LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020823\
Data File : BP013793.D
Acq On : 08 Feb 2023 12:56
Operator : CG/JU
Sample : 01412-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A44M4

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-BP020723.M
Title : SVOA CALIBRATION

Signal : TIC: BP013793.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.440	38	43	45	rBV2	32314	44531	0.64%	0.071%
2	3.469	45	48	58	rVB	33727	52461	0.76%	0.084%
3	3.905	119	122	138	rBV	141118	240758	3.47%	0.385%
4	4.140	157	162	165	rBV	11915	17505	0.25%	0.028%
5	4.240	175	179	186	rVB	14364	19348	0.28%	0.031%
6	4.346	192	197	201	rBV7	6598	10461	0.15%	0.017%
7	4.881	284	288	292	rBV3	6691	9942	0.14%	0.016%
8	5.187	333	340	352	rVB3	13835	29064	0.42%	0.046%
9	5.399	369	376	391	rVB	1207165	1867901	26.95%	2.986%
10	5.581	404	407	414	rBV4	9675	22708	0.33%	0.036%
11	5.810	442	446	449	rBV	6168	7615	0.11%	0.012%
12	6.105	489	496	500	rVB7	3995	7964	0.11%	0.013%
13	6.493	557	562	565	rVB5	6848	10636	0.15%	0.017%
14	6.534	565	569	575	rVB7	4371	7658	0.11%	0.012%
15	7.093	657	664	667	rBV7	4307	8330	0.12%	0.013%
16	7.163	672	676	681	rVV5	4562	7392	0.11%	0.012%
17	7.269	684	694	706	rVV	126583	217224	3.13%	0.347%
18	7.440	715	723	733	rBV	472510	740922	10.69%	1.184%
19	7.646	750	758	770	rBV	421398	702876	10.14%	1.124%
20	8.010	812	820	828	rBV	71265	121167	1.75%	0.194%
21	8.122	831	839	842	rVV	630915	1045074	15.08%	1.671%
22	8.157	842	845	856	rVB	700195	1069255	15.43%	1.709%
23	8.481	892	900	913	rBV	4114312	6931208	100.00%	11.079%
24	8.822	952	958	967	rBV	371205	618417	8.92%	0.989%
25	8.928	970	976	985	rVV	176887	290581	4.19%	0.464%
26	9.028	988	993	1002	rVV2	29694	65905	0.95%	0.105%
27	9.116	1002	1008	1017	rVB	36065	64699	0.93%	0.103%
28	9.293	1032	1038	1051	rBV	490260	811219	11.70%	1.297%
29	9.693	1101	1106	1112	rVB2	13670	22068	0.32%	0.035%
30	9.751	1112	1116	1121	rVB2	13311	19351	0.28%	0.031%
31	9.969	1140	1153	1156	rBV	39372	69159	1.00%	0.111%
32	10.022	1156	1162	1175	rVB	410064	665409	9.60%	1.064%
33	10.563	1244	1254	1262	rBV	716557	1197900	17.28%	1.915%
34	10.628	1263	1265	1271	rVB6	9399	15973	0.23%	0.026%
35	10.722	1275	1281	1288	rVB2	17310	32163	0.46%	0.051%
36	10.810	1288	1296	1312	rVB	378264	624957	9.02%	0.999%

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Operator : CG/JU
Sample : 01412-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A44M4

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-BP020723.M
Title : SVOA CALIBRATION

37	10.951	1312	1320	1327	rBV	939066	1604179	23.14%	2.564%
38	11.081	1334	1342	1360	rVB	548998	1045990	15.09%	1.672%
39	11.334	1379	1385	1395	rVB	72419	130819	1.89%	0.209%
40	11.604	1427	1431	1442	rVV8	4655	12386	0.18%	0.020%
41	12.193	1527	1531	1536	rVB7	5866	10631	0.15%	0.017%
42	12.257	1536	1542	1546	rBV9	3891	7426	0.11%	0.012%
43	12.316	1546	1552	1560	rBV	36868	73162	1.06%	0.117%
44	12.457	1569	1576	1583	rBV	40052	67295	0.97%	0.108%
45	12.522	1583	1587	1591	rBV3	15661	26687	0.39%	0.043%
46	12.710	1614	1619	1623	rBV5	10044	20992	0.30%	0.034%
47	12.898	1645	1651	1657	rBV	50581	87304	1.26%	0.140%
48	12.969	1658	1663	1671	rVB	61103	94892	1.37%	0.152%
49	13.187	1695	1700	1704	rBV8	5016	9321	0.13%	0.015%
50	13.292	1711	1718	1724	rBV4	25917	50441	0.73%	0.081%
51	13.357	1727	1729	1733	rVV5	6418	7407	0.11%	0.012%
52	13.569	1758	1765	1767	rBV7	14206	24647	0.36%	0.039%
53	13.592	1767	1769	1773	rVV	19351	29343	0.42%	0.047%
54	13.639	1774	1777	1789	rVB6	14453	25405	0.37%	0.041%
55	13.763	1794	1798	1802	rVV3	25244	41312	0.60%	0.066%
56	13.810	1802	1806	1815	rVB2	41918	65745	0.95%	0.105%
57	14.157	1858	1865	1876	rBV	1478908	2084052	30.07%	3.331%
58	14.451	1909	1915	1927	rVB	1621019	2416068	34.86%	3.862%
59	14.639	1943	1947	1956	rVB9	9885	19376	0.28%	0.031%
60	14.757	1958	1967	1975	rBV	1612049	2433133	35.10%	3.889%
61	14.863	1983	1985	1993	rVB9	4959	8276	0.12%	0.013%
62	14.945	1993	1999	2010	rBV	96805	185090	2.67%	0.296%
63	15.039	2010	2015	2021	rVB2	23827	40351	0.58%	0.064%
64	15.151	2030	2034	2039	rVB7	10706	18642	0.27%	0.030%
65	15.251	2048	2051	2052	rBV3	9833	9705	0.14%	0.016%
66	15.292	2052	2058	2065	rVB	861223	1184943	17.10%	1.894%
67	15.375	2070	2072	2077	rVB5	8990	12195	0.18%	0.019%
68	15.586	2103	2108	2112	rBV6	7539	10797	0.16%	0.017%
69	15.751	2129	2136	2147	rBV	2208498	3245039	46.82%	5.187%
70	15.857	2148	2154	2164	rBV	409028	571060	8.24%	0.913%
71	16.004	2173	2179	2189	rVB	64583	99882	1.44%	0.160%
72	16.110	2192	2197	2205	rVV2	184090	285333	4.12%	0.456%
73	16.192	2207	2211	2215	rVB	53357	62497	0.90%	0.100%
74	16.251	2215	2221	2230	rBV	306633	453746	6.55%	0.725%
75	16.716	2297	2300	2304	rVB2	18241	23364	0.34%	0.037%
76	16.886	2326	2329	2333	rVB5	11768	14610	0.21%	0.023%

LSC Area Percent Report

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Data File : BP013793.D
Acq On : 08 Feb 2023 12:56
Operator : CG/JU
Sample : 01412-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : A44M4

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_P\Methods\SFAM-EPA-BP020723.M
Title : SVOA CALIBRATION

77	17.157	2370	2375	2385	rVB	84172	143285	2.07%	0.229%
78	17.380	2409	2413	2423	rBV3	26356	48351	0.70%	0.077%
79	17.510	2429	2435	2443	rBV	2263187	3230234	46.60%	5.163%
80	17.610	2446	2452	2459	rBV	2611297	3660932	52.82%	5.852%
81	17.757	2473	2477	2479	rBV5	14847	21523	0.31%	0.034%
82	18.245	2557	2560	2568	rVB10	16641	31021	0.45%	0.050%
83	18.363	2576	2580	2590	rBV	304884	455305	6.57%	0.728%
84	18.457	2592	2596	2602	rVB	54539	90376	1.30%	0.144%
85	18.616	2619	2623	2628	rBV2	97339	139457	2.01%	0.223%
86	18.686	2630	2635	2642	rVB	447494	565433	8.16%	0.904%
87	19.404	2754	2757	2760	rBV2	42512	49825	0.72%	0.080%
88	19.439	2760	2763	2766	rVV	50732	60275	0.87%	0.096%
89	19.474	2766	2769	2772	rVB	43971	52383	0.76%	0.084%
90	19.586	2784	2788	2801	rBV3	145977	254892	3.68%	0.407%
91	19.827	2823	2829	2837	rBV	3740037	5019668	72.42%	8.024%
92	20.216	2892	2895	2903	rVB5	62971	80534	1.16%	0.129%
93	21.004	3026	3029	3034	rVB	192057	217279	3.13%	0.347%
94	21.263	3069	3073	3084	rBV2	638310	1032159	14.89%	1.650%
95	21.586	3123	3128	3137	rBV	3020044	4049227	58.42%	6.473%
96	22.168	3223	3227	3234	rBV2	173771	252900	3.65%	0.404%
97	22.886	3346	3349	3356	rVB	287690	391391	5.65%	0.626%
98	23.986	3529	3536	3549	rVB2	2111882	4633624	66.85%	7.407%
99	24.151	3557	3564	3574	rVB2	1728997	3777011	54.49%	6.037%

Sum of corrected areas: 62560429

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Acq On : 08 Feb 2023 12:56
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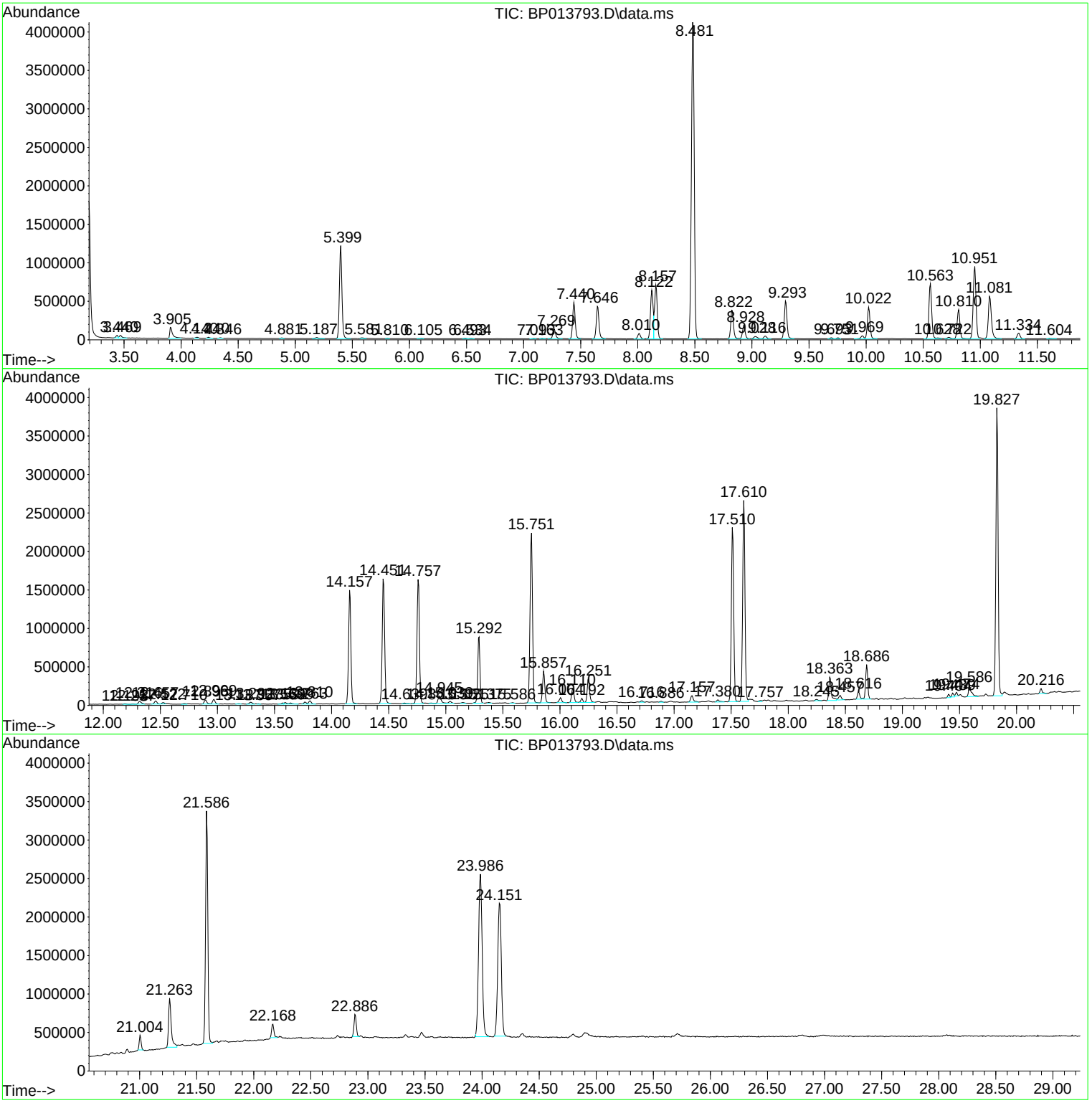
Sample : 01412-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44M4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020723.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020823\
Data File : BP013793.D
Acq On : 08 Feb 2023 12:56
Operator : CG/JU
Sample : 01412-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44M4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020723.M
Quant Title : SVOA CALIBRATION

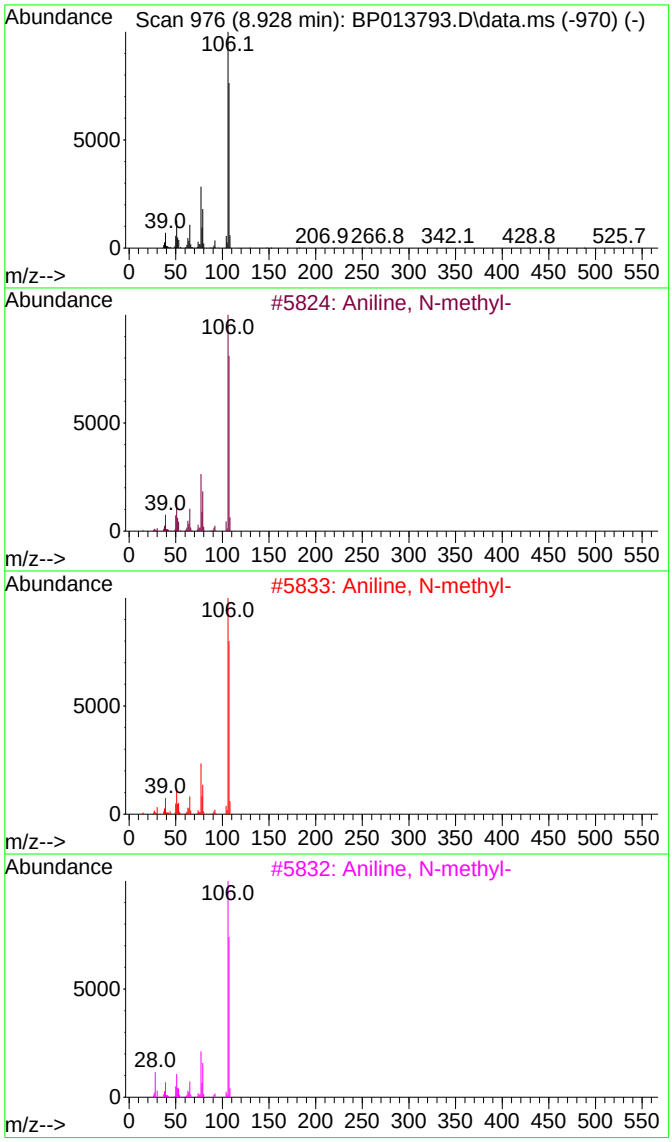
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

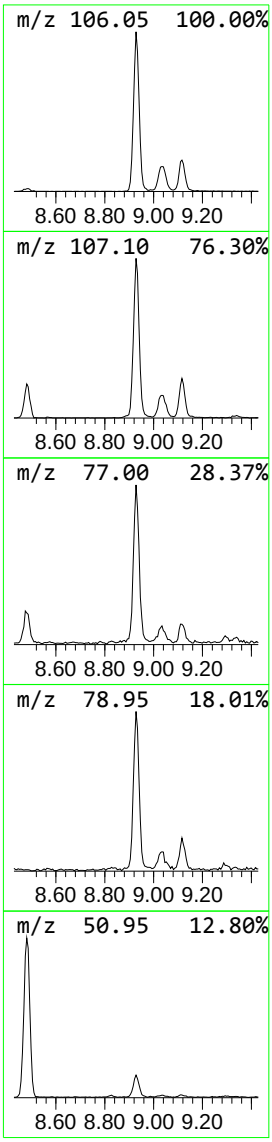
****

Peak Number 4 Aniline, N-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.928	5.56 ng/ul	290581	1,4-Dichlorobenzene-d4	8.122

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Aniline, N-methyl-	107	C7H9N	000100-61-8	96
2	Aniline, N-methyl-	107	C7H9N	000100-61-8	96
3	Aniline, N-methyl-	107	C7H9N	000100-61-8	95
4	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	91
5	Aniline, N-methyl-	107	C7H9N	000100-61-8	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020823\
Data File : BP013793.D
Acq On : 08 Feb 2023 12:56
Operator : CG/JU
Sample : 01412-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44M4

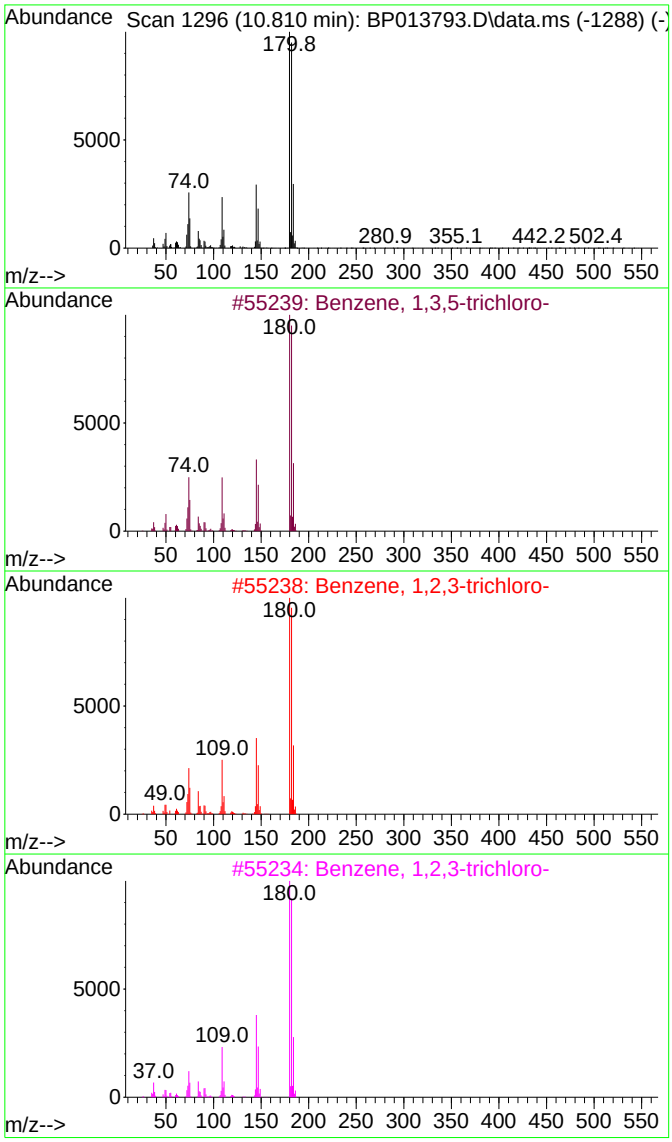
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020723.M
Quant Title : SVOA CALIBRATION

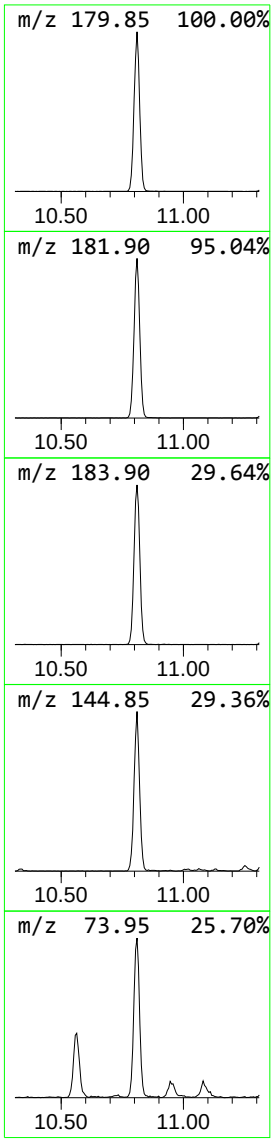
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1,3,5-trichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.810	7.79 ng/ul	624957	Naphthalene-d8	10.951

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
2	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
3	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
4	Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	97
5	Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	97





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ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44M4

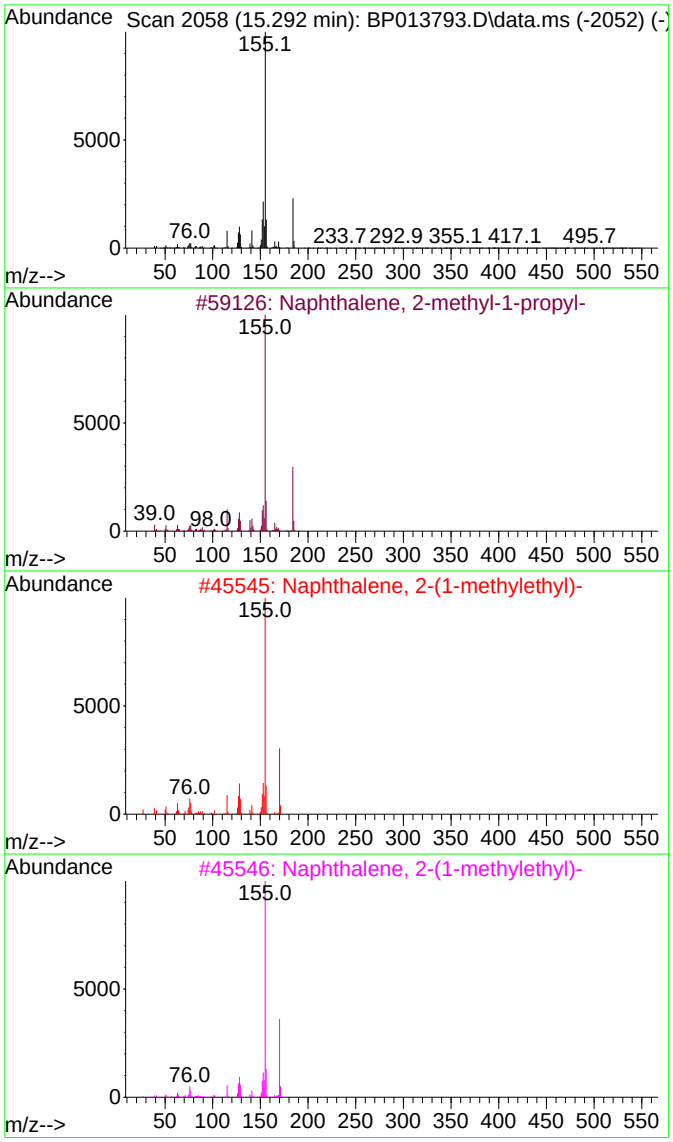
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020723.M
Quant Title : SVOA CALIBRATION

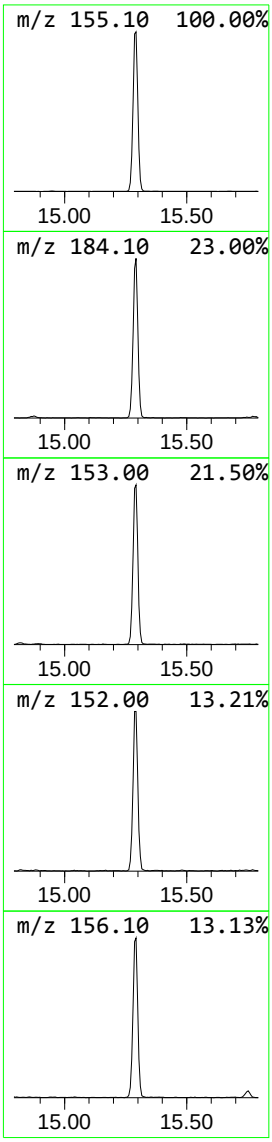
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Naphthalene, 2-methyl-1-pro... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.292	9.74 ng/ul	1184940	Acenaphthene-d10	14.757

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-1-propyl-	184	C14H16	054774-89-9	87
2	Naphthalene, 2-(1-methylethyl)-	170	C13H14	002027-17-0	64
3	Naphthalene, 2-(1-methylethyl)-	170	C13H14	002027-17-0	50
4	Pyridine, 3-phenyl-	155	C11H9N	001008-88-4	43
5	Naphthalene, 2-(1-methylethyl)-	170	C13H14	002027-17-0	43





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BNA_P
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A44M4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020723.M
Quant Title : SVOA CALIBRATION

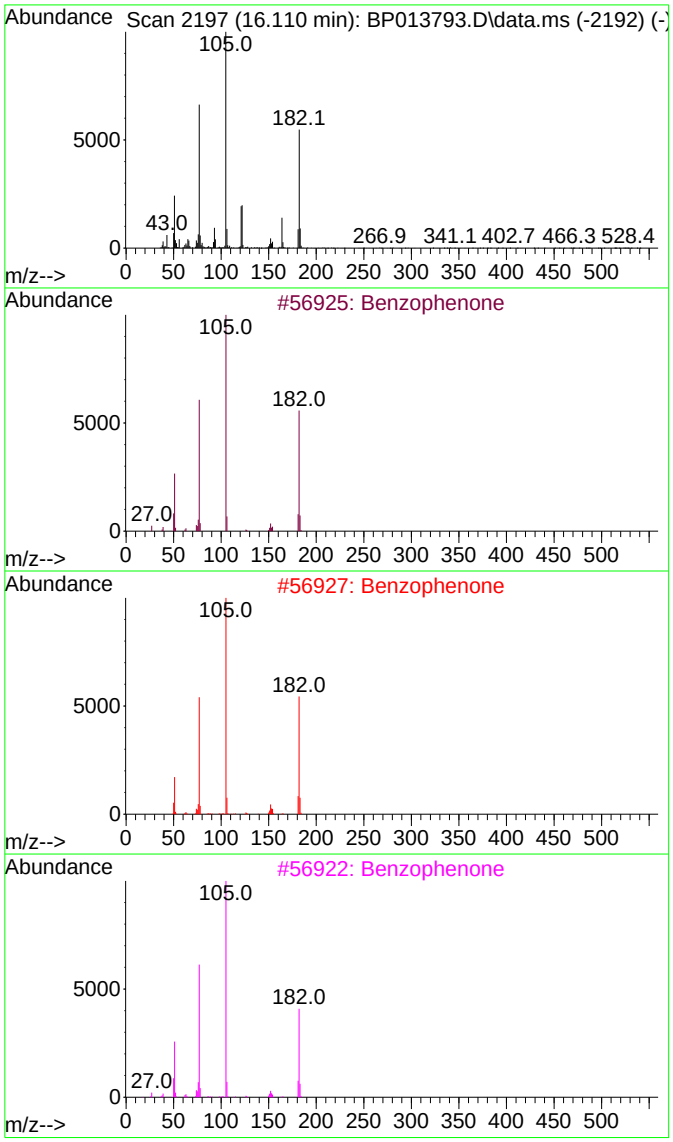
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

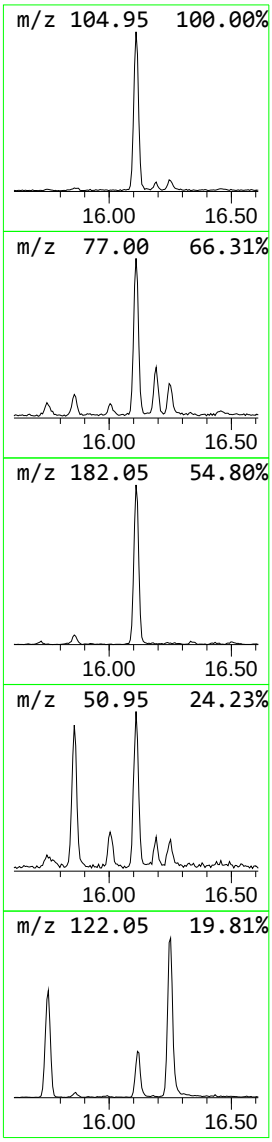
****

Peak Number 7 Benzophenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.110	2.35 ng/ul	285333	Acenaphthene-d10	14.757

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	90
4	Benzophenone	182	C13H10O	000119-61-9	81
5	Azobenzene	182	C12H10N2	000103-33-3	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020823\
Data File : BP013793.D
Acq On : 08 Feb 2023 12:56
Operator : CG/JU
Sample : 01412-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44M4

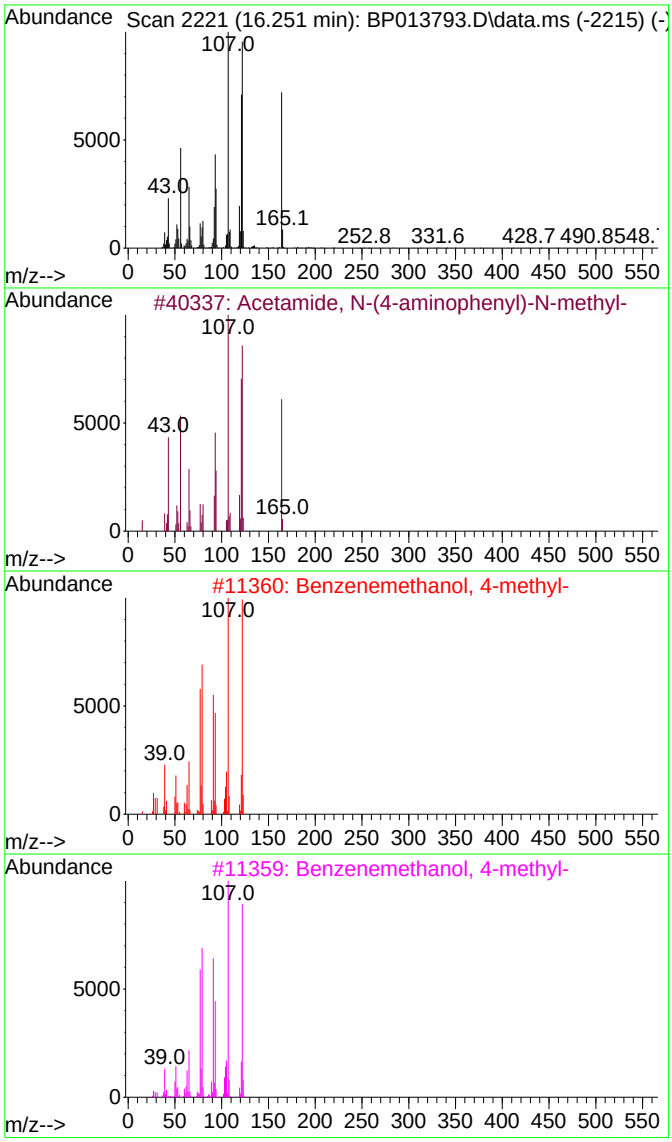
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020723.M
Quant Title : SVOA CALIBRATION

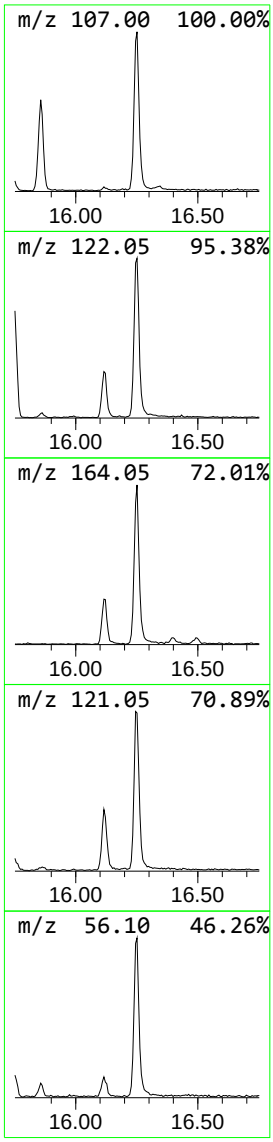
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Acetamide, N-(4-aminophenyl... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.251	2.81 ng/ul	453746	Phenanthrene-d10	17.510

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(4-aminophenyl)-N-m...	164	C9H12N2O	000119-63-1	98
2	Benzenemethanol, 4-methyl-	122	C8H10O	000589-18-4	52
3	Benzenemethanol, 4-methyl-	122	C8H10O	000589-18-4	52
4	2-Ethylphenol, O-acetyl-	164	C10H12O2	1000374-84-9	46
5	2-But-2-enyl-1H-imidazole	122	C7H10N2	1000194-18-8	46





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Data File : BP013793.D
Acq On : 08 Feb 2023 12:56
Operator : CG/JU
Sample : 01412-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44M4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020723.M
Quant Title : SVOA CALIBRATION

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

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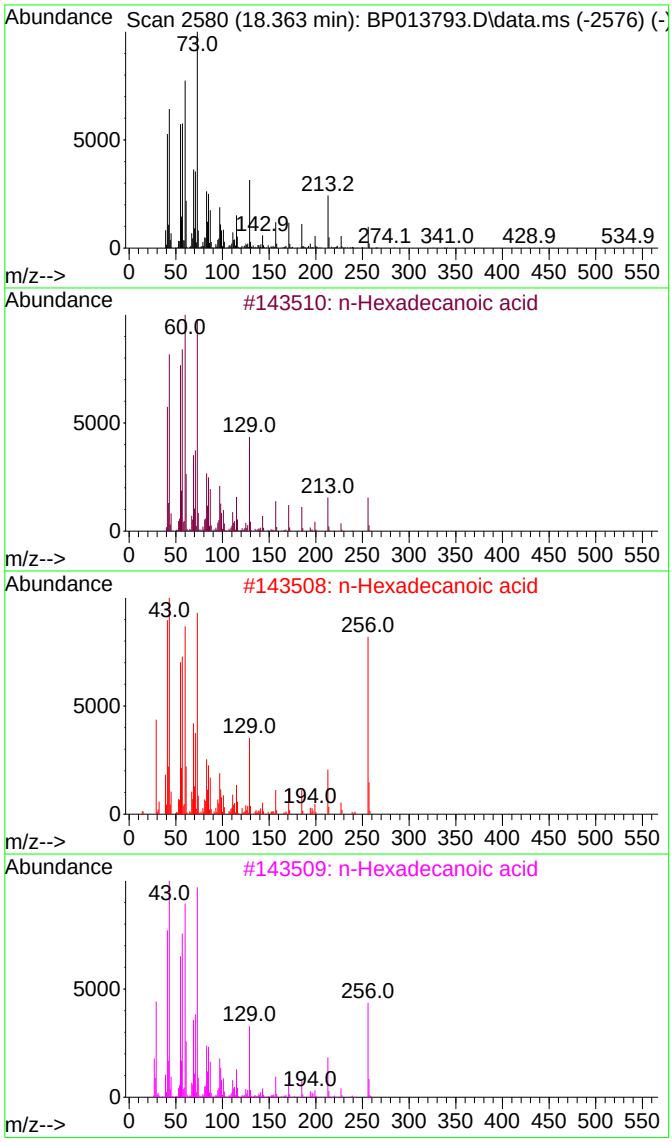
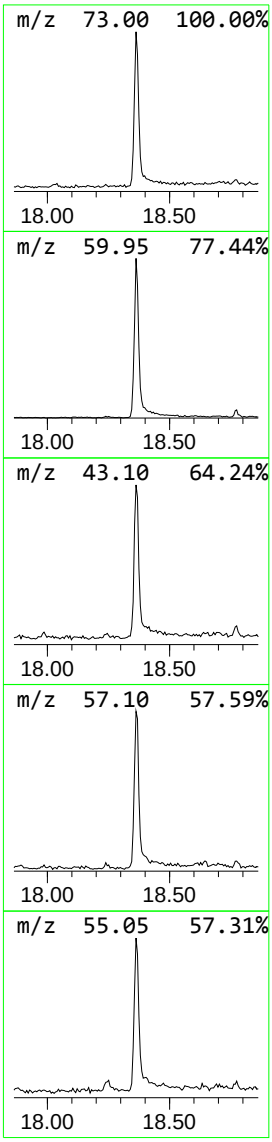
Peak Number 9 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.363	2.82 ng/ul	455305	Phenanthrene-d10	17.510

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	98
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4	Tridecanoic acid	214	C13H26O2	000638-53-9	97
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96





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Acq On : 08 Feb 2023 12:56
Operator : CG/JU
Sample : 01412-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44M4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020723.M
Quant Title : SVOA CALIBRATION

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

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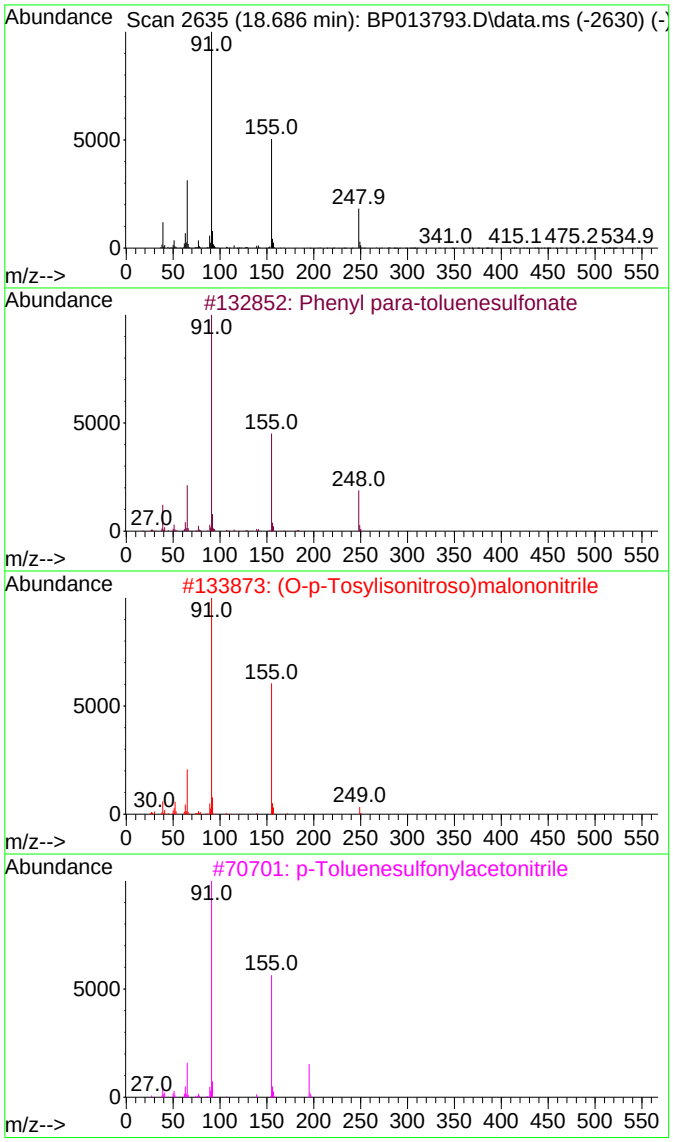
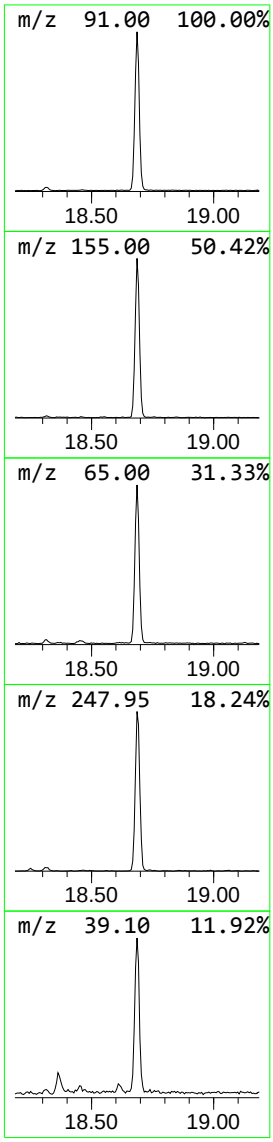
Peak Number 10 Phenyl para-toluenesulfonate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.686	3.50 ng/ul	565433	Phenanthrene-d10	17.510

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenyl para-toluenesulfonate	248	C13H12O3S	000640-60-8	96
2	(O-p-Tosylisonitroso)malononitrile	249	C10H7N3O3S	020893-01-0	76
3	p-Toluenesulfonylacetonitrile	195	C9H9NO2S	005697-44-9	64
4	p-Toluenesulfonylacetonitrile	195	C9H9NO2S	005697-44-9	59
5	Benzene, 1-[(chloromethyl)sulfon...	204	C8H9ClO2S	007569-26-8	56





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Data File : BP013793.D
Acq On : 08 Feb 2023 12:56
Operator : CG/JU
Sample : 01412-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44M4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020723.M
Quant Title : SVOA CALIBRATION

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

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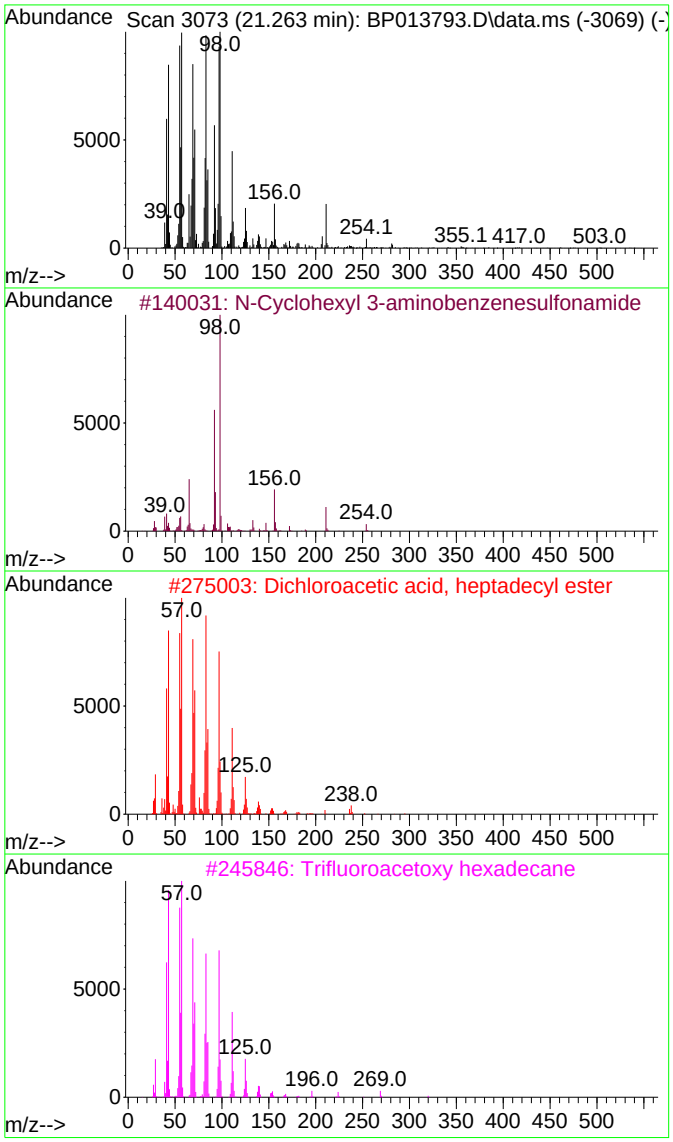
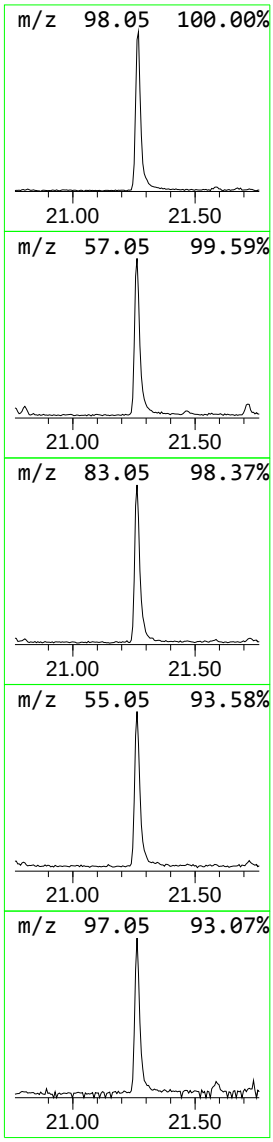
Peak Number 11 N-Cyclohexyl 3-aminobenzene... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.263	5.10 ng/ul	1032160	Chrysene-d12	21.586

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	N-Cyclohexyl 3-aminobenzenesulfo...	254	C12H18N2O2S	061886-26-8	91
2	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	86
3	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	83
4	1-Heneicosanol	312	C21H44O	015594-90-8	83
5	1-Heneicosyl formate	340	C22H44O2	077899-03-7	83





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Data File : BP013793.D
Acq On : 08 Feb 2023 12:56
Operator : CG/JU
Sample : 01412-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44M4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020723.M
Quant Title : SVOA CALIBRATION

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

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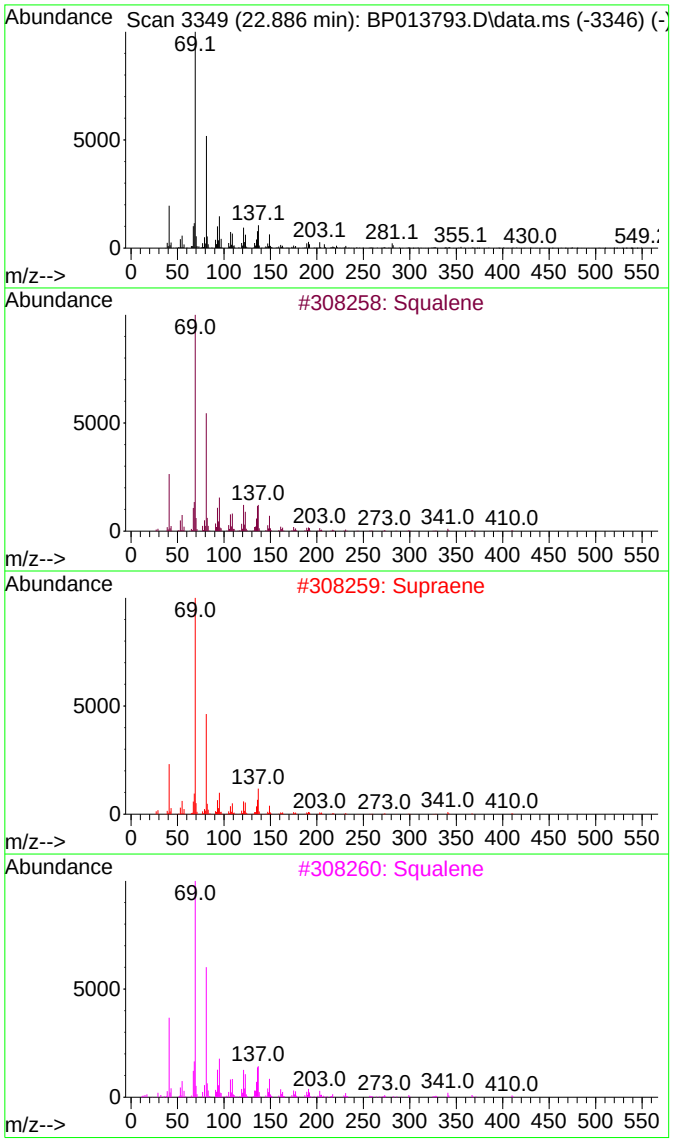
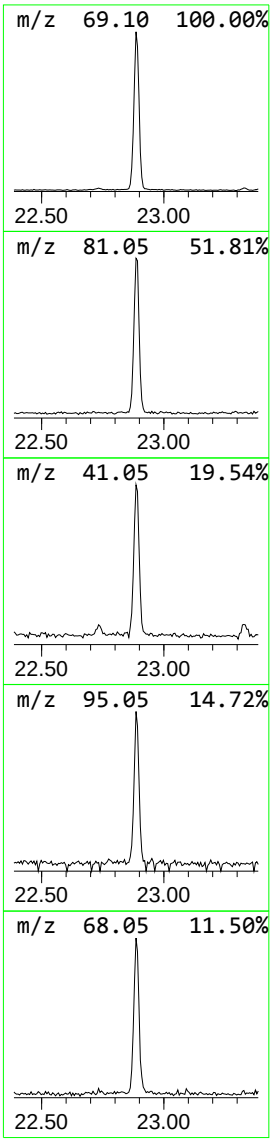
Peak Number 12 Squalene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.886	2.07 ng/ul	391391	Perylene-d12	24.151

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Squalene	410	C30H50	000111-02-4	90
2	Supraene	410	C30H50	007683-64-9	90
3	Squalene	410	C30H50	000111-02-4	86
4	2,6,10,14-Hexadecatetraenoic aci...	332	C22H36O2	024035-35-6	83
5	(2E,6E,10E)-3,7,11,15-Tetramethy...	318	C21H34O2	125456-63-5	72





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020823\
Data File : BP013793.D
Acq On : 08 Feb 2023 12:56
Operator : CG/JU
Sample : 01412-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44M4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020723.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
Aniline, N-methyl-	8.928	5.6	ng/ul	290581	1	8.122	1045070	20.0
Benzene, 1,3,5-...	10.810	7.8	ng/ul	624957	2	10.951	1604180	20.0
Naphthalene, 2-...	15.292	9.7	ng/ul	1184940	3	14.757	2433130	20.0
Benzophenone	16.110	2.4	ng/ul	285333	3	14.757	2433130	20.0
Acetamide, N-(4...	16.251	2.8	ng/ul	453746	4	17.510	3230230	20.0
n-Hexadecanoic ...	18.363	2.8	ng/ul	455305	4	17.510	3230230	20.0
Phenyl para-tol...	18.686	3.5	ng/ul	565433	4	17.510	3230230	20.0
N-Cyclohexyl 3-...	21.263	5.1	ng/ul	1032160	5	21.586	4049230	20.0
Squalene	22.886	2.1	ng/ul	391391	6	24.151	3777010	20.0