LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071421\
Data File : BG049359.D
Acq On : 15 Jul 2021 12:37
Operator : CG/JU
Sample : M2996-03DL 10X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT7DL

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-BG070921.M
Title : SVOA CALIBRATION

Signal : TIC: BG049359.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.109	6	12	22	rBV4	28704	67144	9.87%	1.413%
2	3.185	22	25	32	rVB4	10452	16596	2.44%	0.349%
3	3.549	85	87	89	rBV4	2230	2254	0.33%	0.047%
4	3.691	108	111	114	rVB3	1563	1551	0.23%	0.033%
5	4.213	196	200	210	rVB3	9065	15600	2.29%	0.328%
6	4.501	247	249	252	rBV3	1450	1670	0.25%	0.035%
7	4.719	282	286	290	rVB3	1672	2752	0.40%	0.058%
8	4.772	290	295	297	rBV4	1401	2326	0.34%	0.049%
9	4.901	314	317	318	rBV2	1431	1563	0.23%	0.033%
10	5.347	391	393	396	rBV	1262	1693	0.25%	0.036%
11	5.418	402	405	407	rBV2	1287	1609	0.24%	0.034%
12	5.535	422	425	427	rBV3	1244	1522	0.22%	0.032%
13	7.110	691	693	696	rVB	1321	1552	0.23%	0.033%
14	7.228	708	713	718	rBV3	2264	5117	0.75%	0.108%
15	7.386	734	740	747	rBV2	7689	13818	2.03%	0.291%
16	7.598	769	776	782	rVB3	8944	16649	2.45%	0.350%
17	8.062	848	855	862	rBV	112104	187866	27.61%	3.954%
18	8.185	870	876	879	rBV2	1775	2975	0.44%	0.063%
19	8.450	919	921	926	rVB3	2077	1935	0.28%	0.041%
20	8.773	971	976	982	rVV3	7537	12635	1.86%	0.266%
21	8.896	995	997	1000	rVV2	2017	2574	0.38%	0.054%
22	8.920	1000	1001	1005	rVV	2584	2172	0.32%	0.046%
23	9.231	1048	1054	1063	rVV2	12955	25080	3.69%	0.528%
24	9.725	1135	1138	1141	rBV2	1289	1699	0.25%	0.036%
25	9.960	1172	1178	1185	rBV5	9563	18365	2.70%	0.387%
26	10.277	1228	1232	1234	rBV	1937	1690	0.25%	0.036%
27	10.500	1265	1270	1277	rBV3	14892	28005	4.12%	0.589%
28	10.653	1291	1296	1297	rBV4	1375	1710	0.25%	0.036%
29	10.735	1306	1310	1315	rVB	999	1754	0.26%	0.037%
30	10.794	1315	1320	1323	rVB2	823	1540	0.23%	0.032%
31	10.882	1326	1335	1344	rVB	149523	267670	39.34%	5.634%
32	11.023	1354	1359	1366	rVB5	10128	18581	2.73%	0.391%
33	11.987	1519	1523	1526	rBV3	2399	2889	0.42%	0.061%
34	12.151	1547	1551	1554	rBV2	1328	1528	0.22%	0.032%
35	12.269	1566	1571	1575	rVV2	1709	2081	0.31%	0.044%
36	12.433	1595	1599	1600	rBV2	1307	1649	0.24%	0.035%

LSC Area Percent Report

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Sample : M2996-03DL 10X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT7DL

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-BG070921.M
Title : SVOA CALIBRATION

37	12.750	1651	1653	1658	rVB2	2441	2529	0.37%	0.053%
38	13.144	1716	1720	1728	rVV3	5338	9221	1.36%	0.194%
39	13.226	1730	1734	1740	rVB4	4539	6770	1.00%	0.143%
40	13.303	1743	1747	1748	rBV2	2010	2083	0.31%	0.044%
41	13.414	1762	1766	1773	rBV2	10578	18184	2.67%	0.383%
42	13.538	1785	1787	1793	rVB2	2134	2397	0.35%	0.050%
43	13.726	1815	1819	1825	rBV3	7771	12121	1.78%	0.255%
44	13.843	1832	1839	1841	rBV3	5015	6196	0.91%	0.130%
45	13.884	1844	1846	1850	rVB2	1804	1699	0.25%	0.036%
46	14.108	1878	1884	1890	rVV	35948	51153	7.52%	1.077%
47	14.401	1929	1934	1942	rVB	40950	60596	8.91%	1.275%
48	14.542	1954	1958	1963	rVB3	1884	3083	0.45%	0.065%
49	14.707	1979	1986	1995	rVB	247962	392444	57.68%	8.261%
50	14.842	2004	2009	2011	rBV2	1593	2180	0.32%	0.046%
51	14.913	2017	2021	2023	rBV2	2122	2354	0.35%	0.050%
52	15.013	2033	2038	2041	rBV2	998	1826	0.27%	0.038%
53	15.253	2076	2079	2083	rBV5	3037	3991	0.59%	0.084%
54	15.289	2083	2085	2087	rVV	2244	2048	0.30%	0.043%
55	15.336	2087	2093	2099	rVB3	14164	25961	3.82%	0.546%
56	15.600	2136	2138	2140	rBV	1364	1484	0.22%	0.031%
57	15.700	2148	2155	2160	rBV	50172	76992	11.32%	1.621%
58	15.753	2162	2164	2170	rVB2	3076	3822	0.56%	0.080%
59	15.835	2170	2178	2187	rBV	71221	121970	17.93%	2.567%
60	16.005	2206	2207	2211	rBV2	1522	1617	0.24%	0.034%
61	16.058	2214	2216	2218	rVB2	1866	1750	0.26%	0.037%
62	16.229	2240	2245	2249	rBV2	1715	2269	0.33%	0.048%
63	16.558	2298	2301	2304	rVV3	1114	1558	0.23%	0.033%
64	16.599	2304	2308	2310	rVV2	1675	2121	0.31%	0.045%
65	16.728	2326	2330	2334	rBV	1926	2620	0.39%	0.055%
66	17.110	2389	2395	2405	rBV	430981	680324	100.00%	14.320%
67	17.222	2413	2414	2415	rBV	2624	1606	0.24%	0.034%
68	17.463	2448	2455	2466	rVB	306507	475894	69.95%	10.017%
69	17.557	2466	2471	2477	rBV2	53904	82239	12.09%	1.731%
70	17.621	2480	2482	2483	rBV2	1906	1535	0.23%	0.032%
71	17.651	2483	2487	2489	rVV5	1816	2375	0.35%	0.050%
72	17.809	2509	2514	2527	rBV2	56458	102705	15.10%	2.162%
73	18.056	2554	2556	2558	rBV2	2182	1730	0.25%	0.036%
74	18.115	2563	2566	2571	rVV3	4838	6786	1.00%	0.143%
75	18.303	2595	2598	2602	rBV3	2952	4516	0.66%	0.095%
76	18.338	2602	2604	2605	rBV2	1958	1555	0.23%	0.033%

LSC Area Percent Report

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Acq On : 15 Jul 2021 12:37
Operator : CG/JU
Sample : M2996-03DL 10X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT7DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
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Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Title : SVOA CALIBRATION

77	18.409	2613	2616	2618	rBV2	1587	1479	0.22%	0.031%
78	18.638	2652	2655	2656	rBV	2086	1606	0.24%	0.034%
79	18.855	2689	2692	2695	rBV2	1816	2249	0.33%	0.047%
80	19.090	2730	2732	2736	rVB2	1946	1910	0.28%	0.040%
81	19.466	2793	2796	2797	rBV2	1892	1608	0.24%	0.034%
82	19.543	2804	2809	2813	rBV	98368	115387	16.96%	2.429%
83	19.707	2832	2837	2843	rBV2	86410	139257	20.47%	2.931%
84	19.754	2843	2845	2849	rVB3	20736	22434	3.30%	0.472%
85	19.842	2855	2860	2867	rBV	63496	93855	13.80%	1.976%
86	20.036	2891	2893	2895	rBV2	2194	1670	0.25%	0.035%
87	20.230	2924	2926	2930	rVB4	2574	2742	0.40%	0.058%
88	20.535	2974	2978	2984	rBV7	7606	15621	2.30%	0.329%
89	20.753	3013	3015	3018	rBV3	3089	2623	0.39%	0.055%
90	20.800	3018	3023	3028	rVV8	13299	19144	2.81%	0.403%
91	21.217	3090	3094	3104	rBV	43491	72954	10.72%	1.536%
92	21.740	3177	3183	3190	rBV	290766	493210	72.50%	10.382%
93	22.921	3379	3384	3391	rBV8	12453	29092	4.28%	0.612%
94	24.795	3699	3703	3713	rVB2	30536	78417	11.53%	1.651%
95	25.036	3734	3744	3757	rVB	188355	580321	85.30%	12.215%
96	26.217	3936	3945	3955	rBV6	23980	78072	11.48%	1.643%
97	27.621	4178	4184	4193	rVB5	20516	60893	8.95%	1.282%
98	29.037	4423	4425	4428	rBV4	4432	5185	0.76%	0.109%
99	31.293	4807	4809	4810	rBV2	7056	5292	0.78%	0.111%
100	31.552	4851	4853	4856	rBV4	4004	3690	0.54%	0.078%

Sum of corrected areas: 4750829

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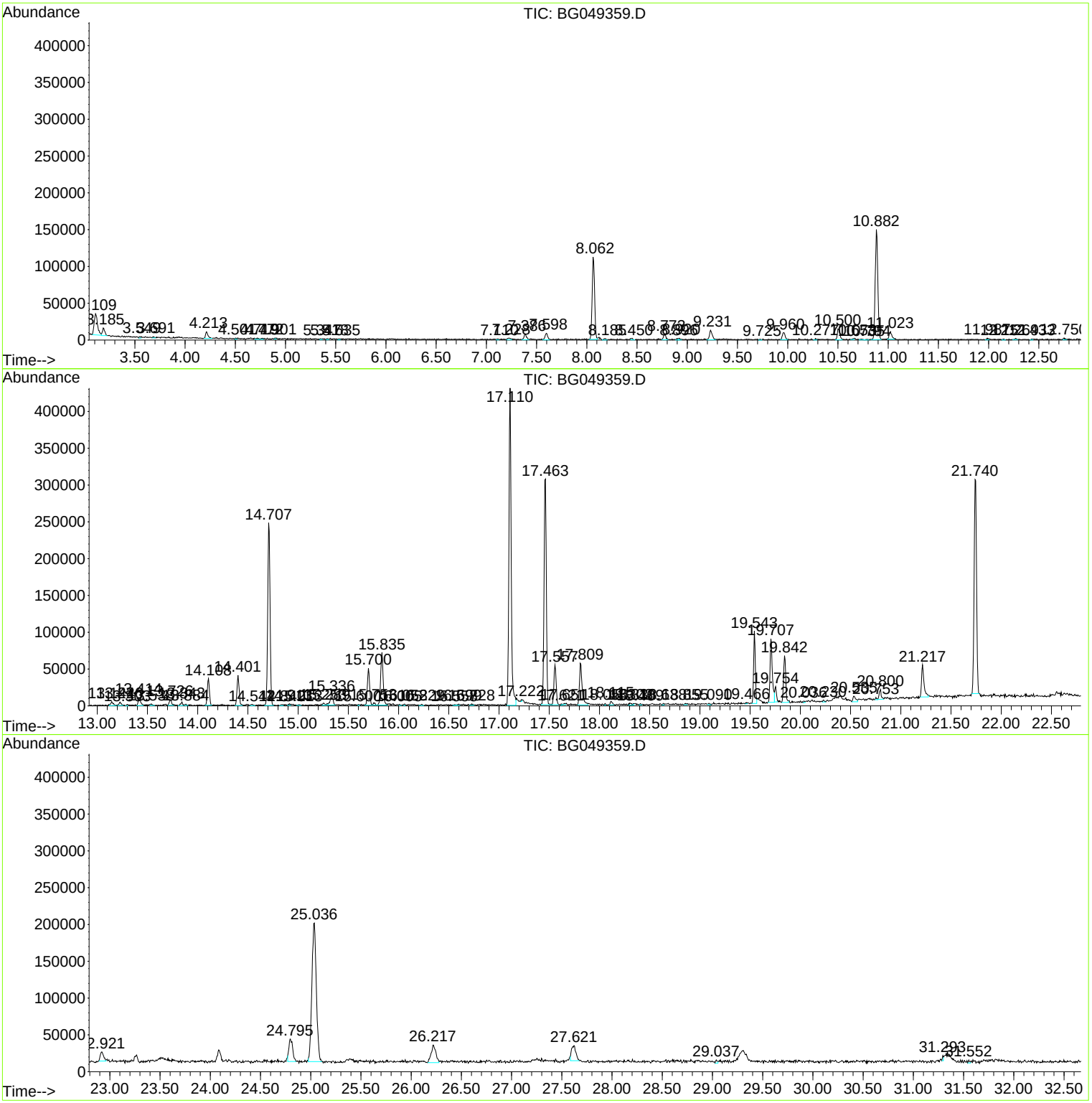
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AT7DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071421\
Data File : BG049359.D
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Instrument : BNA_G
ClientSampleId : C0AT7DL

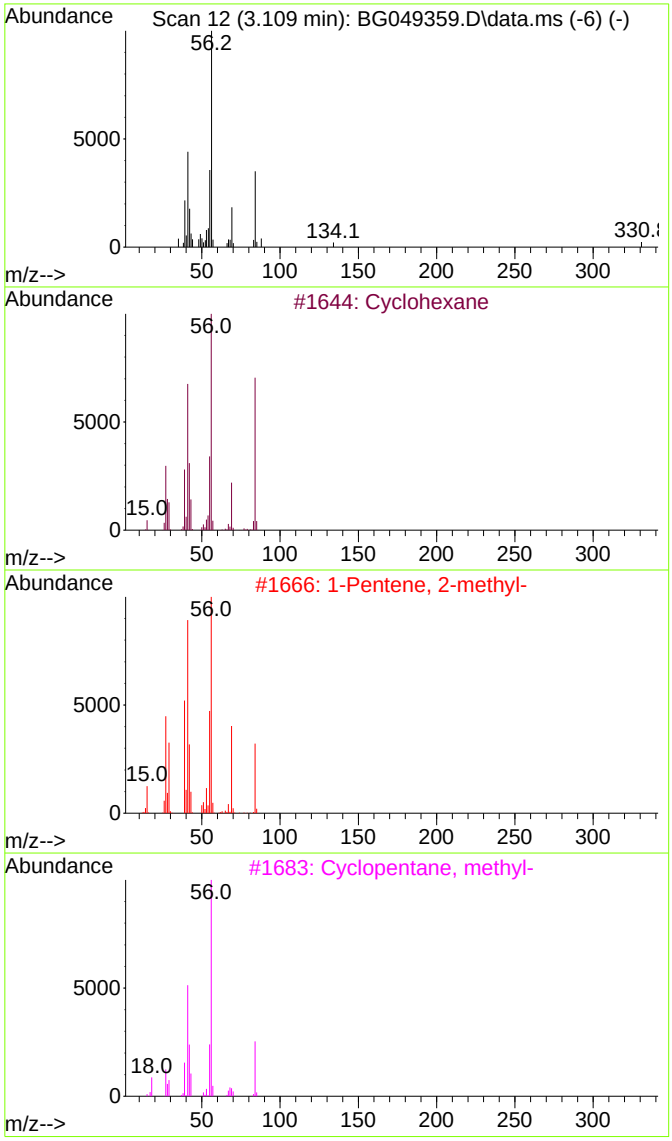
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

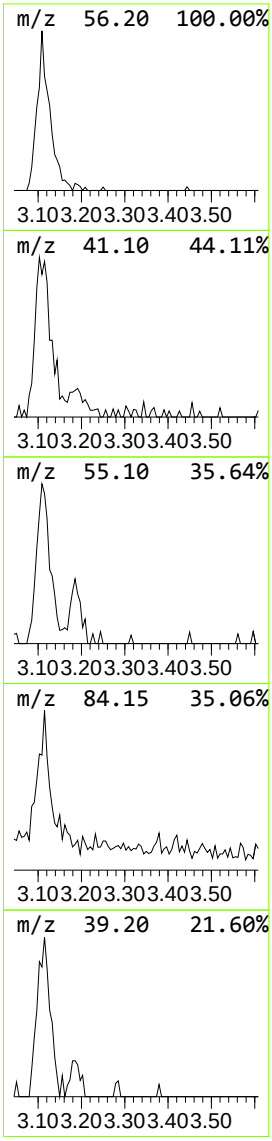
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Cyclic3.11 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.110	7.15 ng/ul	67144	1,4-Dichlorobenzene-d4	8.062

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane	84	C6H12	000110-82-7	74
2	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	64
3	Cyclopentane, methyl-	84	C6H12	000096-37-7	64
4	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	59
5	Cyclohexane	84	C6H12	000110-82-7	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071421\
Data File : BG049359.D
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Sample : M2996-03DL 10X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AT7DL

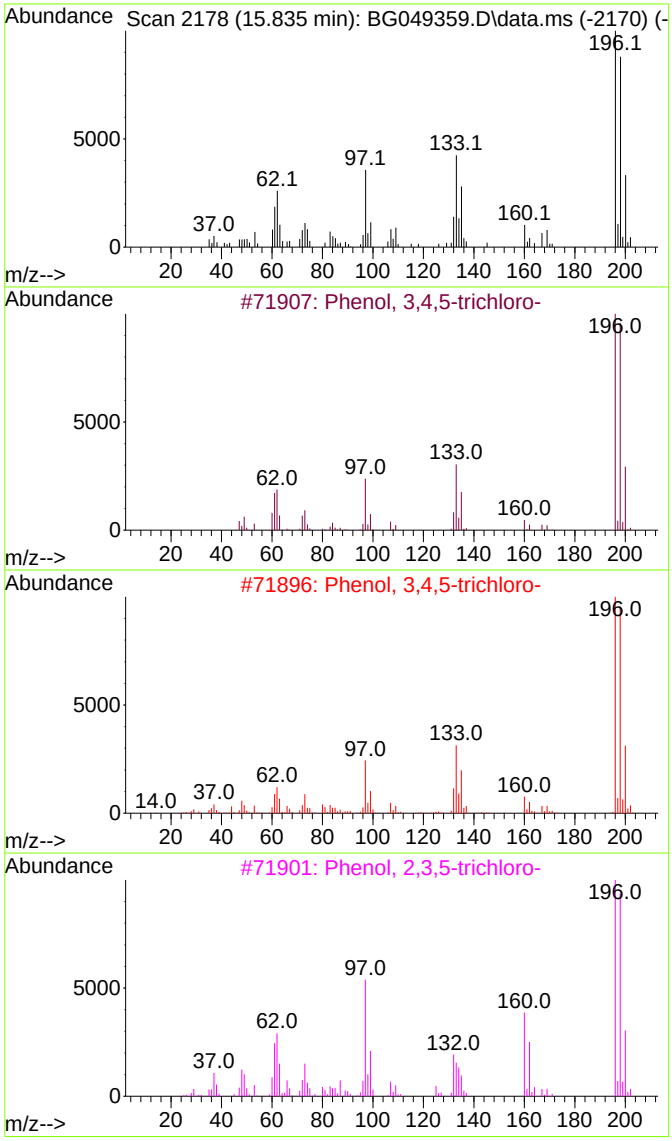
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

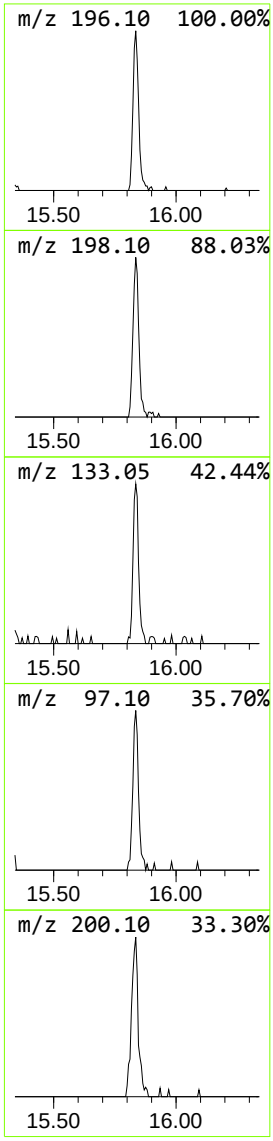
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Phenol, 3,4,5-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.840	6.22 ng/ul	121970	Acenaphthene-d10	14.707

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,4,5-trichloro-	196	C6H3Cl3O	000609-19-8	95
2	Phenol, 3,4,5-trichloro-	196	C6H3Cl3O	000609-19-8	91
3	Phenol, 2,3,5-trichloro-	196	C6H3Cl3O	000933-78-8	90
4	Phenol, 2,3,5-trichloro-	196	C6H3Cl3O	000933-78-8	90
5	Phenol, 2,3,6-trichloro-	196	C6H3Cl3O	000933-75-5	81





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Instrument : BNA_G
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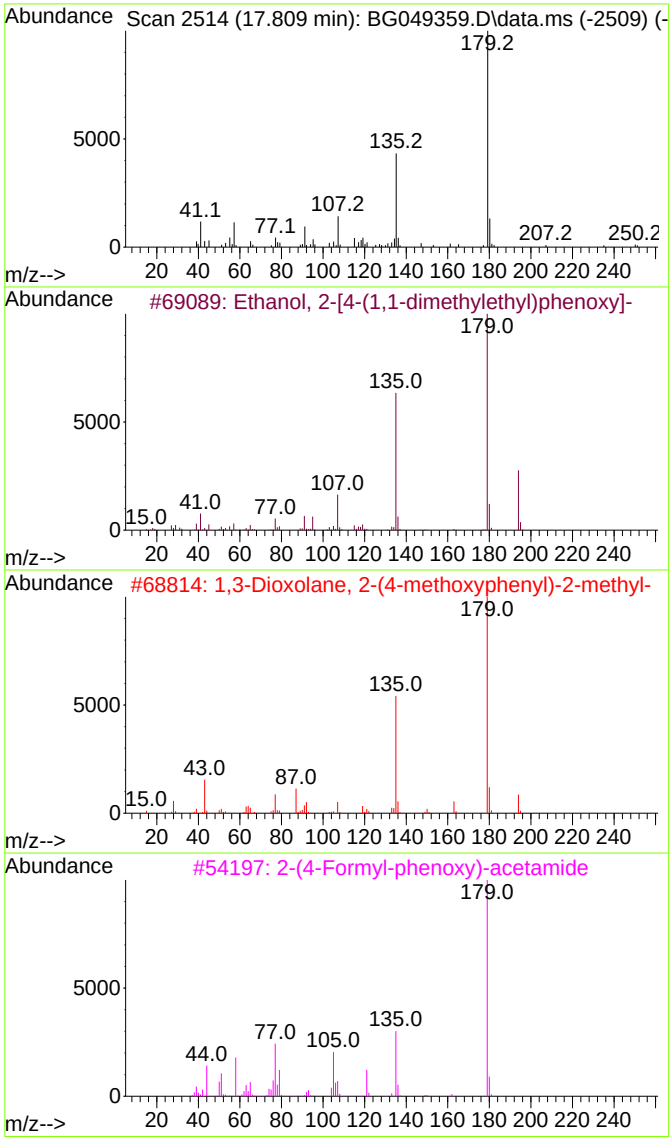
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

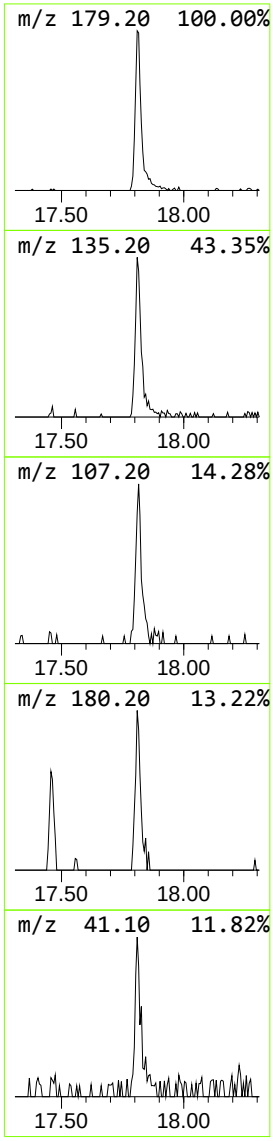
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Ethanol, 2-[4-(1,1-dimethyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.810	4.32 ng/ul	102705	Phenanthrene-d10	17.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-[4-(1,1-dimethylethyl...	194	C12H18O2	000713-46-2	78
2	1,3-Dioxolane, 2-(4-methoxypheny...	194	C11H14O3	036881-00-2	72
3	2-(4-Formyl-phenoxy)-acetamide	179	C9H9NO3	135857-20-4	64
4	Ethanol, 2-[4-(1,1-dimethylpropy...	208	C13H20O2	006382-07-6	64
5	2-Mercapto-3-(trifluoromethyl)py...	221	C8H6F3NOS	1000463-18-0	50





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ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT7DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

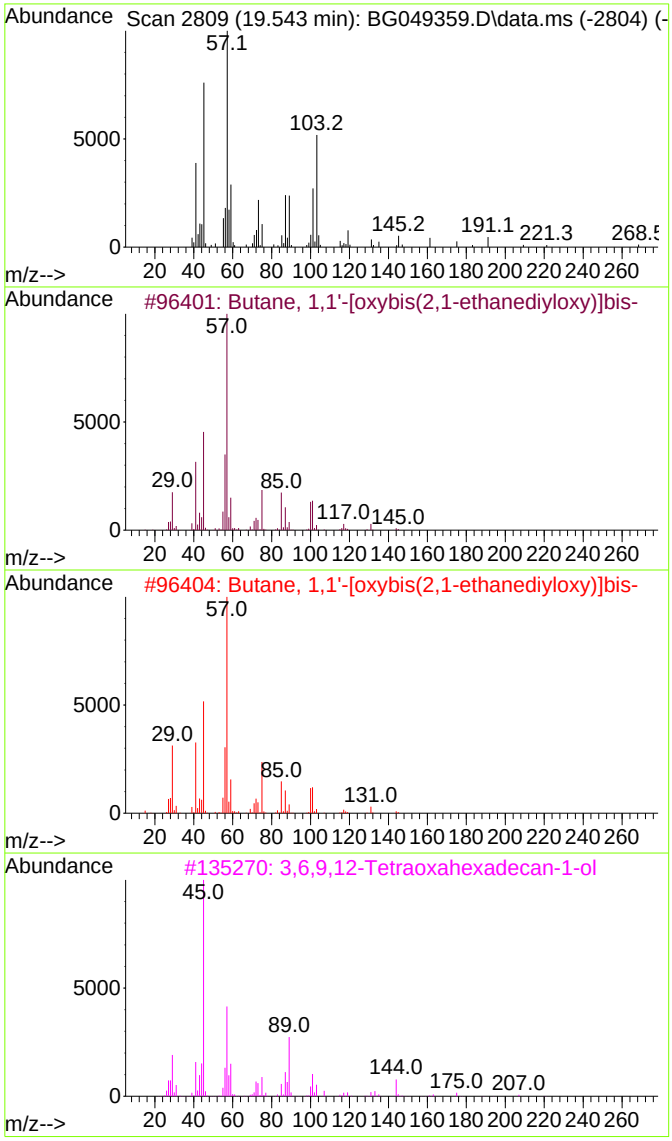
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

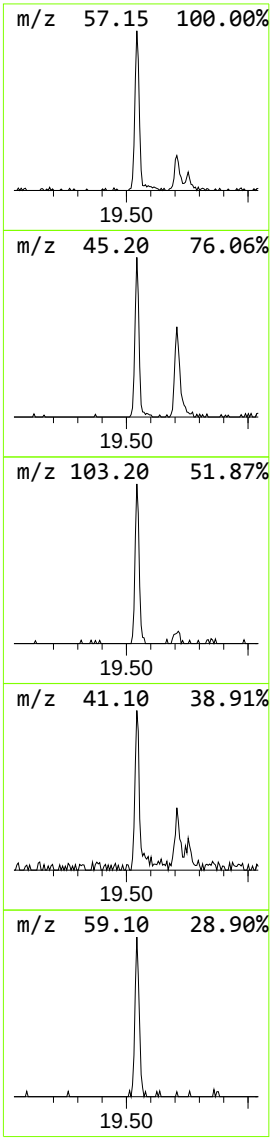
****

Peak Number 4 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.540	4.85 ng/ul	115387	Phenanthrene-d10	17.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 1,1'-[oxybis(2,1-ethaned...	218	C12H26O3	000112-73-2	47
2	Butane, 1,1'-[oxybis(2,1-ethaned...	218	C12H26O3	000112-73-2	47
3	3,6,9,12-Tetraoxahexadecan-1-ol	250	C12H26O5	001559-34-8	43
4	1,4,7,10,13,16-Hexaoxacyclooctad...	264	C12H24O6	017455-13-9	43
5	Di-sec-Butyl ether	130	C8H18O	006863-58-7	43





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Data File : BG049359.D
Acq On : 15 Jul 2021 12:37
Operator : CG/JU
Sample : M2996-03DL 10X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AT7DL

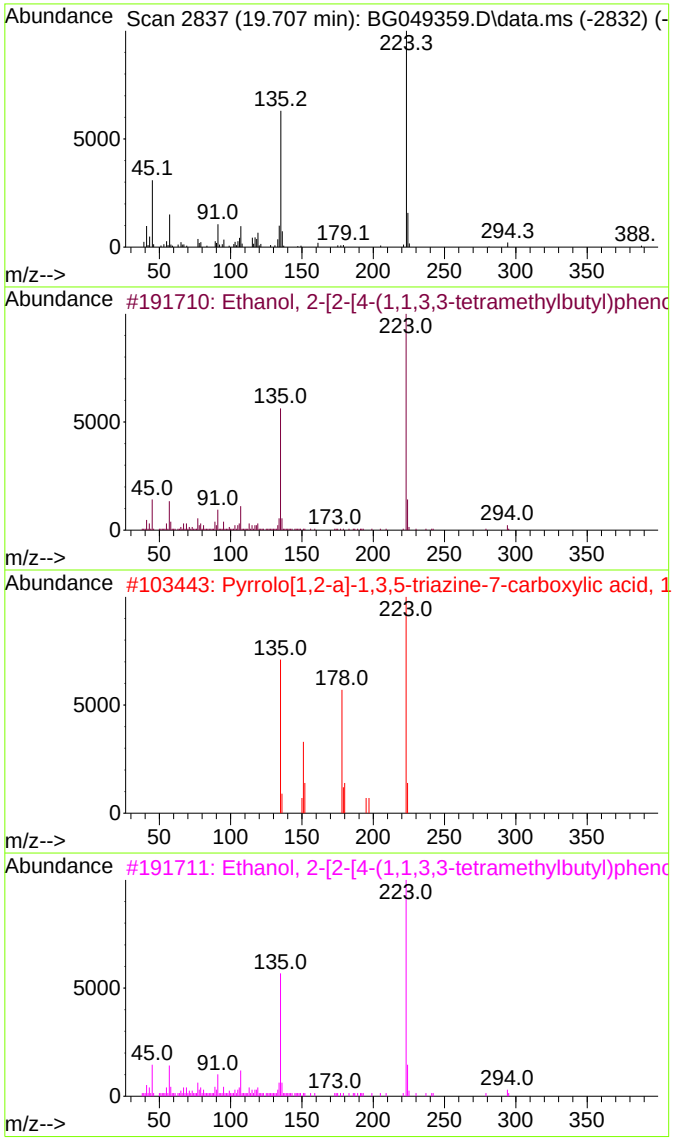
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

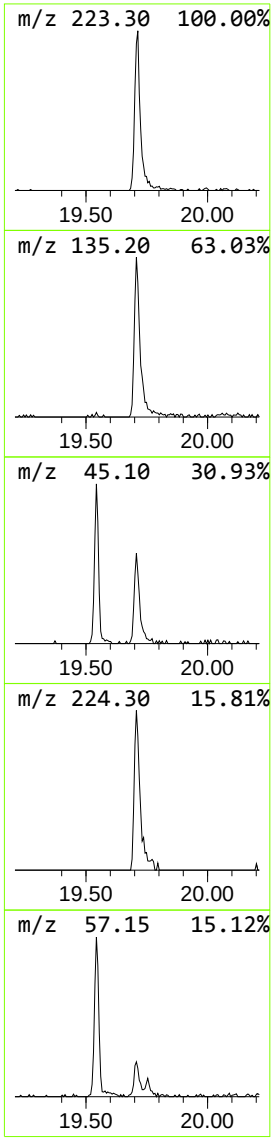
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Ethanol, 2-[2-[4-(1,1,3,3-t... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.710	5.65 ng/ul	139257	Chrysene-d12	21.746

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-[2-[4-(1,1,3,3-tetram...	294	C18H30O3	002315-61-9	93
2	Pyrrolo[1,2-a]-1,3,5-triazine-7-...	223	C9H9N3O4	054449-89-7	90
3	Ethanol, 2-[2-[4-(1,1,3,3-tetram...	294	C18H30O3	002315-61-9	72
4	Ethanol, 2-[2-[4-(1,1,3,3-tetram...	294	C18H30O3	002315-61-9	70
5	5-Methyl-2-(4-methylphenyl)-1,3-...	223	C15H13NO	016155-94-5	46





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Data File : BG049359.D
Acq On : 15 Jul 2021 12:37
Operator : CG/JU
Sample : M2996-03DL 10X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT7DL

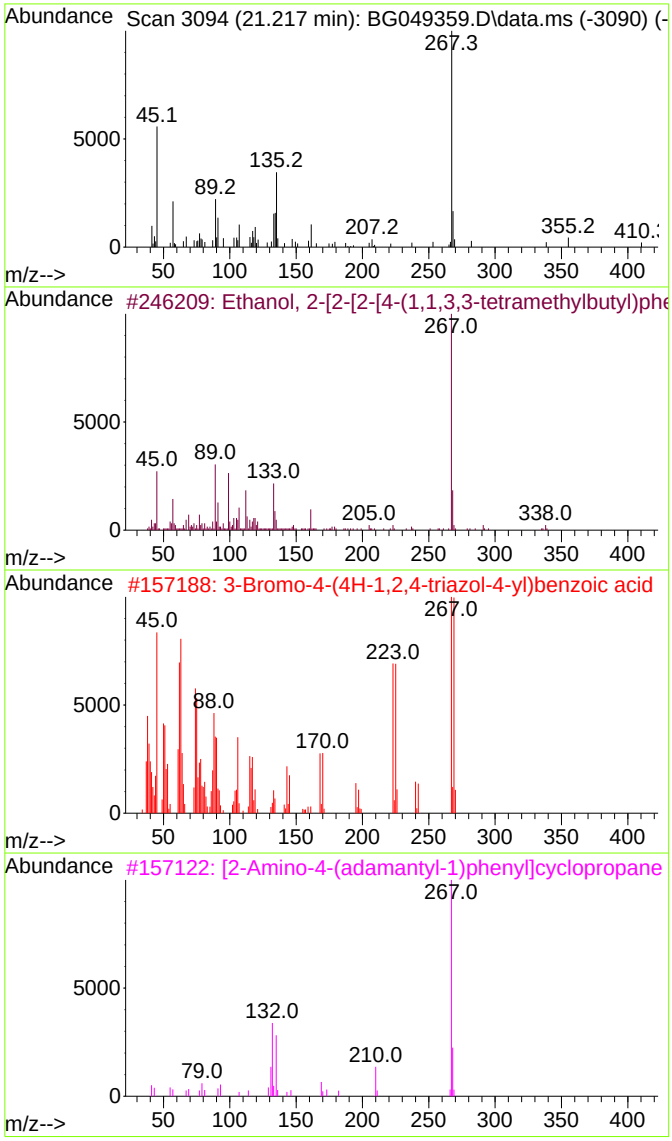
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

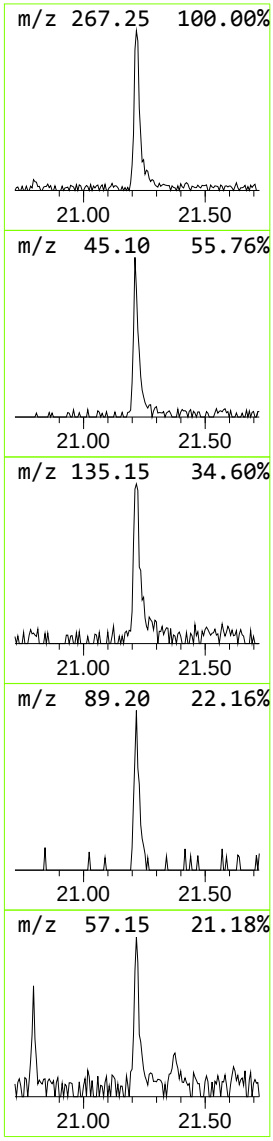
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Ethanol, 2-[2-[2-[4-(1,1,3,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.220	2.96 ng/ul	72954	Chrysene-d12	21.746

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-[2-[2-[4-(1,1,3,3-tet...	338	C20H34O4	002315-62-0	50
2	3-Bromo-4-(4H-1,2,4-triazol-4-yl...	267	C9H6BrN3O2	1000386-73-9	43
3	[2-Amino-4-(adamantyl-1)phenyl]c...	267	C19H25N	1010140-61-3	43
4	13H-Dibenzo[a,i]carbazole	267	C20H13N	000239-64-5	38
5	Propanamide, N-ethyl-N-(3-methyl...	267	C18H21NO	1010308-21-3	38





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Sample : M2996-03DL 10X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT7DL

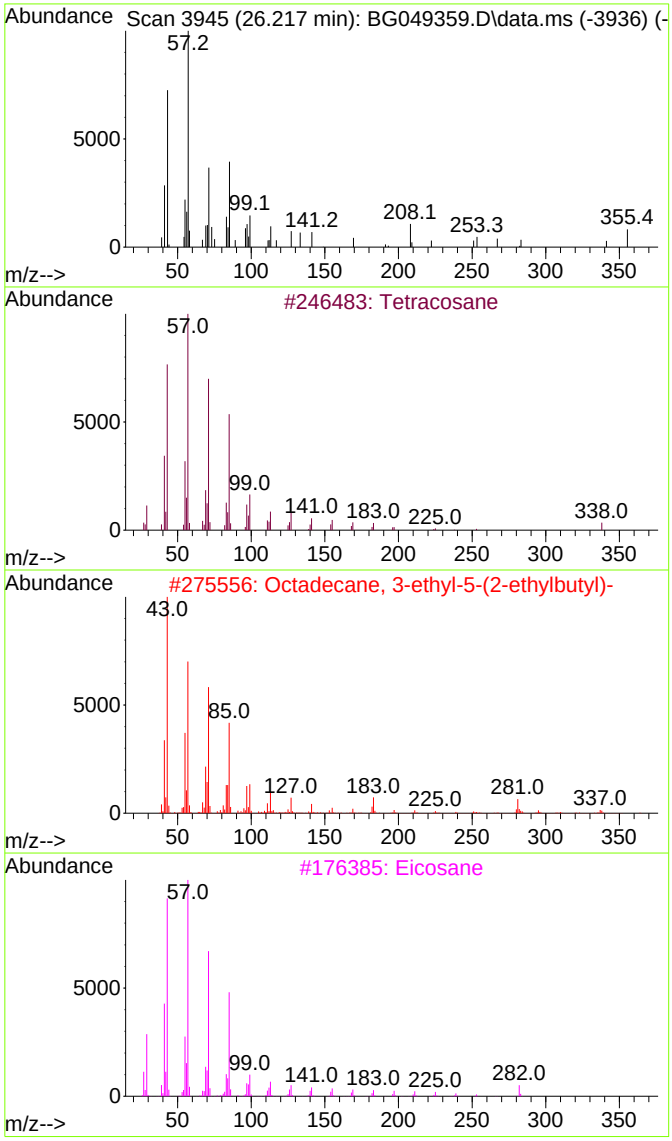
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

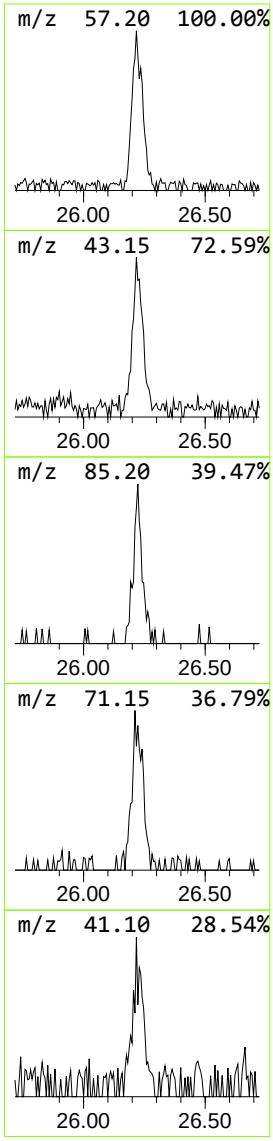
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.220	2.69 ng/ul	78072	Perylene-d12	25.030

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	64
2	Octadecane, 3-ethyl-5-(2-ethylbu...	366	C26H54	055282-12-7	59
3	Eicosane	282	C20H42	000112-95-8	58
4	Heptadecane	240	C17H36	000629-78-7	53
5	Octadecane	254	C18H38	000593-45-3	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071421\
Data File : BG049359.D
Acq On : 15 Jul 2021 12:37
Operator : CG/JU
Sample : M2996-03DL 10X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT7DL

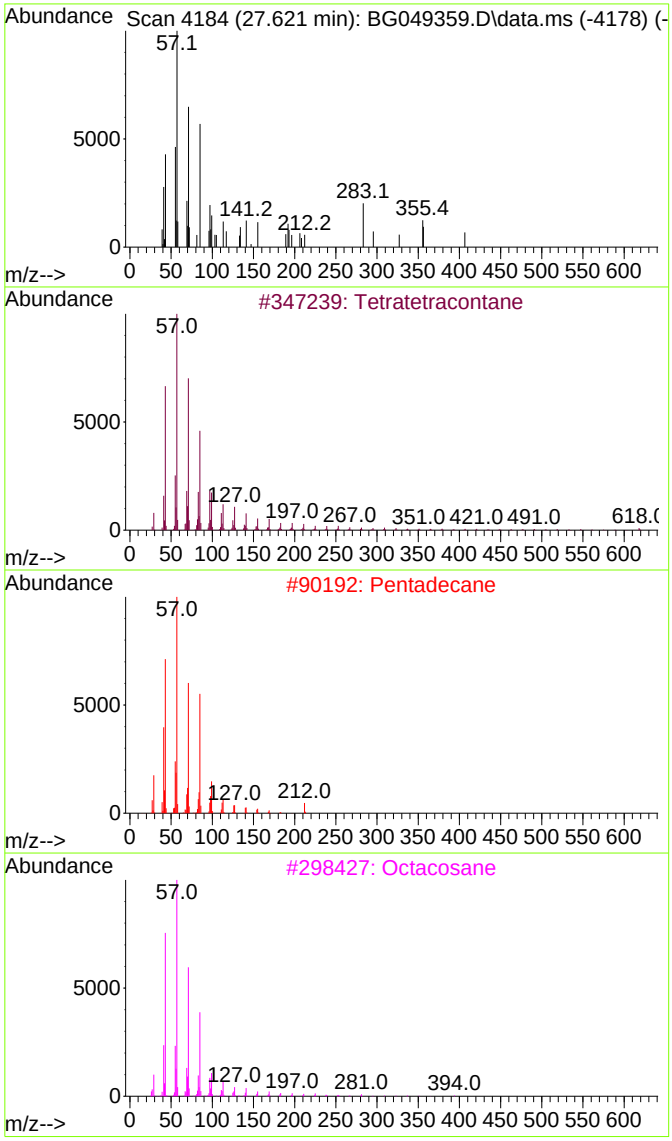
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
Quant Title : SVOA CALIBRATION

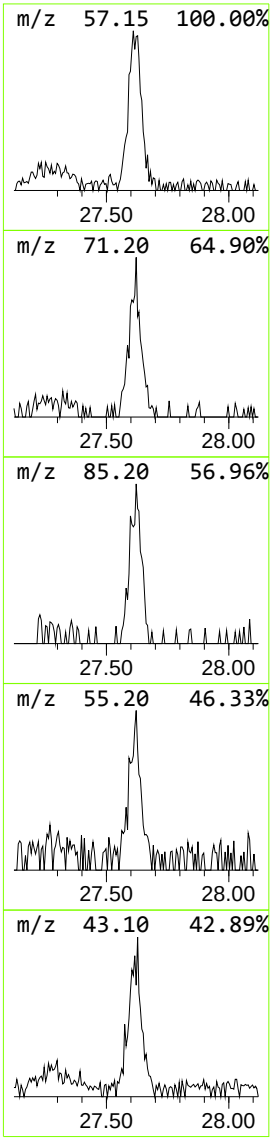
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.620	2.10 ng/ul	60893	Perylene-d12	25.030

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetratetracontane	619	C44H90	007098-22-8	47
2	Pentadecane	212	C15H32	000629-62-9	45
3	Octacosane	394	C28H58	000630-02-4	43
4	Hexatriacontane	507	C36H74	000630-06-8	43
5	Pentadecane	212	C15H32	000629-62-9	41





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071421\
Data File : BG049359.D
Acq On : 15 Jul 2021 12:37
Operator : CG/JU
Sample : M2996-03DL 10X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AT7DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.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...	3.110	7.2	ng/ul	67144	1	8.062	187866	20.0
Phenol, 3,4,5-t...	15.840	6.2	ng/ul	121970	3	14.707	392444	20.0
Ethanol, 2-[4-(...	17.810	4.3	ng/ul	102705	4	17.463	475894	20.0
unknown-01	19.540	4.8	ng/ul	115387	4	17.463	475894	20.0
Ethanol, 2-[2-[...	19.710	5.7	ng/ul	139257	5	21.746	493210	20.0
Ethanol, 2-[2-[...	21.220	3.0	ng/ul	72954	5	21.746	493210	20.0
(DEL) Alkane: S...	26.220	2.7	ng/ul	78072	6	25.030	580321	20.0
(DEL) Alkane: S...	27.620	2.1	ng/ul	60893	6	25.030	580321	20.0