LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071822\
Data File : BP011002.D
Acq On : 18 Jul 2022 20:27
Operator : CG/JU
Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0B09

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-BP071422.M
Title : SVOA CALIBRATION

Signal : TIC: BP011002.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.181	12	16	20	rBV2	94085	127093	1.80%	0.115%
2	3.452	58	62	67	rBV2	82339	116585	1.65%	0.105%
3	3.593	82	86	97	rBV	740823	1099738	15.61%	0.993%
4	4.163	178	183	188	rBV	96488	125946	1.79%	0.114%
5	4.340	207	213	224	rVB	2346893	3333917	47.32%	3.009%
6	4.704	269	275	280	rBV	89836	151498	2.15%	0.137%
7	5.010	322	327	337	rVB	520325	758592	10.77%	0.685%
8	5.157	342	352	356	rBV2	68969	128503	1.82%	0.116%
9	6.081	501	509	516	rBV2	57369	94606	1.34%	0.085%
10	6.722	610	618	625	rVV	298505	656465	9.32%	0.593%
11	6.857	633	641	654	rVV	353249	660241	9.37%	0.596%
12	7.022	662	669	675	rVV	996673	1514475	21.50%	1.367%
13	7.216	696	702	710	rVB	931092	1447382	20.54%	1.307%
14	7.328	717	721	727	rVB	62876	109658	1.56%	0.099%
15	7.622	761	771	773	rBV2	92745	150369	2.13%	0.136%
16	7.681	773	781	786	rBV	1281717	2138801	30.36%	1.931%
17	7.757	791	794	797	rVV2	85162	153591	2.18%	0.139%
18	7.792	797	800	807	rVB4	98110	188448	2.67%	0.170%
19	8.281	876	883	887	rBV	199003	350309	4.97%	0.316%
20	8.392	896	902	908	rBV	866370	1332547	18.91%	1.203%
21	8.451	908	912	922	rVV	152463	277429	3.94%	0.250%
22	8.592	929	936	947	rVB4	51099	144997	2.06%	0.131%
23	8.781	965	968	972	rVV2	85191	126694	1.80%	0.114%
24	8.839	972	978	987	rVV	1028306	1745129	24.77%	1.575%
25	8.928	987	993	997	rVV5	41956	95392	1.35%	0.086%
26	9.045	1009	1013	1020	rVV4	119143	221605	3.15%	0.200%
27	9.198	1034	1039	1043	rBV2	95687	165995	2.36%	0.150%
28	9.298	1051	1056	1064	rVB2	88541	172208	2.44%	0.155%
29	9.557	1094	1100	1114	rVB	670873	1135893	16.12%	1.025%
30	9.722	1122	1128	1139	rBV	769966	1299489	18.45%	1.173%
31	9.845	1139	1149	1154	rBV	772249	1522588	21.61%	1.374%
32	10.022	1173	1179	1185	rVB6	112663	250500	3.56%	0.226%
33	10.098	1185	1192	1200	rBV	1266584	2079629	29.52%	1.877%
34	10.263	1214	1220	1226	rBV8	49305	138964	1.97%	0.125%
35	10.375	1232	1239	1243	rBV4	46767	112747	1.60%	0.102%
36	10.428	1244	1248	1250	rVV2	67208	102591	1.46%	0.093%

LSC Area Percent Report

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Acq On : 18 Jul 2022 20:27
Operator : CG/JU
Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0B09

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-BP071422.M
Title : SVOA CALIBRATION

37	10.475	1250	1256	1261	rVV	1795722	2976606	42.25%	2.687%
38	10.522	1261	1264	1272	rVB	149653	247931	3.52%	0.224%
39	10.622	1272	1281	1295	rBV	2471984	4333070	61.50%	3.911%
40	10.845	1314	1319	1322	rBV2	57393	99438	1.41%	0.090%
41	10.881	1323	1325	1333	rVB8	59177	126823	1.80%	0.114%
42	11.128	1363	1367	1374	rVB	53772	98362	1.40%	0.089%
43	11.292	1383	1395	1405	rBV3	187022	511104	7.25%	0.461%
44	11.463	1418	1424	1429	rBV6	60085	127144	1.80%	0.115%
45	11.833	1478	1487	1493	rBV	280038	478651	6.79%	0.432%
46	11.992	1509	1514	1519	rBV	69325	125538	1.78%	0.113%
47	12.098	1528	1532	1537	rVV3	77545	141761	2.01%	0.128%
48	12.363	1573	1577	1583	rVB4	68996	116126	1.65%	0.105%
49	12.469	1588	1595	1599	rBV2	79589	146701	2.08%	0.132%
50	12.551	1605	1609	1614	rVB	63406	101835	1.45%	0.092%
51	12.622	1615	1621	1626	rBV3	62932	125500	1.78%	0.113%
52	12.928	1666	1673	1682	rVB2	322688	626810	8.90%	0.566%
53	13.133	1703	1708	1713	rBV4	163231	276607	3.93%	0.250%
54	13.392	1749	1752	1757	rVB4	63250	93694	1.33%	0.085%
55	13.686	1797	1802	1805	rBV2	72925	122641	1.74%	0.111%
56	13.763	1806	1815	1821	rVB	2617417	3806235	54.03%	3.436%
57	14.033	1853	1861	1867	rVB	2664619	3830156	54.37%	3.457%
58	14.239	1892	1896	1899	rBV	216332	315535	4.48%	0.285%
59	14.269	1899	1901	1906	rVB2	141053	180002	2.55%	0.162%
60	14.339	1906	1913	1918	rBV	2481406	3673081	52.14%	3.316%
61	14.563	1941	1951	1957	rBV4	327186	626347	8.89%	0.565%
62	14.692	1971	1973	1978	rVB	118302	134051	1.90%	0.121%
63	15.104	2037	2043	2051	rBV2	806996	1441042	20.45%	1.301%
64	15.216	2059	2062	2064	rBV4	87072	101996	1.45%	0.092%
65	15.257	2065	2069	2074	rVV3	230662	401189	5.69%	0.362%
66	15.339	2078	2083	2094	rVB	3714571	5327056	75.61%	4.809%
67	15.427	2094	2098	2100	rBV3	126613	164133	2.33%	0.148%
68	15.463	2100	2104	2109	rVB	955566	1238400	17.58%	1.118%
69	15.592	2121	2126	2132	rBV	661360	1128119	16.01%	1.018%
70	15.639	2132	2134	2139	rVB	162304	173068	2.46%	0.156%
71	15.727	2146	2149	2155	rVB3	81184	127354	1.81%	0.115%
72	15.869	2168	2173	2178	rBV	1449714	1858849	26.38%	1.678%
73	15.945	2183	2186	2191	rVB2	101746	145256	2.06%	0.131%
74	16.051	2198	2204	2216	rBV	2220503	3726696	52.90%	3.364%
75	16.386	2255	2261	2274	rBV	1008055	1577220	22.39%	1.424%
76	16.792	2326	2330	2334	rVB	116151	152646	2.17%	0.138%

LSC Area Percent Report

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Acq On : 18 Jul 2022 20:27
Operator : CG/JU
Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0B09

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-BP071422.M
Title : SVOA CALIBRATION

77	16.921	2348	2352	2357	rVB	508631	622822	8.84%	0.562%
78	16.974	2357	2361	2368	rBV	299468	443637	6.30%	0.400%
79	17.110	2378	2384	2389	rBV	2730916	3986232	56.58%	3.598%
80	17.204	2394	2400	2407	rBV2	4074409	5800456	82.33%	5.236%
81	17.304	2412	2417	2427	rVB	1363204	2038322	28.93%	1.840%
82	17.416	2432	2436	2442	rVB2	386882	539107	7.65%	0.487%
83	17.521	2451	2454	2461	rVB	82124	120234	1.71%	0.109%
84	18.021	2535	2539	2547	rBV	552182	851003	12.08%	0.768%
85	18.463	2610	2614	2618	rBV	198945	265673	3.77%	0.240%
86	18.568	2629	2632	2639	rVB3	137067	175452	2.49%	0.158%
87	18.663	2645	2648	2651	rBV	115893	141403	2.01%	0.128%
88	18.768	2662	2666	2674	rBV3	262922	384432	5.46%	0.347%
89	19.204	2736	2740	2746	rVB	1498339	1742728	24.74%	1.573%
90	19.268	2747	2751	2757	rBV3	158567	241785	3.43%	0.218%
91	19.462	2779	2784	2791	rBV	5352930	6865303	97.45%	6.197%
92	19.521	2791	2794	2799	rVB	451589	531834	7.55%	0.480%
93	19.904	2855	2859	2864	rVB	306476	364622	5.18%	0.329%
94	20.486	2955	2958	2968	rBV2	972430	1116002	15.84%	1.007%
95	20.957	3034	3038	3047	rVB	1547472	2014026	28.59%	1.818%
96	21.156	3069	3072	3075	rBV	214207	254248	3.61%	0.230%
97	21.227	3079	3084	3089	rBV	3676825	4526032	64.24%	4.086%
98	22.333	3269	3272	3280	rVB	443951	685949	9.74%	0.619%
99	23.427	3451	3458	3466	rBV2	3693783	7045134	100.00%	6.359%
100	23.580	3478	3484	3493	rVB2	2454036	4790405	68.00%	4.324%

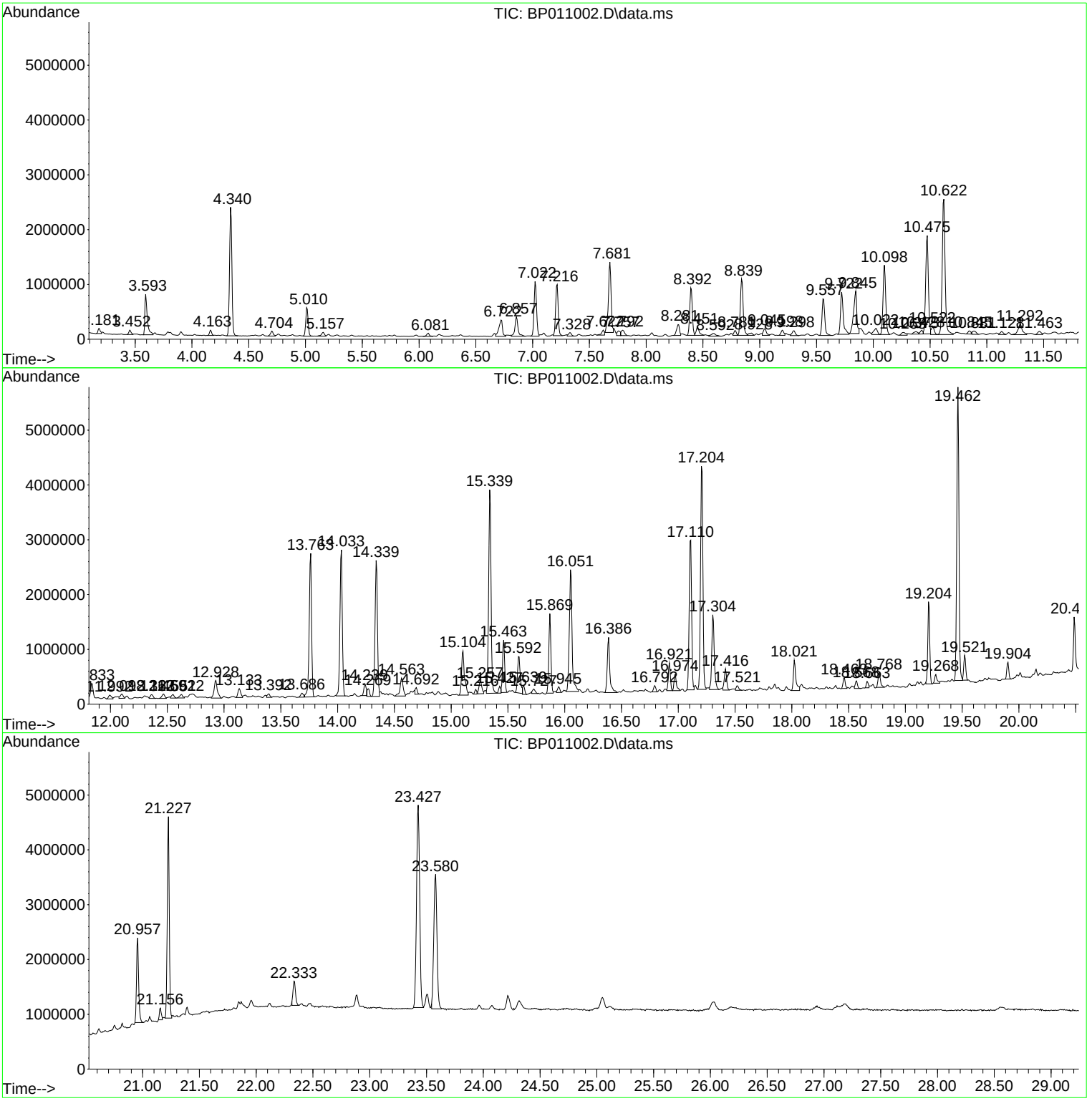
Sum of corrected areas: 110782228

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071822\
Data File : BP011002.D
Acq On : 18 Jul 2022 20:27
Operator : CG/JU
Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0B09

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071822\
Data File : BP011002.D
Acq On : 18 Jul 2022 20:27
Operator : CG/JU
Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0B09

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

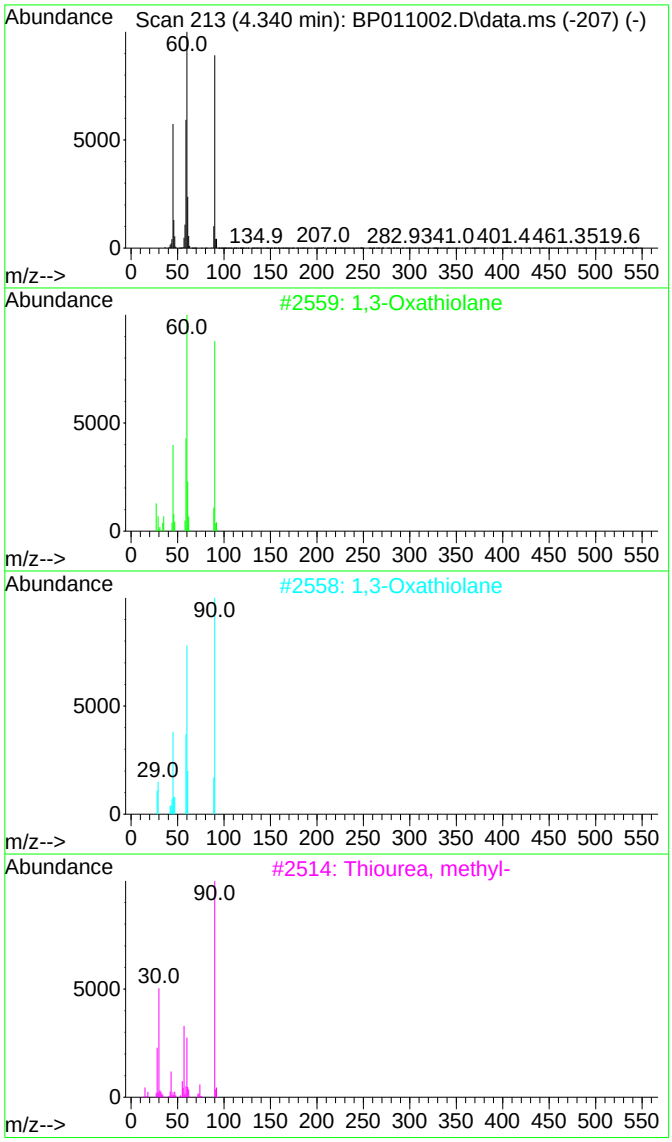
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

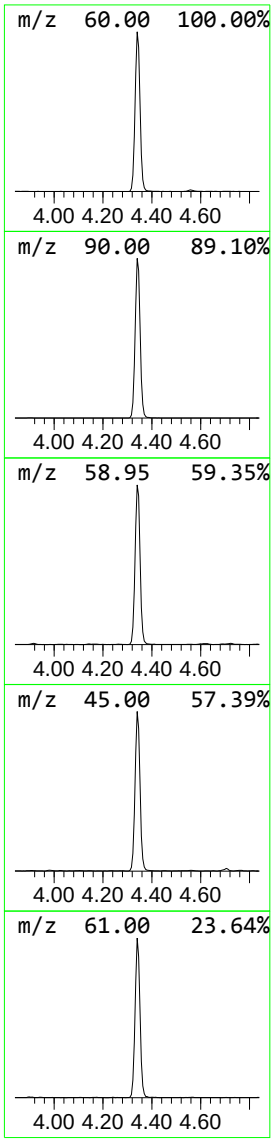
****

Peak Number 1 1,3-Oxathiolane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.340	31.18 ng/ul	3333920	1,4-Dichlorobenzene-d4	7.681

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Oxathiolane	90	C3H6OS	002094-97-5	90
2	1,3-Oxathiolane	90	C3H6OS	002094-97-5	56
3	Thiourea, methyl-	90	C2H6N2S	000598-52-7	50
4	Hydrazinecarboxylic acid, methyl...	90	C2H6N2O2	006294-89-9	45
5	Heptanoic acid, 2-hydroxy-, meth...	160	C8H16O3	054340-91-9	36





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071822\
Data File : BP011002.D
Acq On : 18 Jul 2022 20:27
Operator : CG/JU
Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0B09

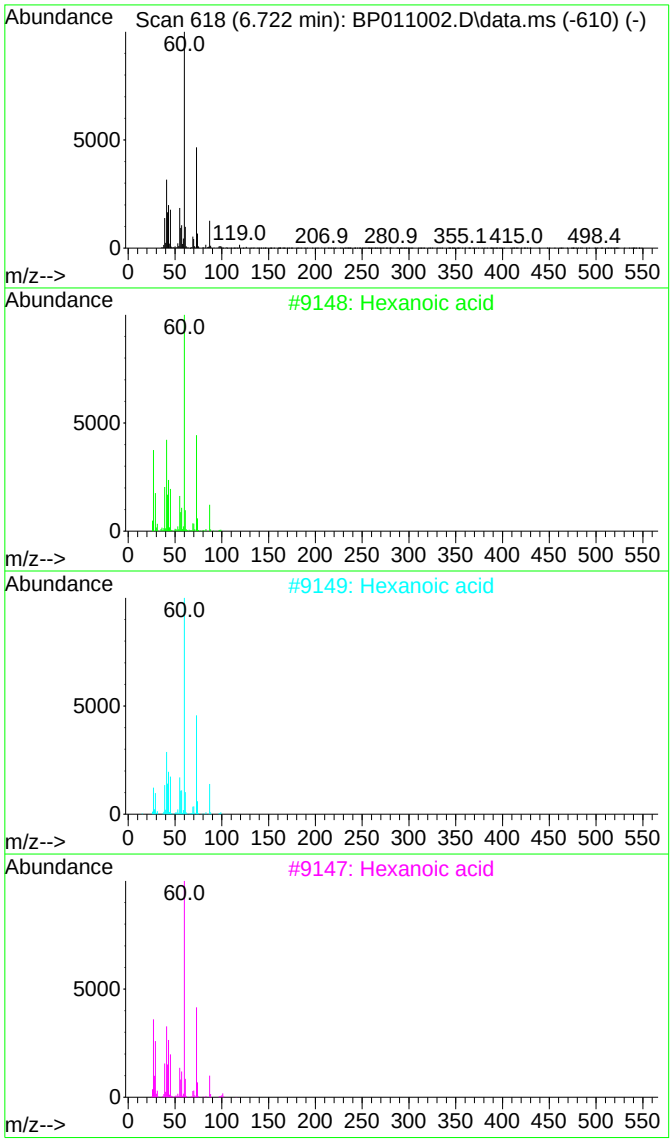
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

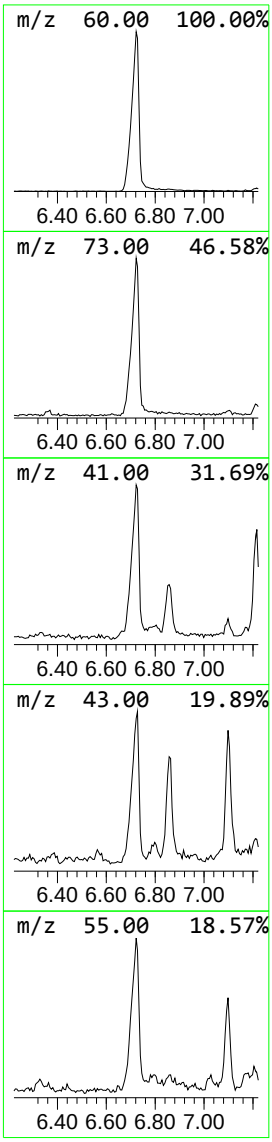
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Hexanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.722	6.14 ng/ul	656465	1,4-Dichlorobenzene-d4	7.681

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid	116	C6H12O2	000142-62-1	90
2	Hexanoic acid	116	C6H12O2	000142-62-1	90
3	Hexanoic acid	116	C6H12O2	000142-62-1	83
4	Hexanoic acid	116	C6H12O2	000142-62-1	78
5	Heptanoic acid	130	C7H14O2	000111-14-8	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071822\
Data File : BP011002.D
Acq On : 18 Jul 2022 20:27
Operator : CG/JU
Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0B09

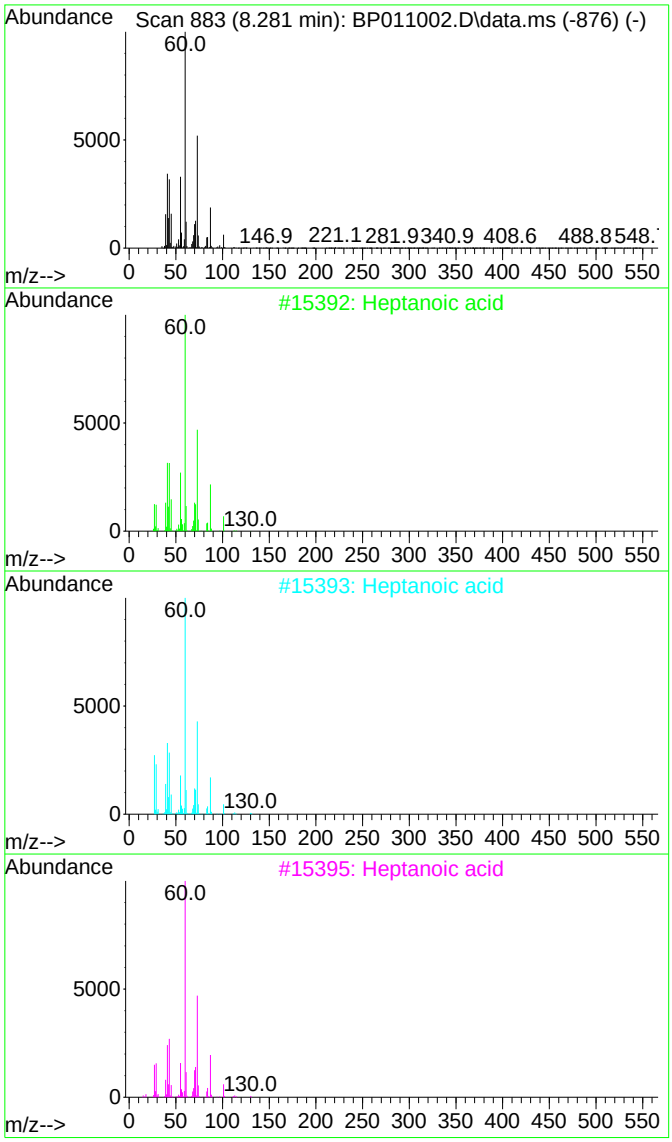
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

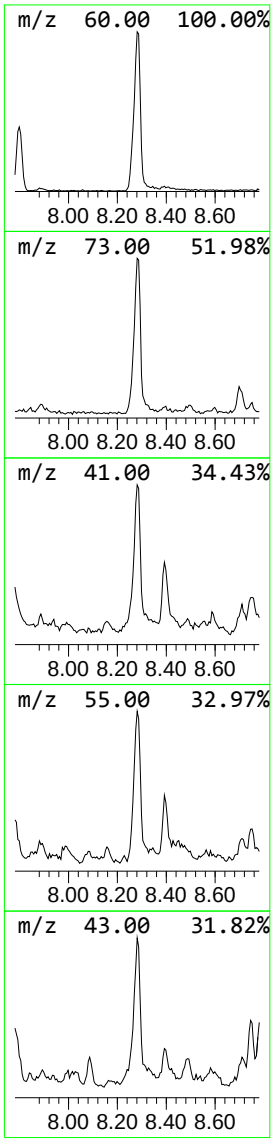
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Heptanoic acid Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.281	3.28 ng/ul	350309	1,4-Dichlorobenzene-d4	7.681

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptanoic acid	130	C7H14O2	000111-14-8	91
2	Heptanoic acid	130	C7H14O2	000111-14-8	90
3	Heptanoic acid	130	C7H14O2	000111-14-8	90
4	Heptanoic acid	130	C7H14O2	000111-14-8	78
5	Hexanoic acid	116	C6H12O2	000142-62-1	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071822\
Data File : BP011002.D
Acq On : 18 Jul 2022 20:27
Operator : CG/JU
Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0B09

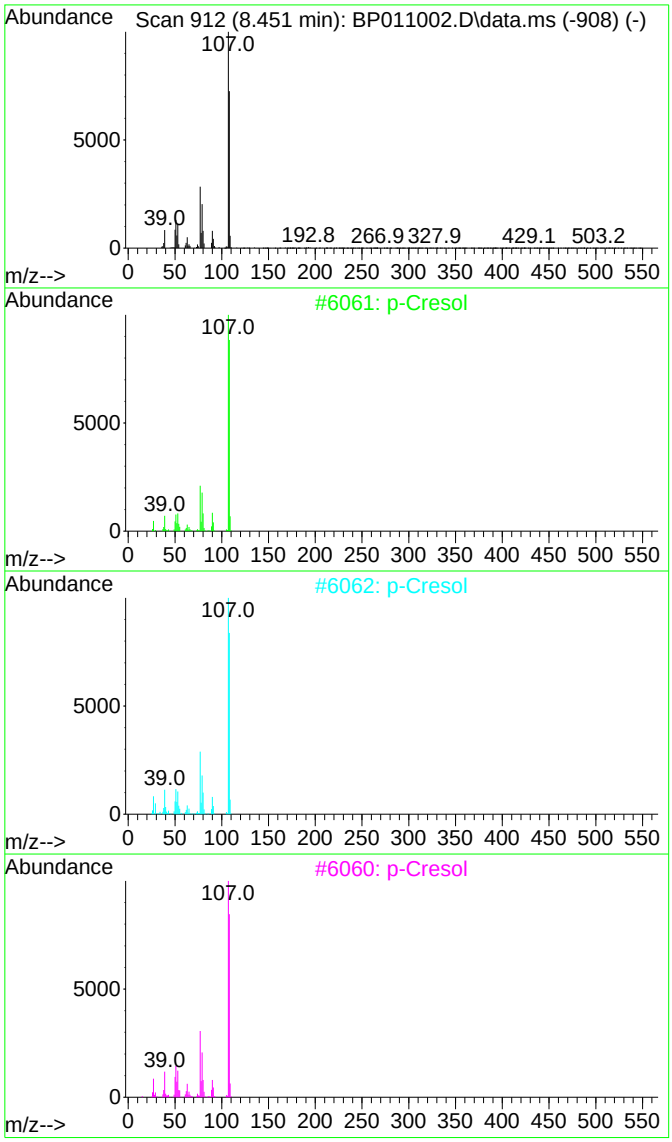
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

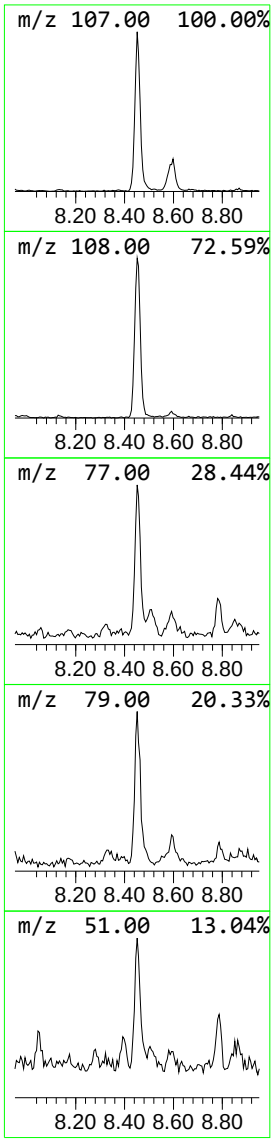
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 p-Cresol Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.451	2.59 ng/ul	277429	1,4-Dichlorobenzene-d4	7.681

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cresol	108	C7H8O	000106-44-5	97
2	p-Cresol	108	C7H8O	000106-44-5	94
3	p-Cresol	108	C7H8O	000106-44-5	94
4	p-Cresol	108	C7H8O	000106-44-5	81
5	Phenol, 3-methyl-	108	C7H8O	000108-39-4	74





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071822\
Data File : BP011002.D
Acq On : 18 Jul 2022 20:27
Operator : CG/JU
Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0B09

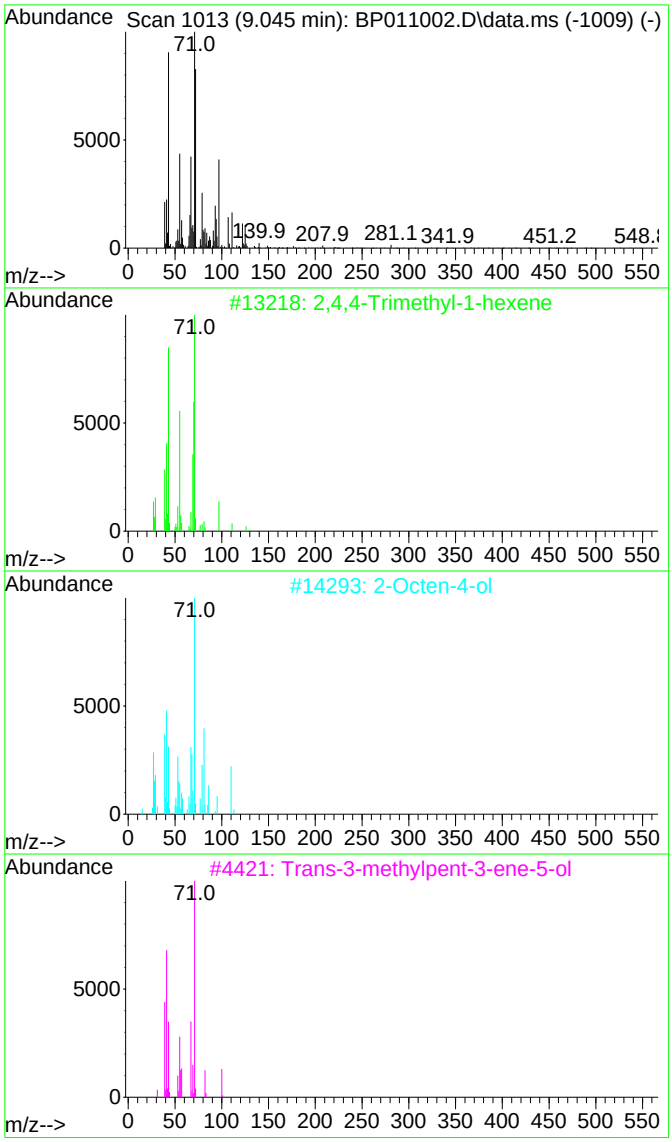
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

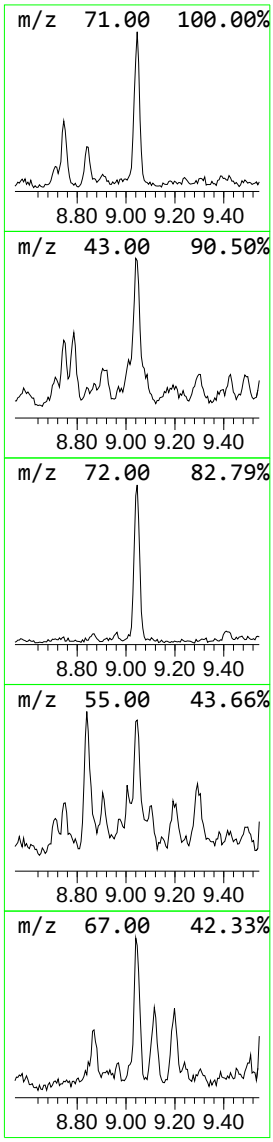
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.045	2.07 ng/ul	221605	1,4-Dichlorobenzene-d4	7.681

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,4-Trimethyl-1-hexene	126	C9H18	051174-12-0	25
2	2-Octen-4-ol	128	C8H16O	004798-61-2	16
3	Trans-3-methylpent-3-ene-5-ol	100	C6H12O	030801-95-7	14
4	2-Propenamide	71	C3H5NO	000079-06-1	11
5	Propafenone	341	C21H27NO3	054063-53-5	10





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Acq On : 18 Jul 2022 20:27
Operator : CG/JU
Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0B09

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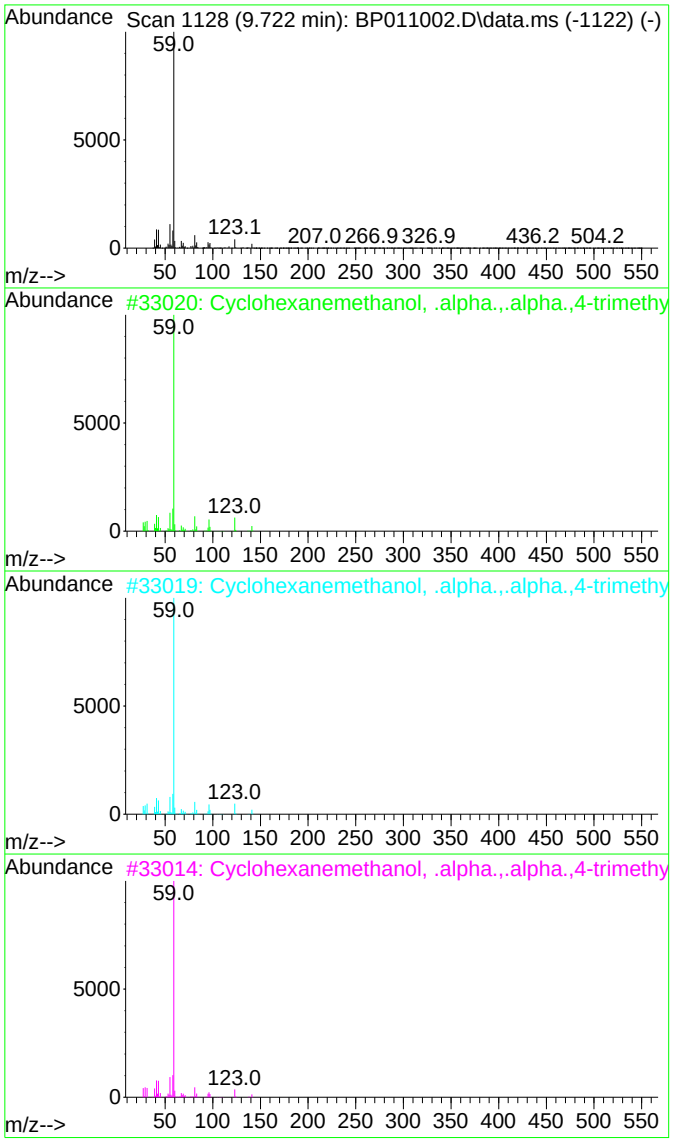
Quant Title : SVOA CALIBRATION

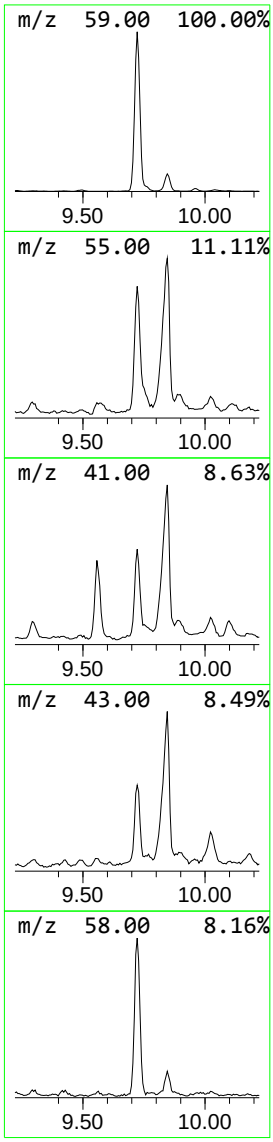
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Cyclohexanemethanol, .alpha... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.722	8.73 ng/ul	1299490	Naphthalene-d8	10.475

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	005114-00-1	78
2	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	007322-63-6	74
3	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	64
4	o-Menthan-8-ol	156	C10H20O	343855-44-7	56
5	2-Hydroxy-2,5-dimethyl-hept-6-en...	156	C9H16O2	1000192-55-8	45





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ClientSampleId :
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Quant Title : SVOA CALIBRATION

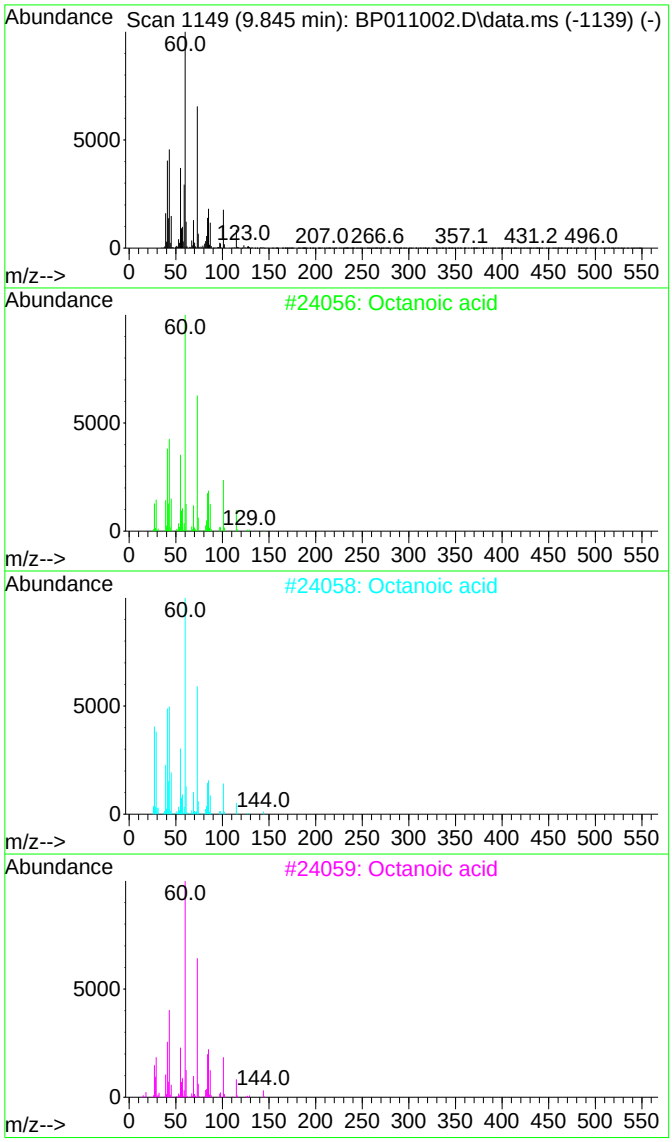
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

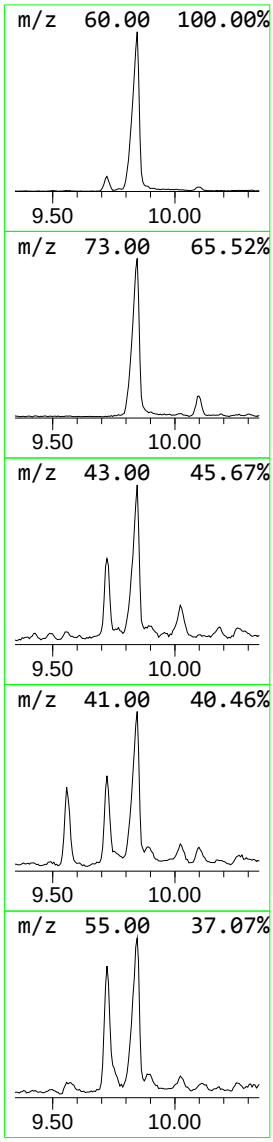
****

Peak Number 8 Octanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.845	10.23 ng/ul	1522590	Naphthalene-d8	10.475

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octanoic acid	144	C8H16O2	000124-07-2	96
2	Octanoic acid	144	C8H16O2	000124-07-2	80
3	Octanoic acid	144	C8H16O2	000124-07-2	72
4	Octanoic acid	144	C8H16O2	000124-07-2	64
5	Hexanoic acid	116	C6H12O2	000142-62-1	58





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Sample : N3710-17
Misc :
ALS Vial : 17 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

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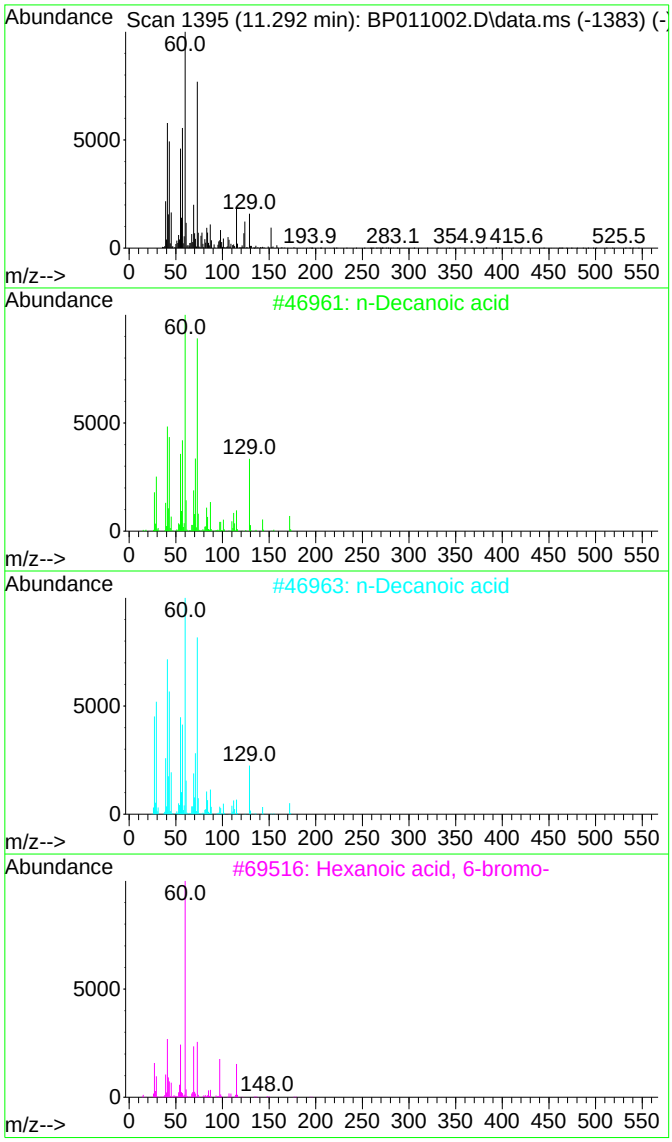
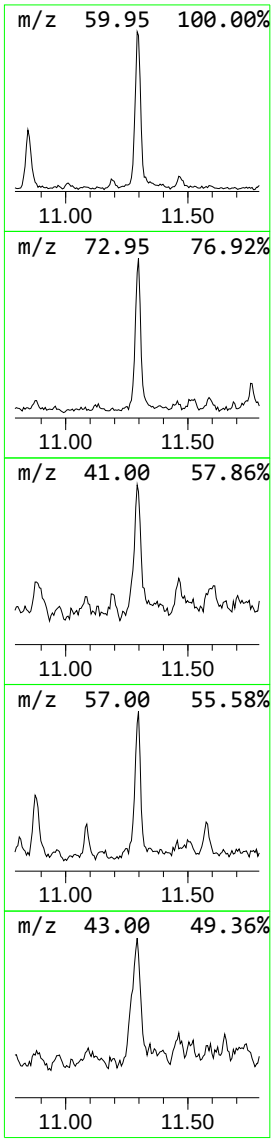
Peak Number 9 n-Decanoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.292	3.43 ng/ul	511104	Naphthalene-d8	10.475

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Decanoic acid	172	C10H20O2	000334-48-5	64
2	n-Decanoic acid	172	C10H20O2	000334-48-5	59
3	Hexanoic acid, 6-bromo-	194	C6H11BrO2	004224-70-8	53
4	Hexanoic acid	116	C6H12O2	000142-62-1	53
5	Hexanoic acid	116	C6H12O2	000142-62-1	53





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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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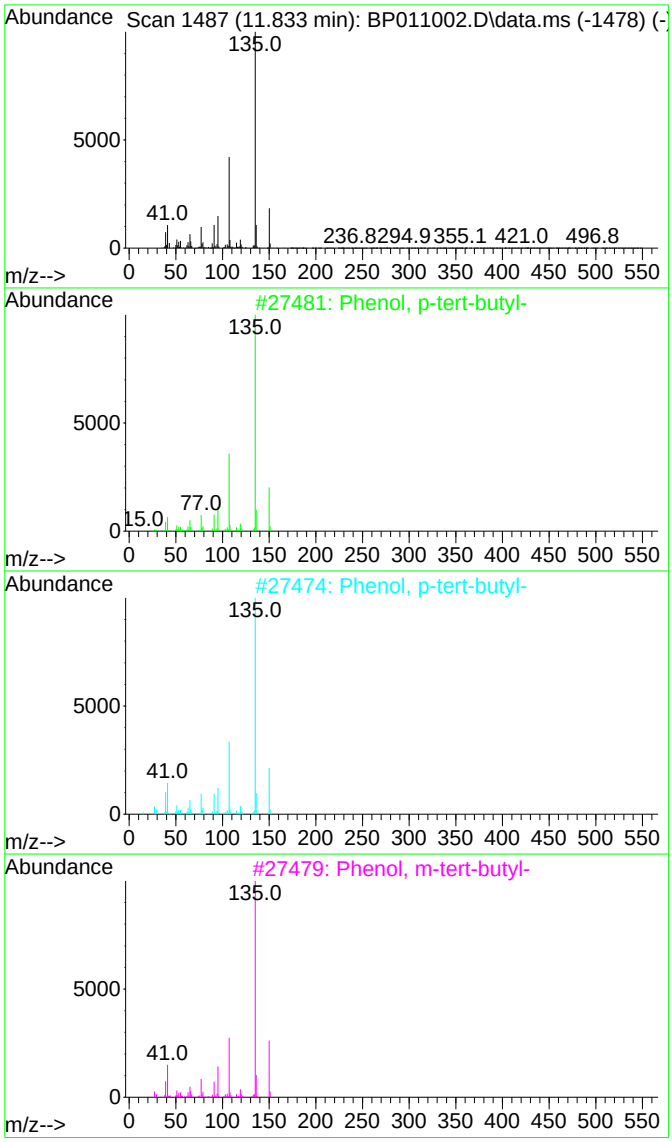
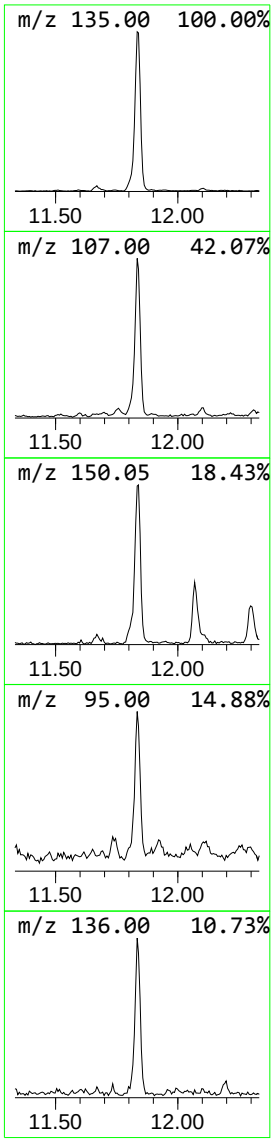
Peak Number 10 Phenol, p-tert-butyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.834	3.22 ng/ul	478651	Naphthalene-d8	10.475

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	95
3	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	94
4	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91
5	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	90





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Operator : CG/JU
Sample : N3710-17
Misc :
ALS Vial : 17 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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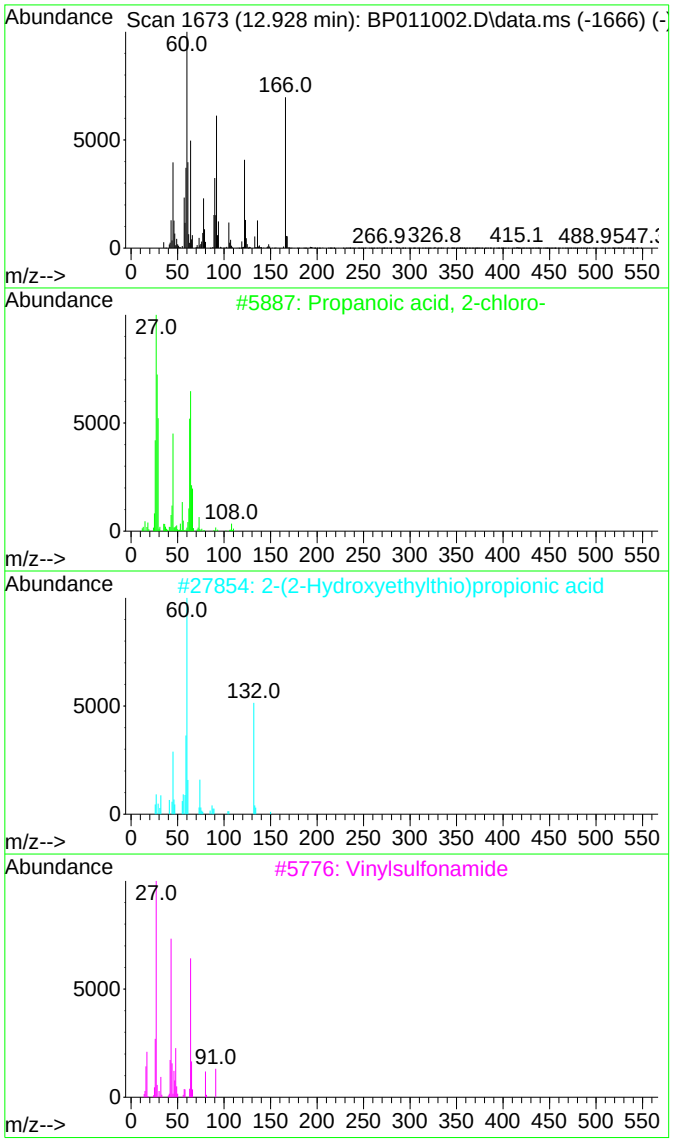
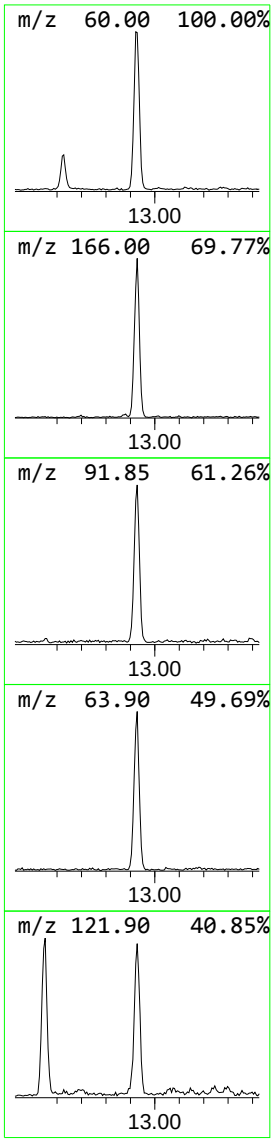
Peak Number 11 unknown-01 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.928	3.41 ng/ul	626810	Acenaphthene-d10	14.339

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, 2-chloro-	108	C3H5ClO2	000598-78-7	16
2	2-(2-Hydroxyethylthio)propionic ...	150	C5H10O3S	209276-28-8	14
3	Vinylsulfonamide	107	C2H5NO2S	002386-58-5	12
4	2-Methyl-1,3-oxathiolane	104	C4H8OS	017642-74-9	12
5	(S)-(-)-2-Chloropropionic acid	108	C3H5ClO2	029617-66-1	12





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Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

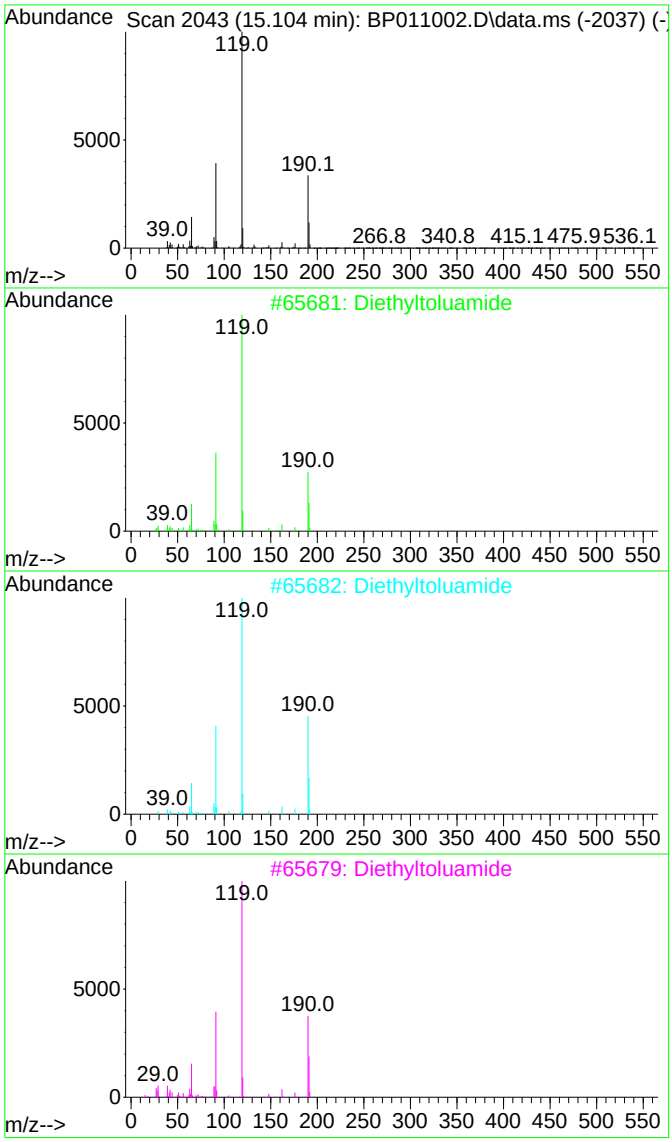
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

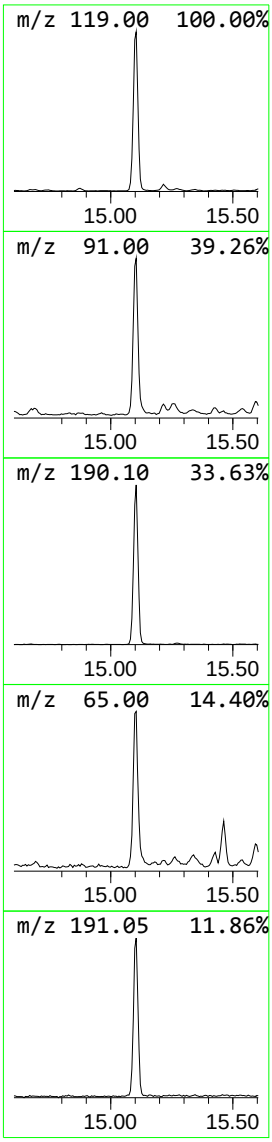
****

Peak Number 12 Diethyltoluamide Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.104	7.85 ng/ul	1441040	Acenaphthene-d10	14.339

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethyltoluamide	191	C12H17NO	000134-62-3	96
2	Diethyltoluamide	191	C12H17NO	000134-62-3	96
3	Diethyltoluamide	191	C12H17NO	000134-62-3	93
4	Benzamide, 3-methyl-N-methyl-N-p...	191	C12H17NO	1000421-46-2	87
5	Benzamide, 3-methyl-N-butyl-	191	C12H17NO	1000407-41-1	87





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Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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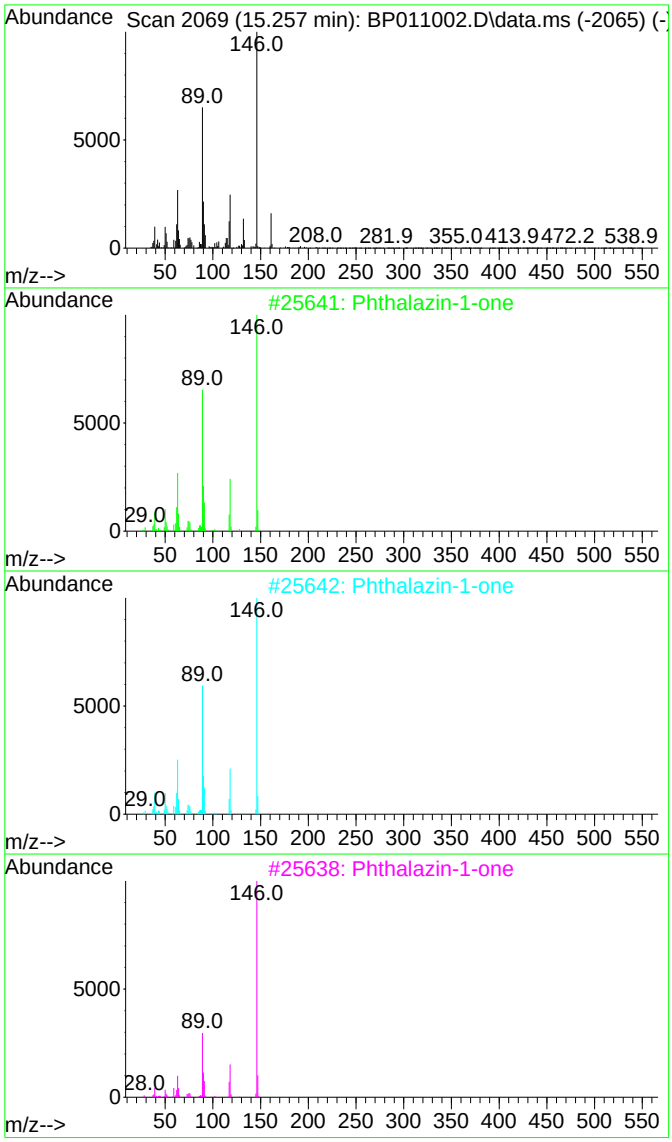
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

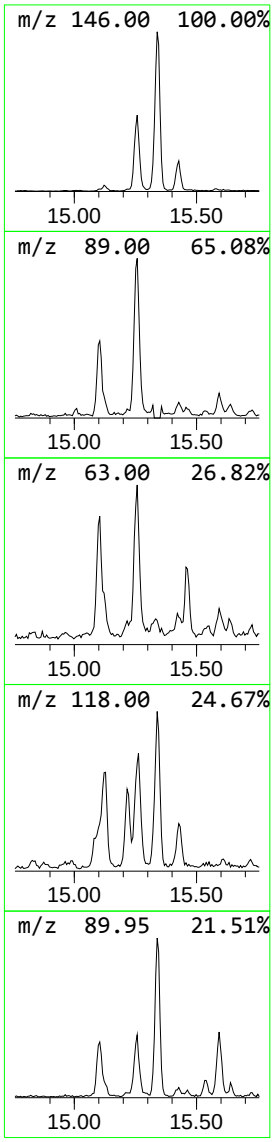
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Phthalazin-1-one Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.257	2.18 ng/ul	401189	Acenaphthene-d10	14.339

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phthalazin-1-one	146	C8H6N2O	000119-39-1	96
2	Phthalazin-1-one	146	C8H6N2O	000119-39-1	93
3	Phthalazin-1-one	146	C8H6N2O	000119-39-1	87
4	Cinnoline, 2-oxide	146	C8H6N2O	004215-44-5	59
5	Coumarin	146	C9H6O2	000091-64-5	50





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ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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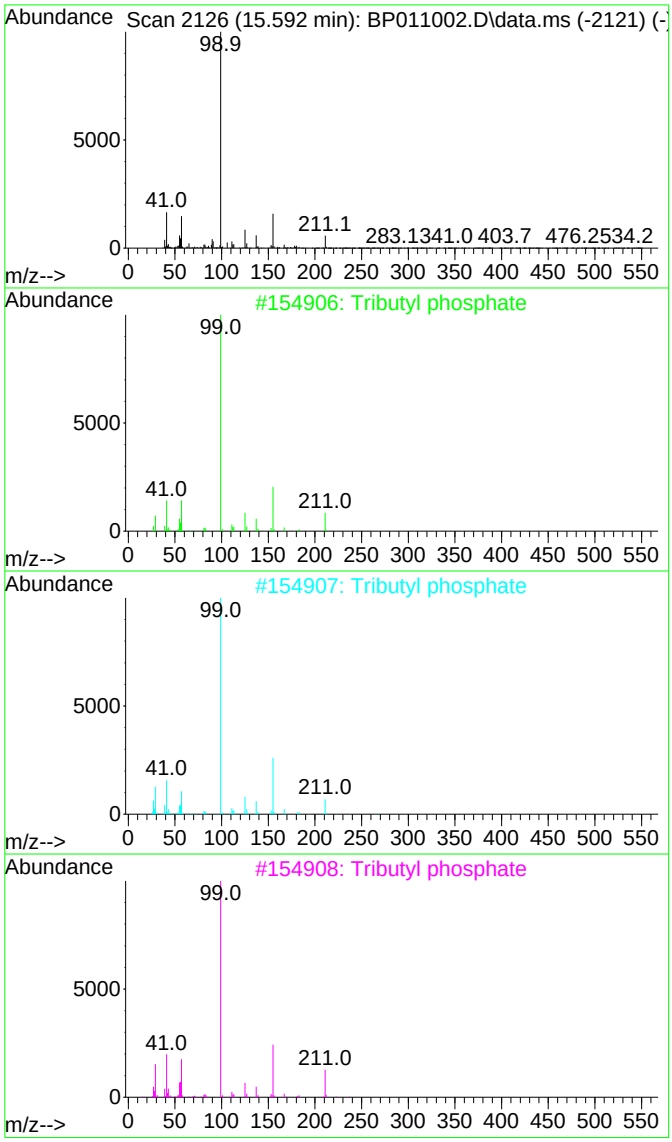
Quant Title : SVOA CALIBRATION

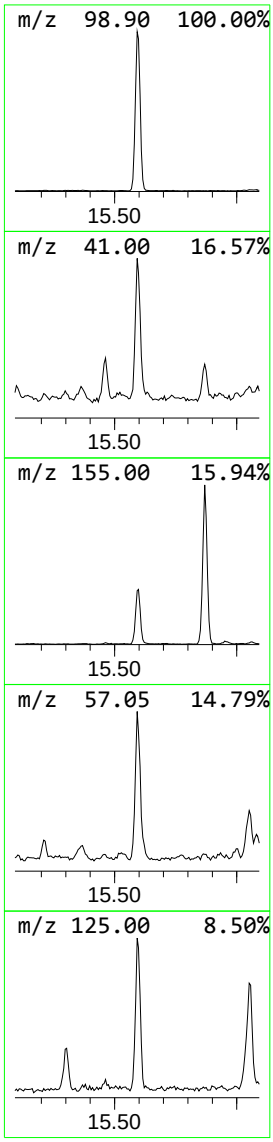
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Tributyl phosphate Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.592	6.14 ng/ul	1128120	Acenaphthene-d10	14.339

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tributyl phosphate	266	C12H27O4P	000126-73-8	87
2	Tributyl phosphate	266	C12H27O4P	000126-73-8	81
3	Tributyl phosphate	266	C12H27O4P	000126-73-8	72
4	Tributyl phosphate	266	C12H27O4P	000126-73-8	72
5	Phosphoric acid, dibutyl 1,1-dim...	352	C13H25F4O4P	1000298-93-9	56





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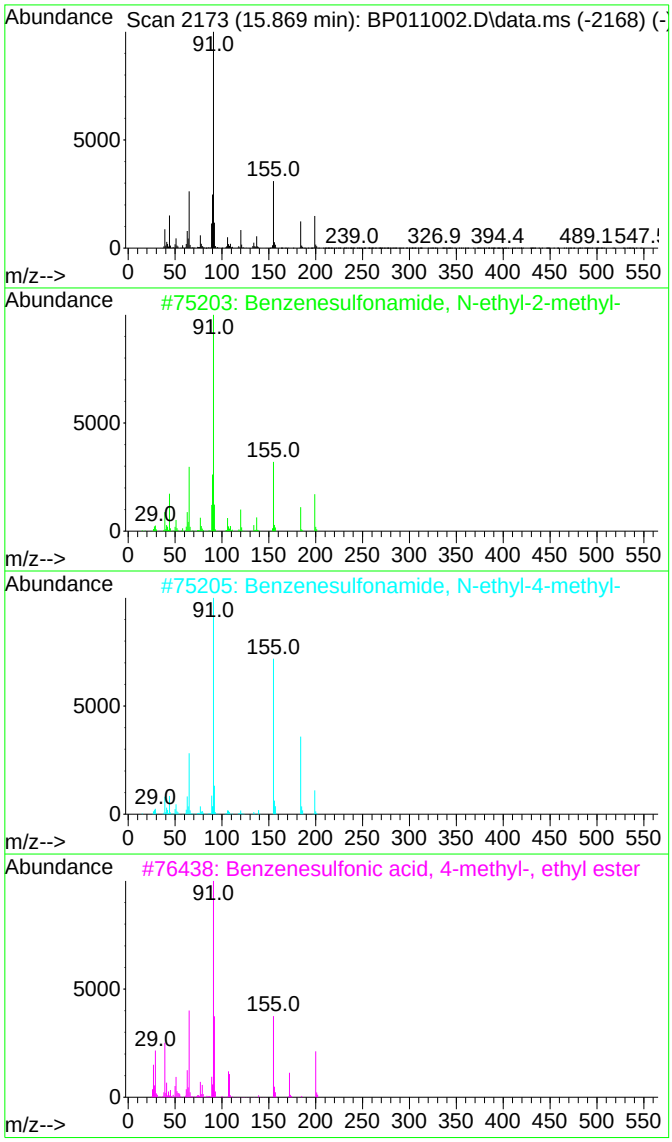
Quant Title : SVOA CALIBRATION

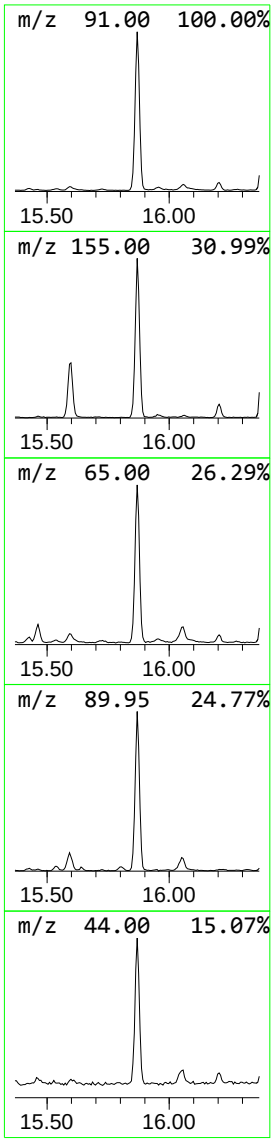
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Benzenesulfonamide, N-ethyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.868	9.33 ng/ul	1858850	Phenanthrene-d10	17.110

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	98
2	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	59
3	Benzenesulfonic acid, 4-methyl-,...	200	C9H12O3S	000080-40-0	53
4	Tolbutamide	270	C12H18N2O3S	000064-77-7	50
5	4-Toluenesulfonylmethyl isocyanide	195	C9H9NO2S	036635-61-7	50





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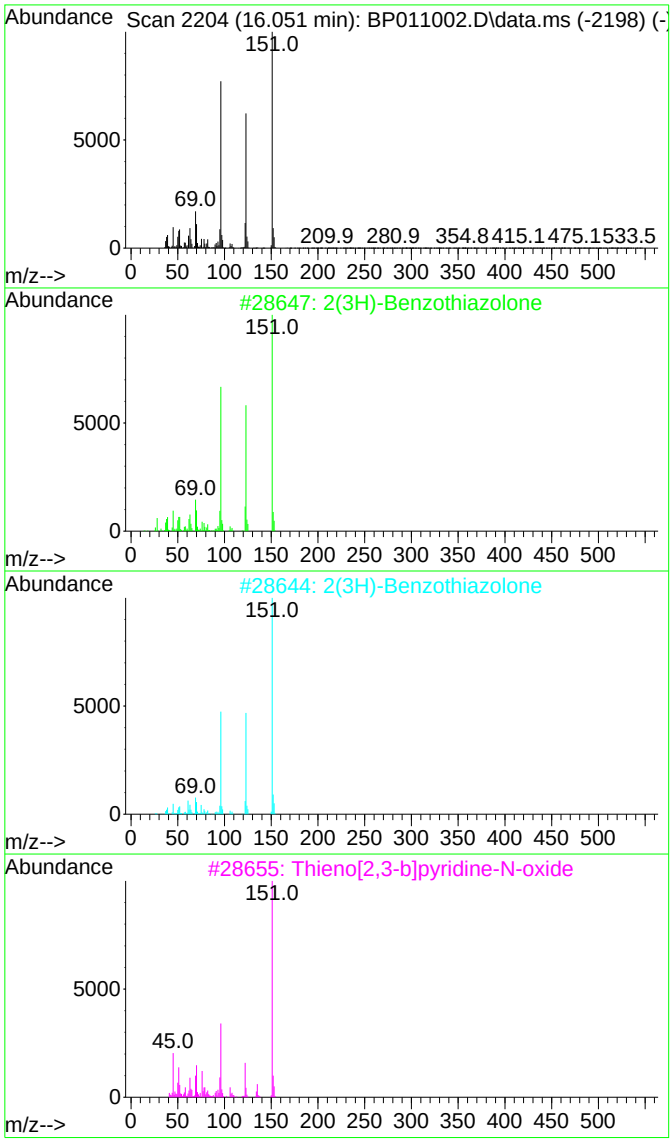
Quant Title : SVOA CALIBRATION

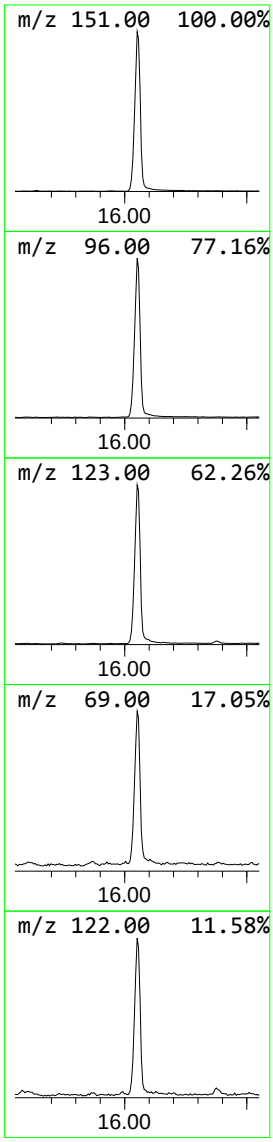
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 2(3H)-Benzothiazolone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.051	18.70 ng/ul	3726700	Phenanthrene-d10	17.110

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	96
2	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	93
3	Thieno[2,3-b]pyridine-N-oxide	151	C7H5NOS	025557-50-0	59
4	Thieno[3,2-b]pyridin-7-ol	151	C7H5NOS	107818-20-2	49
5	1,2-Benzisothiazol-3(2H)-one	151	C7H5NOS	002634-33-5	46





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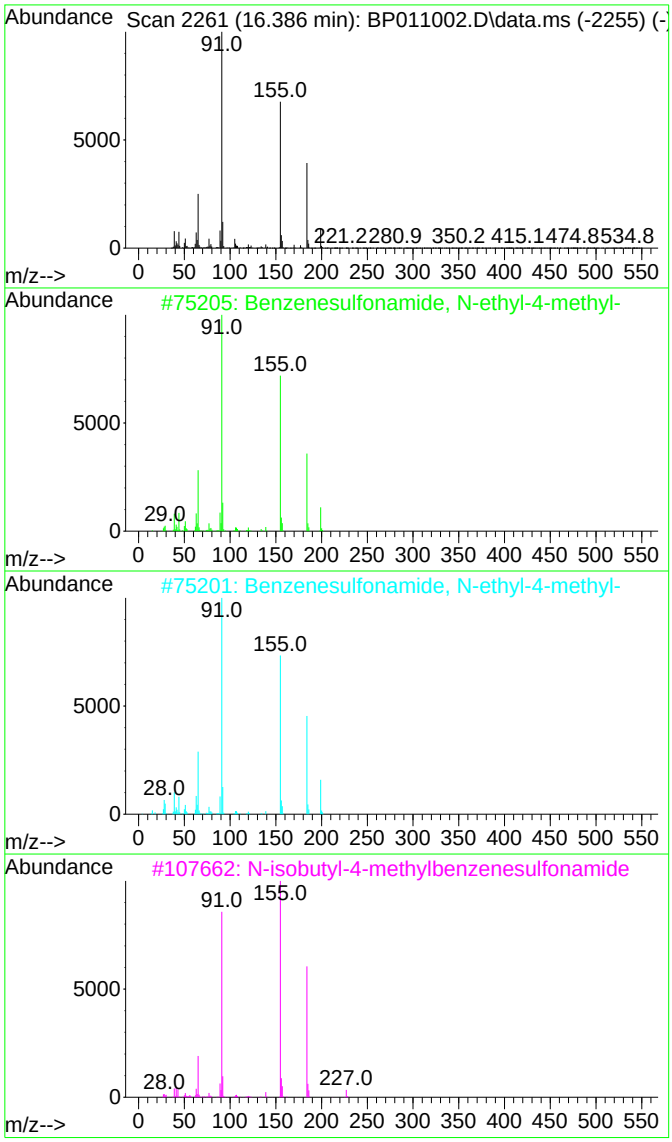
Quant Title : SVOA CALIBRATION

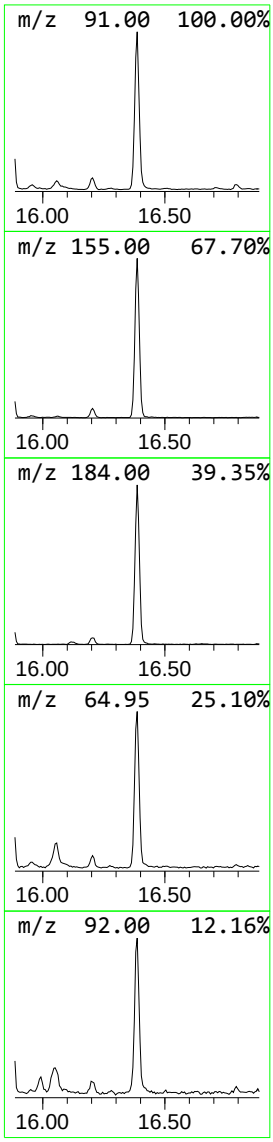
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Benzenesulfonamide, N-ethyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.386	7.91 ng/ul	1577220	Phenanthrene-d10	17.110

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	N-isobutyl-4-methylbenzenesulfon...	227	C11H17NO2S	023705-38-6	86
4	4-Methyl-N-(2-oxobutyl)benzenesu...	241	C11H15NO3S	1000192-14-5	86
5	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	78





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ClientSampleId :
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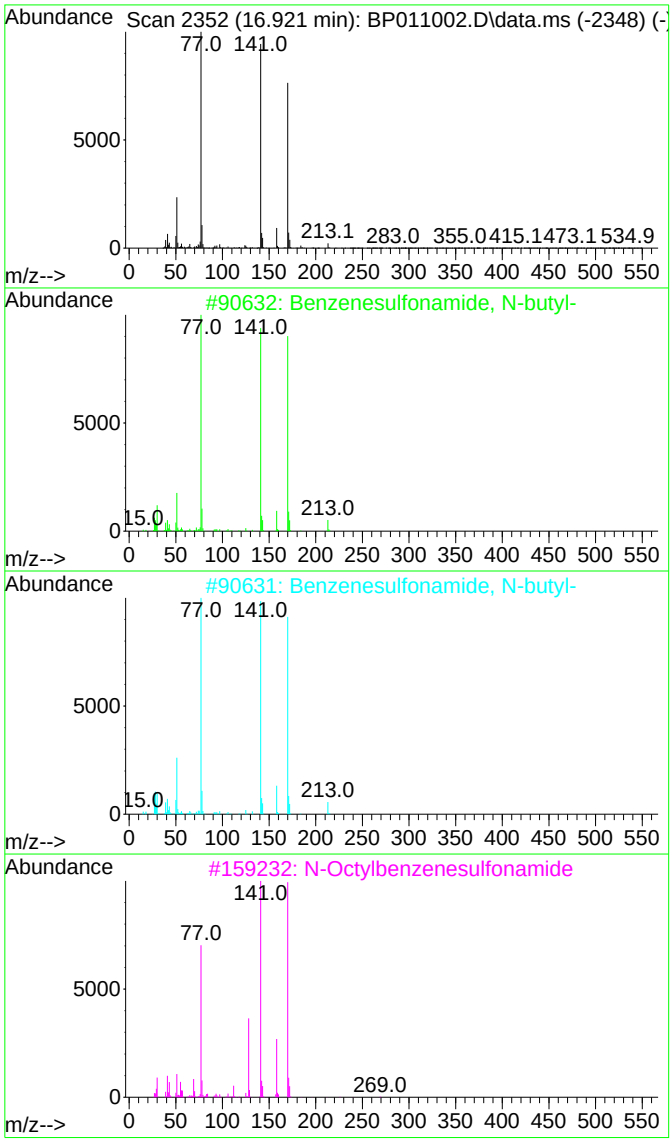
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

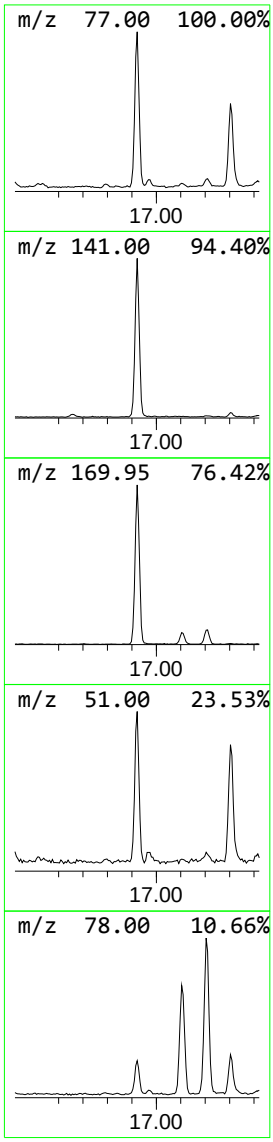
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Benzenesulfonamide, N-butyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.921	3.12 ng/ul	622822	Phenanthrene-d10	17.110

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-butyl-	213	C10H15NO2S	003622-84-2	91
2	Benzenesulfonamide, N-butyl-	213	C10H15NO2S	003622-84-2	91
3	N-Octylbenzenesulfonamide	269	C14H23NO2S	016358-32-0	74
4	N-Benzenesulfonyloxy-2,2-bis(tri...	335	C10H7F6NO3S	055734-41-3	50
5	2-propenoic acid, 2-methyl-, 2-[...	269	C12H15NO4S	1000401-71-8	50





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Data File : BP011002.D
Acq On : 18 Jul 2022 20:27
Operator : CG/JU
Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0B09

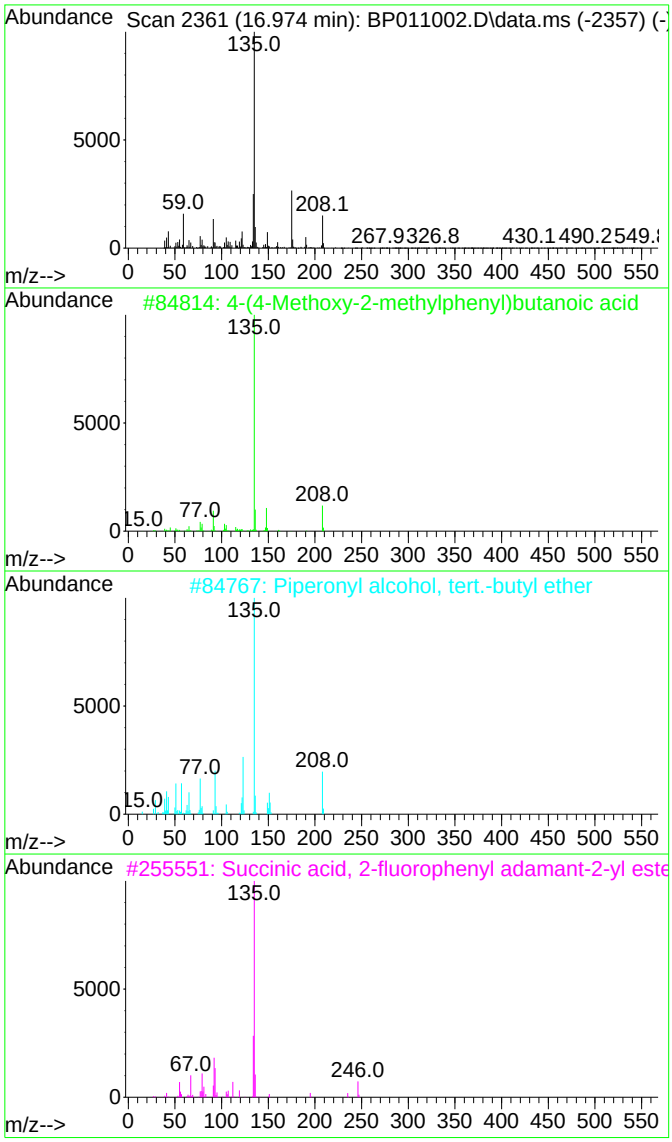
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

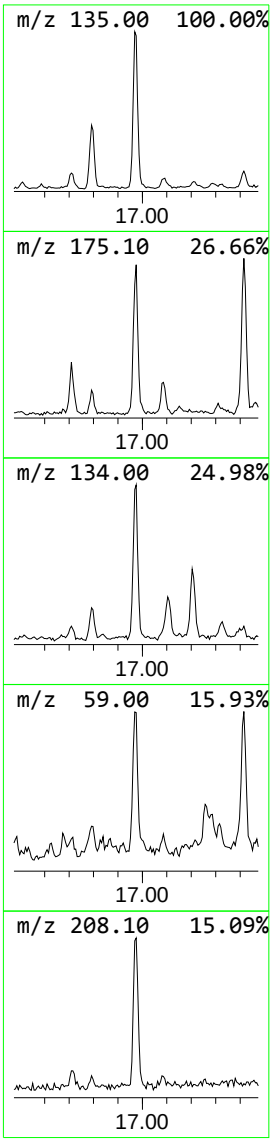
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 4-(4-Methoxy-2-methylphenyl... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.974	2.23 ng/ul	443637	Phenanthrene-d10	17.110

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(4-Methoxy-2-methylphenyl)buta...	208	C12H16O3	006307-30-8	60
2	Piperonyl alcohol, tert.-butyl e...	208	C12H16O3	1000395-11-7	50
3	Succinic acid, 2-fluorophenyl ad...	346	C20H23FO4	1010391-34-3	50
4	Benzenepropanoic acid, 4-methoxy...	208	C11H12O4	1000338-13-0	47
5	Silane, methyldimethoxyethoxy-	150	C5H14O3Si	1010416-18-1	47





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Data File : BP011002.D
Acq On : 18 Jul 2022 20:27
Operator : CG/JU
Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0B09

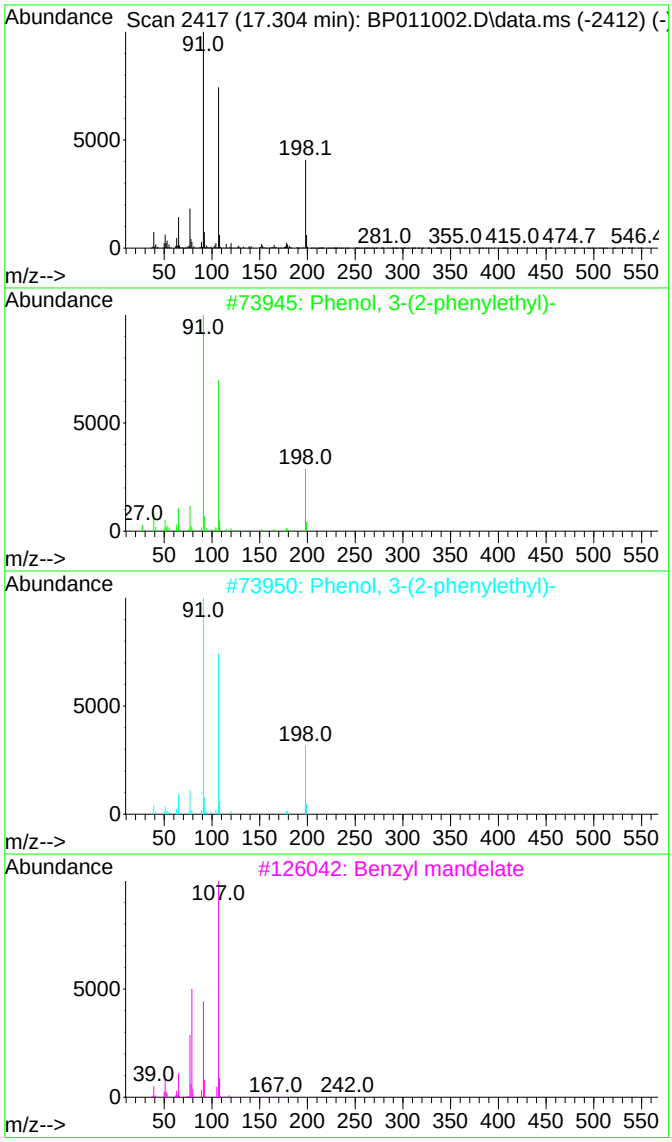
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

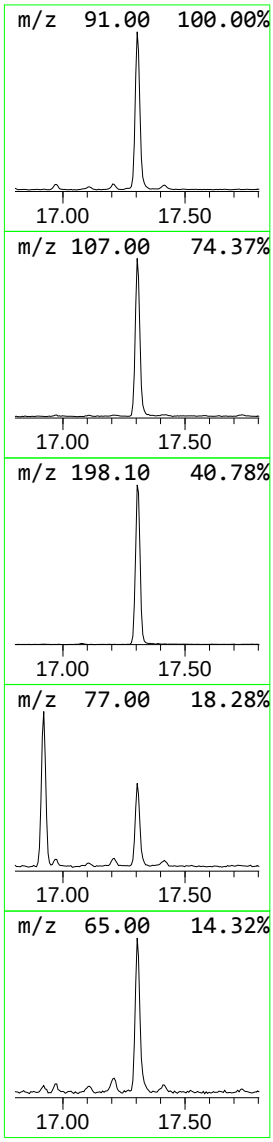
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Phenol, 3-(2-phenylethyl)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.304	10.23 ng/ul	2038320	Phenanthrene-d10	17.110

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3-(2-phenylethyl)-	198	C14H14O	033675-75-1	97
2	Phenol, 3-(2-phenylethyl)-	198	C14H14O	033675-75-1	95
3	Benzyl mandelate	242	C15H14O3	000890-98-2	47
4	p-Hydroxyphenylacetone	150	C9H10O2	000770-39-8	45
5	Pyridine, 2,5-dimethyl-	107	C7H9N	000589-93-5	43





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Data File : BP011002.D
Acq On : 18 Jul 2022 20:27
Operator : CG/JU
Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0B09

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

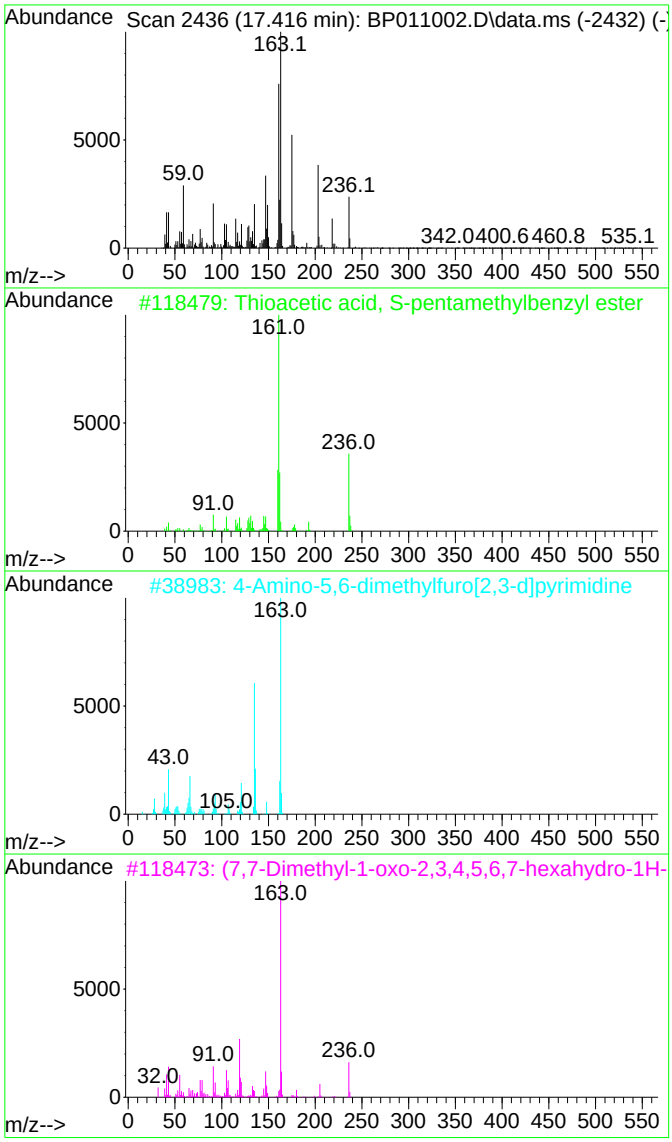
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

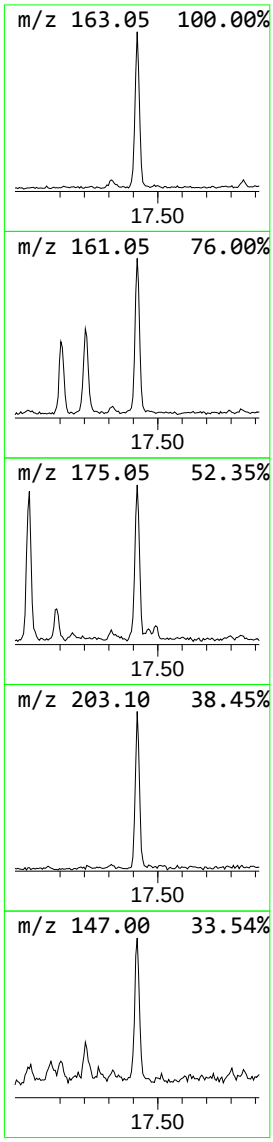
****

Peak Number 21 unknown-02 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.416	2.70 ng/ul	539107	Phenanthrene-d10	17.110

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Thioacetic acid, S-pentamethylbe...	236	C14H20OS	1000267-40-4	22
2	4-Amino-5,6-dimethylfuro[2,3-d]p...	163	C8H9N3O	005117-94-2	18
3	(7,7-Dimethyl-1-oxo-2,3,4,5,6,7-...	236	C14H20O3	055085-50-2	14
4	2(3H)-Naphthalenone, 4,4a,5,6,7,...	218	C15H22O	019598-45-9	14
5	3-Piperidinamine, 1-(2-pyrimidin...	262	C13H18N4O2	1000470-29-7	14





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Data File : BP011002.D
Acq On : 18 Jul 2022 20:27
Operator : CG/JU
Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0B09

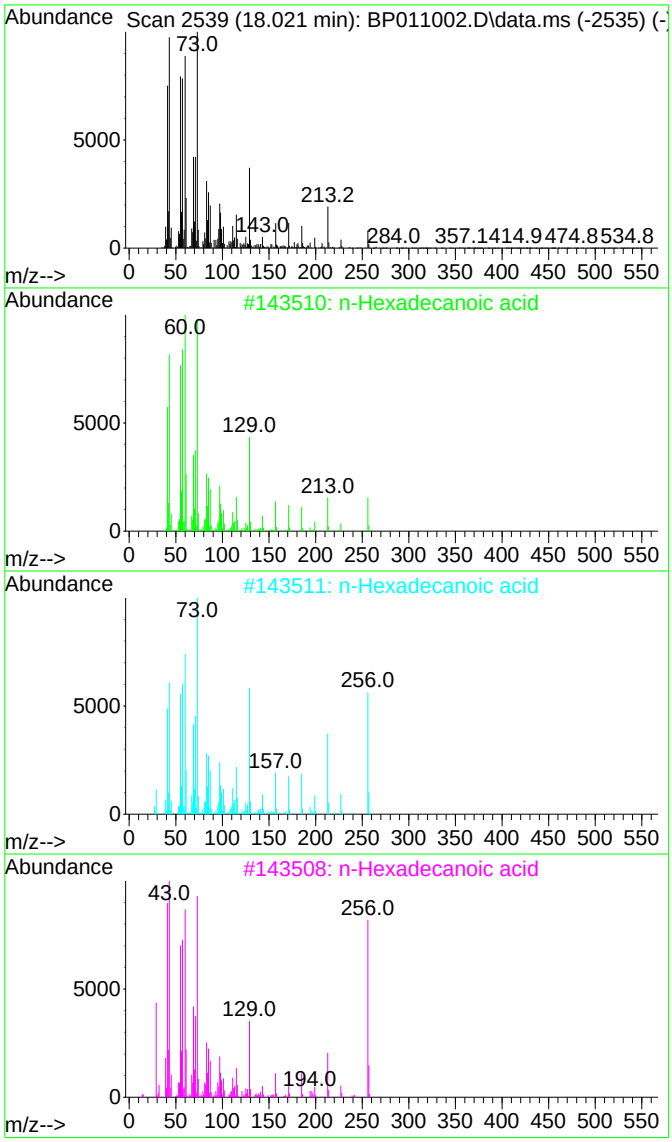
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

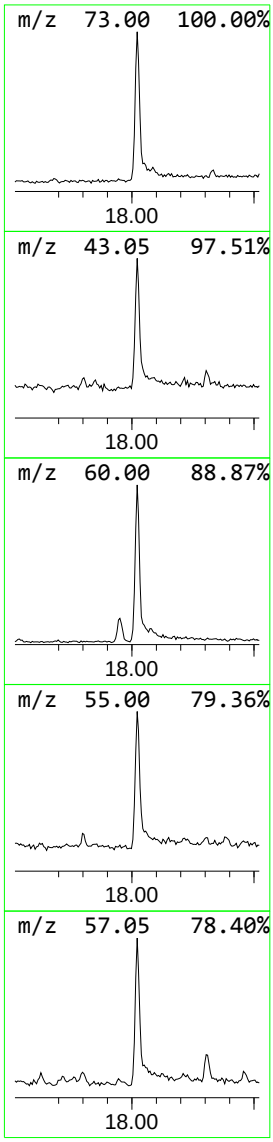
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 n-Hexadecanoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.021	4.27 ng/ul	851003	Phenanthrene-d10	17.110

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	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071822\
Data File : BP011002.D
Acq On : 18 Jul 2022 20:27
Operator : CG/JU
Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0B09

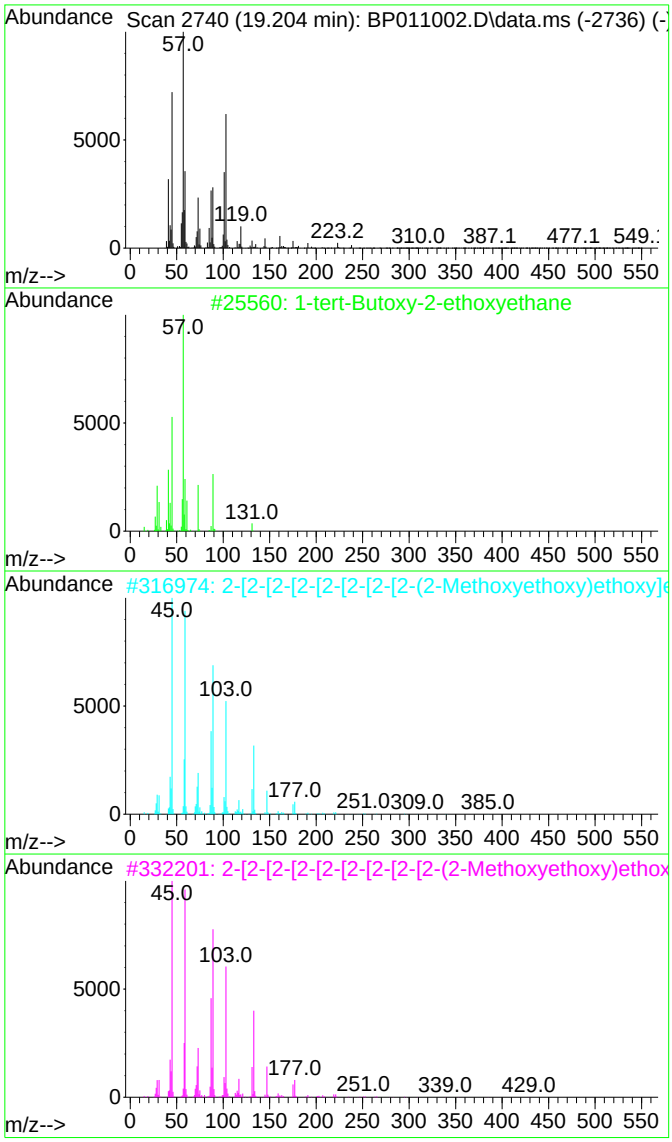
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

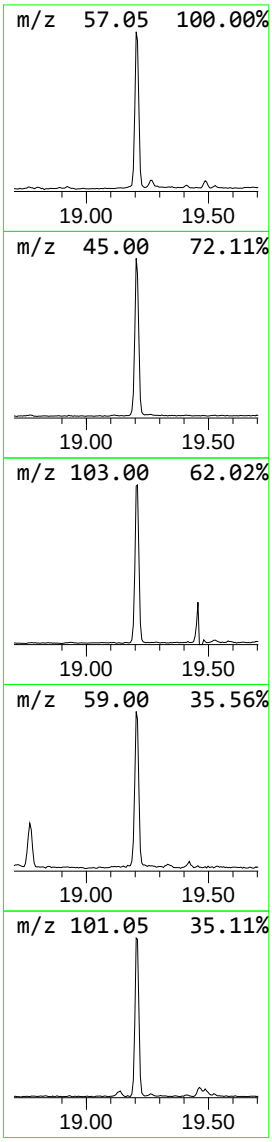
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 unknown-03 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.204	7.70 ng/ul	1742730	Chrysene-d12	21.227

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-tert-Butoxy-2-ethoxyethane	146	C8H18O2	051422-54-9	43
2	2-[2-[2-[2-[2-[2-[2-(2-Methox...	428	C19H40O10	006048-68-6	27
3	2-[2-[2-[2-[2-[2-[2-[2-(2-Met...	472	C21H44O11	027425-92-9	27
4	2-[2-[2-[2-[2-[2-(2-Methoxyet...	384	C17H36O9	025990-96-9	27
5	Butyl 2,5,8,11,14,17,20,23-octao...	454	C21H42O10	1000366-81-8	27





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071822\
Data File : BP011002.D
Acq On : 18 Jul 2022 20:27
Operator : CG/JU
Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0B09

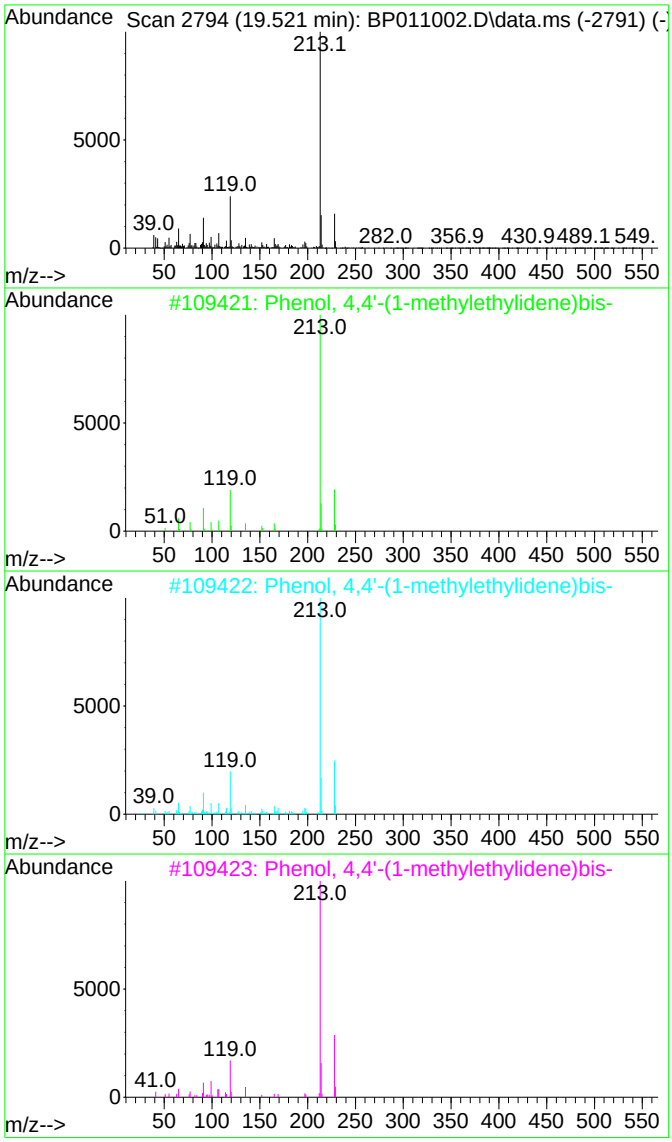
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

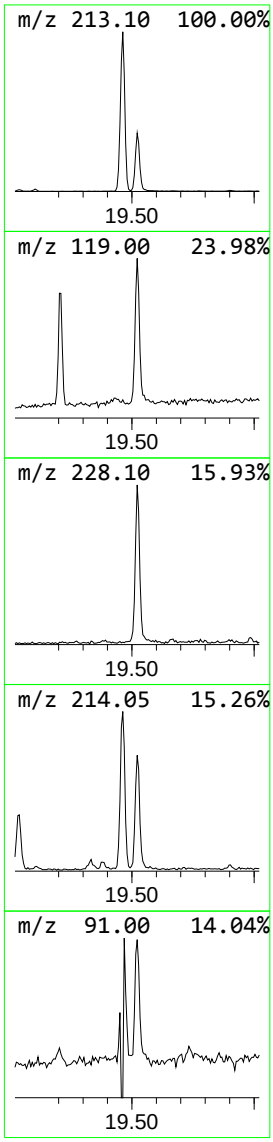
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Phenol, 4,4'-(1-methylethyl... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.521	2.35 ng/ul	531834	Chrysene-d12	21.227

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	97
2	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	92
3	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	91
4	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	90
5	4,4'-(1,3-Dimethylbutylidene)bis...	270	C18H22O2	006807-17-6	72





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Acq On : 18 Jul 2022 20:27
Operator : CG/JU
Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0B09

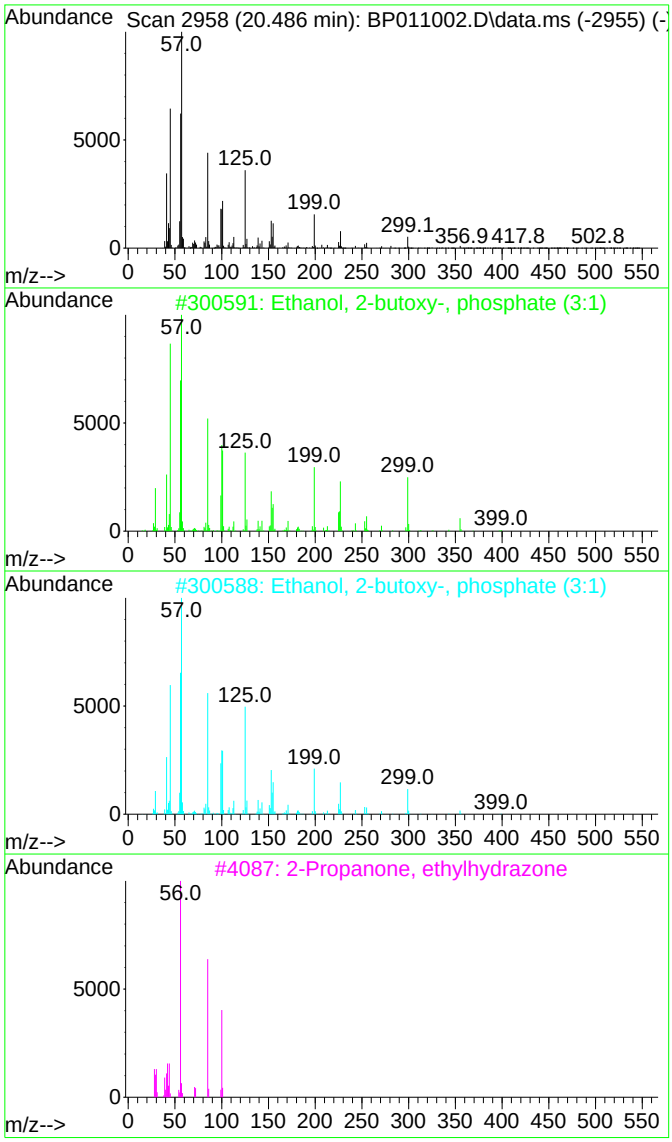
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

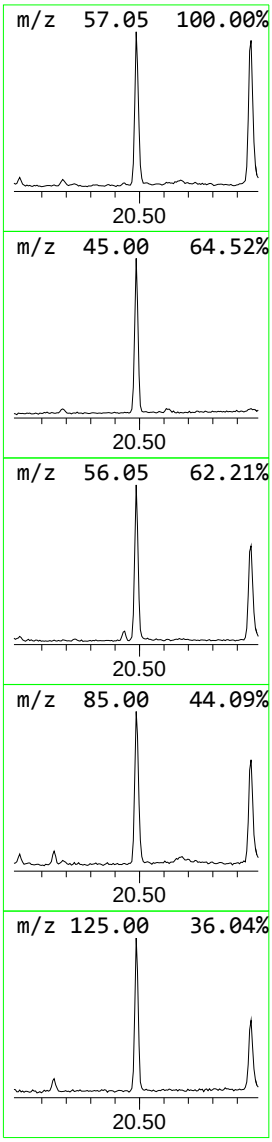
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 Ethanol, 2-butoxy-, phospha... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.486	4.93 ng/ul	1116000	Chrysene-d12	21.227

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-butoxy-, phosphate (3:1)	398	C18H39O7P	000078-51-3	83
2	Ethanol, 2-butoxy-, phosphate (3:1)	398	C18H39O7P	000078-51-3	52
3	2-Propanone, ethylhydrazone	100	C5H12N2	007422-99-3	35
4	Furan, tetrahydro-2,5-dimethyl-	100	C6H12O	001003-38-9	35
5	2(3H)-Furanone, dihydro-5-methyl-	100	C5H8O2	000108-29-2	35





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Acq On : 18 Jul 2022 20:27
Operator : CG/JU
Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0B09

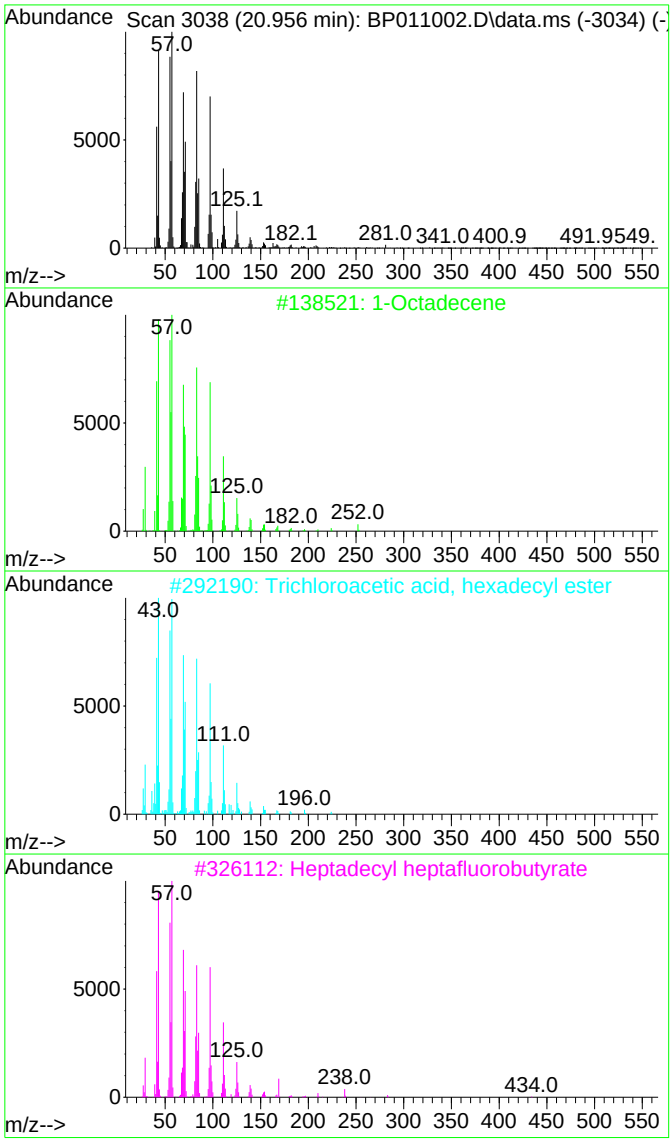
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

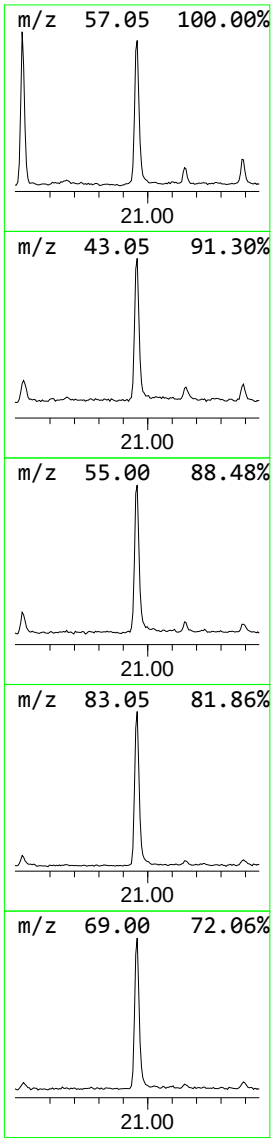
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.956	8.90 ng/ul	2014030	Chrysene-d12	21.227

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Octadecene	252	C18H36	000112-88-9	95
2	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
3	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
4	Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	94
5	n-Tetracosanol-1	354	C24H50O	000506-51-4	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071822\
Data File : BP011002.D
Acq On : 18 Jul 2022 20:27
Operator : CG/JU
Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0B09

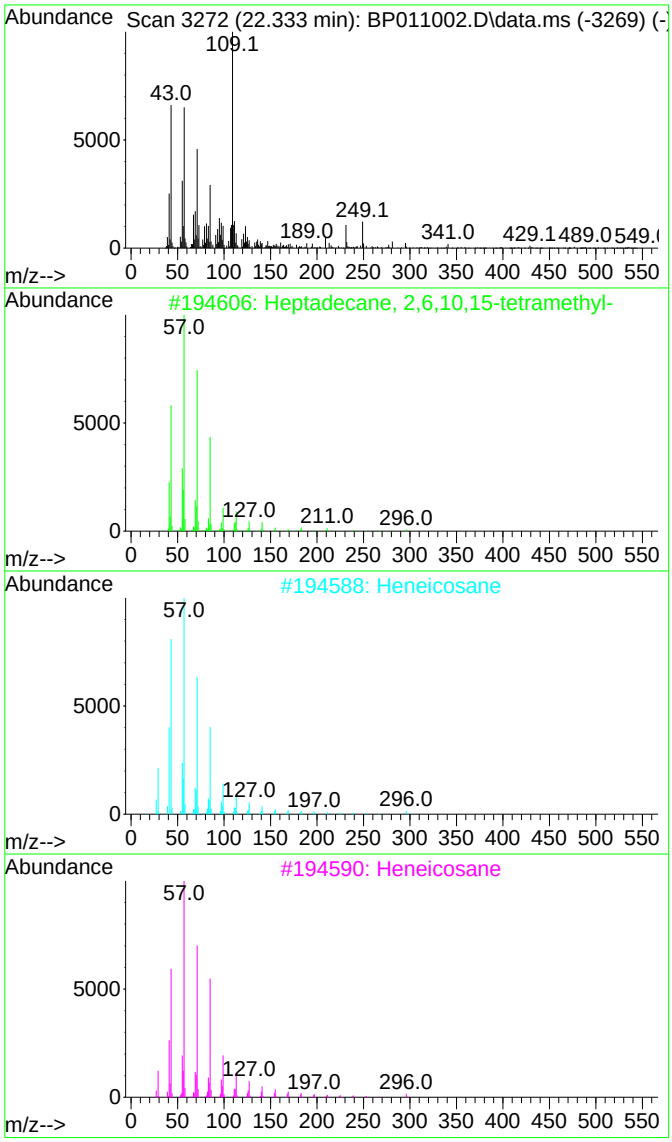
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

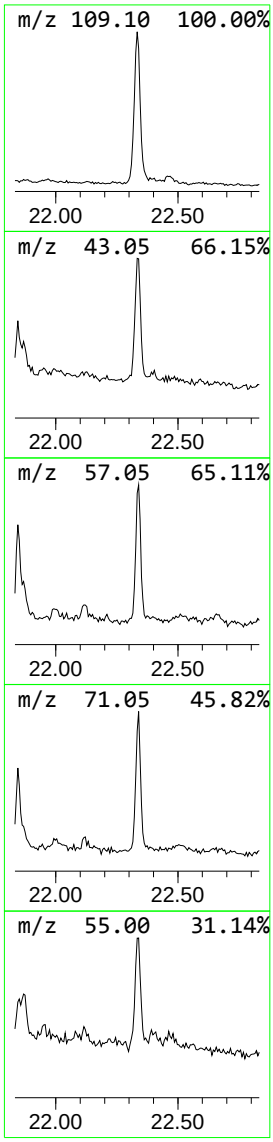
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.333	3.03 ng/ul	685949	Chrysene-d12	21.227

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	64
2	Heneicosane	296	C21H44	000629-94-7	50
3	Heneicosane	296	C21H44	000629-94-7	43
4	8-Hydroxy-7(11)-eremophilen-12,8...	250	C15H22O3	065562-67-6	41
5	p-Decyloxyaniline	249	C16H27NO	039905-47-0	38





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071822\
Data File : BP011002.D
Acq On : 18 Jul 2022 20:27
Operator : CG/JU
Sample : N3710-17
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0B09

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.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	4.340	31.2	ng/ul	3333920	1	7.681	2138800	20.0
Hexanoic acid	6.722	6.1	ng/ul	656465	1	7.681	2138800	20.0
Heptanoic acid	8.281	3.3	ng/ul	350309	1	7.681	2138800	20.0
p-Cresol	8.451	2.6	ng/ul	277429	1	7.681	2138800	20.0
(DEL) Alkane: S...	9.045	2.1	ng/ul	221605	1	7.681	2138800	20.0
Cyclohexanemeth...	9.722	8.7	ng/ul	1299490	2	10.475	2976610	20.0
Octanoic acid	9.845	10.2	ng/ul	1522590	2	10.475	2976610	20.0
n-Decanoic acid	11.292	3.4	ng/ul	511104	2	10.475	2976610	20.0
Phenol, p-tert-...	11.834	3.2	ng/ul	478651	2	10.475	2976610	20.0
unknown-01	12.928	3.4	ng/ul	626810	3	14.339	3673080	20.0
Diethyltoluamide	15.104	7.8	ng/ul	1441040	3	14.339	3673080	20.0
Phthalazin-1-one	15.257	2.2	ng/ul	401189	3	14.339	3673080	20.0
Tributyl phosphate	15.592	6.1	ng/ul	1128120	3	14.339	3673080	20.0
Benzenesulfonam...	15.868	9.3	ng/ul	1858850	4	17.110	3986230	20.0
2(3H)-Benzothia...	16.051	18.7	ng/ul	3726700	4	17.110	3986230	20.0
Benzenesulfonam...	16.386	7.9	ng/ul	1577220	4	17.110	3986230	20.0
Benzenesulfonam...	16.921	3.1	ng/ul	622822	4	17.110	3986230	20.0
4-(4-Methoxy-2-...	16.974	2.2	ng/ul	443637	4	17.110	3986230	20.0
Phenol, 3-(2-ph...	17.304	10.2	ng/ul	2038320	4	17.110	3986230	20.0
unknown-02	17.416	2.7	ng/ul	539107	4	17.110	3986230	20.0
n-Hexadecanoic ...	18.021	4.3	ng/ul	851003	4	17.110	3986230	20.0
unknown-03	19.204	7.7	ng/ul	1742730	5	21.227	4526030	20.0
Phenol, 4,4'-(1...	19.521	2.4	ng/ul	531834	5	21.227	4526030	20.0
Ethanol, 2-buto...	20.486	4.9	ng/ul	1116000	5	21.227	4526030	20.0
(DEL) Alkane: S...	20.956	8.9	ng/ul	2014030	5	21.227	4526030	20.0
(DEL) Alkane: S...	22.333	3.0	ng/ul	685949	5	21.227	4526030	20.0