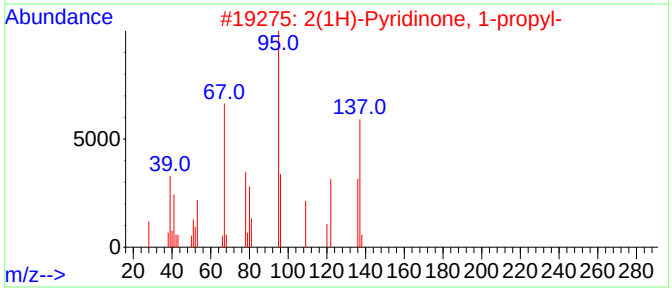
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4,7-Methano-1H-indene, octahydro...	176	C13H20	074793-54-7	44
2	2(1H)-Pyridinone, 1-propyl-	137	C8H11NO	019006-63-4	38
3	Phenanthrene, 7-ethenyl-1,2,3,4,...	272	C20H32	001686-56-2	30
4	1,5,9,11-Tridecatetraene, 12-met...	190	C14H22	062338-27-6	30
5	.beta.-Bisabolene	204	C15H24	000495-61-4	25

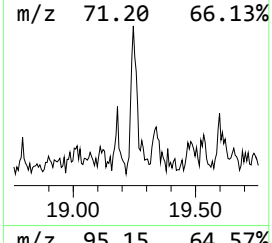


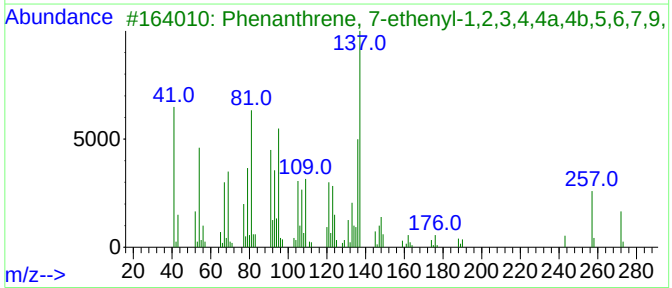


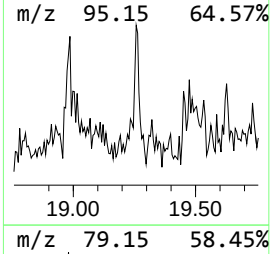


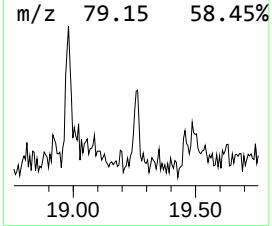












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Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

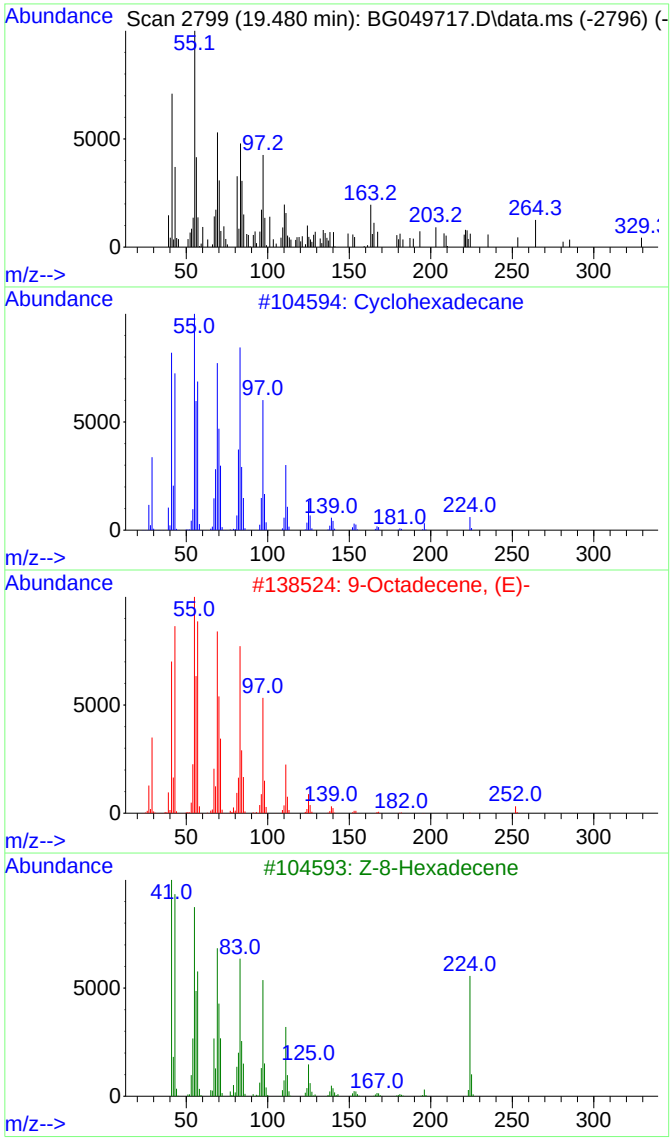
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

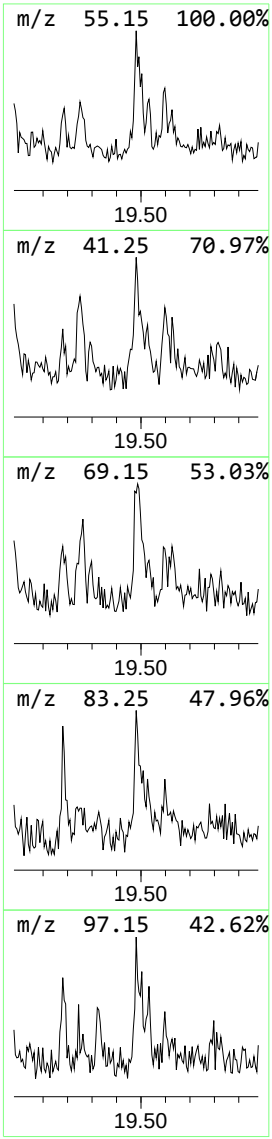
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 (DEL) Alkane: Cyclic19.480 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.480	2.97 ng/ul	63961	Phenanthrene-d10	17.448

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexadecane	224	C16H32	000295-65-8	83
2	9-Octadecene, (E)-	252	C18H36	007206-25-9	52
3	Z-8-Hexadecene	224	C16H32	1000130-87-5	50
4	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	49
5	Cycloeicosane	280	C20H40	000296-56-0	47





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Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
JCE01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

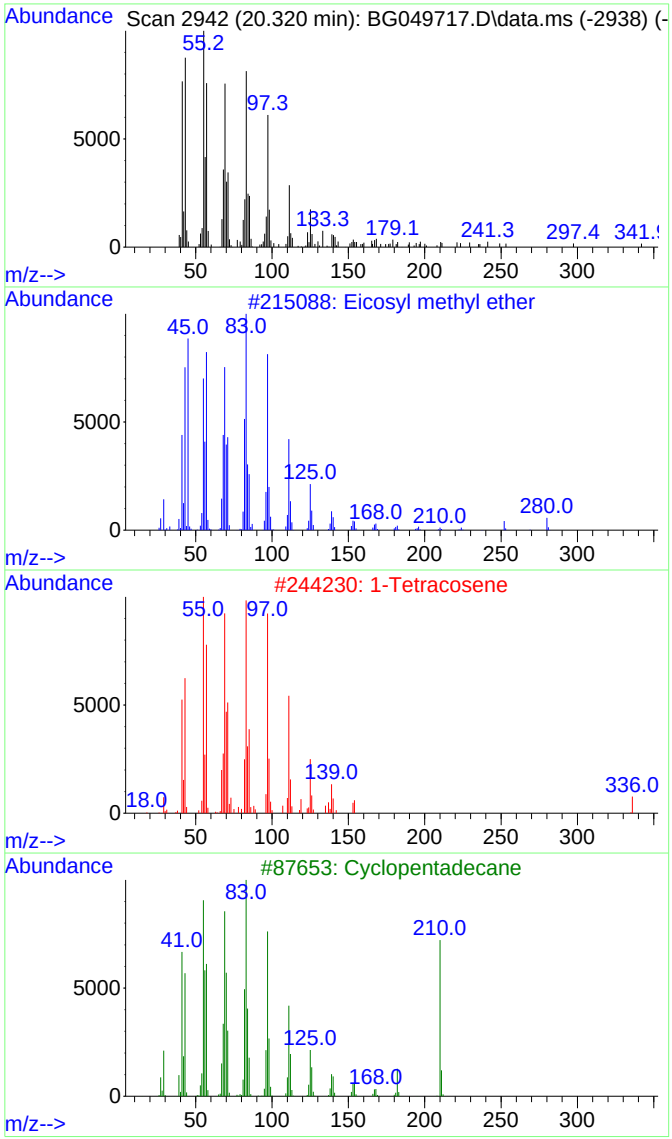
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

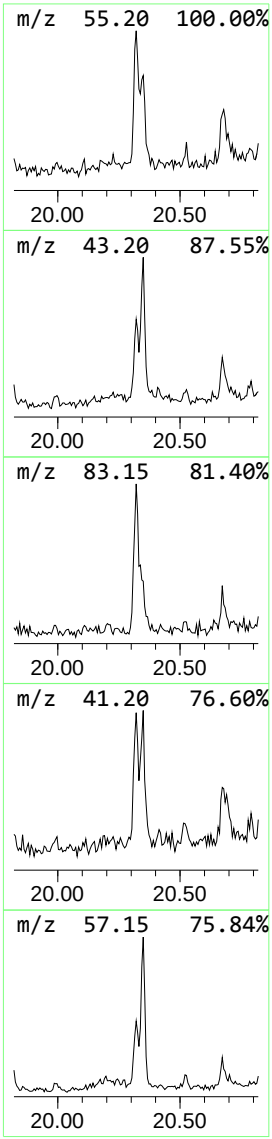
****

Peak Number 14 Eicosyl methyl ether Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.320	6.05 ng/ul	148571	Chrysene-d12	21.730

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosyl methyl ether	312	C21H44O	1000406-29-4	91
2	1-Tetracosene	336	C24H48	010192-32-2	87
3	Cyclopentadecane	210	C15H30	000295-48-7	87
4	1-Tetracosene	336	C24H48	010192-32-2	83
5	Trifluoroacetic acid,n-tridecyl ...	296	C15H27F3O2	053800-02-5	83





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Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

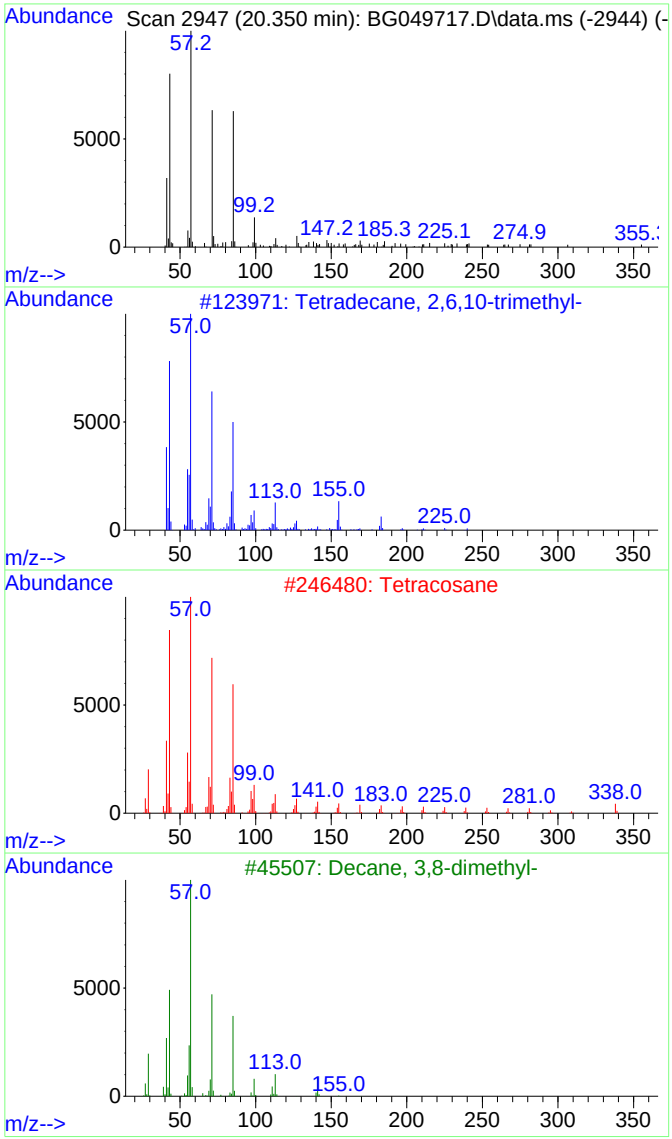
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

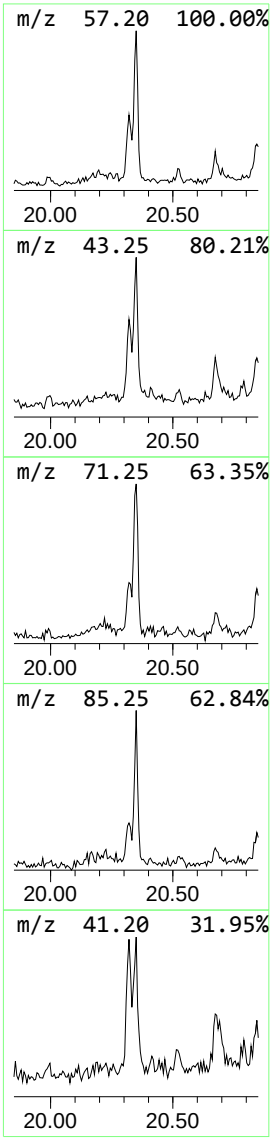
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.350	6.72 ng/ul	165049	Chrysene-d12	21.730

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	80
2	Tetracosane	338	C24H50	000646-31-1	72
3	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	72
4	Dodecane, 2-methyl-	184	C13H28	001560-97-0	72
5	Tridecane	184	C13H28	000629-50-5	72





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Acq On : 7 Aug 2021 23:39
Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

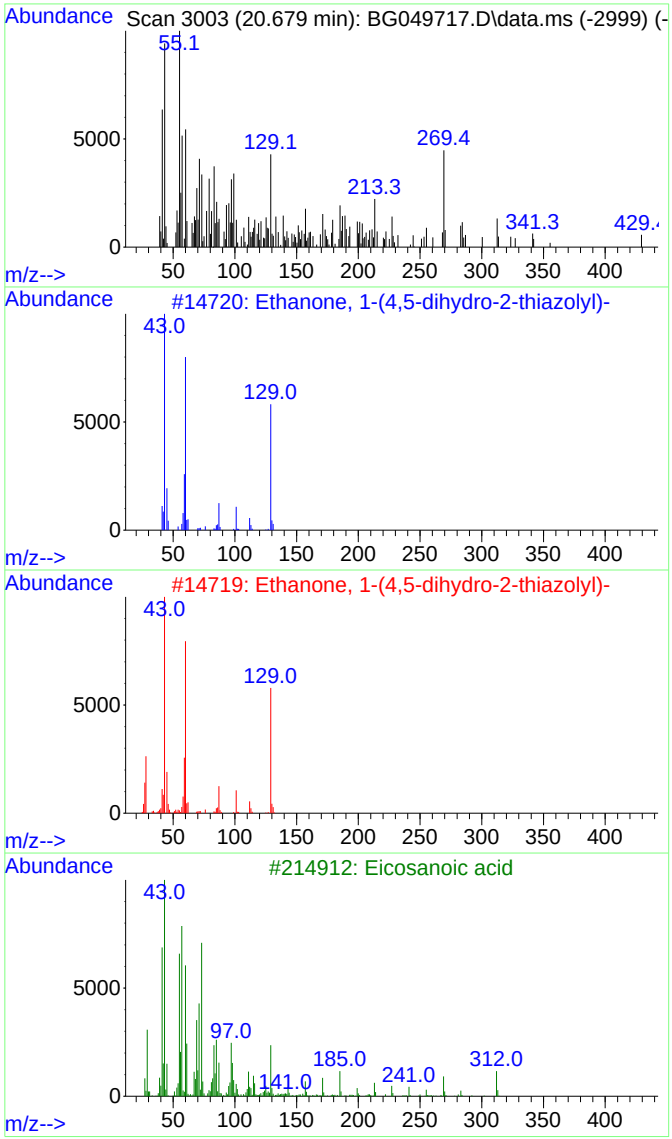
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

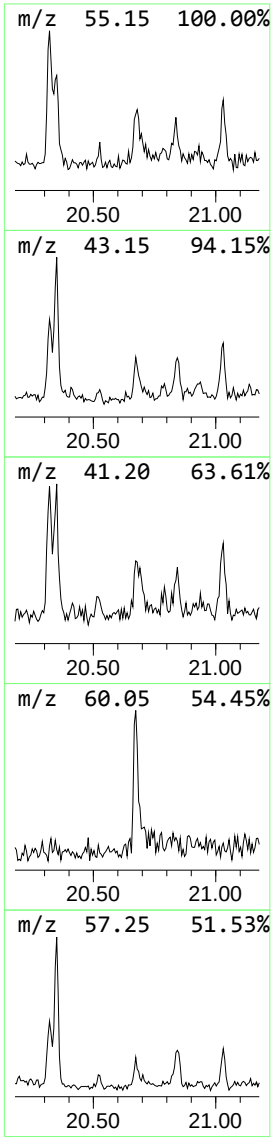
****

Peak Number 16 unknown-05 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.679	5.60 ng/ul	137608	Chrysene-d12	21.730

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanone, 1-(4,5-dihydro-2-thiaz...	129	C5H7NOS	029926-41-8	25
2	Ethanone, 1-(4,5-dihydro-2-thiaz...	129	C5H7NOS	029926-41-8	25
3	Eicosanoic acid	312	C20H40O2	000506-30-9	25
4	2,4-Di-tert-butylphenol mesylate	284	C15H24O3S	1071592-55-6	15
5	Isopropyl stearate	326	C21H42O2	000112-10-7	12





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Acq On : 7 Aug 2021 23:39
Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

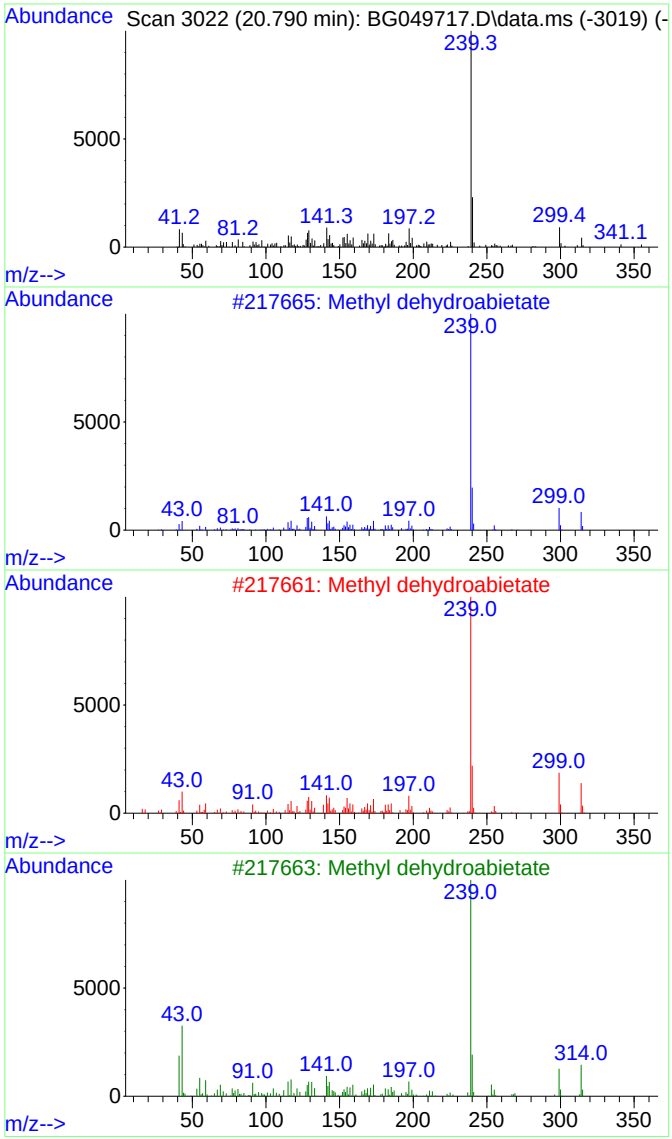
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

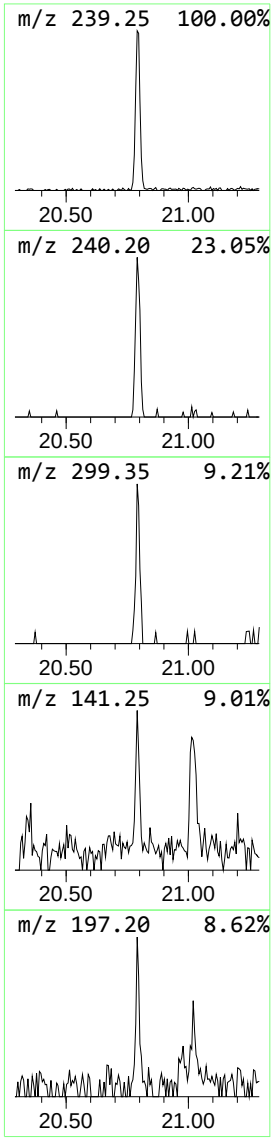
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Methyl dehydroabietate Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.790	4.67 ng/ul	114764	Chrysene-d12	21.730

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl dehydroabietate	314	C21H30O2	001235-74-1	95
2	Methyl dehydroabietate	314	C21H30O2	001235-74-1	91
3	Methyl dehydroabietate	314	C21H30O2	001235-74-1	83
4	9,10-Anthracenedione, 1-amino-4-...	239	C14H9NO3	000116-85-8	53
5	9,10-Anthracenedione, 1-amino-4-...	239	C14H9NO3	000116-85-8	53





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Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

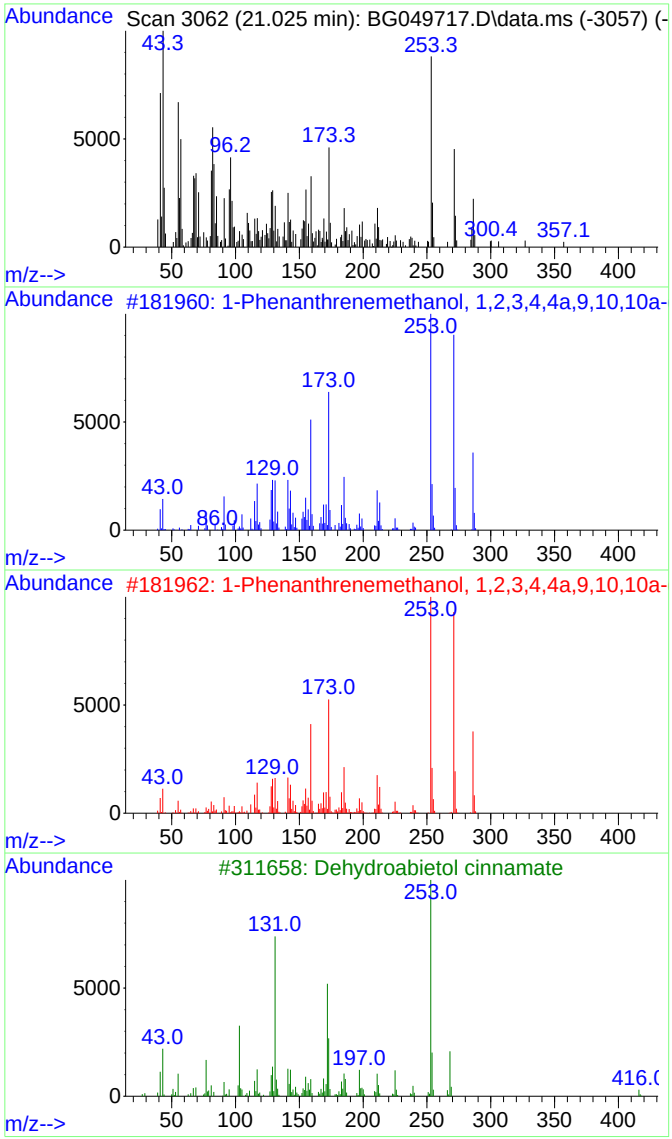
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

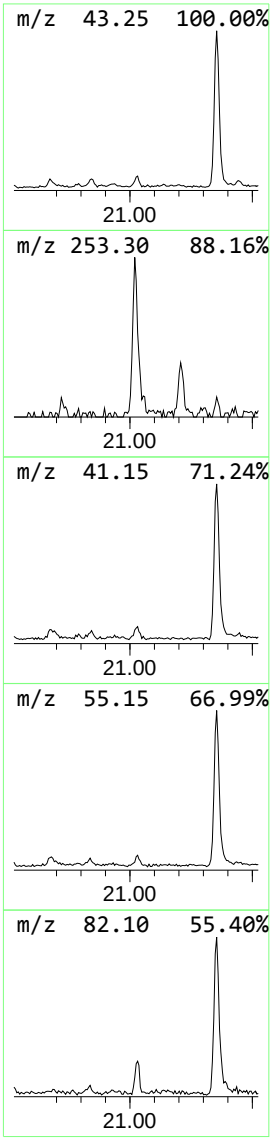
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 1-Phenanthrenemethanol, 1,2... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.025	8.08 ng/ul	198501	Chrysene-d12	21.730

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenanthrenemethanol, 1,2,3,4,...	286	C20H30O	003772-55-2	60
2	1-Phenanthrenemethanol, 1,2,3,4,...	286	C20H30O	003772-55-2	45
3	Dehydroabietol cinnamate	416	C29H36O2	1000414-71-7	27
4	4-epi-Dehydroabietinol acetate	328	C22H32O2	024462-15-5	27
5	8-Chloro-5-methyl-2-phenylquinoline	253	C16H12ClN	025413-16-5	20





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Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

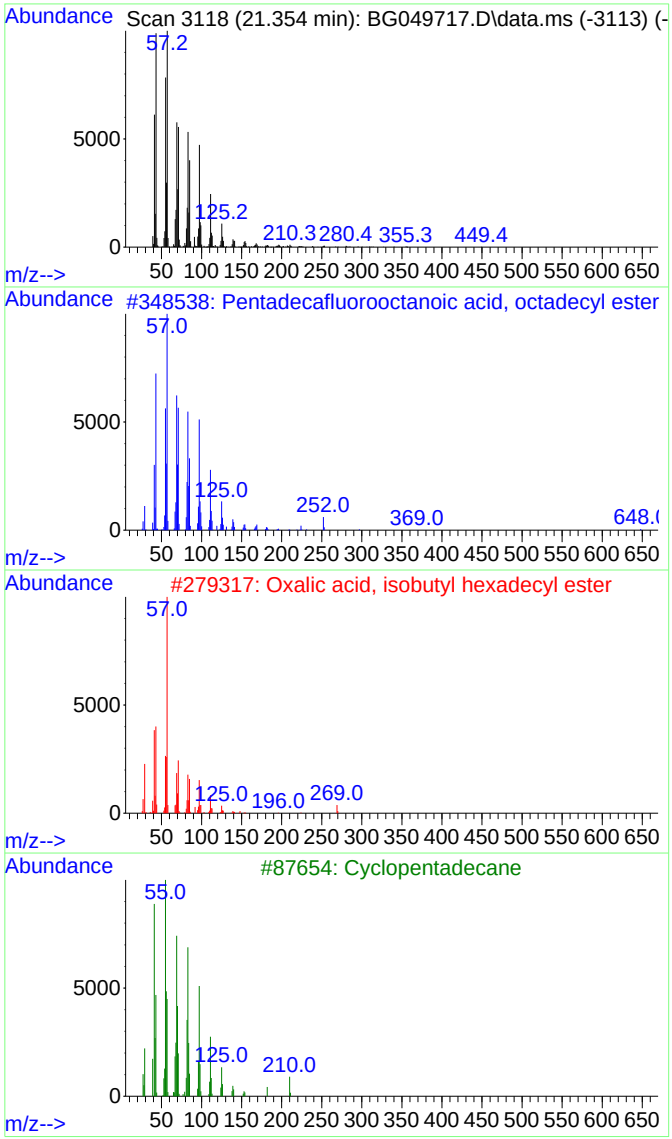
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

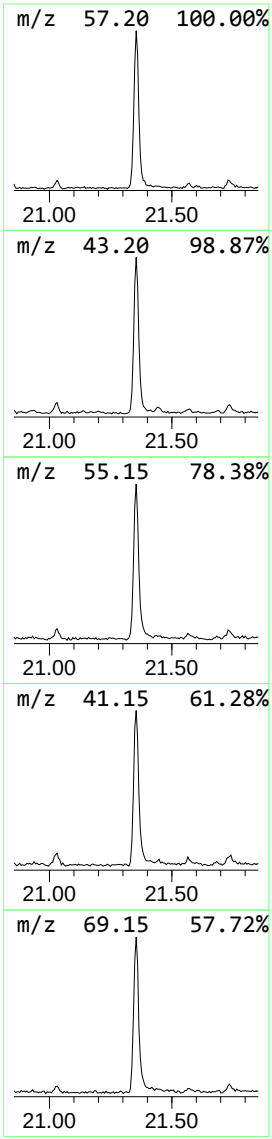
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Pentadecafluorooctanoic aci... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.354	58.40 ng/ul	1434660	Chrysene-d12	21.730

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93
2	Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	91
3	Cyclopentadecane	210	C15H30	000295-48-7	83
4	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	81
5	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	76





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Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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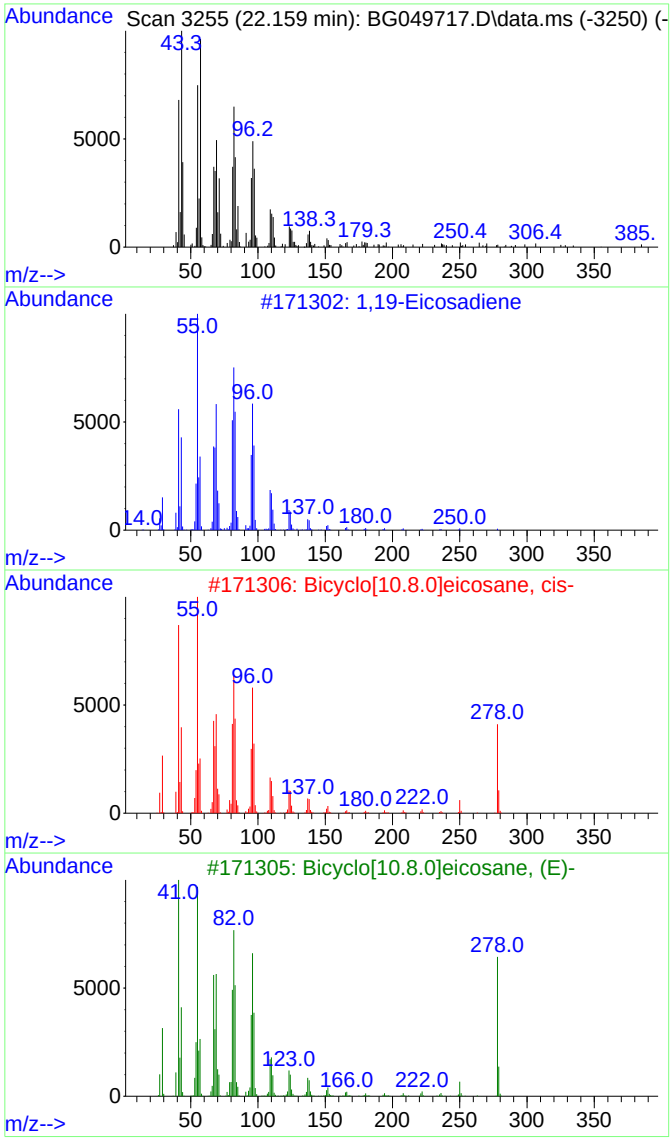
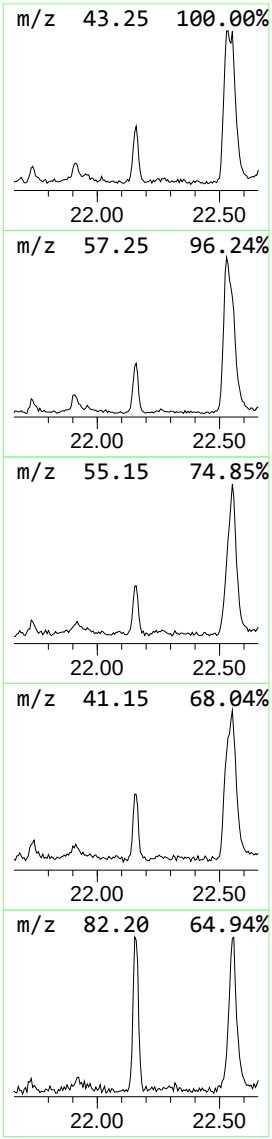
Peak Number 20 1,19-Eicosadiene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.159	12.62 ng/ul	309928	Chrysene-d12	21.730

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,19-Eicosadiene	278	C20H38	014811-95-1	95
2	Bicyclo[10.8.0]eicosane, cis-	278	C20H38	1000155-82-2	93
3	Bicyclo[10.8.0]eicosane, (E)-	278	C20H38	1010155-85-0	93
4	E-9-Tetradecen-1-ol formate	240	C15H28O2	1000130-78-2	90
5	(Z)-14-Tricosenyl formate	366	C24H46O2	077899-10-6	90





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Acq On : 7 Aug 2021 23:39
Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

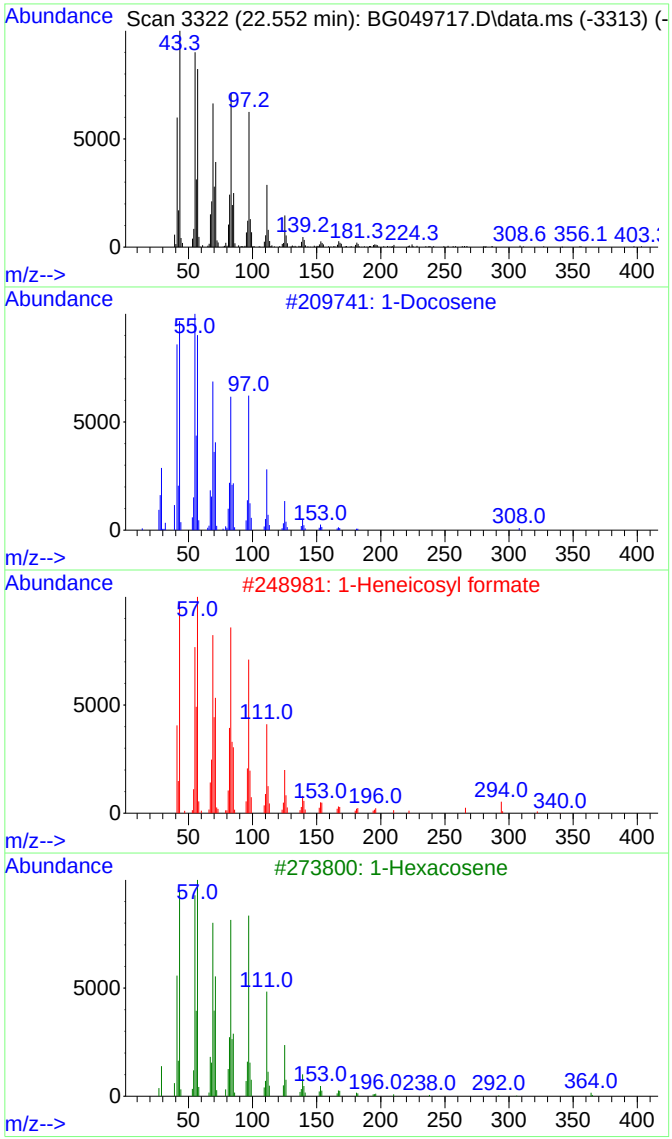
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

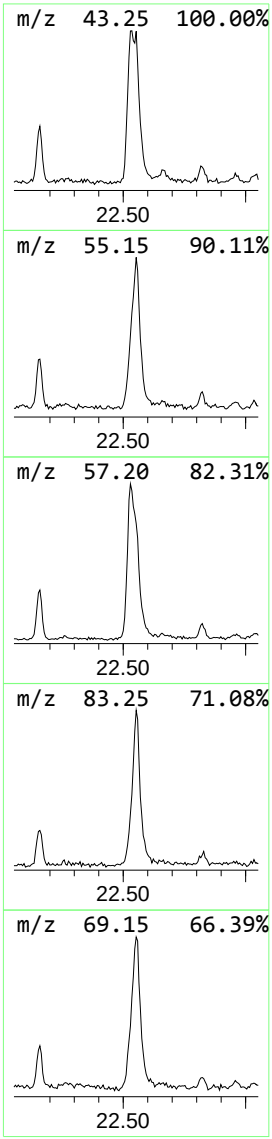
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.552	58.08 ng/ul	1426730	Chrysene-d12	21.730

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	98
2	1-Heneicosyl formate	340	C22H44O2	077899-03-7	95
3	1-Hexacosene	364	C26H52	018835-33-1	94
4	Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91
5	Octacosanol	410	C28H58O	000557-61-9	91





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Acq On : 7 Aug 2021 23:39
Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

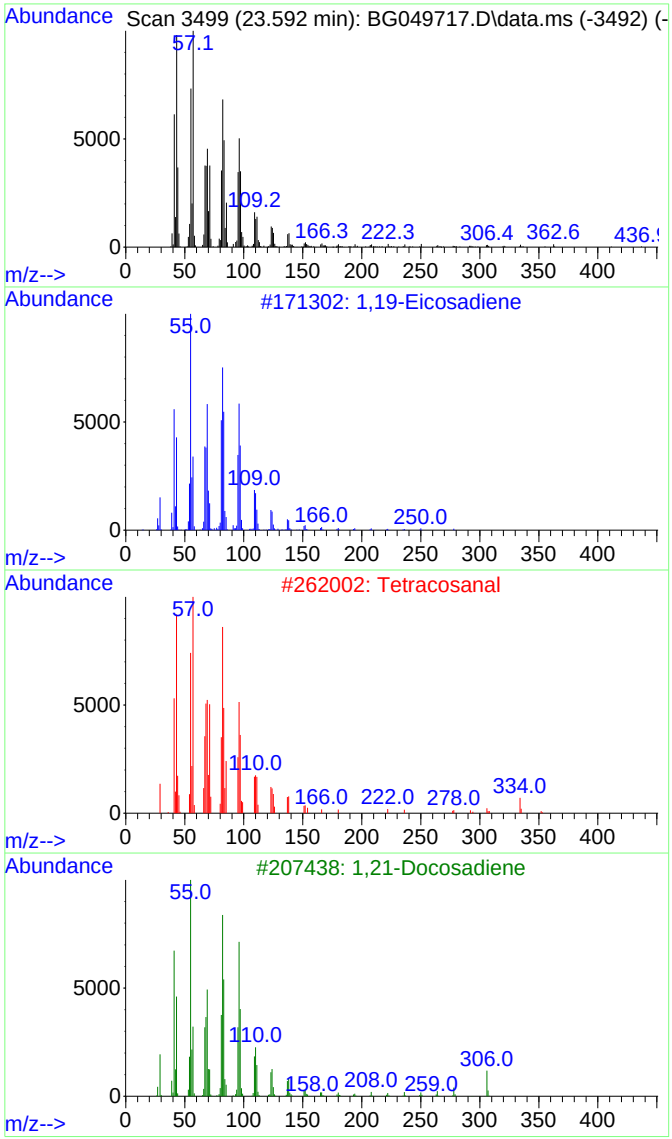
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

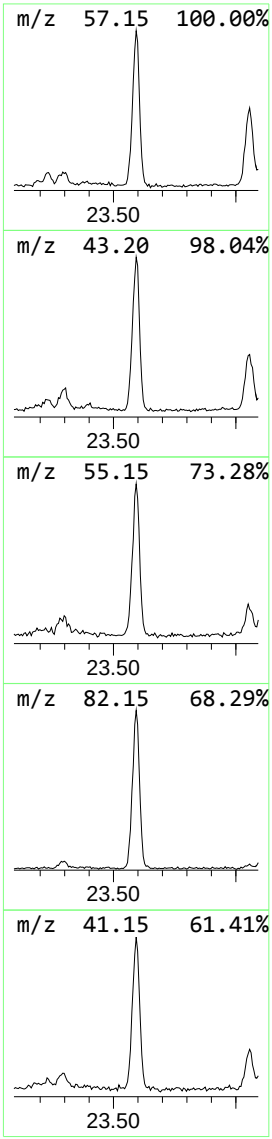
****

Peak Number 22 Tetracosanal Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.592	65.98 ng/ul	1415770	Perylene-d12	25.008

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,19-Eicosadiene	278	C20H38	014811-95-1	95
2	Tetracosanal	352	C24H48O	057866-08-7	94
3	1,21-Docosadiene	306	C22H42	053057-53-7	93
4	Eicosanal-	296	C20H40O	002400-66-0	91
5	(Z)-14-Tricosenyl formate	366	C24H46O2	077899-10-6	91





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Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

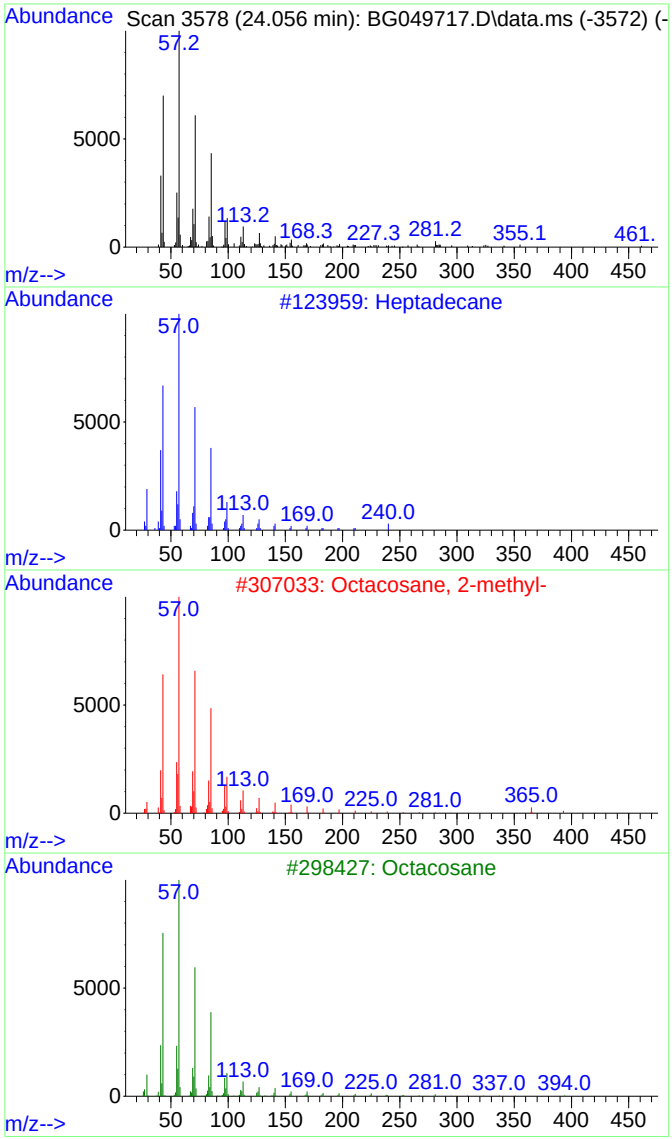
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

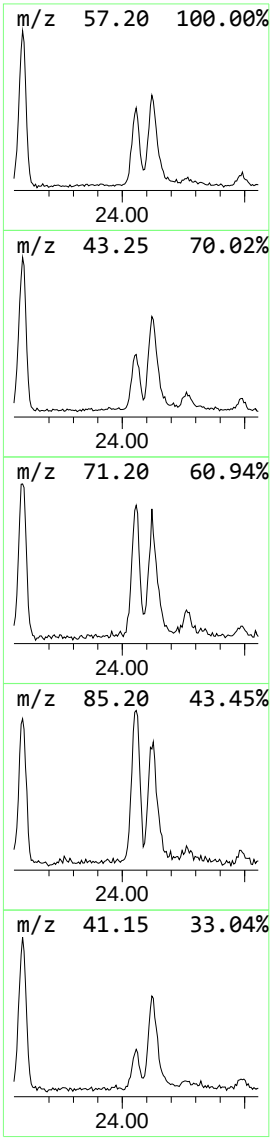
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.056	14.79 ng/ul	317262	Perylene-d12	25.008

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	91
2	Octacosane, 2-methyl-	408	C29H60	001560-98-1	90
3	Octacosane	394	C28H58	000630-02-4	90
4	Hexatriacontane	507	C36H74	000630-06-8	90
5	Tetratetracontane	619	C44H90	007098-22-8	90





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Acq On : 7 Aug 2021 23:39
Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

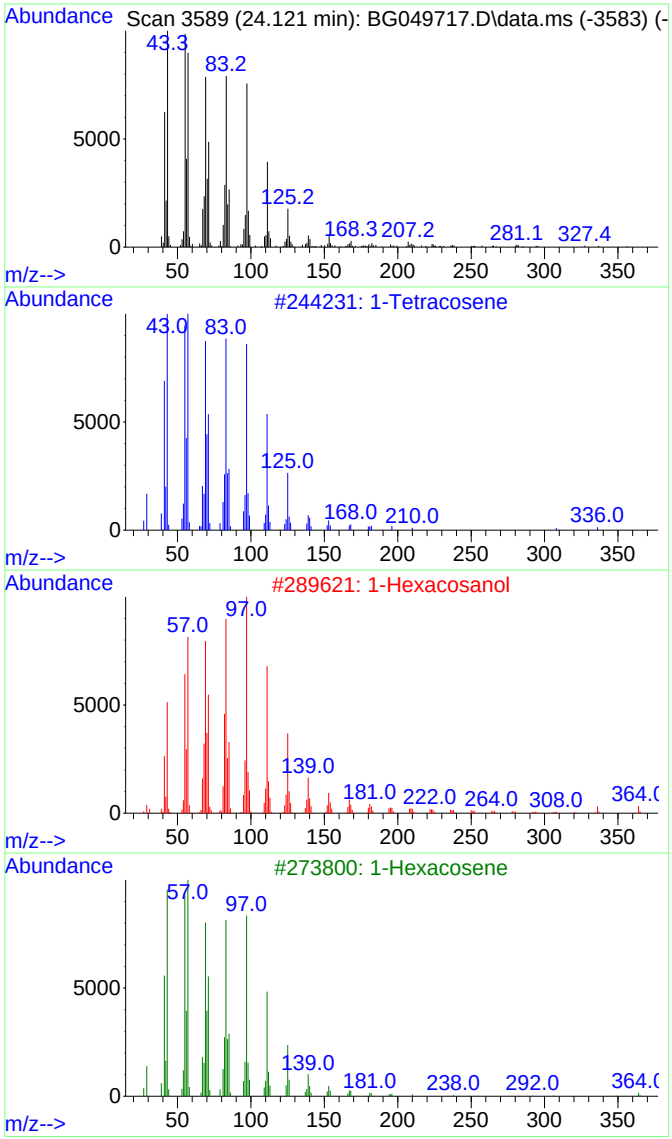
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

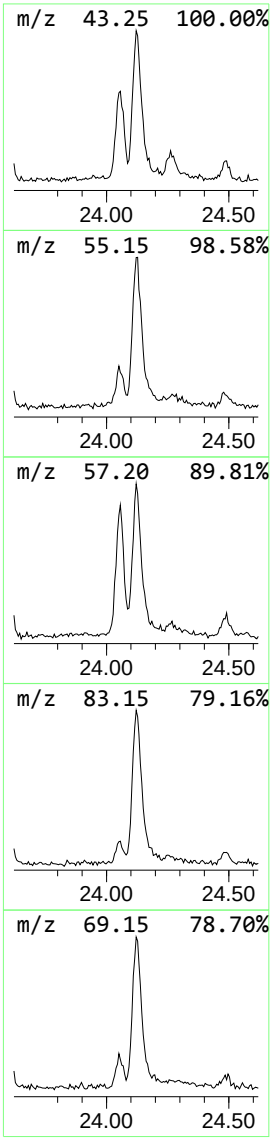
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.121	44.22 ng/ul	948763	Perylene-d12	25.008

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetracosene	336	C24H48	010192-32-2	91
2	1-Hexacosanol	382	C26H54O	000506-52-5	91
3	1-Hexacosene	364	C26H52	018835-33-1	91
4	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	91
5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	86





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Acq On : 7 Aug 2021 23:39
Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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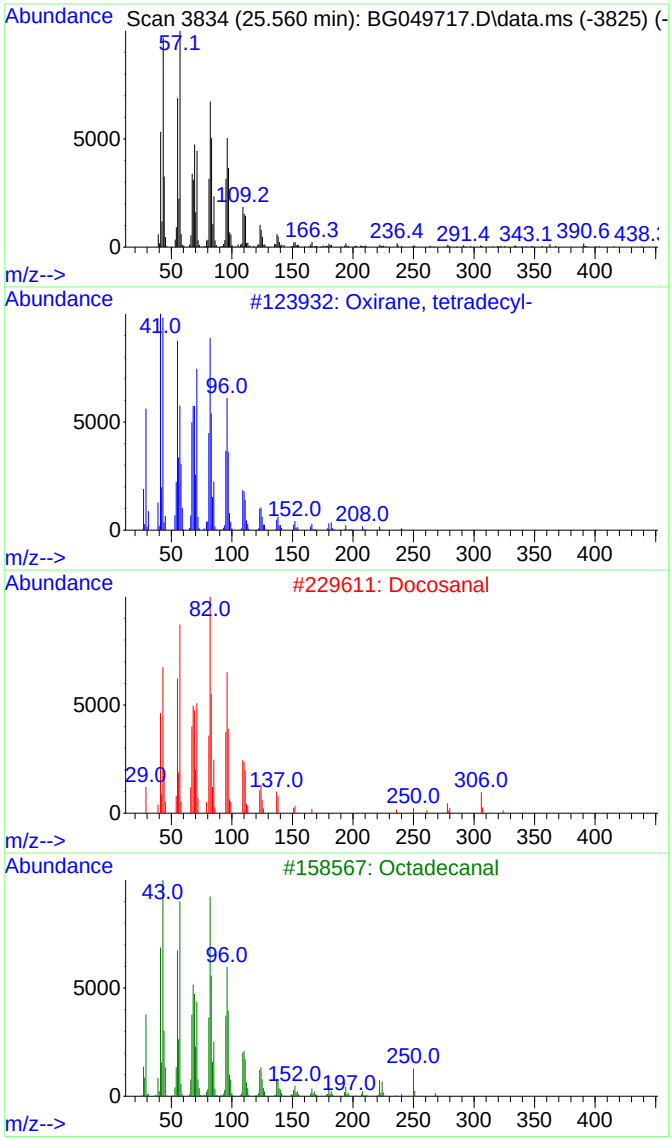
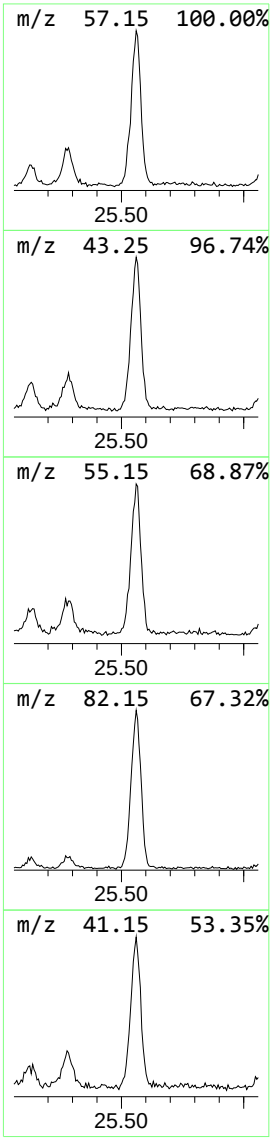
Peak Number 25 Oxirane, tetradecyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.560	64.52 ng/ul	1384540	Perylene-d12	25.008

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, tetradecyl-	240	C16H32O	007320-37-8	95
2	Docosanal	324	C22H44O	057402-36-5	91
3	Octadecanal	268	C18H36O	000638-66-4	90
4	1,19-Eicosadiene	278	C20H38	014811-95-1	89
5	E-9-Tetradecen-1-ol formate	240	C15H28O2	1000130-78-2	87





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Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

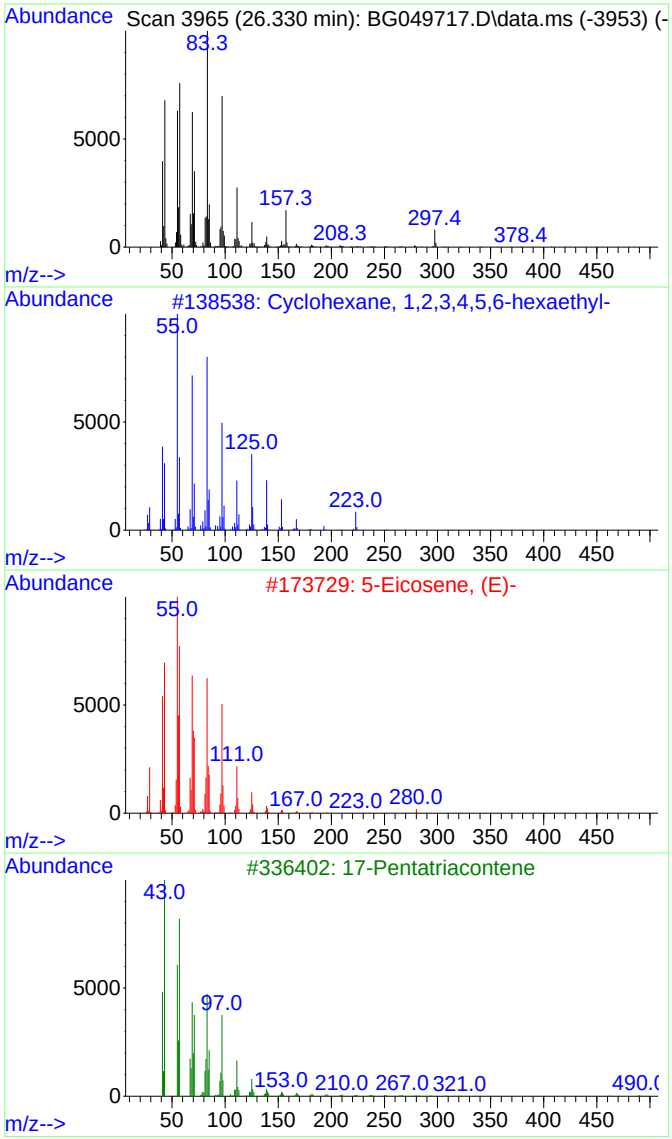
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

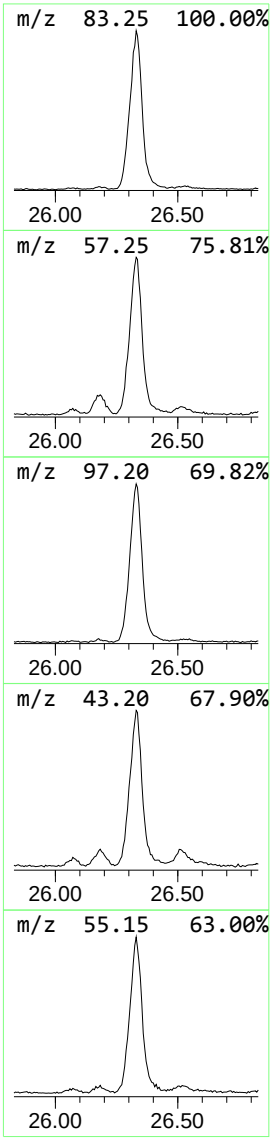
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 (DEL) Alkane: Cyclic26.330 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.330	170.67 ng/ul	3662160	Perylene-d12	25.008

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,2,3,4,5,6-hexaethyl-	252	C18H36	001795-14-8	72
2	5-Eicosene, (E)-	280	C20H40	074685-30-6	58
3	17-Pentatriacontene	491	C35H70	006971-40-0	50
4	Tetracosyl heptafluorobutyrate	550	C28H49F7O2	1000351-83-7	49
5	Hexacosyl heptafluorobutyrate	578	C30H53F7O2	1000351-83-3	49





Tentatively Identified Compound (LSC) summary

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Data File : BG049717.D
Acq On : 7 Aug 2021 23:39
Operator : CG/JU
Sample : M3244-01
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JCE01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: C...	3.079	12.5	ng/ul	138316	1	8.043	221161	20.0
Butane, 2-metho...	3.161	48.0	ng/ul	531036	1	8.043	221161	20.0
(DEL) Alkane: S...	4.412	6.2	ng/ul	68084	1	8.043	221161	20.0
unknown-01	4.829	4.3	ng/ul	47844	1	8.043	221161	20.0
Diethyltoluamide	15.427	2.6	ng/ul	54472	3	14.693	413233	20.0
unknown-02	15.497	3.5	ng/ul	73132	3	14.693	413233	20.0
Benzoic acid, 2...	16.572	2.1	ng/ul	44397	4	17.448	431010	20.0
Carbonic acid, ...	17.195	2.2	ng/ul	48278	4	17.448	431010	20.0
n-Hexadecanoic ...	18.329	8.5	ng/ul	183970	4	17.448	431010	20.0
Succinic acid, ...	18.646	3.7	ng/ul	79657	4	17.448	431010	20.0
unknown-03	18.981	2.6	ng/ul	56914	4	17.448	431010	20.0
unknown-04	19.257	4.3	ng/ul	93509	4	17.448	431010	20.0
(DEL) Alkane: C...	19.480	3.0	ng/ul	63961	4	17.448	431010	20.0
Eicosyl methyl ...	20.320	6.0	ng/ul	148571	5	21.730	491282	20.0
(DEL) Alkane: S...	20.350	6.7	ng/ul	165049	5	21.730	491282	20.0
unknown-05	20.679	5.6	ng/ul	137608	5	21.730	491282	20.0
Methyl dehydroa...	20.790	4.7	ng/ul	114764	5	21.730	491282	20.0
1-Phenanthrenem...	21.025	8.1	ng/ul	198501	5	21.730	491282	20.0
Pentadecafluoro...	21.354	58.4	ng/ul	1434660	5	21.730	491282	20.0
1,19-Eicosadiene	22.159	12.6	ng/ul	309928	5	21.730	491282	20.0
(DEL) Alkane: S...	22.552	58.1	ng/ul	1426730	5	21.730	491282	20.0
Tetracosanal	23.592	66.0	ng/ul	1415770	6	25.008	429158	20.0
(DEL) Alkane: S...	24.056	14.8	ng/ul	317262	6	25.008	429158	20.0
(DEL) Alkane: S...	24.121	44.2	ng/ul	948763	6	25.008	429158	20.0
Oxirane, tetrad...	25.560	64.5	ng/ul	1384540	6	25.008	429158	20.0
(DEL) Alkane: C...	26.330	170.7	ng/ul	3662160	6	25.008	429158	20.0