LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031623\
Data File : BP014109.D
Acq On : 16 Mar 2023 18:47
Operator : CG/JU
Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB925

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_P\Methods\SFAM-EPA-BP030723.M
Title : SVOA CALIBRATION

Signal : TIC: BP014109.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.234	4	8	22	rVV	52244	96382	2.00%	0.123%
2	3.423	36	40	43	rBV	34197	43363	0.90%	0.056%
3	3.458	43	46	58	rVB	83195	121827	2.52%	0.156%
4	3.864	112	115	131	rBV	247450	449756	9.31%	0.576%
5	4.646	240	248	261	rVB	337453	527802	10.93%	0.676%
6	5.340	358	366	374	rBV	172620	273357	5.66%	0.350%
7	5.787	438	442	447	rBV2	26633	43935	0.91%	0.056%
8	6.528	563	568	576	rBV2	17057	34252	0.71%	0.044%
9	7.222	673	686	699	rBV	128864	276277	5.72%	0.354%
10	7.387	707	714	732	rBV	468714	777383	16.10%	0.996%
11	7.593	742	749	756	rVV	435797	739502	15.31%	0.947%
12	7.658	756	760	765	rVV6	18244	36043	0.75%	0.046%
13	7.916	796	804	808	rBV5	24500	49503	1.02%	0.063%
14	7.958	808	811	817	rVB	21223	31704	0.66%	0.041%
15	8.064	817	829	841	rBV	682342	1151018	23.83%	1.475%
16	8.399	884	886	895	rVV9	11635	32741	0.68%	0.042%
17	8.546	903	911	916	rVV7	16749	41964	0.87%	0.054%
18	8.775	943	950	960	rVV	400182	670667	13.89%	0.859%
19	8.975	980	984	991	rVB3	15911	41324	0.86%	0.053%
20	9.058	991	998	1007	rBV	540808	892075	18.47%	1.143%
21	9.175	1013	1018	1022	rBV	67126	107441	2.22%	0.138%
22	9.240	1022	1029	1038	rVB	574342	1001024	20.73%	1.282%
23	9.440	1060	1063	1070	rVB7	26395	44661	0.92%	0.057%
24	9.593	1082	1089	1093	rBV2	142543	254120	5.26%	0.326%
25	9.693	1099	1106	1112	rVB9	19915	48544	1.01%	0.062%
26	9.822	1117	1128	1132	rBV5	37406	72122	1.49%	0.092%
27	9.963	1146	1152	1165	rVB	462381	788204	16.32%	1.010%
28	10.099	1171	1175	1180	rVV6	20672	37540	0.78%	0.048%
29	10.375	1216	1222	1224	rBV3	29260	59375	1.23%	0.076%
30	10.405	1224	1227	1237	rVB5	43998	103721	2.15%	0.133%
31	10.505	1238	1244	1254	rBV	685735	1307128	27.06%	1.675%
32	10.681	1265	1274	1281	rBV9	25838	79968	1.66%	0.102%
33	10.816	1290	1297	1302	rBV2	57437	106258	2.20%	0.136%
34	10.887	1302	1309	1315	rVV	949932	1570437	32.51%	2.012%
35	11.028	1324	1333	1347	rVV	690734	1309590	27.11%	1.678%
36	11.140	1347	1352	1356	rVB3	21092	38425	0.80%	0.049%

LSC Area Percent Report

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Data File : BP014109.D
Acq On : 16 Mar 2023 18:47
Operator : CG/JU
Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB925

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_P\Methods\SFAM-EPA-BP030723.M
Title : SVOA CALIBRATION

37	11.246	1365	1370	1376	rVB9	14147	33094	0.69%	0.042%
38	11.469	1405	1408	1415	rVB9	20521	40334	0.84%	0.052%
39	11.693	1440	1446	1448	rBV6	26937	55861	1.16%	0.072%
40	12.087	1508	1513	1518	rBV6	22879	41486	0.86%	0.053%
41	12.134	1518	1521	1526	rVB3	18293	30271	0.63%	0.039%
42	12.216	1526	1535	1540	rBV	226664	396401	8.21%	0.508%
43	12.263	1540	1543	1552	rVB6	23892	48410	1.00%	0.062%
44	12.387	1558	1564	1572	rBV3	121933	244854	5.07%	0.314%
45	12.487	1577	1581	1590	rVB5	73415	165804	3.43%	0.212%
46	12.646	1604	1608	1614	rBV6	20005	36993	0.77%	0.047%
47	12.869	1642	1646	1651	rBV4	31159	53382	1.11%	0.068%
48	12.963	1658	1662	1673	rVB8	34874	81348	1.68%	0.104%
49	13.510	1750	1755	1758	rBV3	80374	131057	2.71%	0.168%
50	13.757	1793	1797	1805	rVB7	37174	63959	1.32%	0.082%
51	13.928	1823	1826	1831	rBV6	19475	31163	0.65%	0.040%
52	14.110	1850	1857	1865	rVB	1565354	2163830	44.80%	2.772%
53	14.263	1879	1883	1885	rBV3	40949	65234	1.35%	0.084%
54	14.398	1900	1906	1913	rBV	1543170	2335369	48.35%	2.992%
55	14.510	1921	1925	1932	rBV2	41394	96053	1.99%	0.123%
56	14.628	1940	1945	1949	rBV	163807	237246	4.91%	0.304%
57	14.704	1952	1958	1964	rVV	1504519	2286068	47.33%	2.929%
58	14.769	1965	1969	1976	rVB	137604	213100	4.41%	0.273%
59	14.893	1987	1990	1993	rBV5	42634	64733	1.34%	0.083%
60	14.934	1993	1997	2006	rVB2	77798	172290	3.57%	0.221%
61	15.110	2025	2027	2032	rVB4	43144	54911	1.14%	0.070%
62	15.440	2078	2083	2091	rBV	1733071	2434761	50.41%	3.119%
63	15.698	2121	2127	2133	rBV	2138823	3010172	62.32%	3.857%
64	15.816	2142	2147	2153	rBV	641005	864074	17.89%	1.107%
65	15.910	2160	2163	2167	rBV	119512	138722	2.87%	0.178%
66	15.957	2167	2171	2178	rVB2	133669	213355	4.42%	0.273%
67	16.063	2184	2189	2201	rVB	582040	847785	17.55%	1.086%
68	16.222	2209	2216	2225	rBV	3270379	4829910	100.00%	6.188%
69	16.363	2237	2240	2243	rBV2	55109	77311	1.60%	0.099%
70	16.404	2244	2247	2251	rVB2	74674	100458	2.08%	0.129%
71	16.545	2267	2271	2275	rBV	189801	241520	5.00%	0.309%
72	16.592	2275	2279	2282	rVB2	59556	76873	1.59%	0.098%
73	16.734	2296	2303	2307	rBV	2173110	3141354	65.04%	4.025%
74	16.775	2307	2310	2314	rVB	926475	1098473	22.74%	1.407%
75	16.851	2319	2323	2330	rVB6	70813	126665	2.62%	0.162%
76	16.992	2344	2347	2354	rVB	131591	176129	3.65%	0.226%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031623\
Data File : BP014109.D
Acq On : 16 Mar 2023 18:47
Operator : CG/JU
Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB925

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_P\Methods\SFAM-EPA-BP030723.M
Title : SVOA CALIBRATION

77	17.463	2422	2427	2435	rVB	1996506	2819636	58.38%	3.612%
78	17.563	2438	2444	2451	rBV	2327061	3260563	67.51%	4.177%
79	18.028	2519	2523	2527	rBV	109502	139319	2.88%	0.178%
80	18.204	2550	2553	2558	rBV2	108056	195665	4.05%	0.251%
81	18.328	2570	2574	2583	rBV	1376578	2104440	43.57%	2.696%
82	18.728	2638	2642	2647	rVB	1256525	1452994	30.08%	1.862%
83	19.086	2701	2703	2707	rVB2	172988	184863	3.83%	0.237%
84	19.439	2756	2763	2768	rBV4	177669	440728	9.12%	0.565%
85	19.486	2768	2771	2778	rVV	232281	304770	6.31%	0.390%
86	19.551	2778	2782	2798	rVV2	401026	742399	15.37%	0.951%
87	19.786	2815	2822	2826	rBV	3199041	4571297	94.65%	5.857%
88	19.828	2826	2829	2838	rVB	1732530	2196880	45.48%	2.815%
89	20.281	2903	2906	2911	rVB	376987	407867	8.44%	0.523%
90	20.645	2965	2968	2978	rBV5	155328	302224	6.26%	0.387%
91	20.763	2984	2988	2992	rBV	682238	792034	16.40%	1.015%
92	21.222	3062	3066	3077	rBV	1684230	2269025	46.98%	2.907%
93	21.545	3116	3121	3127	rBV	2345654	3172608	65.69%	4.065%
94	21.669	3138	3142	3149	rVB	663729	754236	15.62%	0.966%
95	22.151	3220	3224	3234	rVB	549616	819095	16.96%	1.049%
96	22.675	3309	3313	3326	rVB3	538863	1238945	25.65%	1.587%
97	22.827	3334	3339	3347	rVB	1911048	2780585	57.57%	3.562%
98	23.263	3409	3413	3418	rVB	320035	481967	9.98%	0.617%
99	23.916	3517	3524	3535	rVB2	1848931	3927495	81.32%	5.032%
100	24.080	3546	3552	3565	rVB2	1386815	2913678	60.33%	3.733%

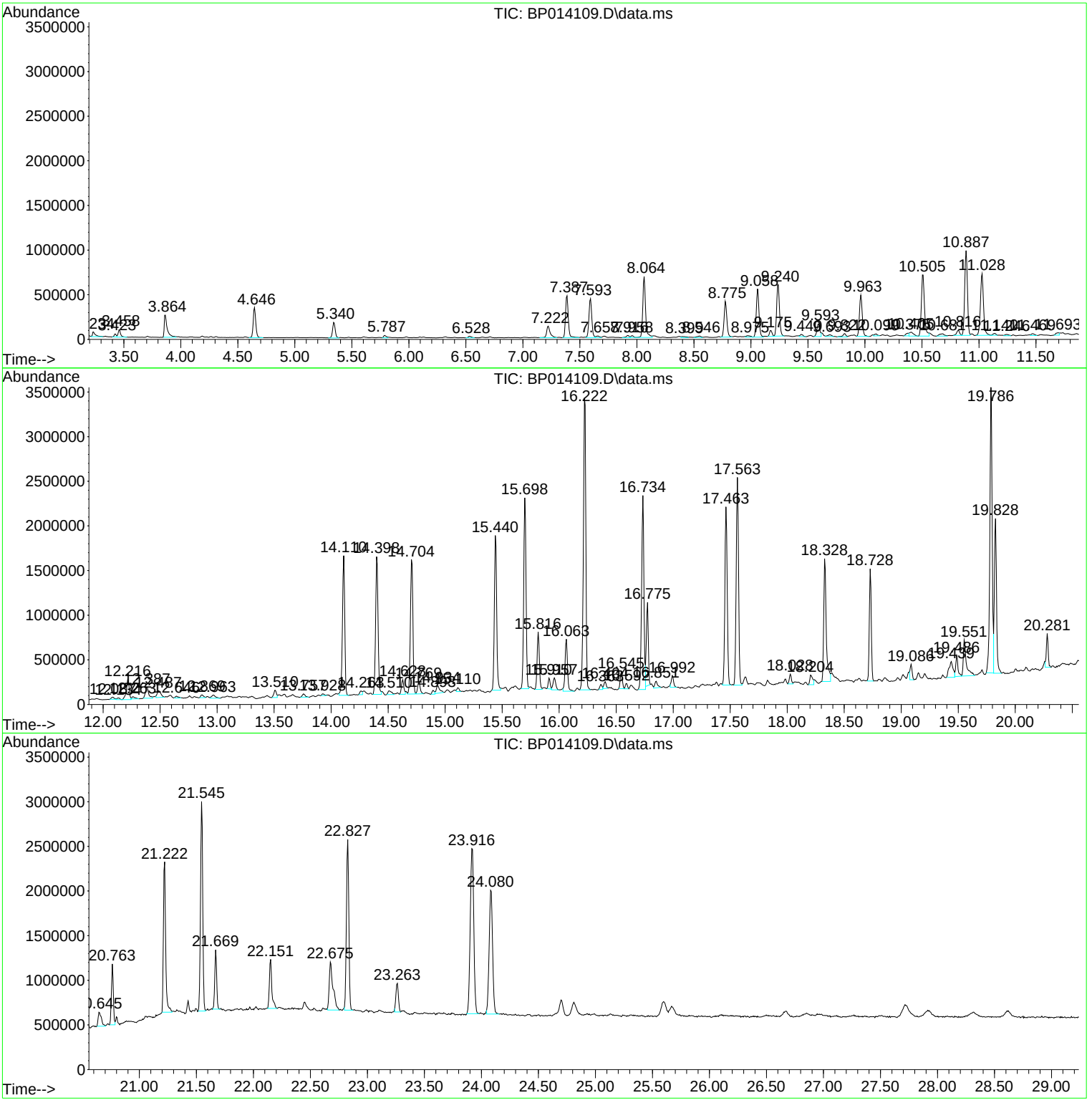
Sum of corrected areas: 78053559

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031623\
Data File : BP014109.D
Acq On : 16 Mar 2023 18:47
Operator : CG/JU
Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB925

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031623\
Data File : BP014109.D
Acq On : 16 Mar 2023 18:47
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Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB925

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

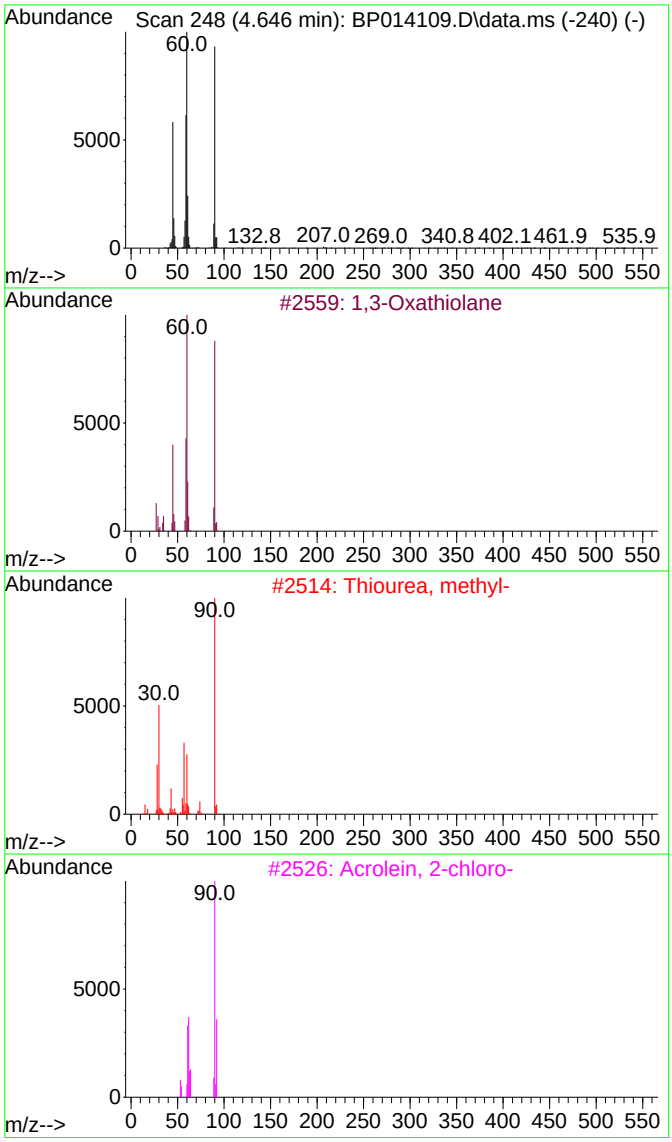
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

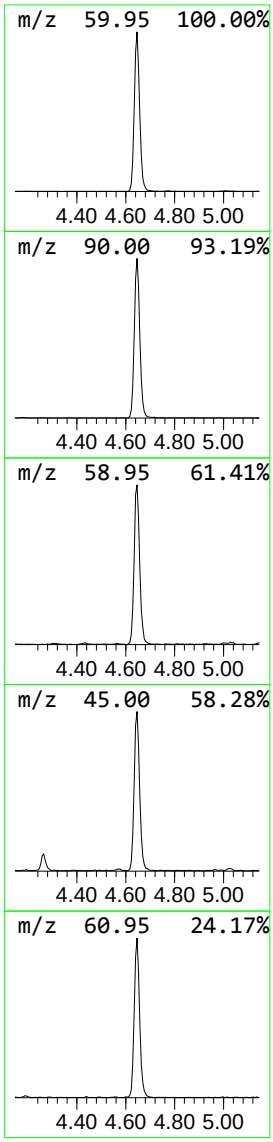
****

Peak Number 1 1,3-Oxathiolane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.646	9.17 ng/ul	527802	1,4-Dichlorobenzene-d4	8.064

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Oxathiolane	90	C3H6OS	002094-97-5	94
2	Thiourea, methyl-	90	C2H6N2S	000598-52-7	50
3	Acrolein, 2-chloro-	90	C3H3ClO	000683-51-2	42
4	Thiourea, methyl-	90	C2H6N2S	000598-52-7	36
5	Heptanoic acid, 2-hydroxy-, meth...	160	C8H16O3	054340-91-9	36





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031623\
Data File : BP014109.D
Acq On : 16 Mar 2023 18:47
Operator : CG/JU
Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB925

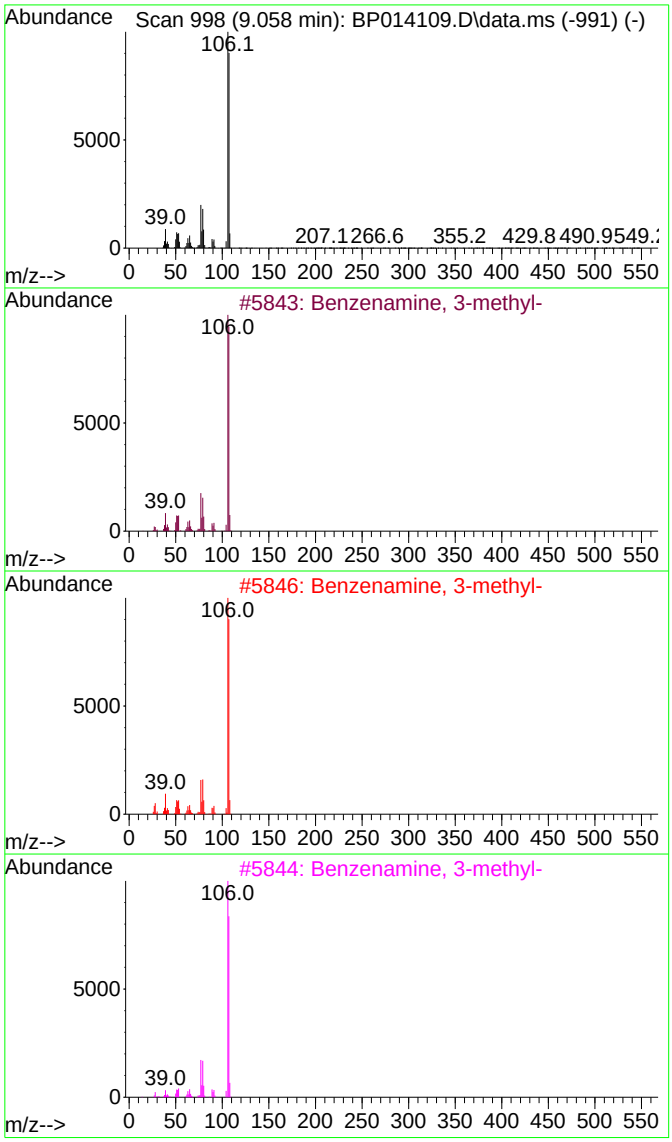
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

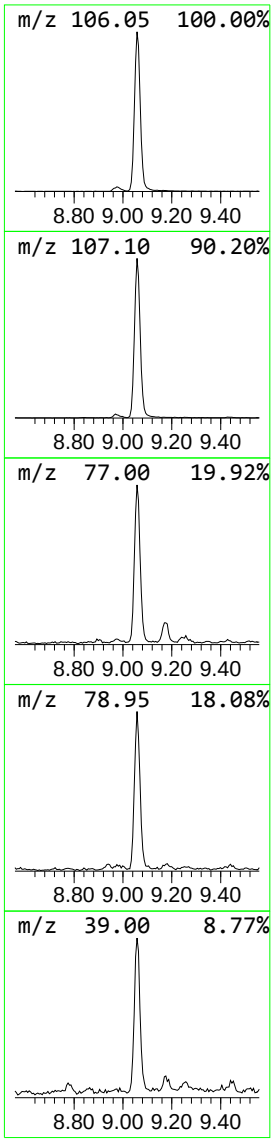
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzenamine, 3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.058	15.50 ng/ul	892075	1,4-Dichlorobenzene-d4	8.064

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
2	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
3	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
4	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	96
5	o-Toluidine	107	C7H9N	000095-53-4	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031623\
Data File : BP014109.D
Acq On : 16 Mar 2023 18:47
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Sample : 01884-13
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ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB925

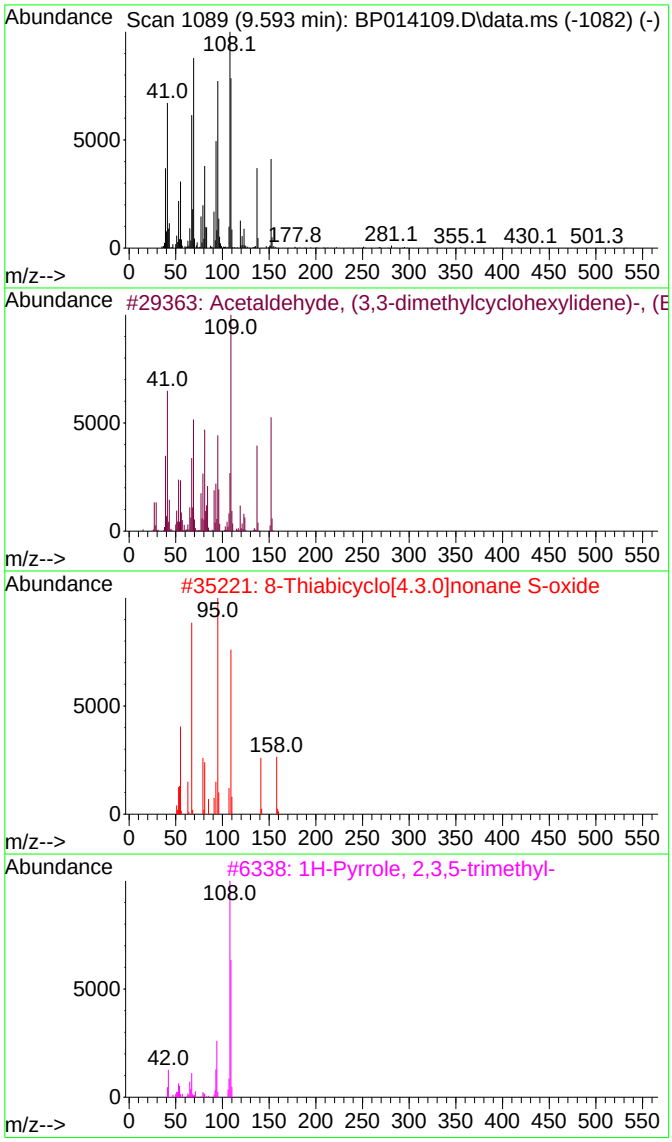
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

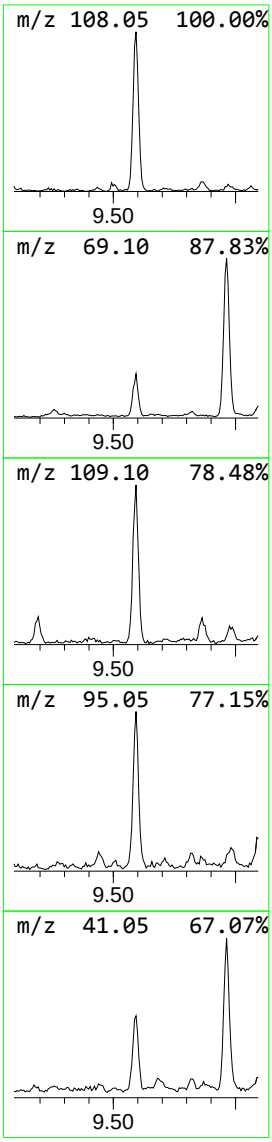
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Acetaldehyde, (3,3-dimethyl... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.593	3.24 ng/ul	254120	Naphthalene-d8	10.887

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetaldehyde, (3,3-dimethylcyclo...	152	C10H16O	026532-25-2	80
2	8-Thiabicyclo[4.3.0]nonane S-oxide	158	C8H14OS	1010215-10-0	35
3	1H-Pyrrole, 2,3,5-trimethyl-	109	C7H11N	002199-41-9	30
4	.alpha.-Campholenal	152	C10H16O	004501-58-0	27
5	1,2,5-Trimethylpyrrole	109	C7H11N	000930-87-0	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031623\
Data File : BP014109.D
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Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB925

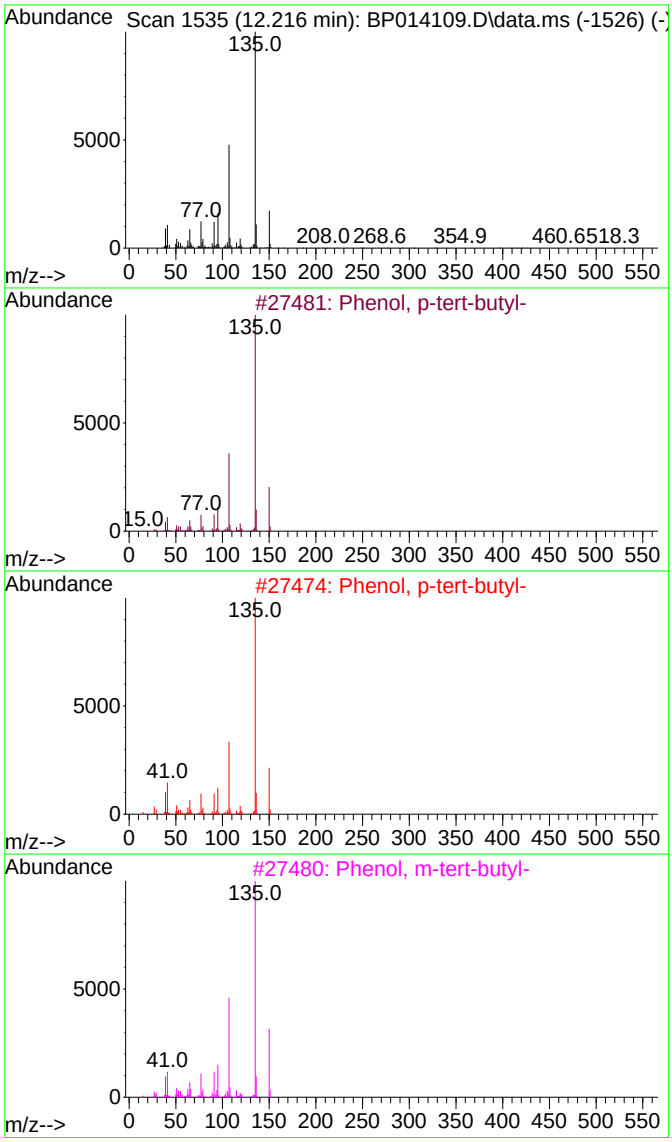
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

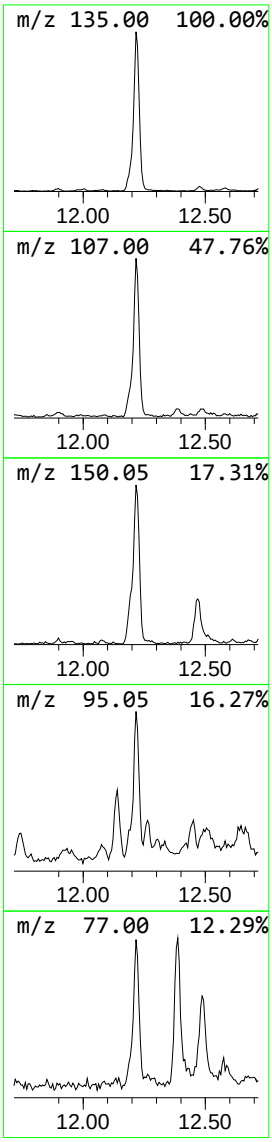
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Phenol, p-tert-butyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.216	5.05 ng/ul	396401	Naphthalene-d8	10.887

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	95
2	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	91
3	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91
4	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	90
5	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031623\
Data File : BP014109.D
Acq On : 16 Mar 2023 18:47
Operator : CG/JU
Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB925

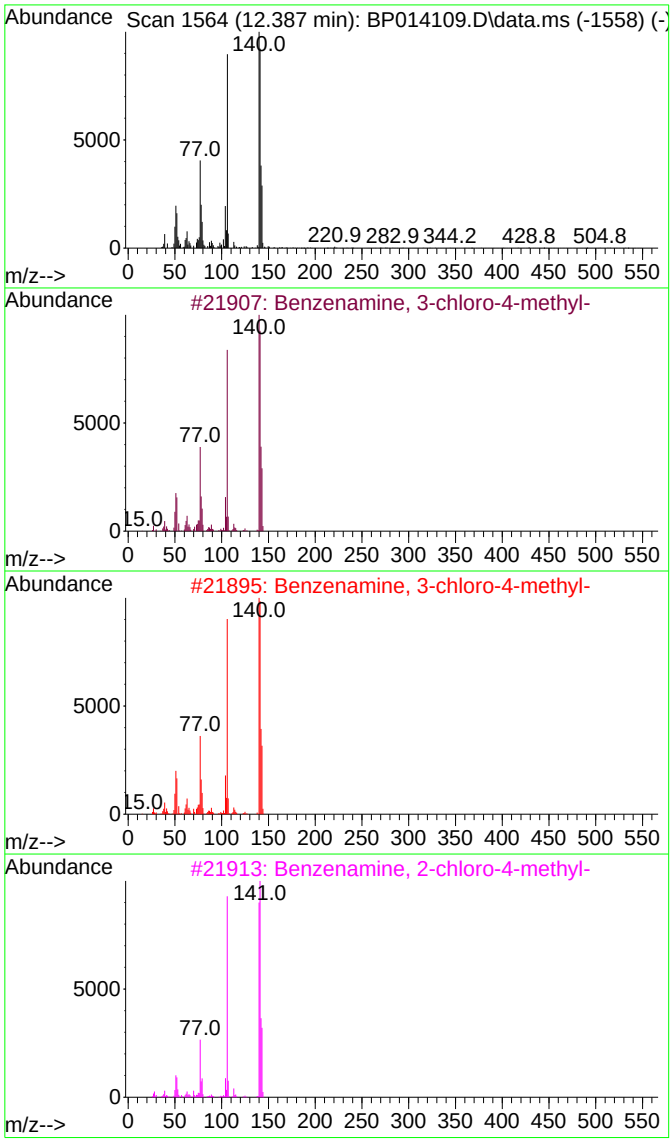
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

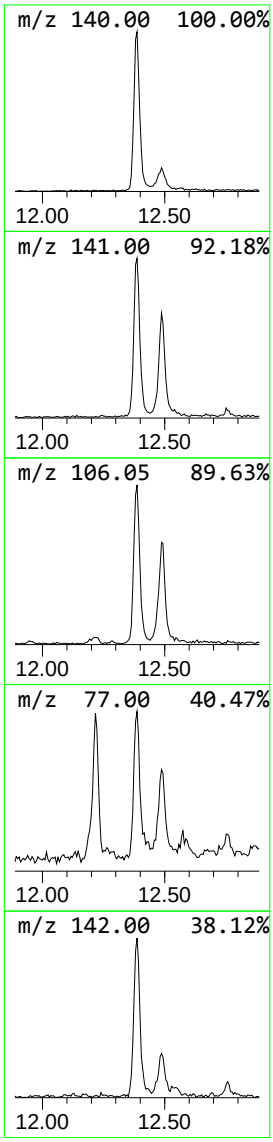
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzenamine, 3-chloro-4-met... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.387	3.12 ng/ul	244854	Naphthalene-d8	10.887

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	96
2	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	94
3	Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	93
4	Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	76
5	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	76





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Acq On : 16 Mar 2023 18:47
Operator : CG/JU
Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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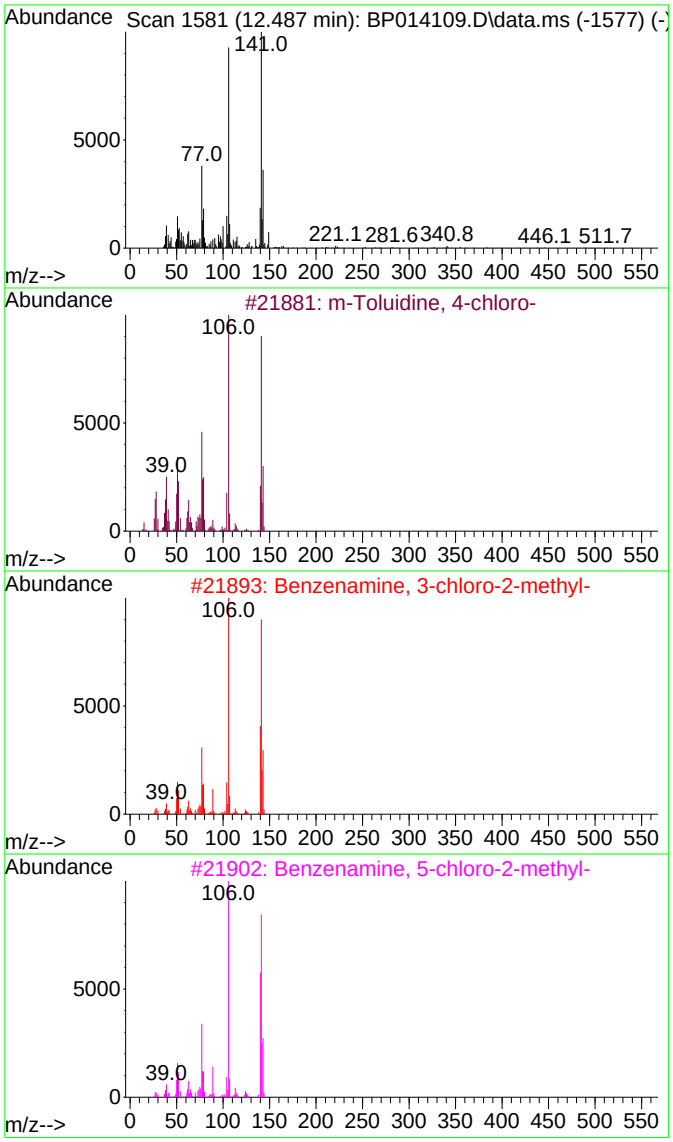
Quant Title : SVOA CALIBRATION

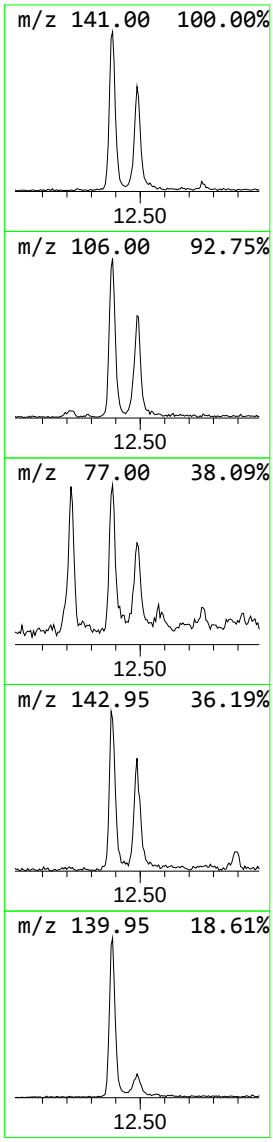
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 m-Toluidine, 4-chloro- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.487	2.11 ng/ul	165804	Naphthalene-d8	10.887

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	m-Toluidine, 4-chloro-	141	C7H8ClN	007149-75-9	93
2	Benzenamine, 3-chloro-2-methyl-	141	C7H8ClN	000087-60-5	87
3	Benzenamine, 5-chloro-2-methyl-	141	C7H8ClN	000095-79-4	87
4	Benzenamine, 3-chloro-2-methyl-	141	C7H8ClN	000087-60-5	83
5	Benzenamine, 4-chloro-2-methyl-	141	C7H8ClN	000095-69-2	80





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Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB925

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Quant Title : SVOA CALIBRATION

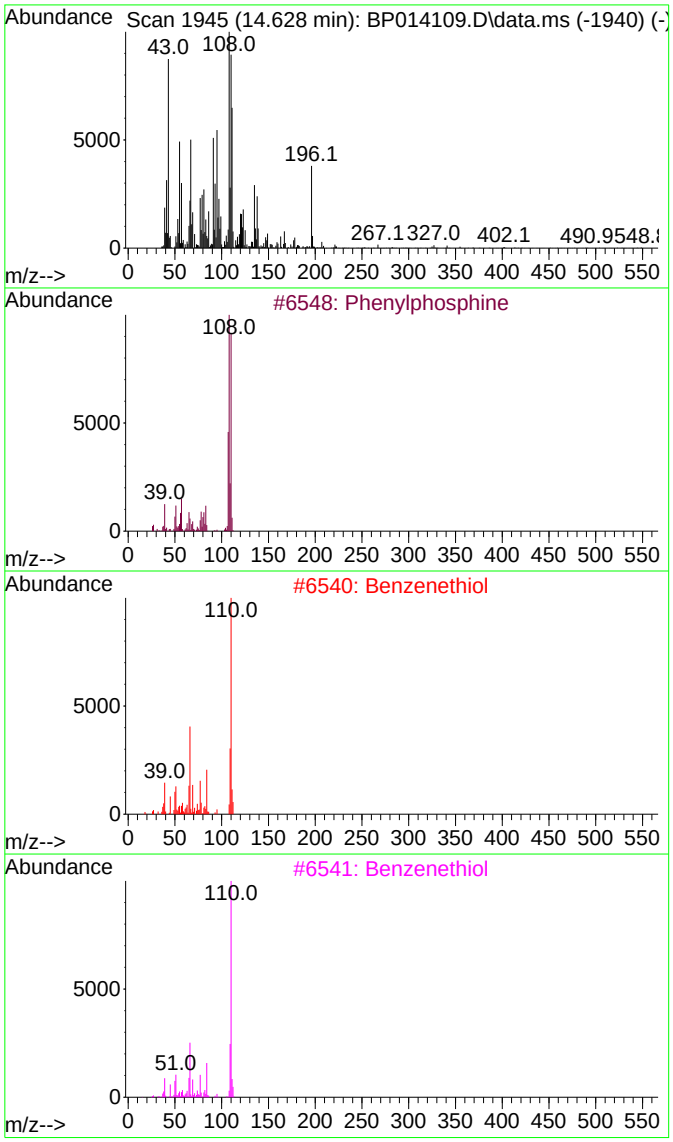
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

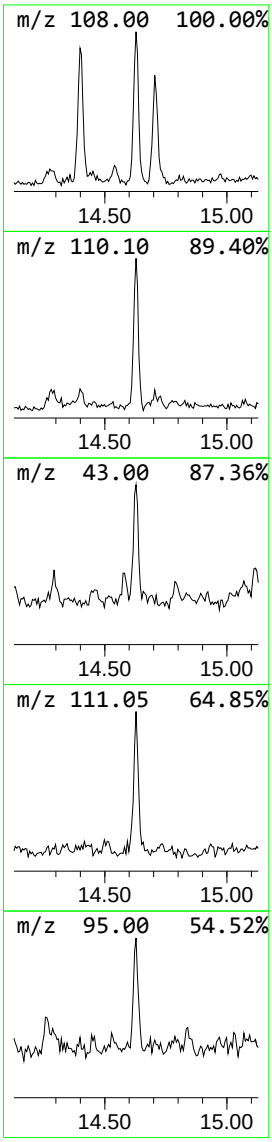
****

Peak Number 8 unknown-01 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.628	2.08 ng/ul	237246	Acenaphthene-d10	14.704

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenylphosphine	110	C6H7P	000638-21-1	38
2	Benzenethiol	110	C6H6S	000108-98-5	35
3	Benzenethiol	110	C6H6S	000108-98-5	35
4	2-Pyridinamine, 1-oxide	110	C5H6N2O	014150-95-9	27
5	Bicyclo(6.1.0)nonane-9-carboxyli...	168	C10H16O2	077841-63-5	25





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Operator : CG/JU
Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

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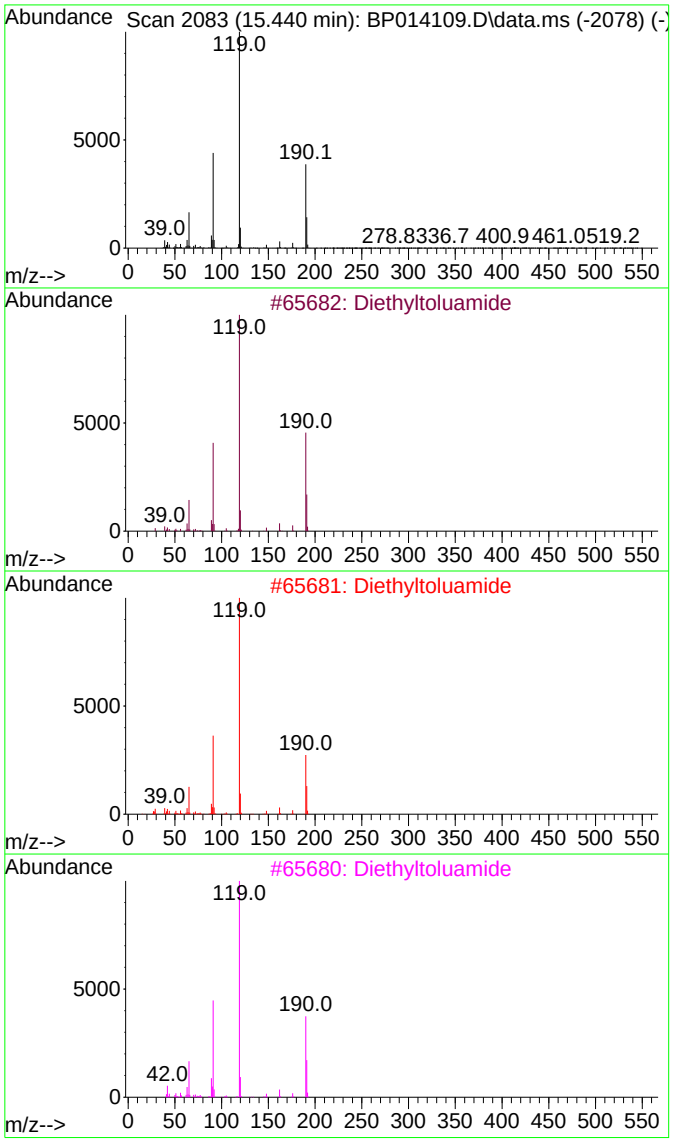
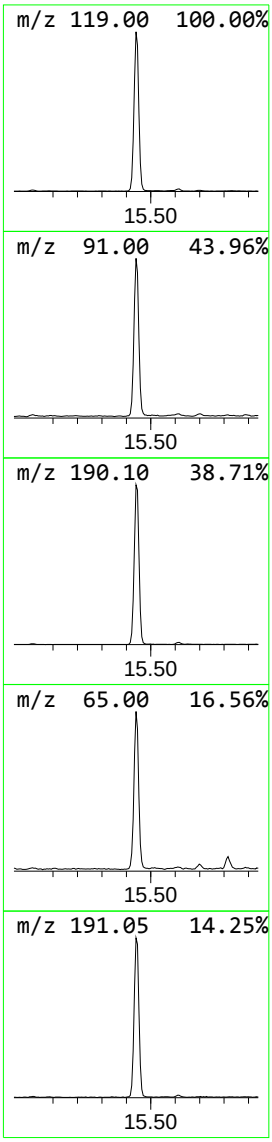
Peak Number 9 Diethyltoluamide Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.440	21.30 ng/ul	2434760	Acenaphthene-d10	14.704

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethyltoluamide	191	C12H17NO	000134-62-3	97
2	Diethyltoluamide	191	C12H17NO	000134-62-3	96
3	Diethyltoluamide	191	C12H17NO	000134-62-3	91
4	Benzamide, N,N-diethyl-4-methyl-	191	C12H17NO	002728-05-4	87
5	Benzamide, 3-methyl-N-methyl-N-p...	191	C12H17NO	1000421-46-2	81





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Operator : CG/JU
Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

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ClientSampleId :
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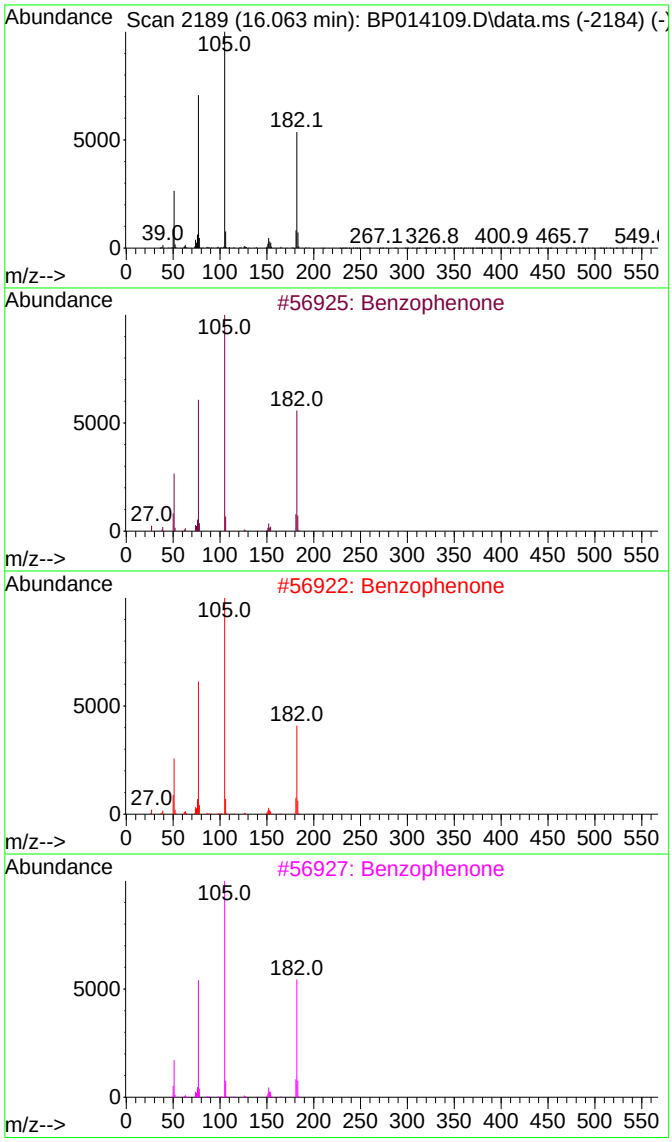
Quant Title : SVOA CALIBRATION

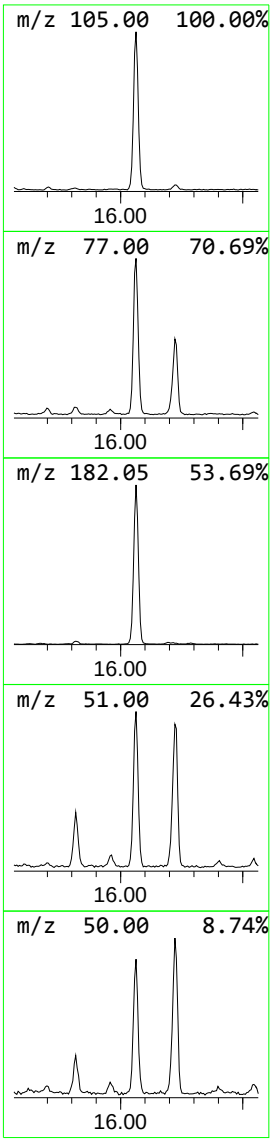
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzophenone Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.063	7.42 ng/ul	847785	Acenaphthene-d10	14.704

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	96
2	Benzophenone	182	C13H10O	000119-61-9	96
3	Benzophenone	182	C13H10O	000119-61-9	95
4	Benzophenone	182	C13H10O	000119-61-9	91
5	Benzophenone	182	C13H10O	000119-61-9	91





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Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

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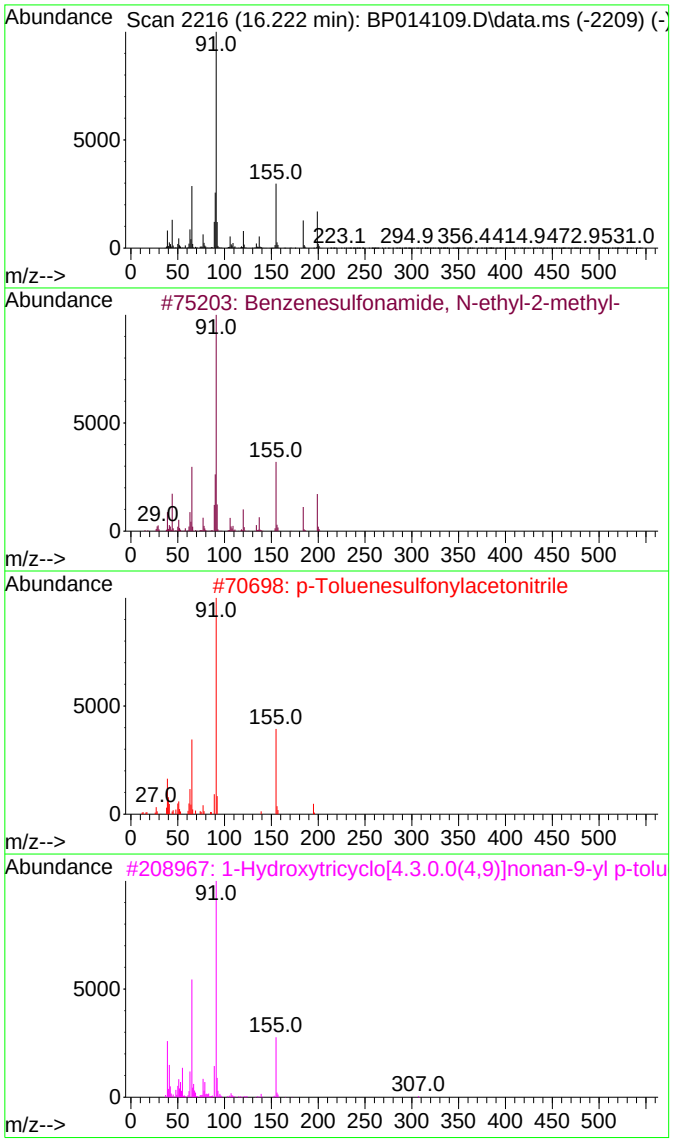
Quant Title : SVOA CALIBRATION

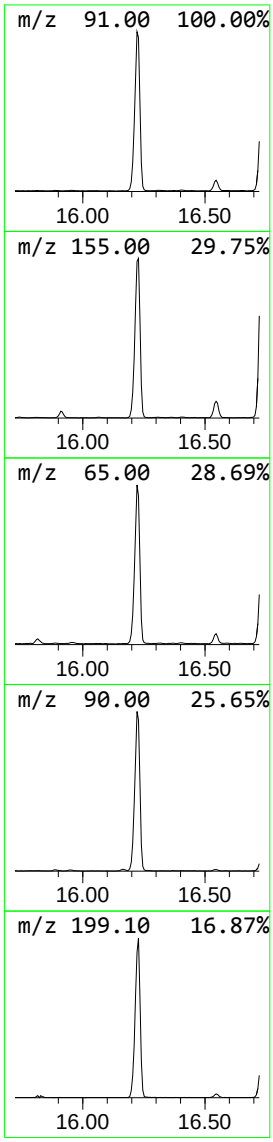
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzenesulfonamide, N-ethyl... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.222	34.26 ng/ul	4829910	Phenanthrene-d10	17.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	98
2	p-Toluenesulfonylacetonitrile	195	C9H9NO2S	005697-44-9	50
3	1-Hydroxytricyclo[4.3.0.0(4,9)]n...	308	C16H20O4S	201992-83-8	50
4	Tolbutamide	270	C12H18N2O3S	000064-77-7	50
5	(O-p-Tosylisonitroso)malononitrile	249	C10H7N3O3S	020893-01-0	47





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Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

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ClientSampleId :
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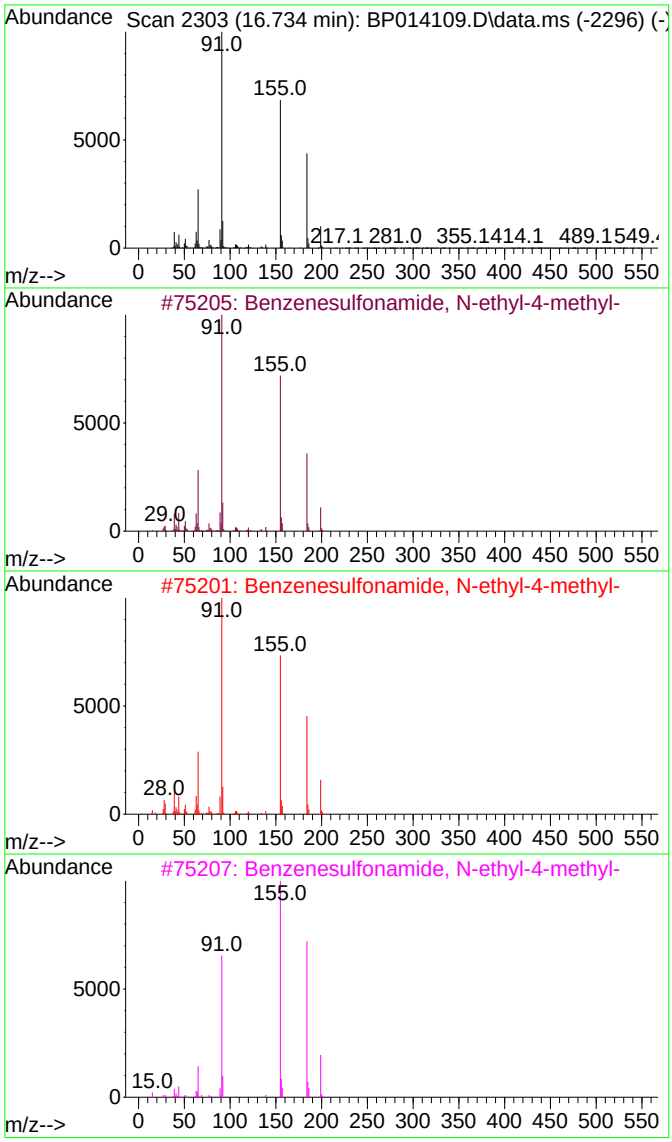
Quant Title : SVOA CALIBRATION

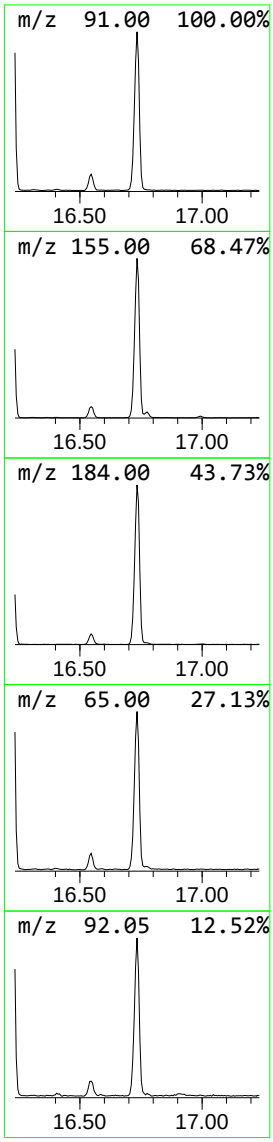
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Benzenesulfonamide, N-ethyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.734	22.28 ng/ul	3141350	Phenanthrene-d10	17.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	97
2	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	96
3	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	91
4	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	87
5	N-isobutyl-4-methylbenzenesulfon...	227	C11H17NO2S	023705-38-6	86





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Acq On : 16 Mar 2023 18:47
Operator : CG/JU
Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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Quant Title : SVOA CALIBRATION

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

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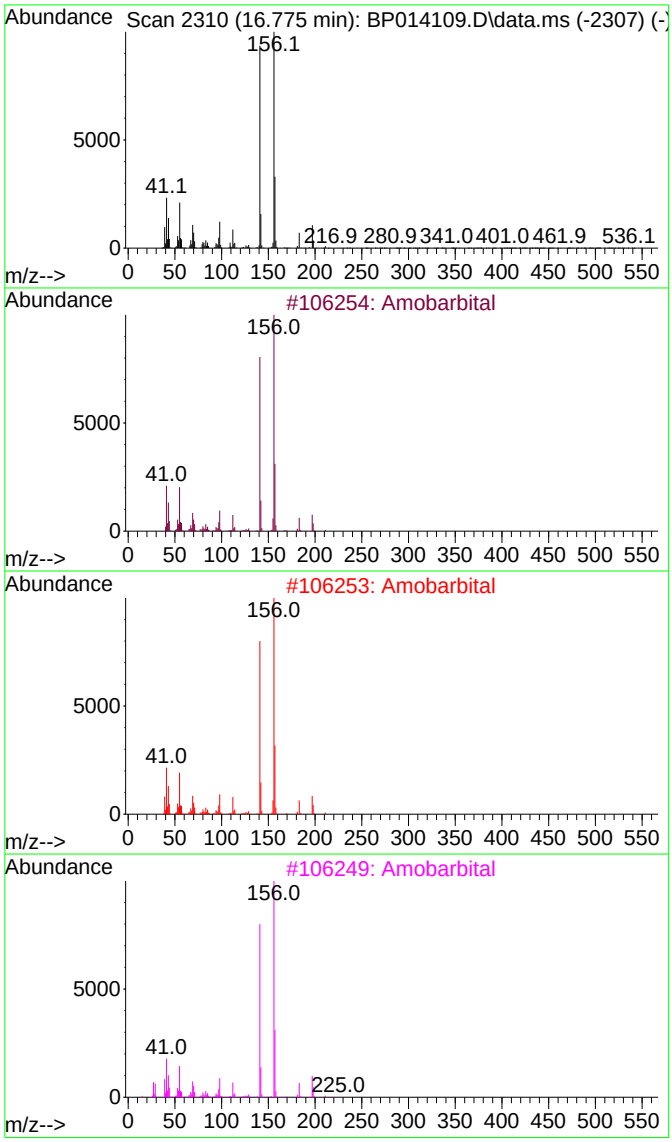
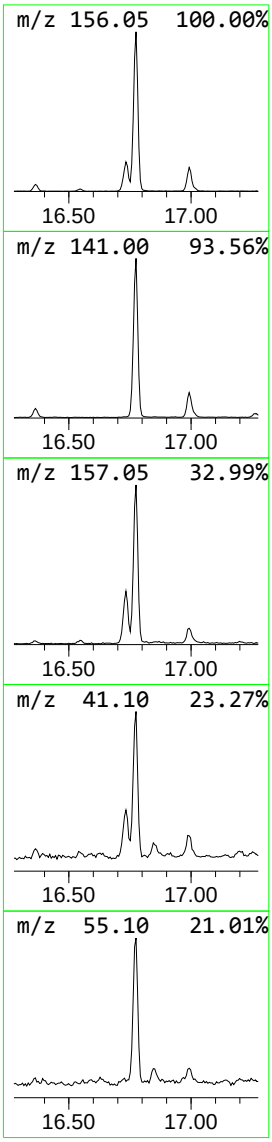
Peak Number 13 Amobarbital Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.775	7.79 ng/ul	1098470	Phenanthrene-d10	17.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Amobarbital	226	C11H18N2O3	000057-43-2	91
2	Amobarbital	226	C11H18N2O3	000057-43-2	91
3	Amobarbital	226	C11H18N2O3	000057-43-2	91
4	Amobarbital	226	C11H18N2O3	000057-43-2	91
5	Amobarbital	226	C11H18N2O3	000057-43-2	90





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Quant Title : SVOA CALIBRATION

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

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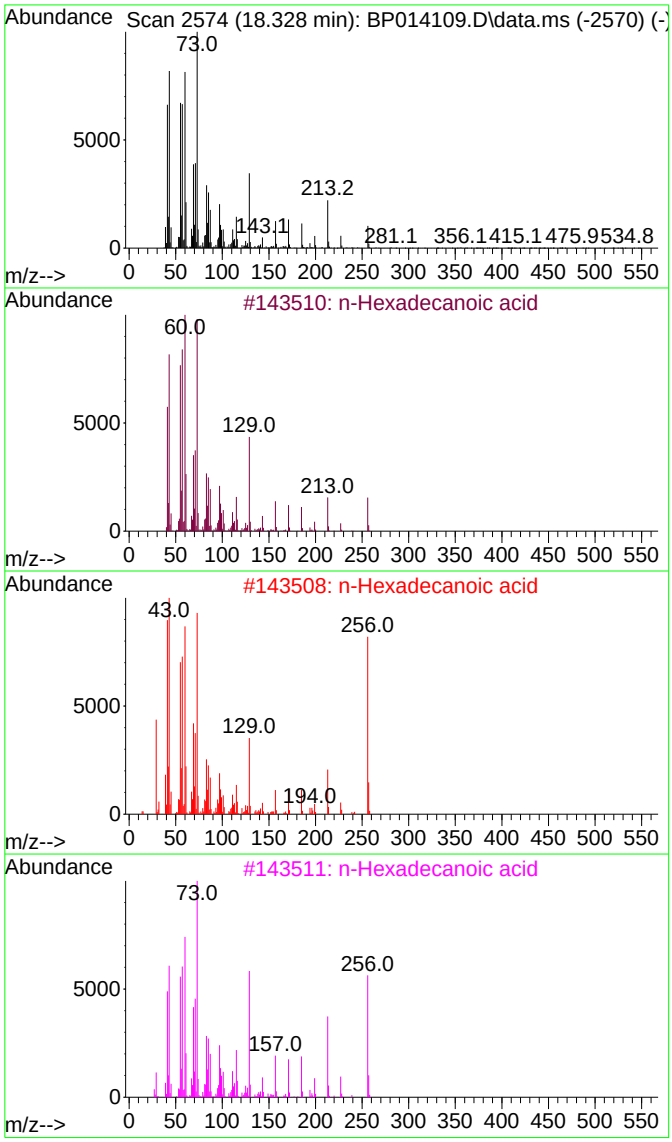
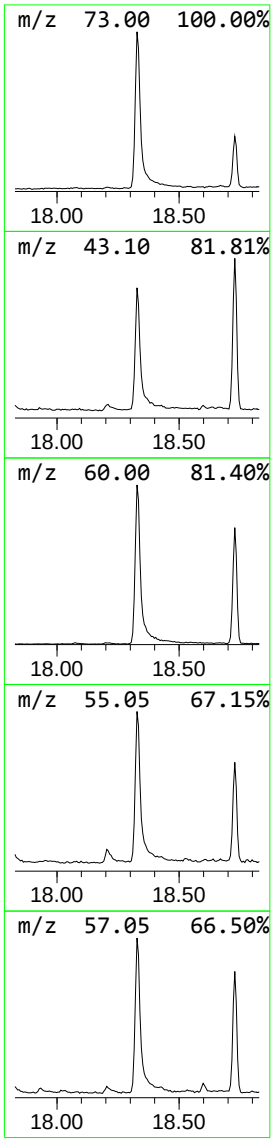
Peak Number 14 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.328	14.93 ng/ul	2104440	Phenanthrene-d10	17.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5	Tridecanoic acid	214	C13H26O2	000638-53-9	93





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Sample : 01884-13
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ALS Vial : 13 Sample Multiplier: 1

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Quant Title : SVOA CALIBRATION

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

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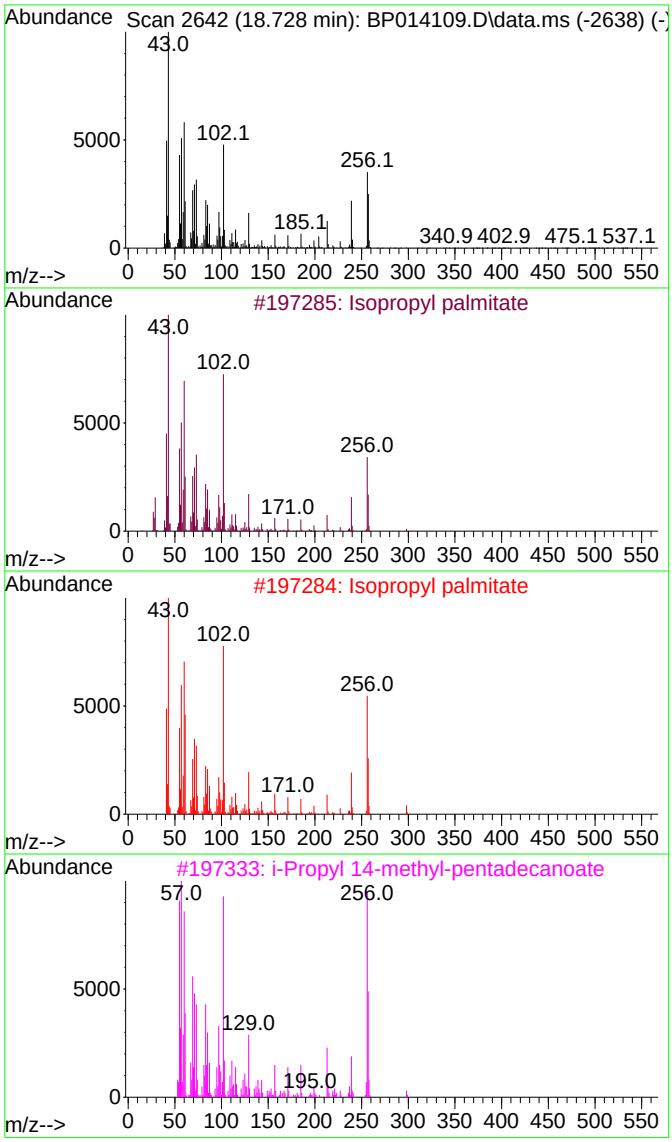
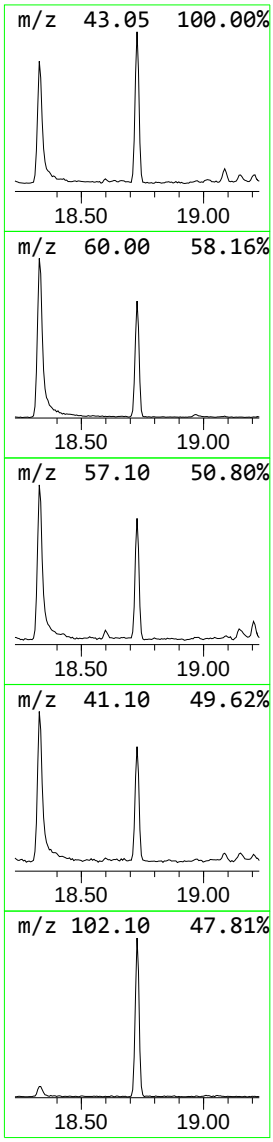
Peak Number 15 Isopropyl palmitate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.728	10.31 ng/ul	1452990	Phenanthrene-d10	17.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isopropyl palmitate	298	C19H38O2	000142-91-6	91
2	Isopropyl palmitate	298	C19H38O2	000142-91-6	59
3	i-Propyl 14-methyl-pentadecanoate	298	C19H38O2	1000336-62-4	53
4	Isopropyl palmitate	298	C19H38O2	000142-91-6	46
5	Nonanoic acid, 1-methylethyl ester	200	C12H24O2	028267-32-5	43





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ALS Vial : 13 Sample Multiplier: 1

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Quant Title : SVOA CALIBRATION

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

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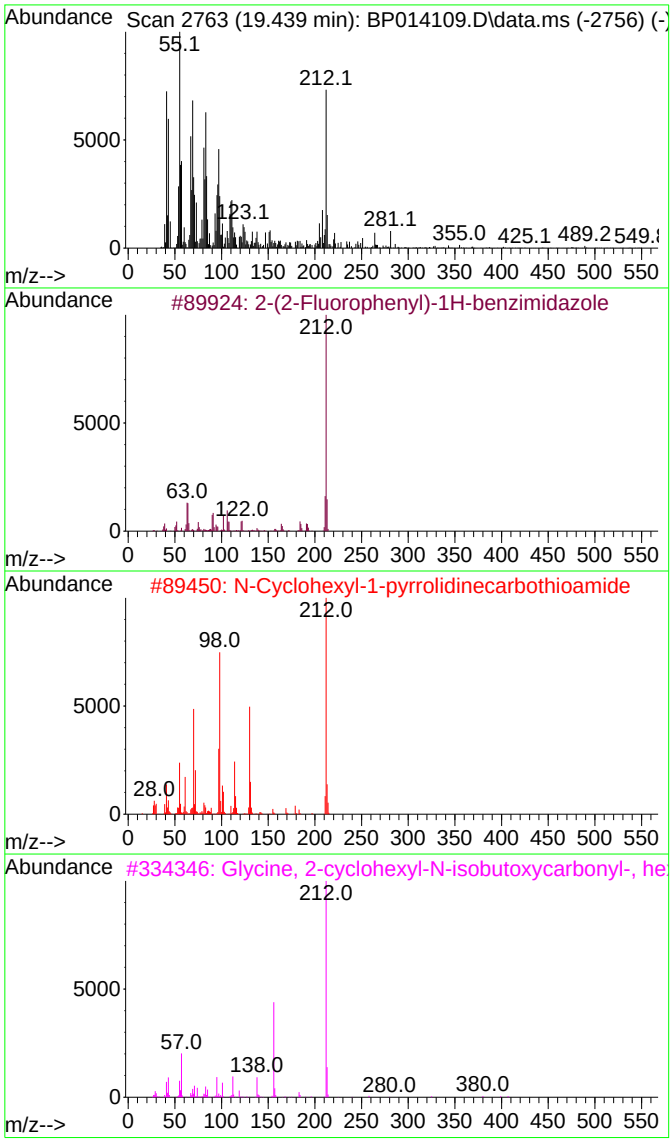
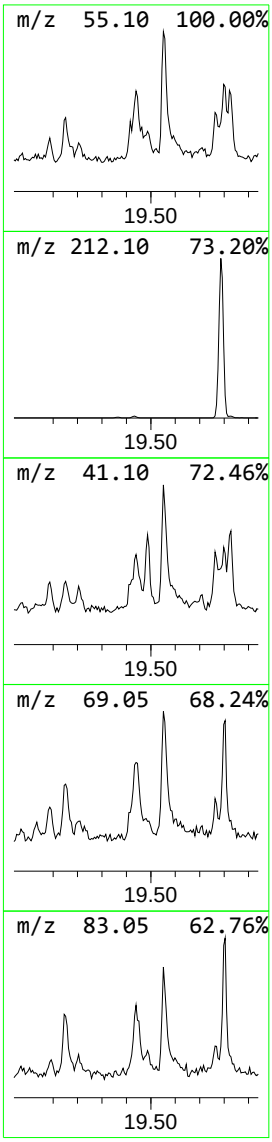
Peak Number 16 unknown-02 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.439	3.13 ng/ul	440728	Phenanthrene-d10	17.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(2-Fluorophenyl)-1H-benzimidazole	212	C13H9FN2	000321-51-7	44
2	N-Cyclohexyl-1-pyrrolidinecarbot...	212	C11H20N2S	015260-45-4	22
3	Glycine, 2-cyclohexyl-N-isobutox...	481	C29H55NO4	1000383-10-6	22
4	5-(Methylthio)-Salicylic acid, O...	212	C10H12O3S	1000374-80-1	22
5	2,6,6-Trimethyl-9-undecen-1-ol	212	C14H28O	1010131-31-6	22





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ALS Vial : 13 Sample Multiplier: 1

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Quant Title : SVOA CALIBRATION

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

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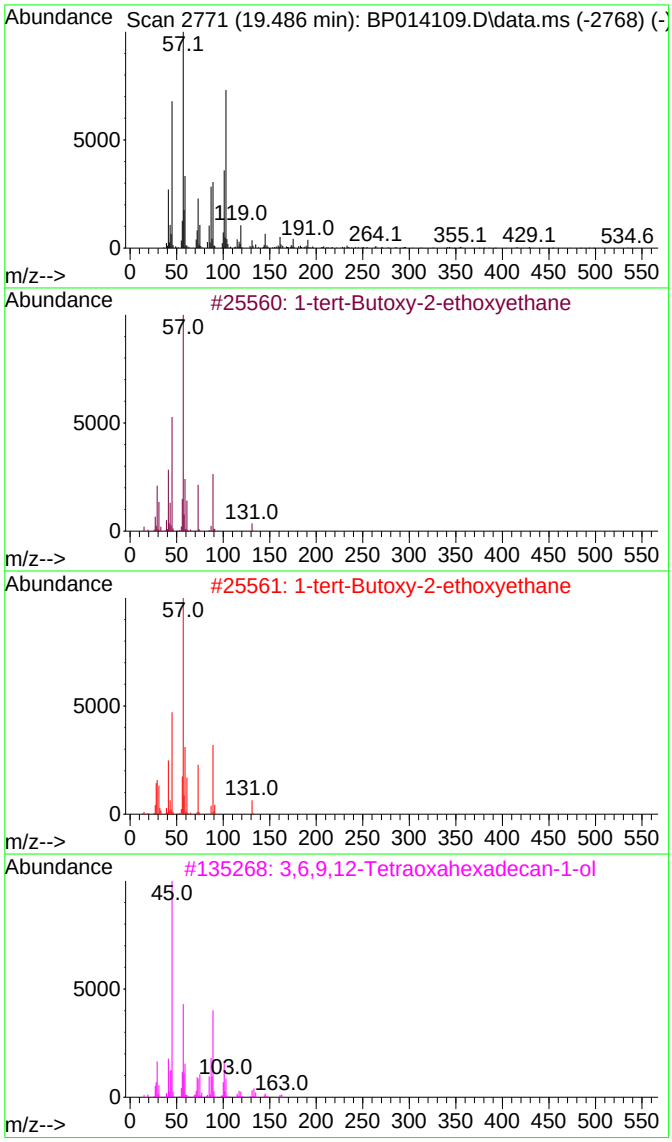
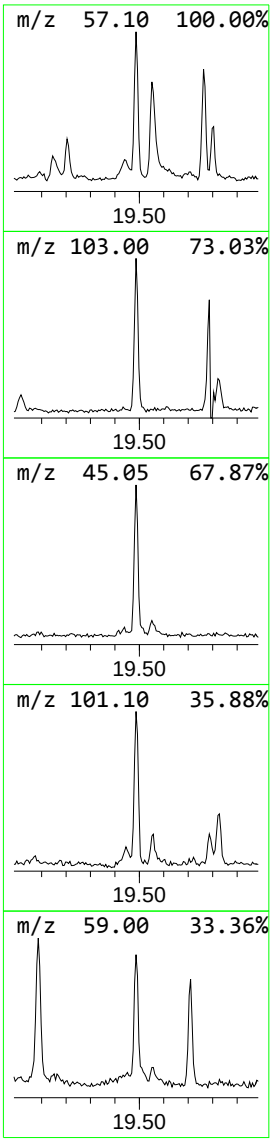
Peak Number 17 unknown-03 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.486	2.16 ng/ul	304770	Phenanthrene-d10	17.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-tert-Butoxy-2-ethoxyethane	146	C8H18O2	051422-54-9	43
2	1-tert-Butoxy-2-ethoxyethane	146	C8H18O2	051422-54-9	37
3	3,6,9,12-Tetraoxahexadecan-1-ol	250	C12H26O5	001559-34-8	35
4	Carbonic acid, bis(1-methylpropy...	174	C9H18O3	000623-63-2	35
5	Di-sec-Butyl ether	130	C8H18O	006863-58-7	35





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Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

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ClientSampleId :
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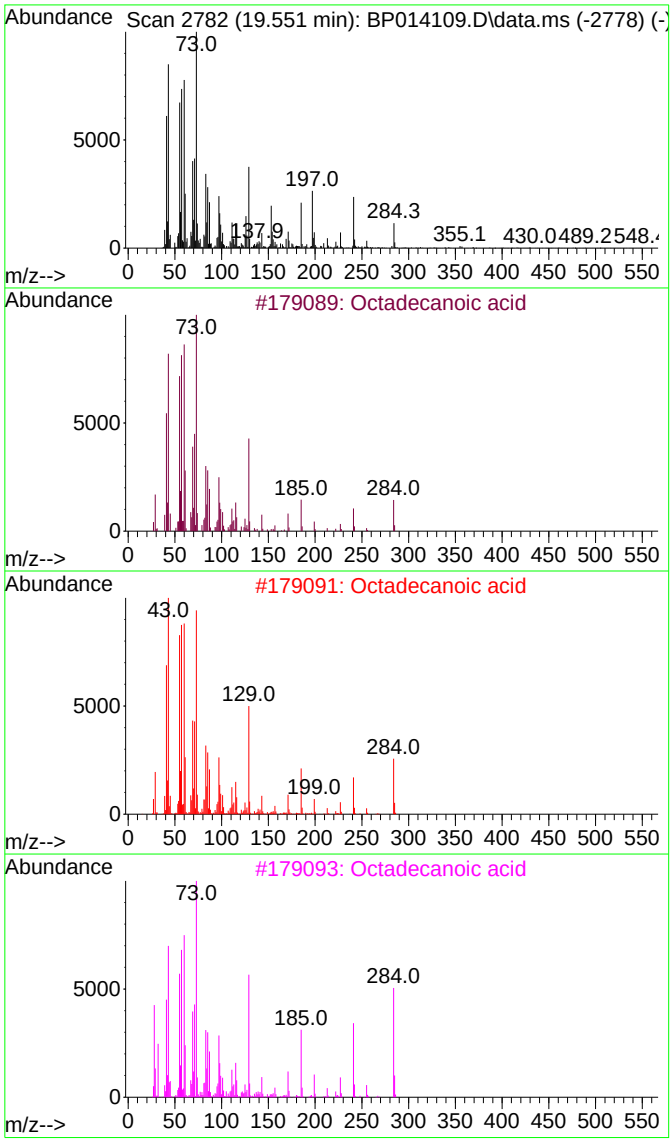
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

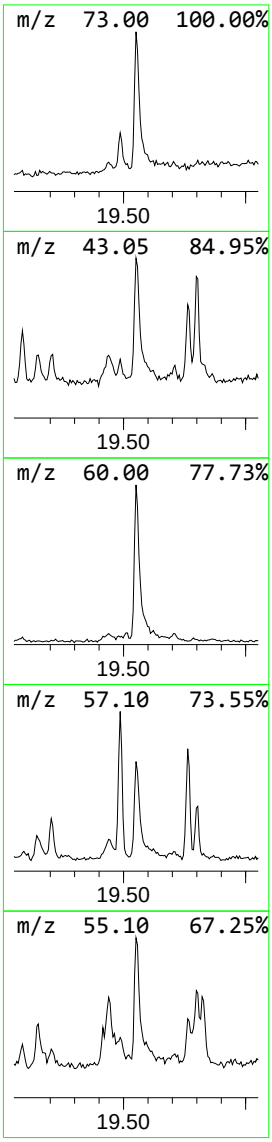
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Octadecanoic acid Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.551	4.68 ng/ul	742399	Chrysene-d12	21.545

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	99
3	Octadecanoic acid	284	C18H36O2	000057-11-4	96
4	Octadecanoic acid	284	C18H36O2	000057-11-4	94
5	Octadecanoic acid	284	C18H36O2	000057-11-4	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031623\
Data File : BP014109.D
Acq On : 16 Mar 2023 18:47
Operator : CG/JU
Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB925

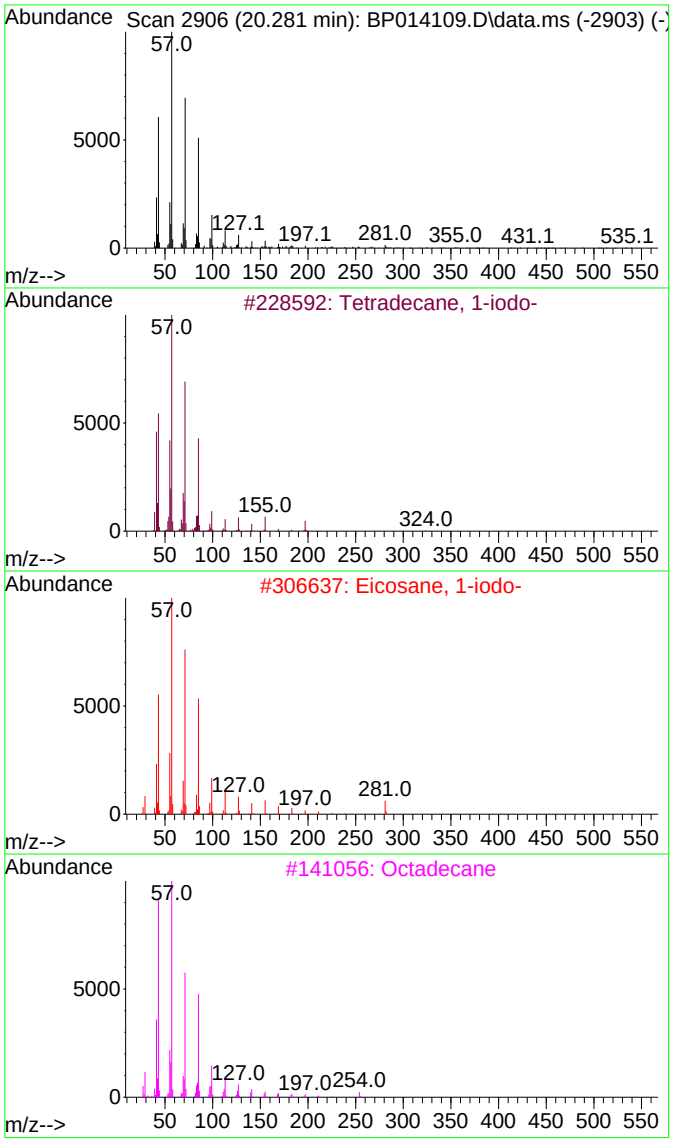
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

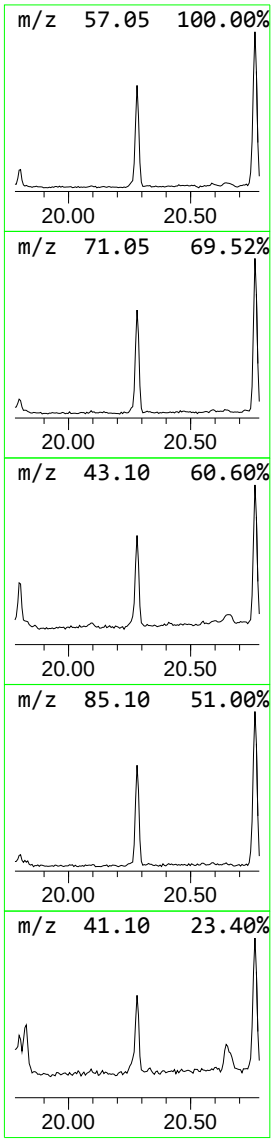
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Tetradecane, 1-iodo- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.281	2.57 ng/ul	407867	Chrysene-d12	21.545

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	91
2	Eicosane, 1-iodo-	408	C20H41I	1000406-31-8	91
3	Octadecane	254	C18H38	000593-45-3	90
4	Nonadecane	268	C19H40	000629-92-5	90
5	Docosane, 1-iodo-	436	C22H45I	1000406-31-9	87





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Acq On : 16 Mar 2023 18:47
Operator : CG/JU
Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB925

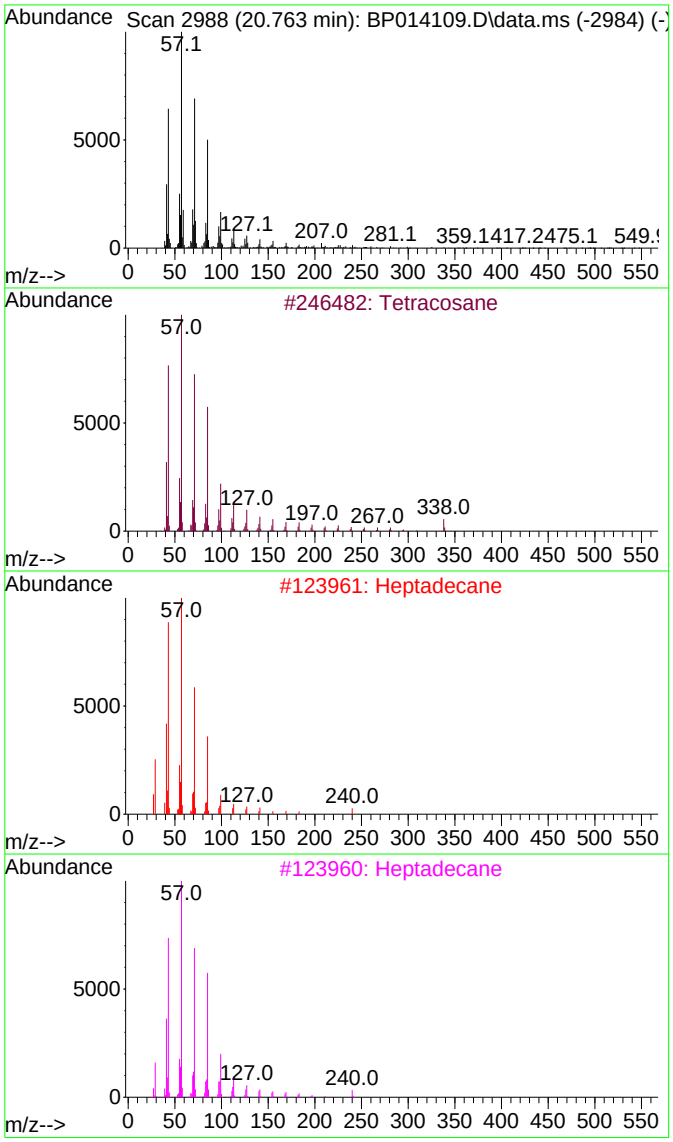
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

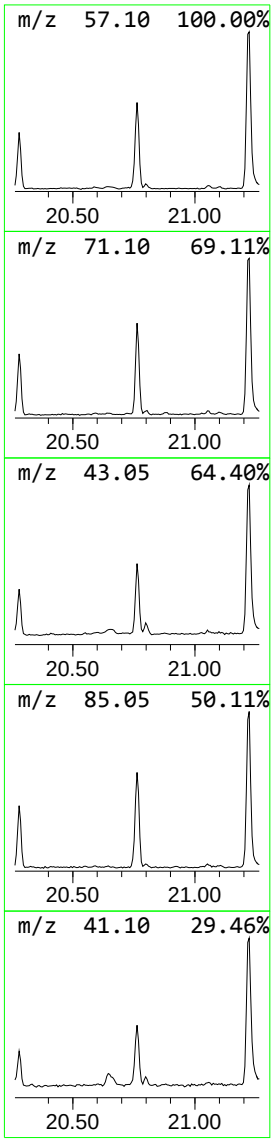
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.763	4.99 ng/ul	792034	Chrysene-d12	21.545

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	96
2	Heptadecane	240	C17H36	000629-78-7	94
3	Heptadecane	240	C17H36	000629-78-7	91
4	2-Bromo dodecane	248	C12H25Br	013187-99-0	90
5	11-Methylpentacosane	366	C26H54	015689-71-1	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031623\
Data File : BP014109.D
Acq On : 16 Mar 2023 18:47
Operator : CG/JU
Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB925

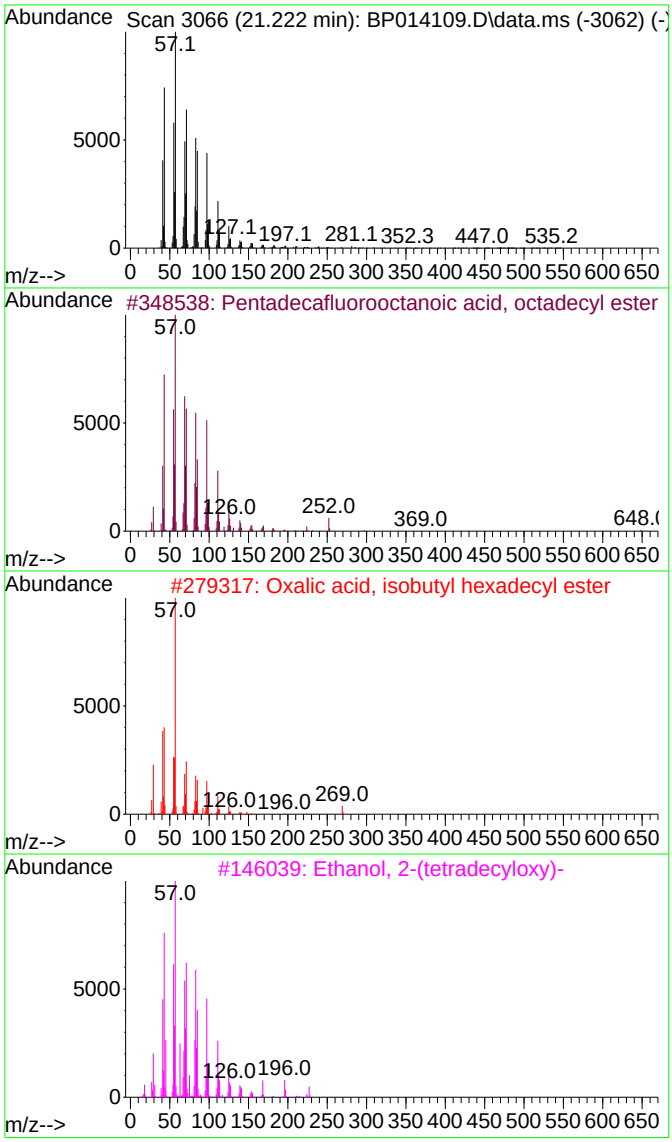
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

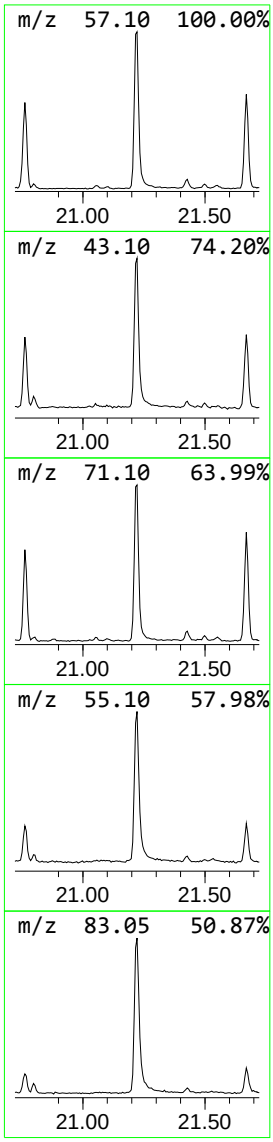
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 Pentadecafluorooctanoic aci... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.222	14.30 ng/ul	2269030	Chrysene-d12	21.545

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	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	90
4	Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	76
5	Pentafluoropropionic acid, tride...	346	C16H27F5O2	959261-22-4	76





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031623\
Data File : BP014109.D
Acq On : 16 Mar 2023 18:47
Operator : CG/JU
Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB925

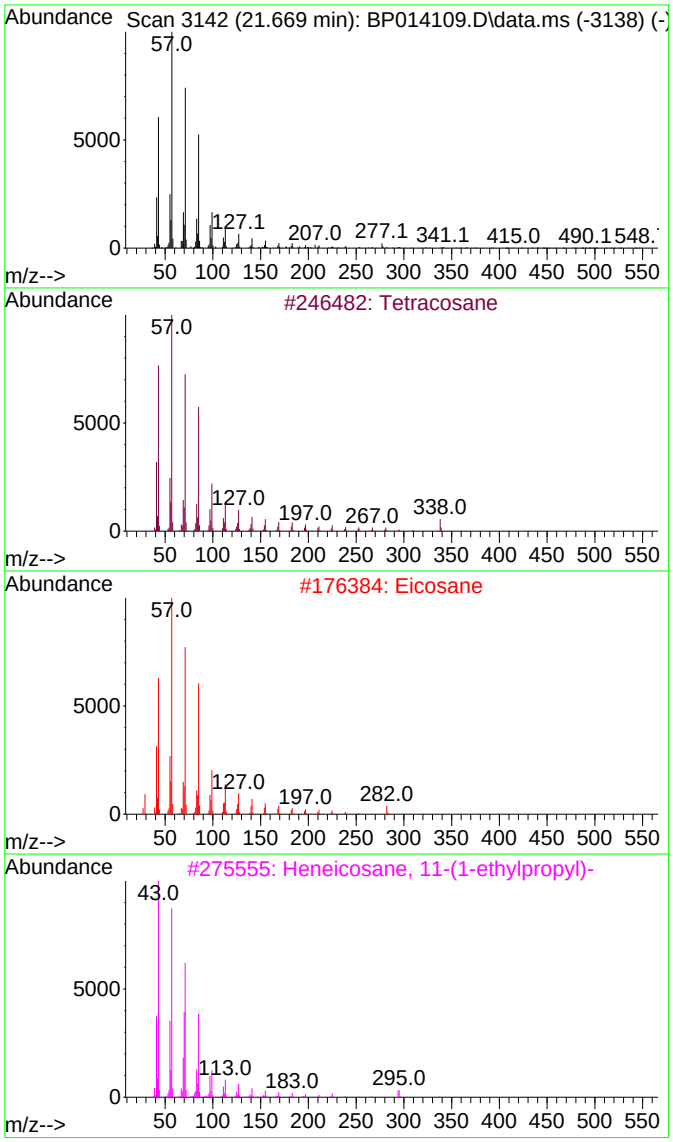
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

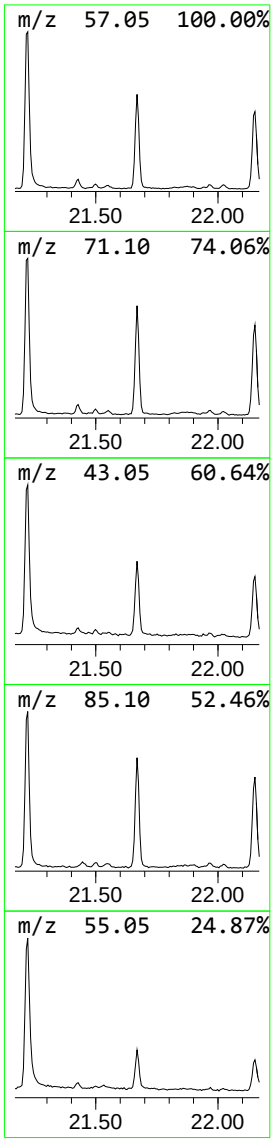
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.669	4.75 ng/ul	754236	Chrysene-d12	21.545

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	96
2	Eicosane	282	C20H42	000112-95-8	94
3	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
4	Heneicosane	296	C21H44	000629-94-7	91
5	Octacosane	394	C28H58	000630-02-4	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031623\
Data File : BP014109.D
Acq On : 16 Mar 2023 18:47
Operator : CG/JU
Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB925

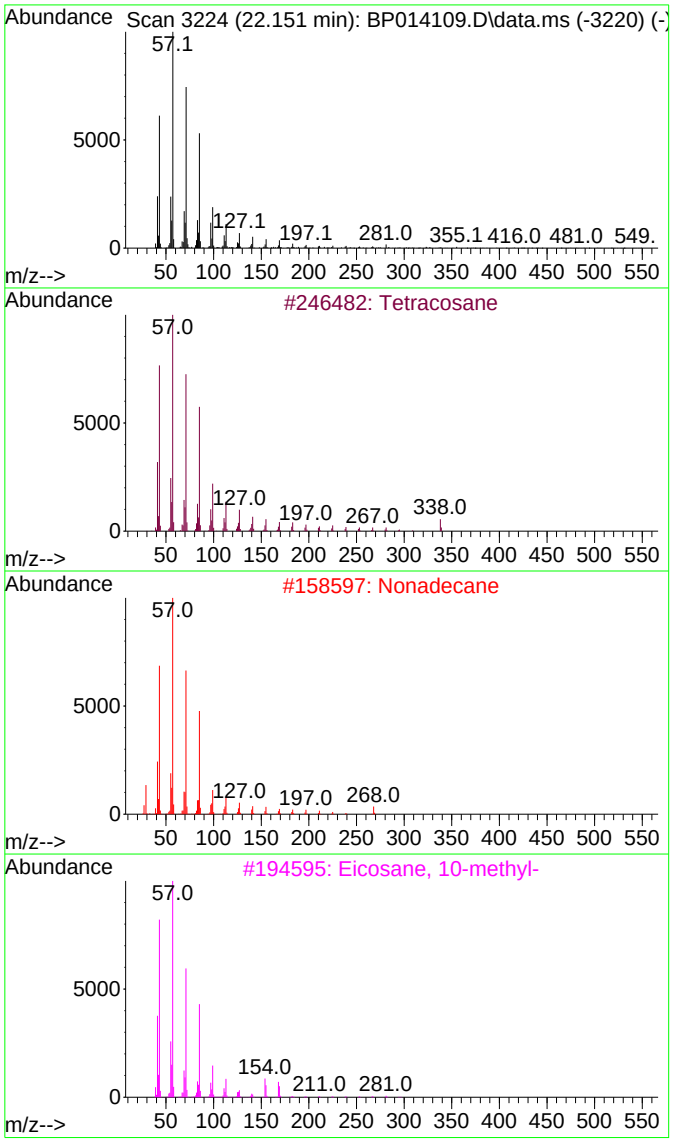
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

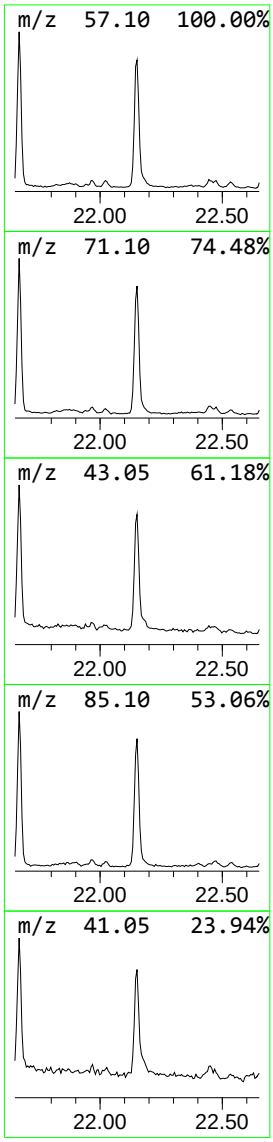
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.151	5.16 ng/ul	819095	Chrysene-d12	21.545

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	97
2	Nonadecane	268	C19H40	000629-92-5	94
3	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93
5	Heptacosane	380	C27H56	000593-49-7	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031623\
Data File : BP014109.D
Acq On : 16 Mar 2023 18:47
Operator : CG/JU
Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB925

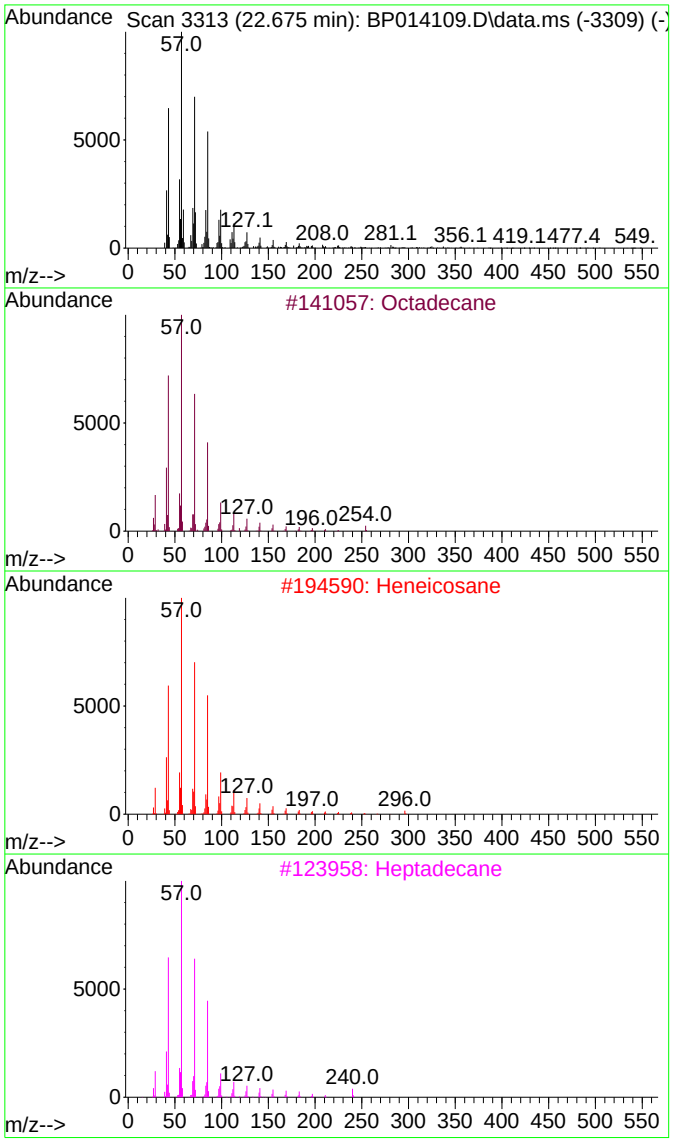
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

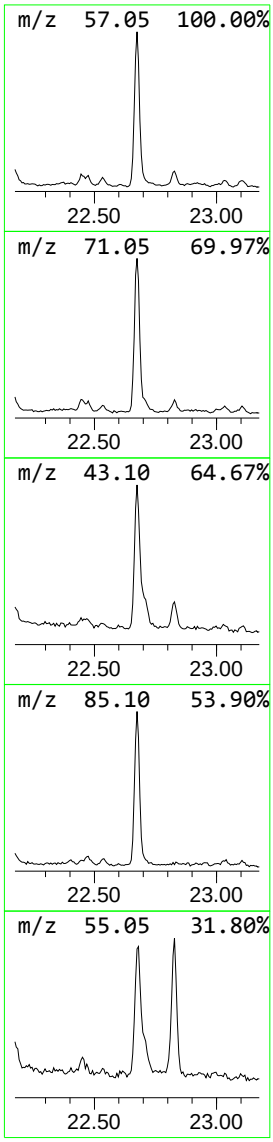
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.674	7.81 ng/ul	1238950	Chrysene-d12	21.545

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	96
2	Heneicosane	296	C21H44	000629-94-7	94
3	Heptadecane	240	C17H36	000629-78-7	93
4	Octacosane	394	C28H58	000630-02-4	91
5	2-Bromo dodecane	248	C12H25Br	013187-99-0	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031623\
Data File : BP014109.D
Acq On : 16 Mar 2023 18:47
Operator : CG/JU
Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB925

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

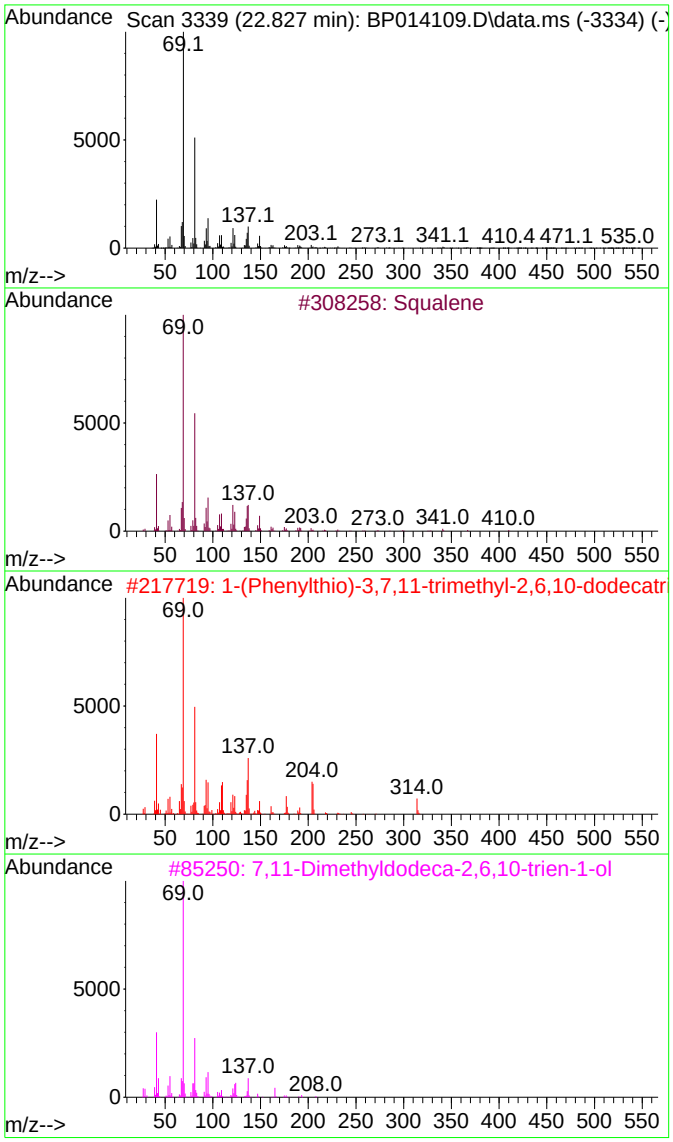
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

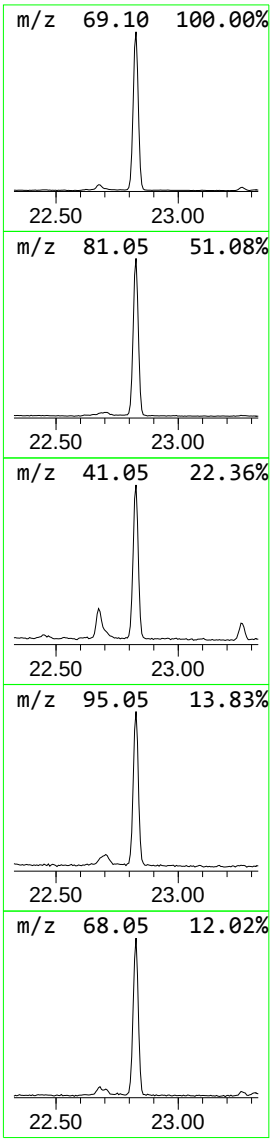
****

Peak Number 25 Squalene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.828	19.09 ng/ul	2780590	Perylene-d12	24.080

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Squalene	410	C30H50	000111-02-4	96
2	1-(Phenylthio)-3,7,11-trimethyl-...	314	C21H30S	028413-57-2	78
3	7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	72
4	(.+/-.)-Lavandulol, pentafluorop...	300	C13H17F5O2	1000352-63-1	64
5	Farnesol isomer a	222	C15H26O	1000108-92-4	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031623\
Data File : BP014109.D
Acq On : 16 Mar 2023 18:47
Operator : CG/JU
Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB925

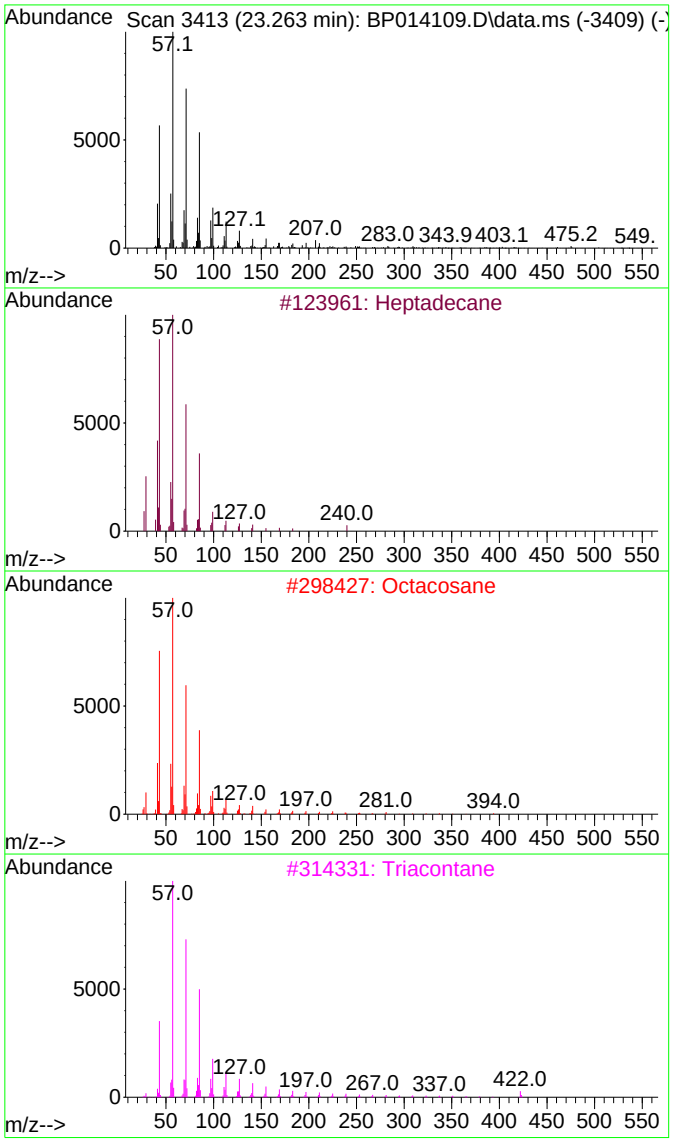
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

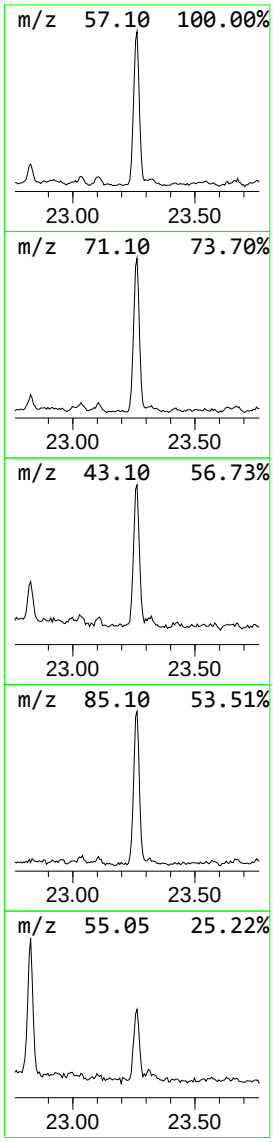
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.263	3.31 ng/ul	481967	Perylene-d12	24.080

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	93
2	Octacosane	394	C28H58	000630-02-4	91
3	Triacontane	422	C30H62	000638-68-6	91
4	Hexacosane	366	C26H54	000630-01-3	91
5	Heneicosane	296	C21H44	000629-94-7	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031623\
Data File : BP014109.D
Acq On : 16 Mar 2023 18:47
Operator : CG/JU
Sample : 01884-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB925

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.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
1,3-Oxathiolane	4.646	9.2	ng/ul	527802	1	8.064	1151020	20.0
Benzenamine, 3-...	9.058	15.5	ng/ul	892075	1	8.064	1151020	20.0
Acetaldehyde, (...	9.593	3.2	ng/ul	254120	2	10.887	1570440	20.0
Phenol, p-tert-...	12.216	5.0	ng/ul	396401	2	10.887	1570440	20.0
Benzenamine, 3-...	12.387	3.1	ng/ul	244854	2	10.887	1570440	20.0
m-Toluidine, 4-...	12.487	2.1	ng/ul	165804	2	10.887	1570440	20.0
unknown-01	14.628	2.1	ng/ul	237246	3	14.704	2286070	20.0
Diethyltoluamide	15.440	21.3	ng/ul	2434760	3	14.704	2286070	20.0
Benzophenone	16.063	7.4	ng/ul	847785	3	14.704	2286070	20.0
Benzenesulfonam...	16.222	34.3	ng/ul	4829910	4	17.463	2819640	20.0
Benzenesulfonam...	16.734	22.3	ng/ul	3141350	4	17.463	2819640	20.0
Amobarbital	16.775	7.8	ng/ul	1098470	4	17.463	2819640	20.0
n-Hexadecanoic ...	18.328	14.9	ng/ul	2104440	4	17.463	2819640	20.0
Isopropyl palmi...	18.728	10.3	ng/ul	1452990	4	17.463	2819640	20.0
unknown-02	19.439	3.1	ng/ul	440728	4	17.463	2819640	20.0
unknown-03	19.486	2.2	ng/ul	304770	4	17.463	2819640	20.0
Octadecanoic acid	19.551	4.7	ng/ul	742399	5	21.545	3172610	20.0
Tetradecane, 1-...	20.281	2.6	ng/ul	407867	5	21.545	3172610	20.0
(DEL) Alkane: S...	20.763	5.0	ng/ul	792034	5	21.545	3172610	20.0
Pentadecafluoro...	21.222	14.3	ng/ul	2269030	5	21.545	3172610	20.0
(DEL) Alkane: S...	21.669	4.8	ng/ul	754236	5	21.545	3172610	20.0
(DEL) Alkane: S...	22.151	5.2	ng/ul	819095	5	21.545	3172610	20.0
(DEL) Alkane: S...	22.674	7.8	ng/ul	1238950	5	21.545	3172610	20.0
Squalene	22.828	19.1	ng/ul	2780590	6	24.080	2913680	20.0
(DEL) Alkane: S...	23.263	3.3	ng/ul	481967	6	24.080	2913680	20.0