LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082124\
Data File : BG062514.D
Acq On : 21 Aug 2024 18:46
Operator : MA/JU
Sample : P3626-21
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DCYC6

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-BG081424.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BG062514.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.407	16	20	25	rBV	14300	19901	1.20%	0.089%
2	3.665	60	64	74	rVB	94530	140940	8.49%	0.633%
3	3.812	84	89	101	rBV	189347	298670	17.99%	1.342%
4	5.058	295	301	307	rVB5	3796	7550	0.45%	0.034%
5	6.027	461	466	470	rBV	7247	9896	0.60%	0.044%
6	6.926	614	619	624	rBV4	3403	5689	0.34%	0.026%
7	7.102	641	649	659	rBV	258895	444032	26.74%	1.995%
8	7.266	669	677	685	rBV	182585	298143	17.96%	1.340%
9	7.472	704	712	718	rBV	246280	406900	24.51%	1.829%
10	7.525	718	721	728	rVB	30627	45127	2.72%	0.203%
11	7.942	785	792	798	rBV	248820	419460	25.26%	1.885%
12	7.995	798	801	809	rVB4	6500	11454	0.69%	0.051%
13	8.635	903	910	920	rBV	281987	482053	29.03%	2.166%
14	9.087	980	987	1000	rVB	227283	407816	24.56%	1.833%
15	9.646	1076	1082	1090	rBV2	30238	51314	3.09%	0.231%
16	9.810	1102	1110	1122	rBV	198239	339509	20.45%	1.526%
17	10.350	1194	1202	1216	rBV	320133	573193	34.52%	2.576%
18	10.732	1259	1267	1278	rBV	399511	692085	41.68%	3.110%
19	10.862	1278	1289	1297	rBV	240738	424194	25.55%	1.906%
20	11.255	1350	1356	1364	rVB2	14803	24605	1.48%	0.111%
21	11.467	1385	1392	1409	rBV	79024	157480	9.48%	0.708%
22	12.313	1531	1536	1541	rVB2	6086	9691	0.58%	0.044%
23	13.393	1714	1720	1725	rBV3	3974	6660	0.40%	0.030%
24	13.693	1765	1771	1779	rBV	9863	27229	1.64%	0.122%
25	13.963	1811	1817	1825	rVV	589336	836052	50.35%	3.757%
26	14.204	1853	1858	1859	rBV2	9952	11130	0.67%	0.050%
27	14.251	1860	1866	1878	rVB2	673603	1066083	64.21%	4.791%
28	14.463	1899	1902	1906	rVV4	4156	5250	0.32%	0.024%
29	14.557	1910	1918	1926	rVV	657925	1039442	62.60%	4.671%
30	14.627	1927	1930	1934	rVB4	3850	5366	0.32%	0.024%
31	14.745	1941	1950	1960	rBV3	244486	468682	28.23%	2.106%
32	14.833	1960	1965	1970	rVV4	14831	29823	1.80%	0.134%
33	14.891	1971	1975	1985	rVV3	21064	33415	2.01%	0.150%
34	14.968	1985	1988	1993	rVV2	7995	11150	0.67%	0.050%
35	15.068	2001	2005	2008	rBV4	5806	8495	0.51%	0.038%
36	15.350	2049	2053	2059	rVB5	4113	6858	0.41%	0.031%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082124\
Data File : BG062514.D
Acq On : 21 Aug 2024 18:46
Operator : MA/JU
Sample : P3626-21
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCYC6

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-BG081424.MA.M
Title : SVOA CALIBRATION

37	15.414	2059	2064	2068	rBV3	5305	7227	0.44%	0.032%
38	15.549	2080	2087	2096	rVB	880476	1337152	80.53%	6.009%
39	15.655	2099	2105	2116	rBV	212598	322457	19.42%	1.449%
40	15.784	2121	2127	2133	rBV8	7113	16488	0.99%	0.074%
41	15.867	2137	2141	2147	rVB7	3452	7269	0.44%	0.033%
42	15.990	2158	2162	2166	rVV6	3900	6127	0.37%	0.028%
43	16.107	2178	2182	2187	rBV6	6079	13112	0.79%	0.059%
44	16.272	2206	2210	2213	rBV4	6413	8125	0.49%	0.037%
45	16.407	2229	2233	2235	rBV3	5217	7238	0.44%	0.033%
46	16.442	2237	2239	2241	rVV2	5110	4559	0.27%	0.020%
47	16.695	2279	2282	2287	rBV5	4919	7110	0.43%	0.032%
48	16.859	2306	2310	2312	rVV2	6291	7688	0.46%	0.035%
49	16.971	2323	2329	2331	rBV5	4636	6614	0.40%	0.030%
50	17.006	2331	2335	2340	rVB4	30539	43121	2.60%	0.194%
51	17.059	2340	2344	2353	rBV7	9729	21707	1.31%	0.098%
52	17.194	2363	2367	2375	rBV3	22707	40814	2.46%	0.183%
53	17.300	2375	2385	2395	rVB	831920	1266168	76.26%	5.690%
54	17.394	2395	2401	2411	rBV	886370	1368538	82.42%	6.150%
55	17.482	2411	2416	2420	rVB5	7144	12238	0.74%	0.055%
56	17.582	2430	2433	2437	rBV4	2860	5089	0.31%	0.023%
57	17.805	2467	2471	2475	rBV5	4213	7831	0.47%	0.035%
58	17.876	2475	2483	2492	rVV	227325	406269	24.47%	1.826%
59	17.970	2496	2499	2506	rVB3	18982	30167	1.82%	0.136%
60	18.069	2509	2516	2522	rBV8	13932	29017	1.75%	0.130%
61	18.134	2522	2527	2532	rBV6	16015	26266	1.58%	0.118%
62	18.193	2532	2537	2545	rBV	194124	268521	16.17%	1.207%
63	18.375	2565	2568	2572	rVB5	7104	8775	0.53%	0.039%
64	18.434	2574	2578	2584	rVB5	14671	21978	1.32%	0.099%
65	18.492	2584	2588	2590	rBV4	11604	14853	0.89%	0.067%
66	18.622	2607	2610	2613	rBV4	4821	6071	0.37%	0.027%
67	18.862	2646	2651	2653	rBV5	3765	5113	0.31%	0.023%
68	18.992	2668	2673	2674	rBV4	3377	4869	0.29%	0.022%
69	19.086	2680	2689	2691	rBV10	6503	18199	1.10%	0.082%
70	19.121	2692	2695	2700	rVV3	10472	13889	0.84%	0.062%
71	19.250	2713	2717	2723	rVB3	15230	23000	1.39%	0.103%
72	19.315	2723	2728	2731	rBV3	35482	50263	3.03%	0.226%
73	19.350	2731	2734	2737	rVB4	8018	10763	0.65%	0.048%
74	19.462	2749	2753	2761	rBV2	46701	75894	4.57%	0.341%
75	19.532	2761	2765	2771	rVB	28183	38528	2.32%	0.173%
76	19.614	2774	2779	2784	rVB7	9219	16732	1.01%	0.075%

LSC Area Percent Report

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Operator : MA/JU
Sample : P3626-21
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCYC6

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

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Title : SVOA CALIBRATION

77	19.679	2784	2790	2796	rBV	1148773	1660362	100.00%	7.461%
78	20.061	2853	2855	2859	rBV5	5789	7518	0.45%	0.034%
79	20.125	2863	2866	2872	rVB2	22580	27384	1.65%	0.123%
80	20.196	2875	2878	2881	rVV4	11630	16080	0.97%	0.072%
81	20.225	2881	2883	2887	rVB3	20340	23234	1.40%	0.104%
82	20.390	2908	2911	2917	rVB6	11087	15509	0.93%	0.070%
83	20.554	2935	2939	2943	rBV	72681	88759	5.35%	0.399%
84	20.601	2943	2947	2952	rVV7	15503	32306	1.95%	0.145%
85	20.654	2953	2956	2960	rVB2	34967	43528	2.62%	0.196%
86	21.024	3016	3019	3024	rBV5	9864	12845	0.77%	0.058%
87	21.206	3045	3050	3061	rBV2	171626	300953	18.13%	1.352%
88	21.535	3097	3106	3118	rVB	819864	1328702	80.02%	5.971%
89	21.741	3138	3141	3148	rBV5	19416	30621	1.84%	0.138%
90	22.346	3236	3244	3258	rBV3	87314	287649	17.32%	1.293%
91	22.992	3350	3354	3360	rVB2	30759	46102	2.78%	0.207%
92	23.174	3381	3385	3391	rVB5	15271	25389	1.53%	0.114%
93	23.386	3417	3421	3425	rBV5	11415	18296	1.10%	0.082%
94	23.762	3478	3485	3491	rBV	153866	307434	18.52%	1.382%
95	24.402	3583	3594	3606	rBV	615536	1596857	96.18%	7.176%
96	24.613	3620	3630	3646	rVB	518080	1460384	87.96%	6.563%
97	25.154	3716	3722	3729	rVB6	38154	91992	5.54%	0.413%
98	25.741	3814	3822	3831	rVB2	85903	239609	14.43%	1.077%
99	25.865	3838	3843	3850	rVB5	21168	53624	3.23%	0.241%
100	28.585	4295	4306	4312	rBV7	32360	126530	7.62%	0.569%

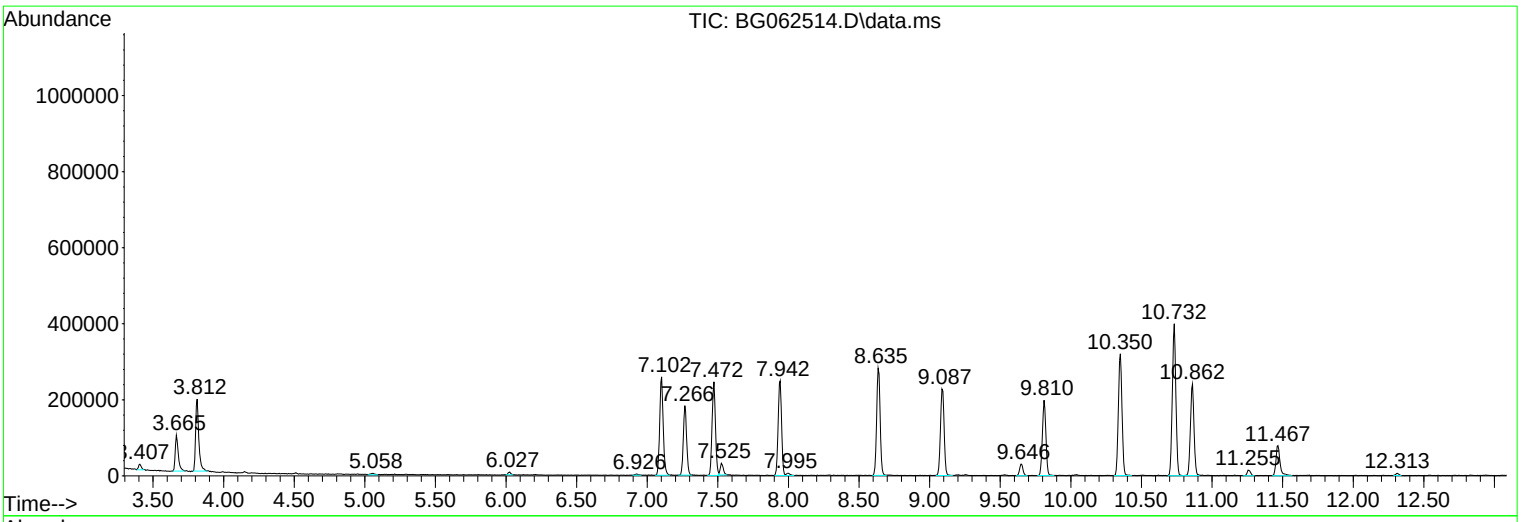
Sum of corrected areas: 22252843

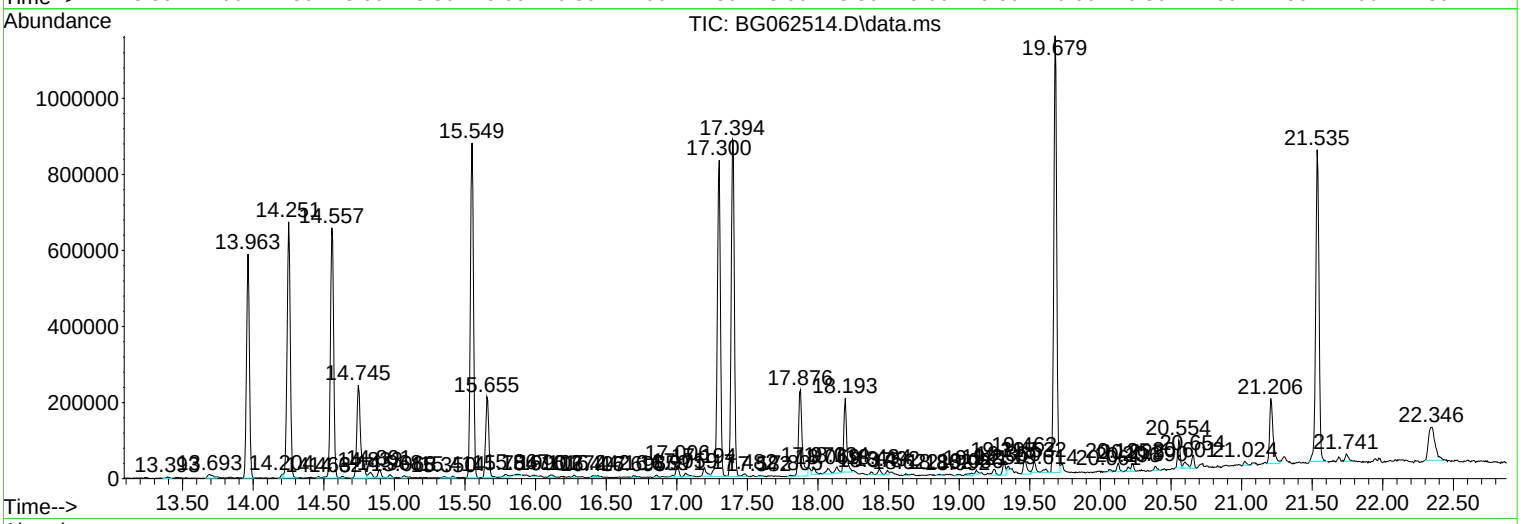
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082124\
Data File : BG062514.D
Acq On : 21 Aug 2024 18:46
Operator : MA/JU
Sample : P3626-21
Misc :
ALS Vial : 15 Sample Multiplier: 1

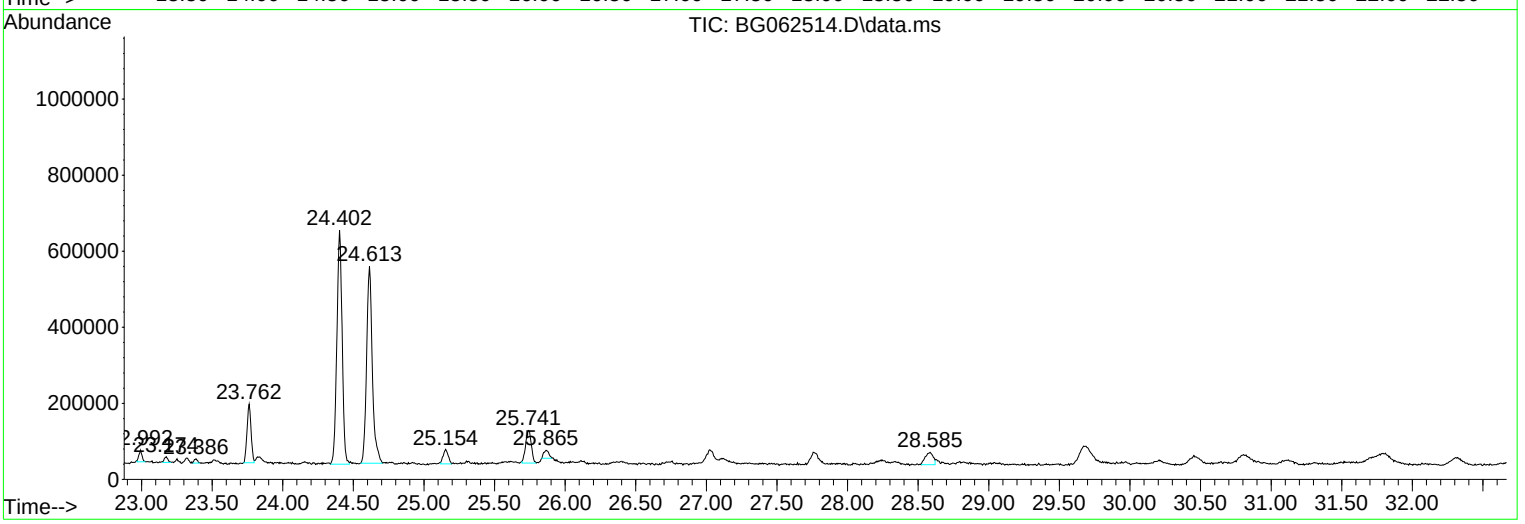
Instrument :
BNA_G
ClientSampleId :
DCYC6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

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







Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082124\
Data File : BG062514.D
Acq On : 21 Aug 2024 18:46
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Sample : P3626-21
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ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCYC6

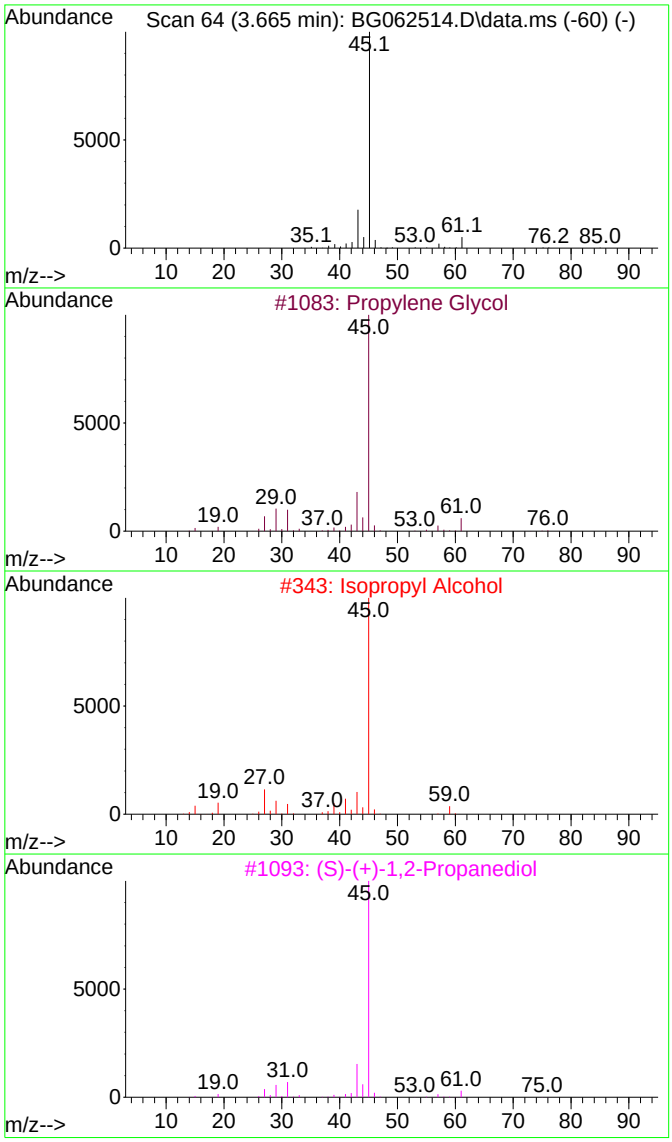
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

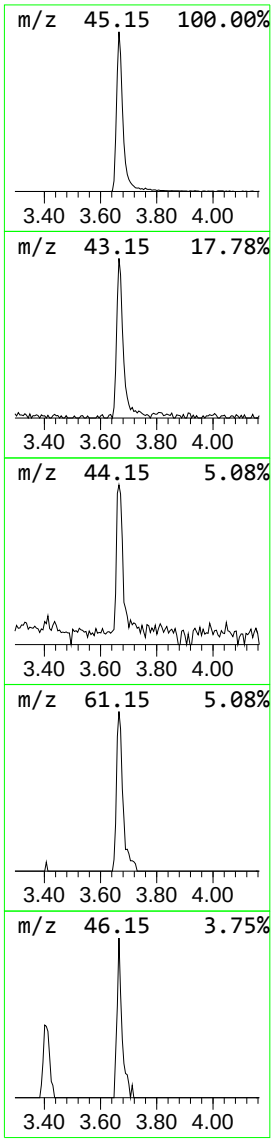
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Propylene Glycol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.665	6.72 ng/ul	140940	1,4-Dichlorobenzene-d4	7.942

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propylene Glycol	76	C3H8O2	000057-55-6	86
2	Isopropyl Alcohol	60	C3H8O	000067-63-0	56
3	(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	56
4	(R)-(-)-3-Methyl-2-butanol	88	C5H12O	001572-93-6	9
5	R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082124\
Data File : BG062514.D
Acq On : 21 Aug 2024 18:46
Operator : MA/JU
Sample : P3626-21
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCYC6

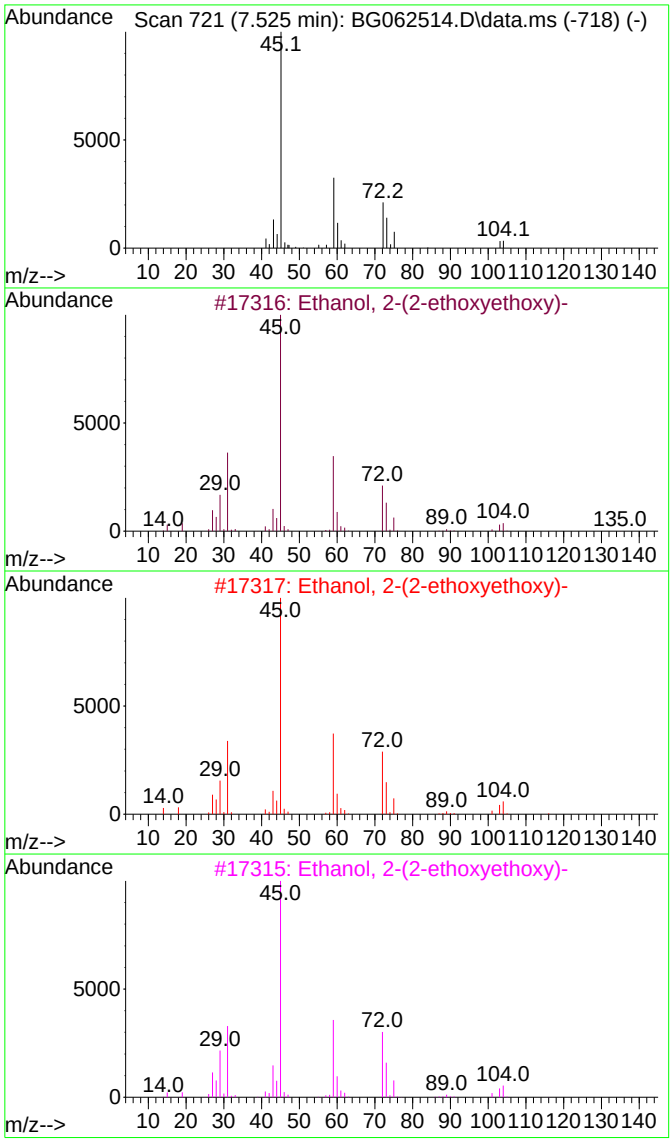
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

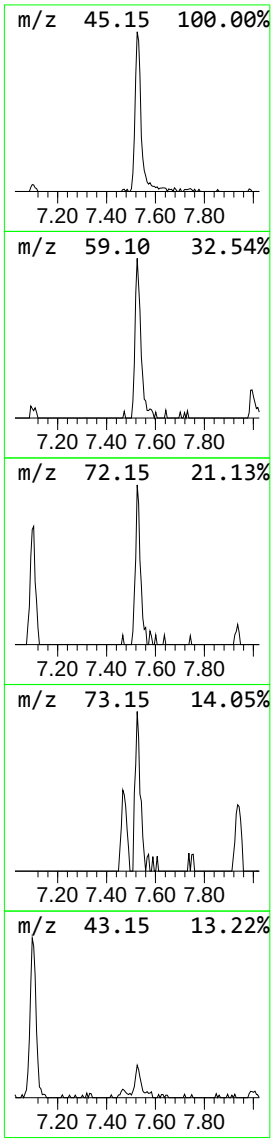
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.525	2.15 ng/ul	45127	1,4-Dichlorobenzene-d4	7.942

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	83
2	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	78
3	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	78
4	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	50
5	Ethanol, 2-[2-(2-ethoxyethoxy)et...	178	C8H18O4	000112-50-5	50





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Data File : BG062514.D
Acq On : 21 Aug 2024 18:46
Operator : MA/JU
Sample : P3626-21
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCYC6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

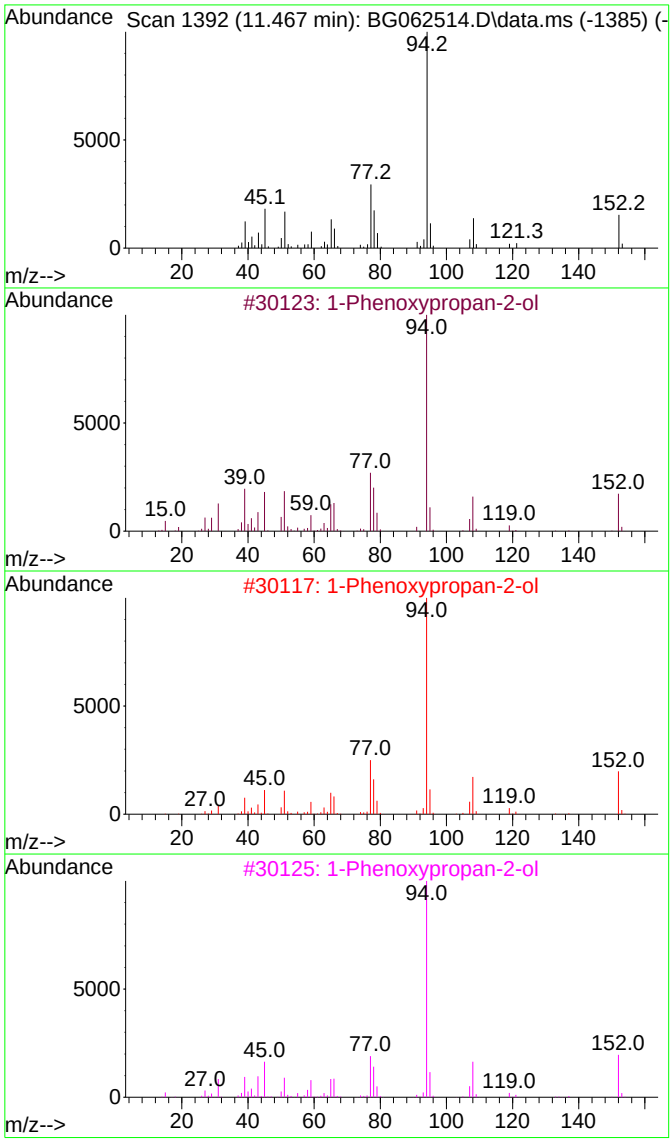
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

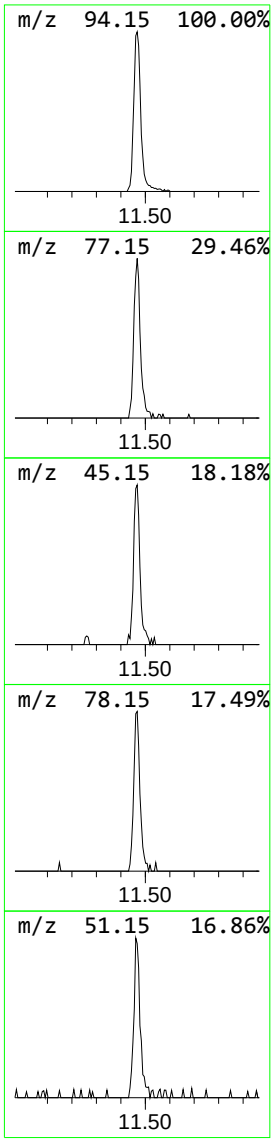
****

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.467	4.55 ng/ul	157480	Naphthalene-d8	10.732

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
2	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
3	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	93
4	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
5	1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082124\
Data File : BG062514.D
Acq On : 21 Aug 2024 18:46
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Sample : P3626-21
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCYC6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

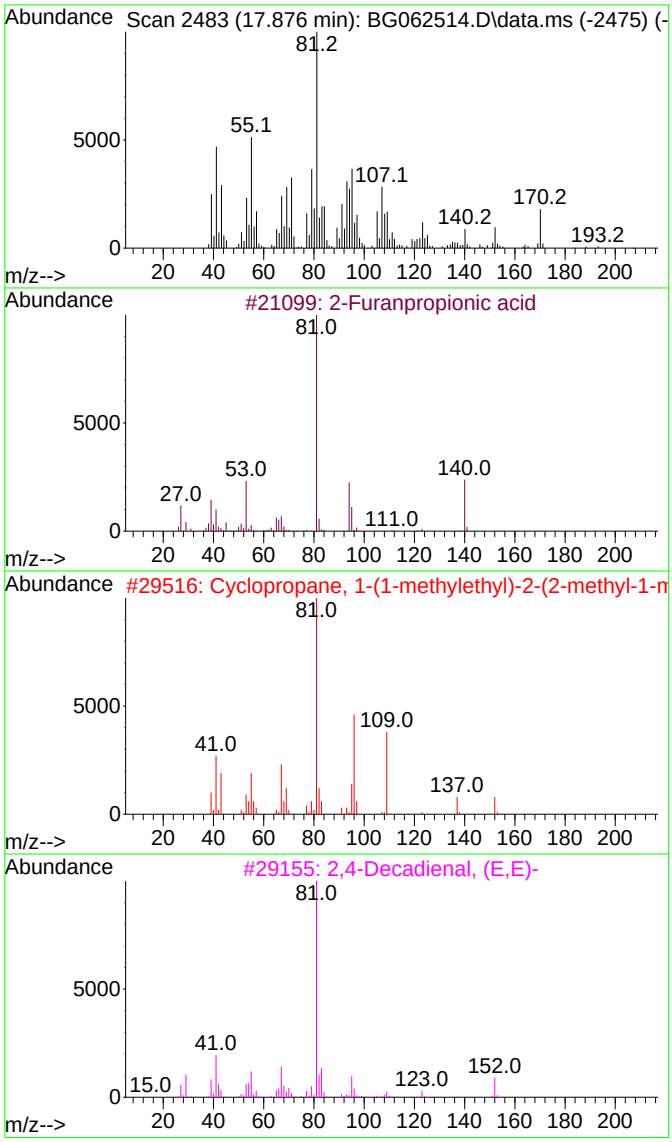
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

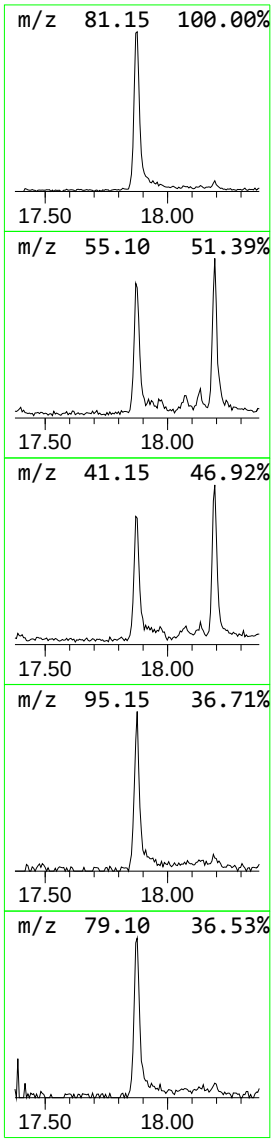
****

Peak Number 4 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.876	6.42 ng/ul	406269	Phenanthrene-d10	17.300

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Furanpropionic acid	140	C7H8O3	000935-13-7	38
2	Cyclopropane, 1-(1-methylethyl)-...	152	C11H20	056259-17-7	38
3	2,4-Decadienal, (E,E)-	152	C10H16O	025152-84-5	38
4	L-Fenchone	152	C10H16O	007787-20-4	35
5	Fenchone	152	C10H16O	001195-79-5	35





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082124\
Data File : BG062514.D
Acq On : 21 Aug 2024 18:46
Operator : MA/JU
Sample : P3626-21
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCYC6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

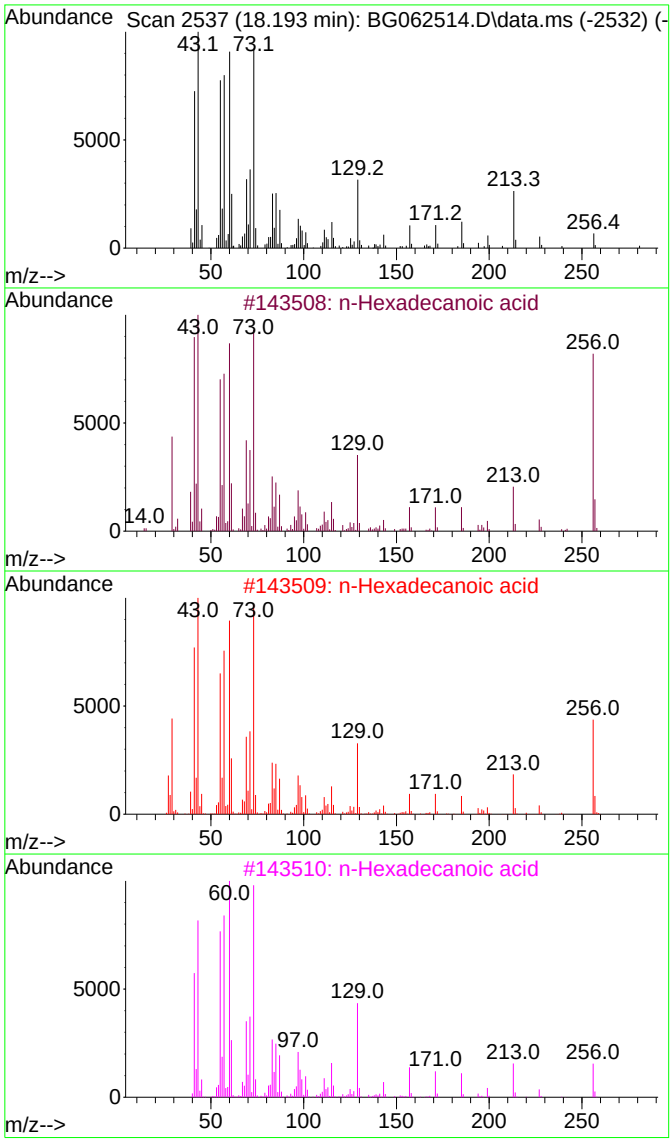
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

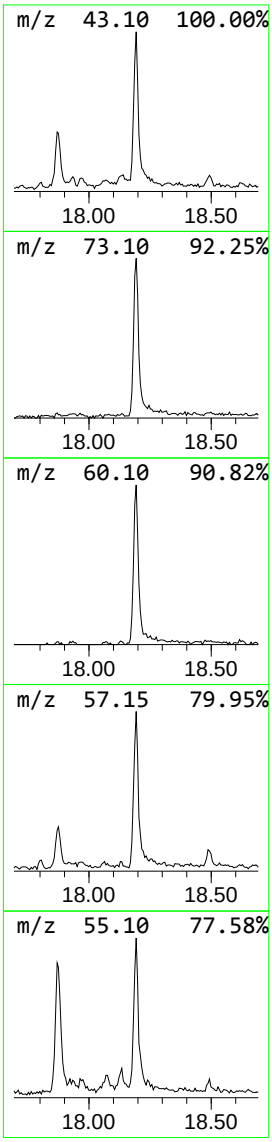
****

Peak Number 5 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.193	4.24 ng/ul	268521	Phenanthrene-d10	17.300

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	81
5	Tridecanoic acid	214	C13H26O2	000638-53-9	81





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082124\
Data File : BG062514.D
Acq On : 21 Aug 2024 18:46
Operator : MA/JU
Sample : P3626-21
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCYC6

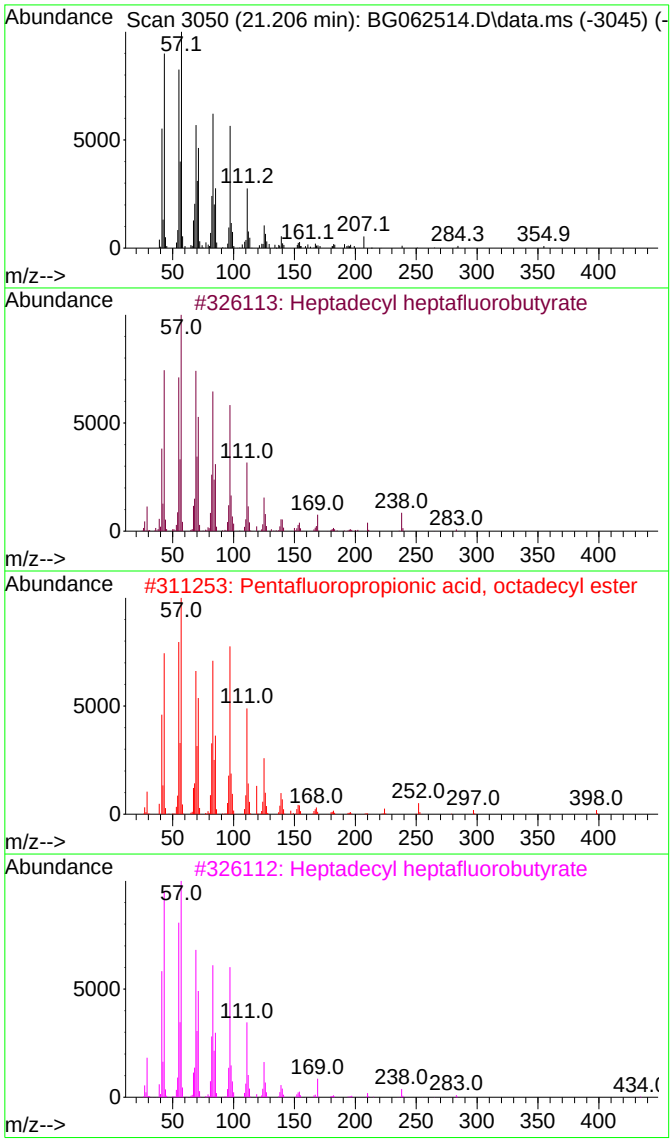
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

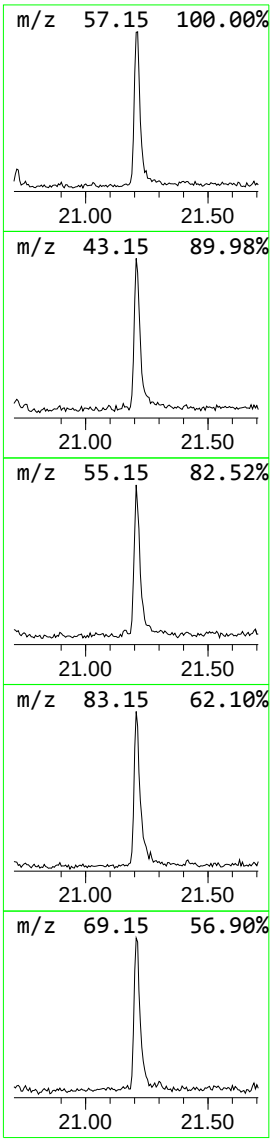
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Heptadecyl heptafluorobutyrate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.206	4.53 ng/ul	300953	Chrysene-d12	21.535

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	93
2	Pentafluoropropionic acid, octad...	416	C21H37F5O2	959261-25-7	91
3	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91
4	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	91
5	Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	91





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Data File : BG062514.D
Acq On : 21 Aug 2024 18:46
Operator : MA/JU
Sample : P3626-21
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCYC6

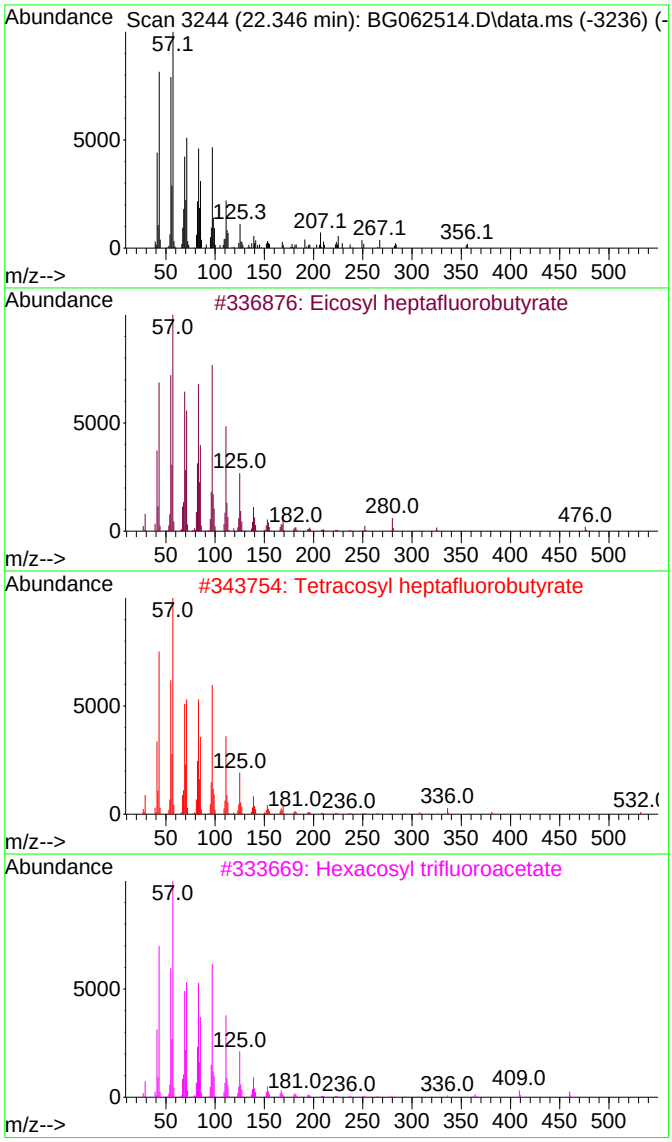
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

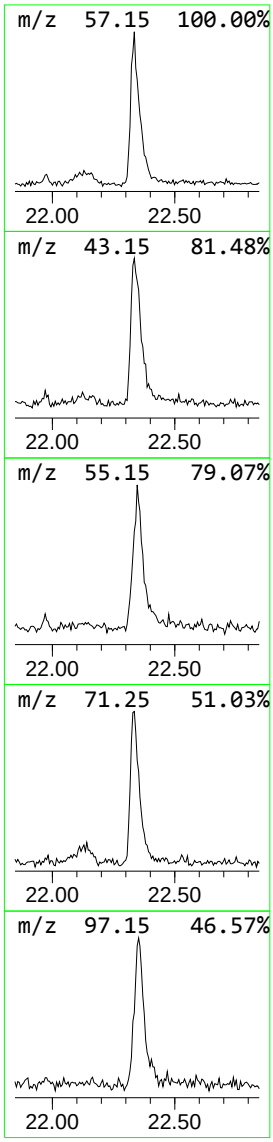
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Eicosyl heptafluorobutyrate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.346	4.33 ng/ul	287649	Chrysene-d12	21.535

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosyl heptafluorobutyrate	494	C24H41F7O2	1000351-83-5	91
2	Tetracosyl heptafluorobutyrate	550	C28H49F7O2	1000351-83-7	91
3	Hexacosyl trifluoroacetate	478	C28H53F3O2	1000351-75-0	91
4	Hexacosyl heptafluorobutyrate	578	C30H53F7O2	1000351-83-3	91
5	Eicosyl trifluoroacetate	394	C22H41F3O2	1000351-75-2	91





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Data File : BG062514.D
Acq On : 21 Aug 2024 18:46
Operator : MA/JU
Sample : P3626-21
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCYC6

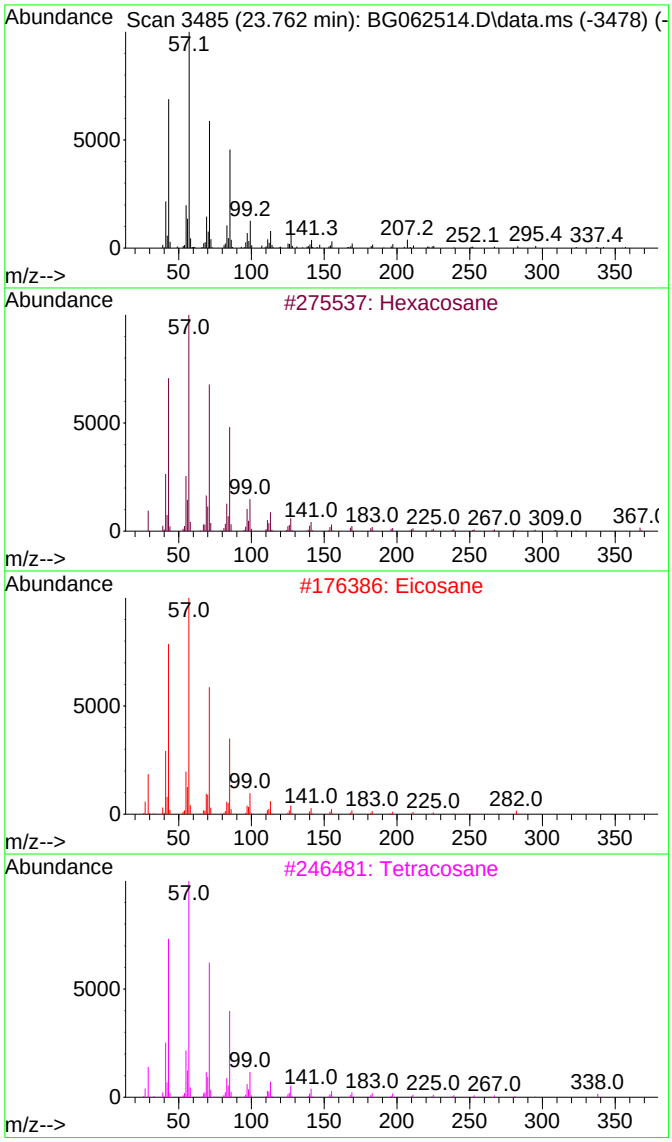
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

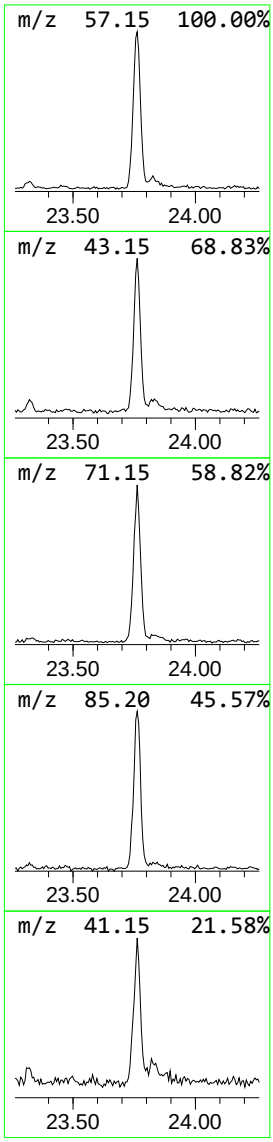
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.762	4.21 ng/ul	307434	Perylene-d12	24.613

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	93
2	Eicosane	282	C20H42	000112-95-8	90
3	Tetracosane	338	C24H50	000646-31-1	90
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
5	Pentacosane	352	C25H52	000629-99-2	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082124\
Data File : BG062514.D
Acq On : 21 Aug 2024 18:46
Operator : MA/JU
Sample : P3626-21
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCYC6

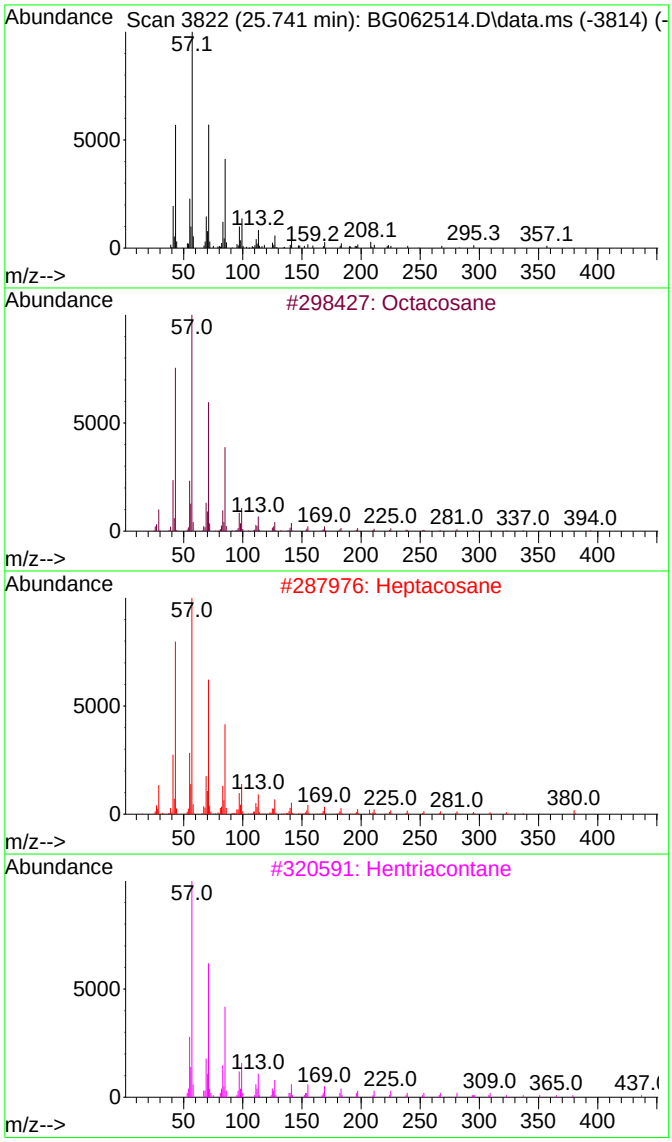
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

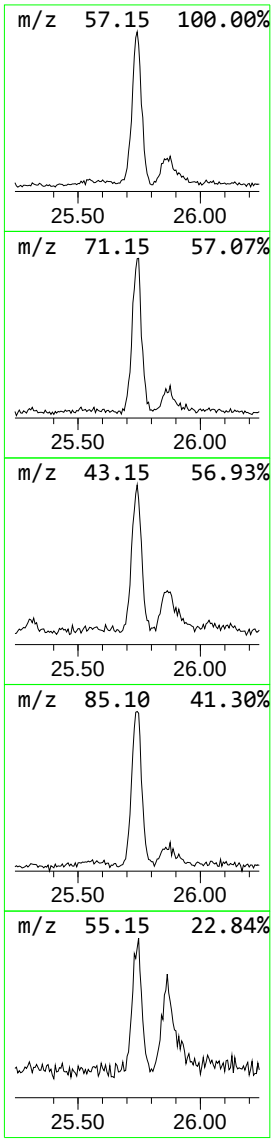
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.741	3.28 ng/ul	239609	Perylene-d12	24.613

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	90
2	Heptacosane	380	C27H56	000593-49-7	90
3	Hentriacontane	437	C31H64	000630-04-6	90
4	Tetracosane	338	C24H50	000646-31-1	87
5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082124\
Data File : BG062514.D
Acq On : 21 Aug 2024 18:46
Operator : MA/JU
Sample : P3626-21
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DCYC6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.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
Propylene Glycol	3.665	6.7	ng/ul	140940	1	7.942	419460	20.0
Ethanol, 2-(2-e...	7.525	2.1	ng/ul	45127	1	7.942	419460	20.0
1-Phenoxypropan...	11.467	4.5	ng/ul	157480	2	10.732	692085	20.0
unknown-01	17.876	6.4	ng/ul	406269	4	17.300	1266170	20.0
n-Hexadecanoic ...	18.193	4.2	ng/ul	268521	4	17.300	1266170	20.0
Heptadecyl hept...	21.206	4.5	ng/ul	300953	5	21.535	1328700	20.0
Eicosyl heptafl...	22.346	4.3	ng/ul	287649	5	21.535	1328700	20.0
(DEL) Alkane: S...	23.762	4.2	ng/ul	307434	6	24.613	1460380	20.0
(DEL) Alkane: S...	25.741	3.3	ng/ul	239609	6	24.613	1460380	20.0