LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012093.D
Acq On : 12 Oct 2022 22:41
Operator : CG/JU
Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8Y7

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-BP101122.M
Title : SVOA CALIBRATION

Signal : TIC: BP012093.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.075	12	15	31	rVB	212342	321460	5.01%	0.385%
2	3.275	45	49	51	rBV	51335	58544	0.91%	0.070%
3	3.310	51	55	67	rVB	416105	537322	8.38%	0.644%
4	3.704	118	122	139	rBV	460618	742752	11.59%	0.890%
5	3.922	155	159	170	rVV	27275	50668	0.79%	0.061%
6	4.469	245	252	261	rVB	316815	473608	7.39%	0.568%
7	4.963	328	336	342	rBV2	16180	38696	0.60%	0.046%
8	5.157	363	369	376	rBV	303073	454856	7.09%	0.545%
9	5.593	439	443	449	rBV	41484	64148	1.00%	0.077%
10	6.128	523	534	542	rBV	26523	65045	1.01%	0.078%
11	6.445	582	588	593	rBV2	34264	56862	0.89%	0.068%
12	7.022	679	686	699	rBV	245769	467770	7.30%	0.561%
13	7.187	707	714	732	rBV	752037	1247284	19.46%	1.495%
14	7.381	741	747	755	rVV	794943	1307840	20.40%	1.567%
15	7.451	755	759	763	rVV6	17947	44838	0.70%	0.054%
16	7.710	794	803	805	rVV9	23491	47674	0.74%	0.057%
17	7.739	806	808	815	rVV7	24755	56027	0.87%	0.067%
18	7.851	816	827	836	rVV	965314	1631219	25.44%	1.955%
19	7.934	837	841	852	rVV3	59988	154687	2.41%	0.185%
20	8.222	882	890	903	rVV	207598	377843	5.89%	0.453%
21	8.339	903	910	915	rVV5	24349	49849	0.78%	0.060%
22	8.510	933	939	942	rBV2	38111	67578	1.05%	0.081%
23	8.569	943	949	958	rVV	696568	1150574	17.95%	1.379%
24	8.769	977	983	989	rVB7	15438	39185	0.61%	0.047%
25	8.845	989	996	1005	rBV	309039	527429	8.23%	0.632%
26	8.957	1010	1015	1020	rBV	87171	146156	2.28%	0.175%
27	9.022	1020	1026	1037	rVB	797337	1379503	21.52%	1.653%
28	9.228	1055	1061	1068	rVB9	27648	64863	1.01%	0.078%
29	9.375	1080	1086	1091	rBV	92269	163523	2.55%	0.196%
30	9.481	1098	1104	1114	rVB2	106280	178575	2.79%	0.214%
31	9.604	1120	1125	1130	rBV3	36941	59265	0.92%	0.071%
32	9.739	1142	1148	1162	rVB	626940	1098213	17.13%	1.316%
33	9.875	1162	1171	1179	rBV8	62497	197198	3.08%	0.236%
34	10.063	1195	1203	1207	rBV8	24443	55896	0.87%	0.067%
35	10.181	1213	1223	1234	rVB5	113932	352703	5.50%	0.423%
36	10.281	1234	1240	1249	rBV	1079194	1870311	29.17%	2.241%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012093.D
Acq On : 12 Oct 2022 22:41
Operator : CG/JU
Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y7

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-BP101122.M
Title : SVOA CALIBRATION

37	10.439	1261	1267	1281	rBV	526381	903003	14.09%	1.082%
38	10.598	1286	1294	1298	rVV	129287	229722	3.58%	0.275%
39	10.657	1298	1304	1312	rVB	1230281	2085510	32.53%	2.499%
40	10.804	1318	1329	1343	rBV	783701	1454031	22.68%	1.742%
41	10.916	1344	1348	1352	rVV4	39758	68257	1.06%	0.082%
42	10.963	1352	1356	1363	rVV5	44710	91395	1.43%	0.110%
43	11.451	1435	1439	1441	rBV	28725	43517	0.68%	0.052%
44	11.869	1506	1510	1516	rVB2	40431	63731	0.99%	0.076%
45	12.004	1525	1533	1538	rBV	281305	479580	7.48%	0.575%
46	12.051	1539	1541	1548	rVB6	33280	62525	0.98%	0.075%
47	12.169	1555	1561	1565	rBV2	134519	232958	3.63%	0.279%
48	12.275	1576	1579	1589	rVB9	51221	135009	2.11%	0.162%
49	12.433	1603	1606	1613	rBV4	30829	49521	0.77%	0.059%
50	12.545	1622	1625	1629	rVB5	31231	44543	0.69%	0.053%
51	12.657	1639	1644	1650	rBV3	70019	119929	1.87%	0.144%
52	12.757	1656	1661	1670	rVB8	39700	94141	1.47%	0.113%
53	13.304	1750	1754	1758	rBV5	44671	81949	1.28%	0.098%
54	13.398	1765	1770	1777	rVB	207227	313584	4.89%	0.376%
55	13.557	1793	1797	1801	rBV2	56909	73307	1.14%	0.088%
56	13.727	1823	1826	1830	rBV3	29049	43074	0.67%	0.052%
57	13.827	1839	1843	1845	rBV4	37232	57779	0.90%	0.069%
58	13.916	1852	1858	1867	rVB	1738653	2534300	39.53%	3.037%
59	14.057	1878	1882	1885	rBV3	65714	103439	1.61%	0.124%
60	14.192	1899	1905	1912	rBV2	2180393	3280344	51.17%	3.931%
61	14.316	1921	1926	1937	rBV	141733	258196	4.03%	0.309%
62	14.427	1940	1945	1949	rVV	298615	395641	6.17%	0.474%
63	14.498	1951	1957	1963	rVV	1796539	2747046	42.85%	3.292%
64	14.563	1963	1968	1975	rVB	611892	882114	13.76%	1.057%
65	14.698	1988	1991	1992	rBV3	47340	50928	0.79%	0.061%
66	14.727	1993	1996	2000	rVB	108847	151242	2.36%	0.181%
67	15.251	2080	2085	2092	rBV	508057	832961	12.99%	0.998%
68	15.374	2103	2106	2109	rBV4	58906	85343	1.33%	0.102%
69	15.498	2119	2127	2132	rBV	2795938	4361818	68.04%	5.227%
70	15.545	2133	2135	2142	rVB2	105300	146062	2.28%	0.175%
71	15.621	2143	2148	2153	rBV	450076	632234	9.86%	0.758%
72	15.727	2163	2166	2167	rBV	70976	76363	1.19%	0.092%
73	15.757	2168	2171	2180	rVB2	230032	337513	5.26%	0.404%
74	15.957	2201	2205	2209	rBV2	179510	249342	3.89%	0.299%
75	16.021	2210	2216	2228	rVB	1724619	2505895	39.09%	3.003%
76	16.198	2239	2246	2251	rBV2	216654	419177	6.54%	0.502%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012093.D
Acq On : 12 Oct 2022 22:41
Operator : CG/JU
Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y7

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-BP101122.M
Title : SVOA CALIBRATION

77	16.351	2269	2272	2275	rVB	72813	79120	1.23%	0.095%
78	16.533	2297	2303	2308	rBV	1034047	1438382	22.44%	1.724%
79	16.804	2345	2349	2358	rVB	169947	291549	4.55%	0.349%
80	17.257	2421	2426	2434	rVB	2251180	3217516	50.19%	3.855%
81	17.357	2438	2443	2452	rVB2	3131240	4568988	71.27%	5.475%
82	18.145	2573	2577	2591	rBV	983287	1574754	24.56%	1.887%
83	18.898	2703	2705	2711	rVB	186620	219936	3.43%	0.264%
84	19.057	2730	2732	2737	rVB2	134201	126832	1.98%	0.152%
85	19.239	2759	2763	2766	rBV	400465	565201	8.82%	0.677%
86	19.315	2774	2776	2781	rVB	125086	131503	2.05%	0.158%
87	19.380	2783	2787	2801	rVB	682958	1008540	15.73%	1.208%
88	19.592	2818	2823	2828	rBV	4002495	5632359	87.85%	6.749%
89	19.645	2828	2832	2840	rVB	2836690	3658631	57.07%	4.384%
90	20.080	2904	2906	2909	rBV	133206	167054	2.61%	0.200%
91	20.109	2909	2911	2915	rVB	203768	205860	3.21%	0.247%
92	20.439	2964	2967	2970	rBV2	239631	264098	4.12%	0.316%
93	20.592	2990	2993	2997	rBV	502916	565440	8.82%	0.678%
94	21.051	3067	3071	3075	rBV	806843	887305	13.84%	1.063%
95	21.351	3117	3122	3128	rVB	2898175	3823003	59.63%	4.581%
96	21.486	3142	3145	3151	rVB	498225	544309	8.49%	0.652%
97	21.951	3221	3224	3233	rVB	439622	569794	8.89%	0.683%
98	22.450	3306	3309	3322	rVB2	449119	862175	13.45%	1.033%
99	23.609	3500	3506	3518	rVB2	3099778	6411029	100.00%	7.682%
100	23.768	3526	3533	3542	rVB2	2065135	4264288	66.51%	5.110%

Sum of corrected areas: 83454653

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
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Acq On : 12 Oct 2022 22:41
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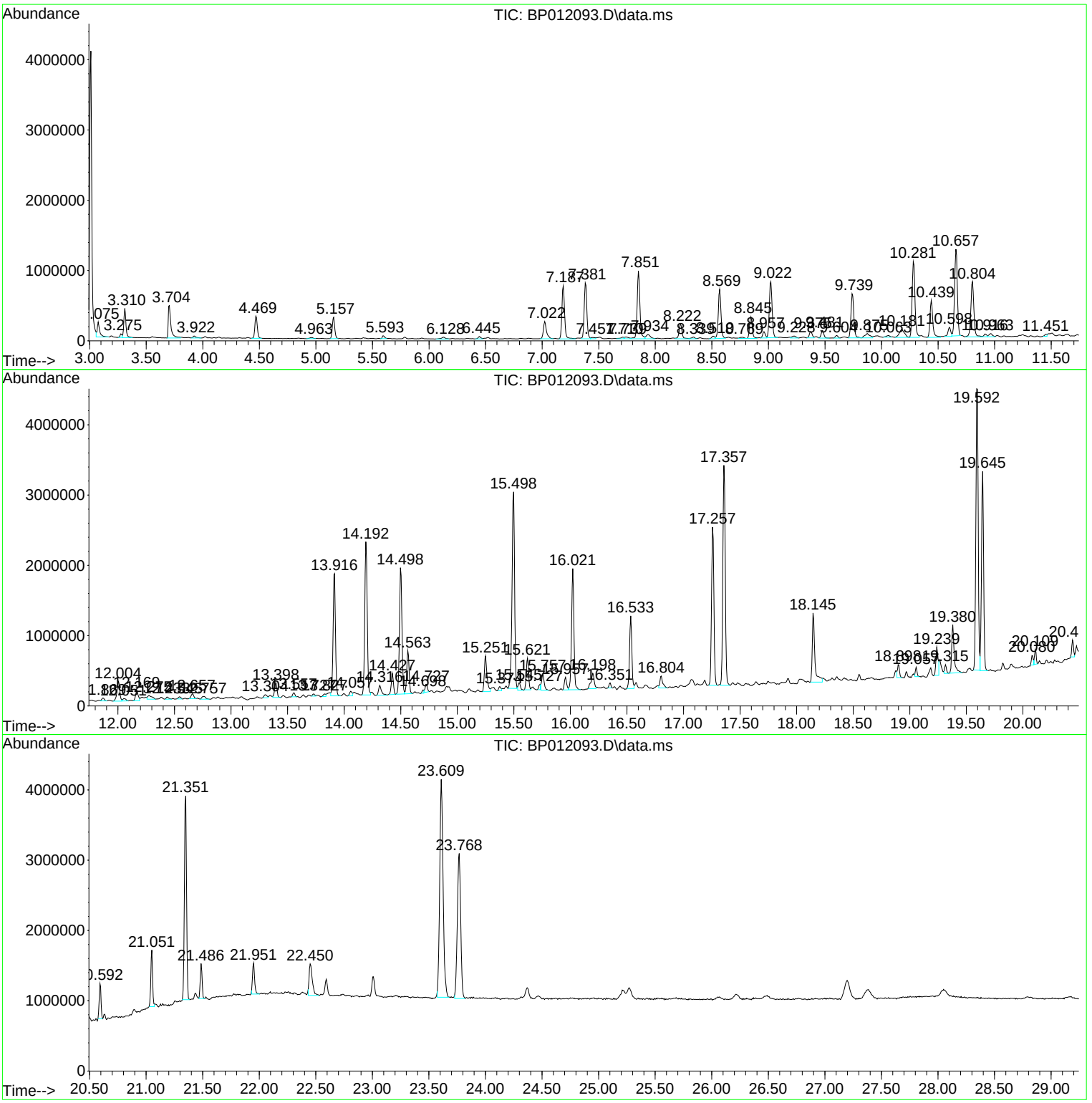
Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012093.D
Acq On : 12 Oct 2022 22:41
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Sample : N4976-11
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

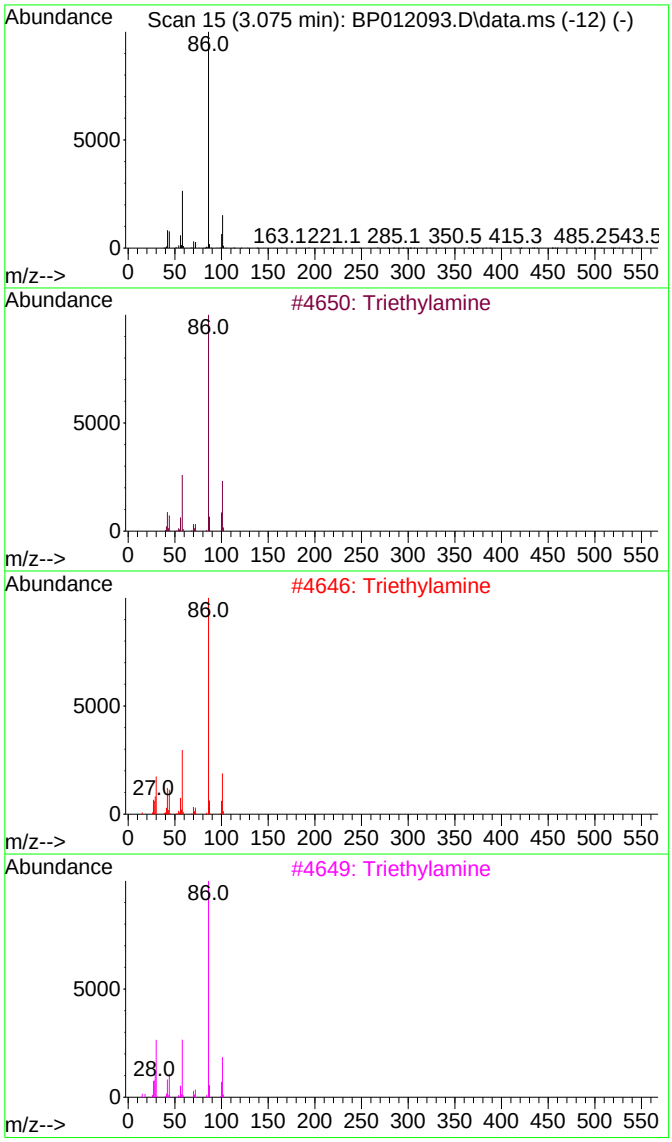
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

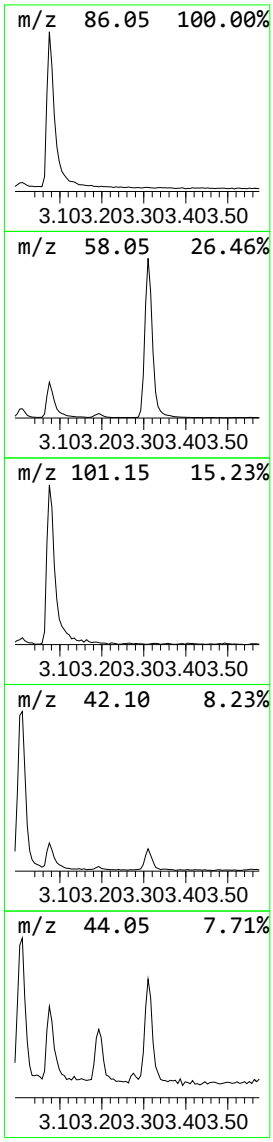
****

Peak Number 1 Triethylamine Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.075	3.94 ng/ul	321460	1,4-Dichlorobenzene-d4	7.851

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triethylamine	101	C6H15N	000121-44-8	91
2	Triethylamine	101	C6H15N	000121-44-8	91
3	Triethylamine	101	C6H15N	000121-44-8	91
4	Triethylamine	101	C6H15N	000121-44-8	86
5	Tetraethyl ammonium fluoride	149	C8H20FN	000665-46-3	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012093.D
Acq On : 12 Oct 2022 22:41
Operator : CG/JU
Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

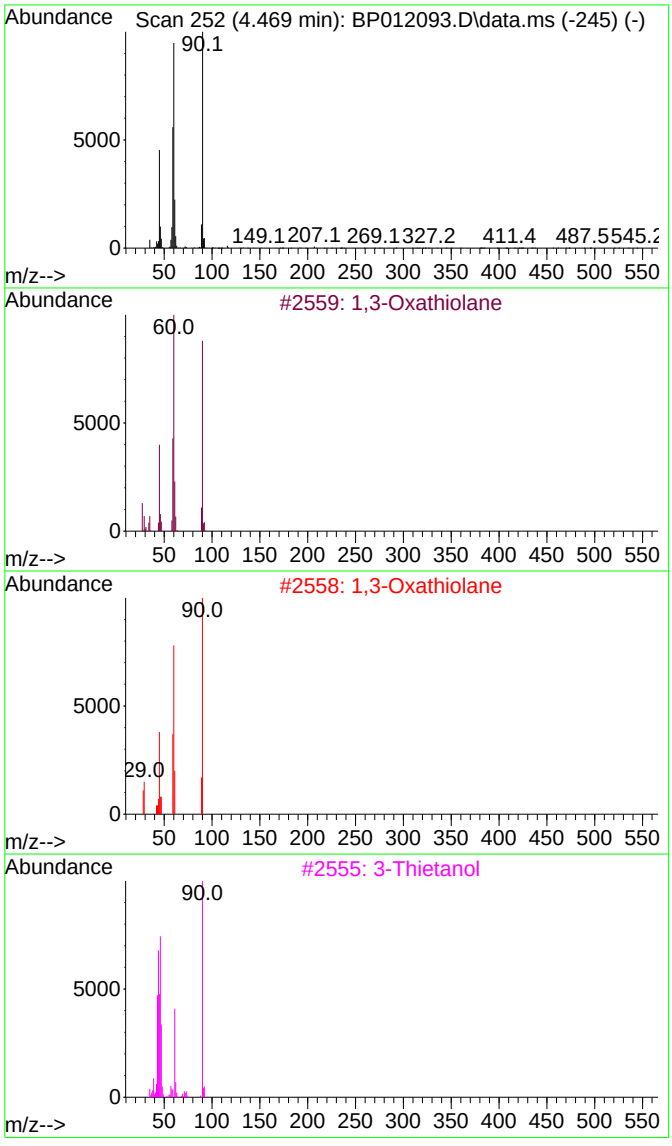
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

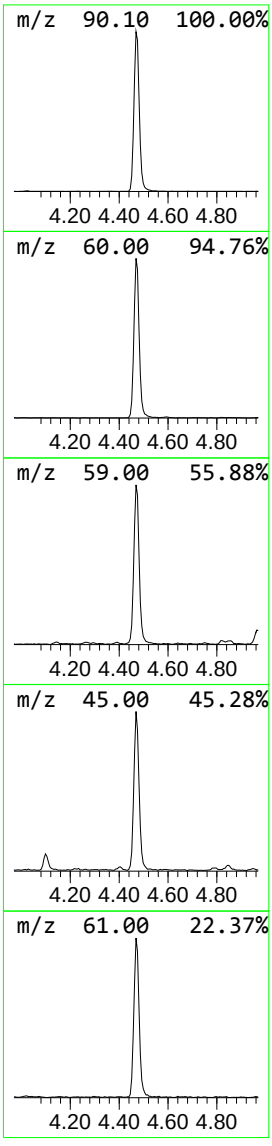
****

Peak Number 2 1,3-Oxathiolane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.469	5.81 ng/ul	473608	1,4-Dichlorobenzene-d4	7.851

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Oxathiolane	90	C3H6OS	002094-97-5	94
2	1,3-Oxathiolane	90	C3H6OS	002094-97-5	86
3	3-Thietanol	90	C3H6OS	010304-16-2	53
4	Heptanoic acid, 2-hydroxy-, meth...	160	C8H16O3	054340-91-9	53
5	Thiourea, methyl-	90	C2H6N2S	000598-52-7	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012093.D
Acq On : 12 Oct 2022 22:41
Operator : CG/JU
Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y7

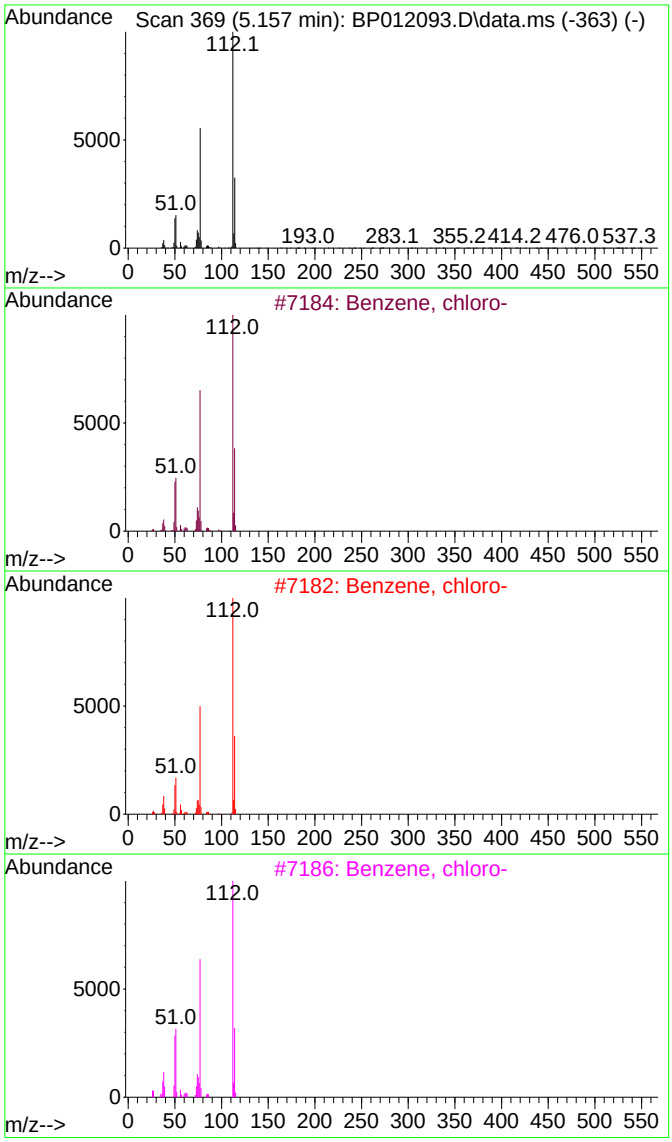
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

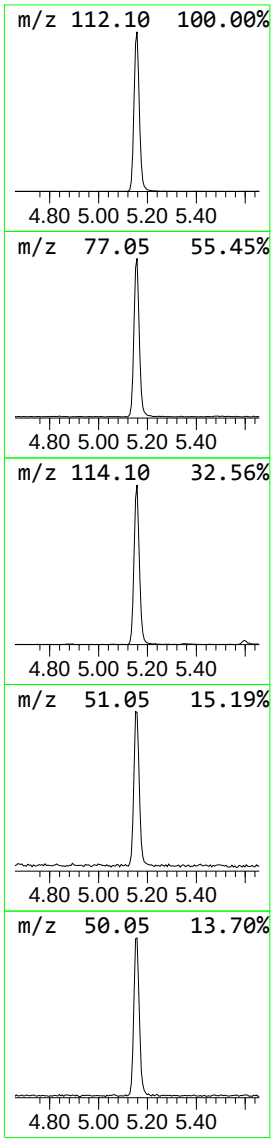
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, chloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.157	5.58 ng/ul	454856	1,4-Dichlorobenzene-d4	7.851

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, chloro-	112	C6H5Cl	000108-90-7	95
2	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
3	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
4	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
5	Benzene, chloro-	112	C6H5Cl	000108-90-7	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
Data File : BP012093.D
Acq On : 12 Oct 2022 22:41
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Instrument :
BNA_P
ClientSampleId :
CB8Y7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

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

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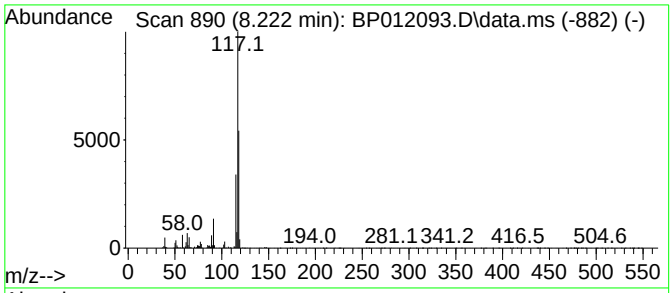
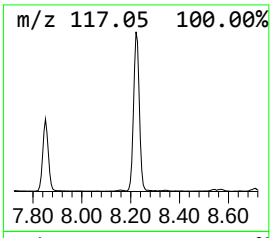
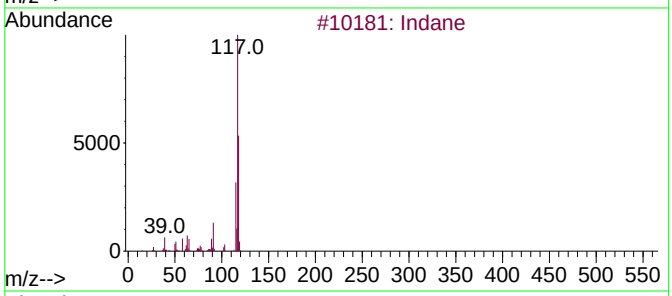
Peak Number 4 Indane Concentration Rank 12

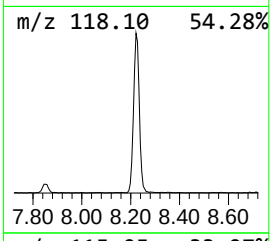
R.T.	EstConc	Area	Relative to ISTD	R.T.
8.222	4.63 ng/ul	377843	1,4-Dichlorobenzene-d4	7.851

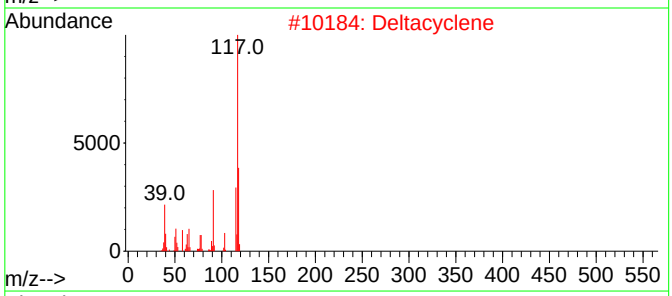
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	95
2	Deltacyclene	118	C9H10	007785-10-6	80
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	72
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	72
5	Benzene, 1-propenyl-	118	C9H10	000637-50-3	72

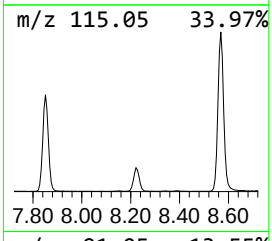


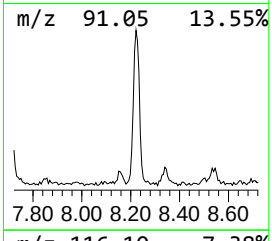


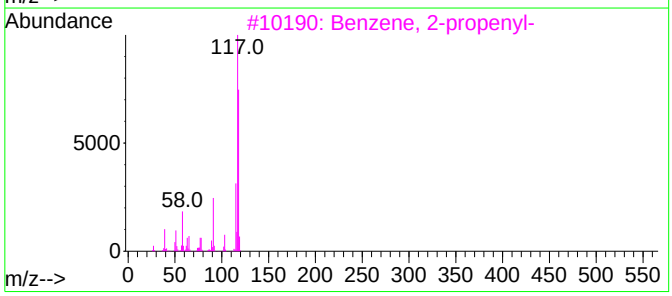


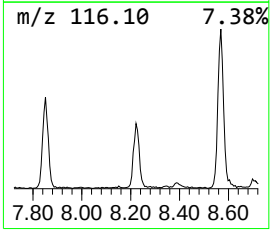












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Data File : BP012093.D
Acq On : 12 Oct 2022 22:41
Operator : CG/JU
Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y7

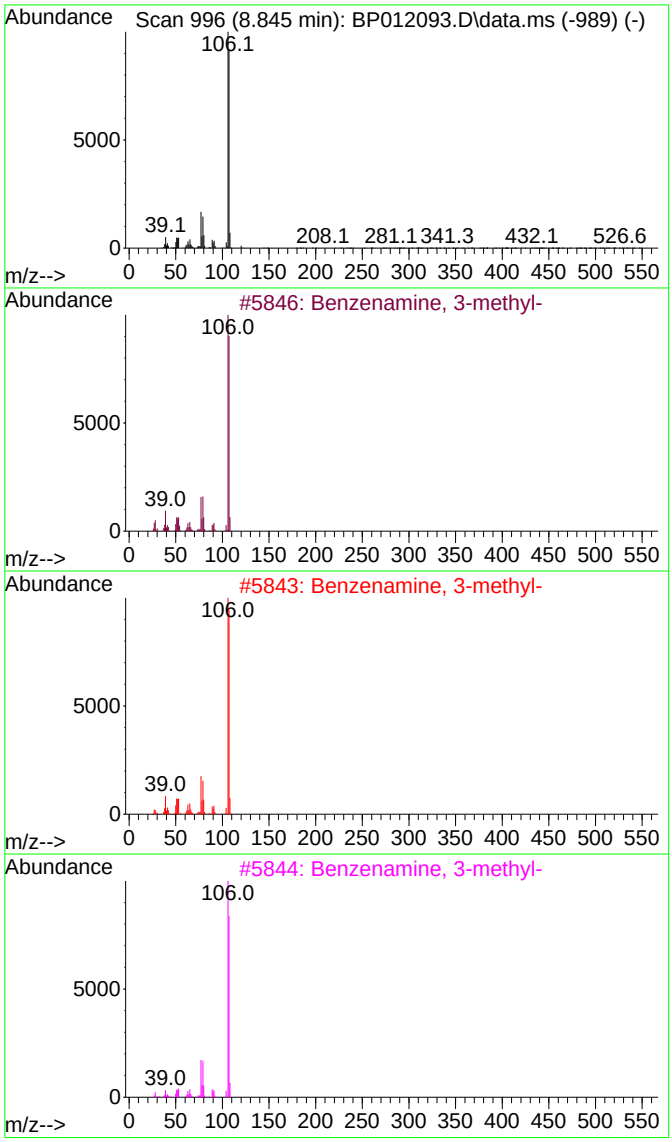
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

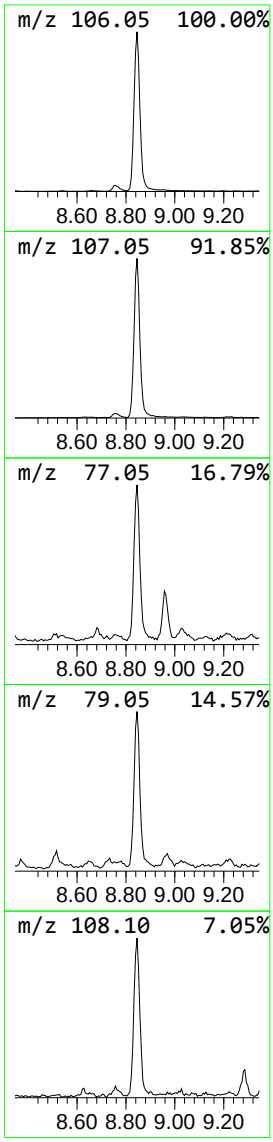
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzenamine, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.845	6.47 ng/ul	527429	1,4-Dichlorobenzene-d4	7.851

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	95
5	o-Toluidine	107	C7H9N	000095-53-4	94





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Operator : CG/JU
Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CB8Y7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

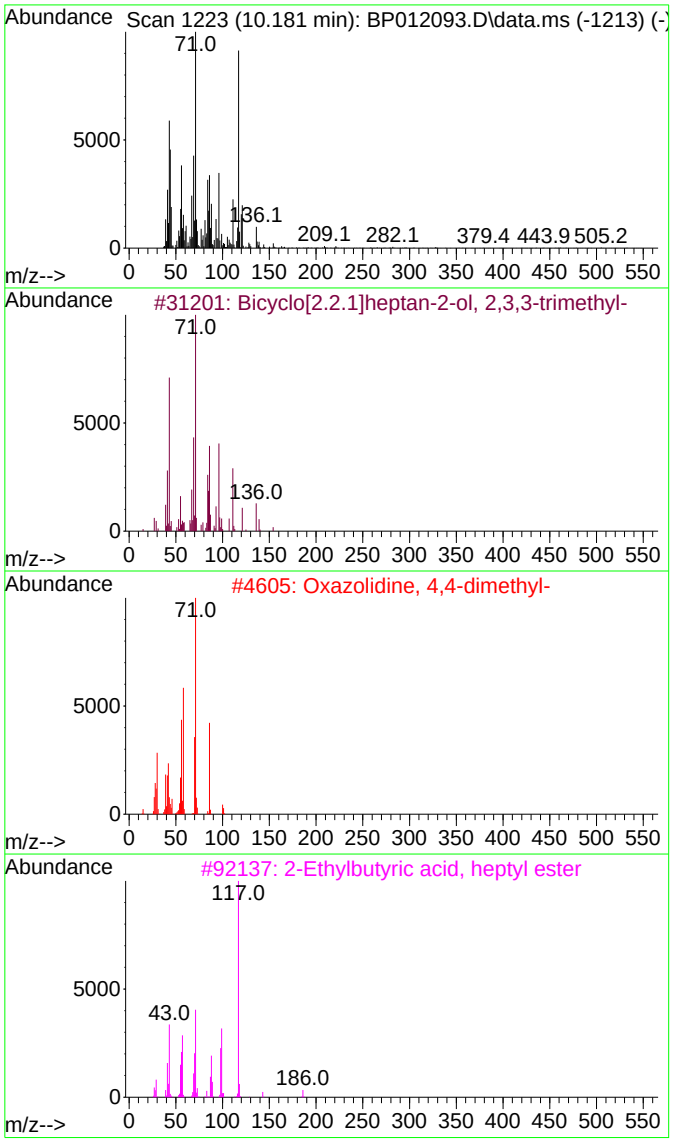
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

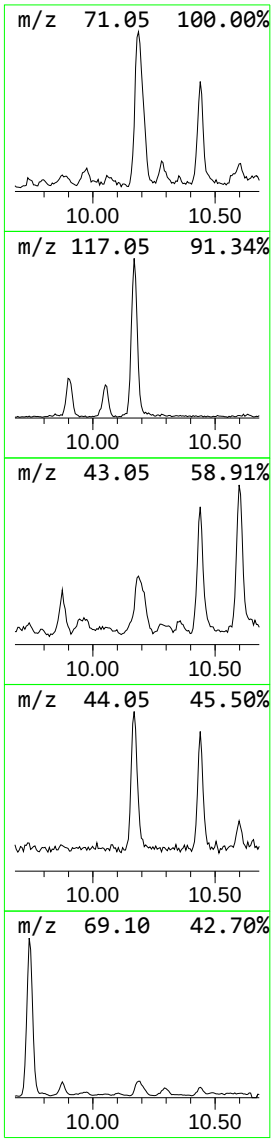
****

Peak Number 6 unknown-01 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.181	3.38 ng/ul	352703	Naphthalene-d8	10.657

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptan-2-ol, 2,3,3...	154	C10H18O	000465-31-6	30
2	Oxazolidine, 4,4-dimethyl-	101	C5H11NO	051200-87-4	27
3	2-Ethylbutyric acid, heptyl ester	214	C13H26O2	1000369-52-8	27
4	.+/-.-Tetrahydro-3-furanmethanol	102	C5H10O2	015833-61-1	18
5	N'-Hydroxypropimidamide	88	C3H8N2O	029335-36-2	14





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Operator : CG/JU
Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CB8Y7

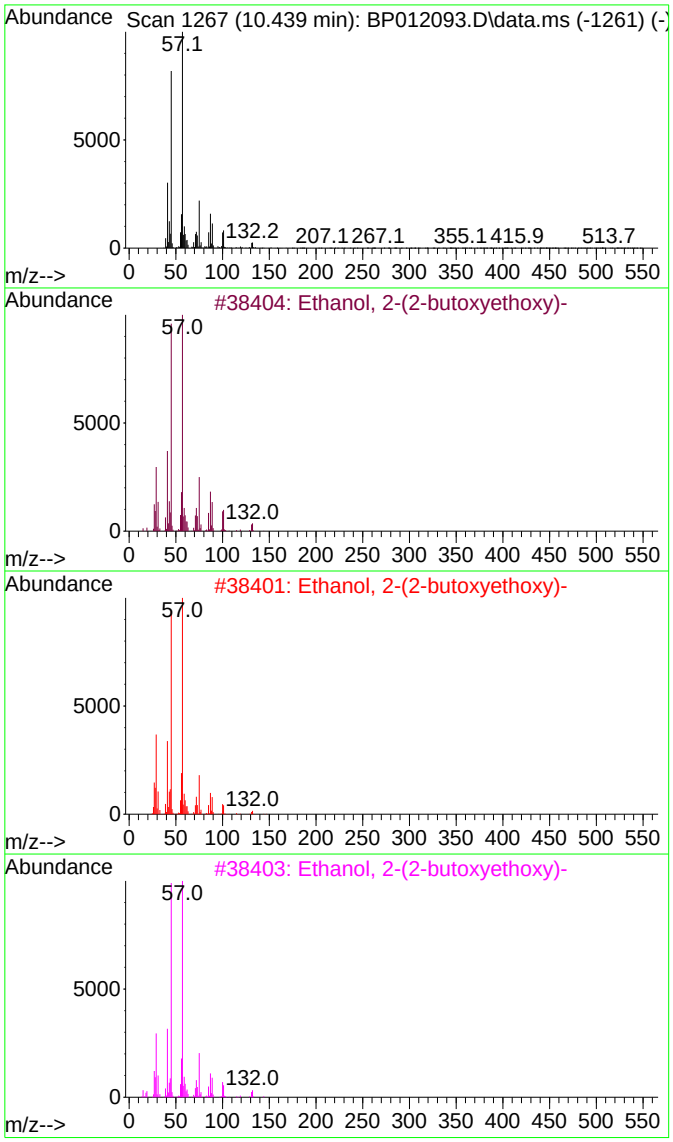
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

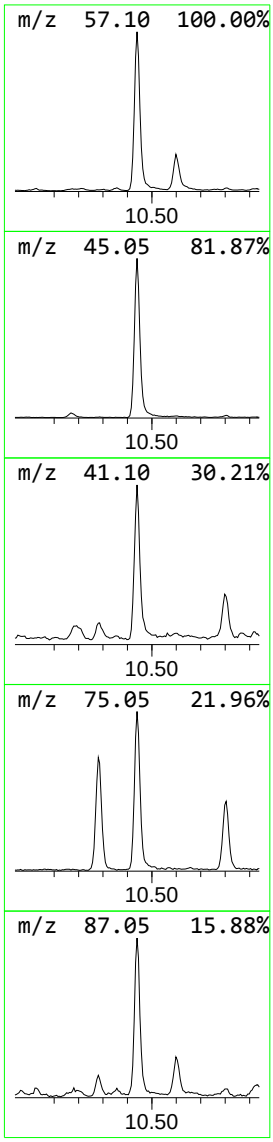
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.439	8.66 ng/ul	903003	Naphthalene-d8	10.657

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	91
2	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
4	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
5	Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	72





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Acq On : 12 Oct 2022 22:41
Operator : CG/JU
Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y7

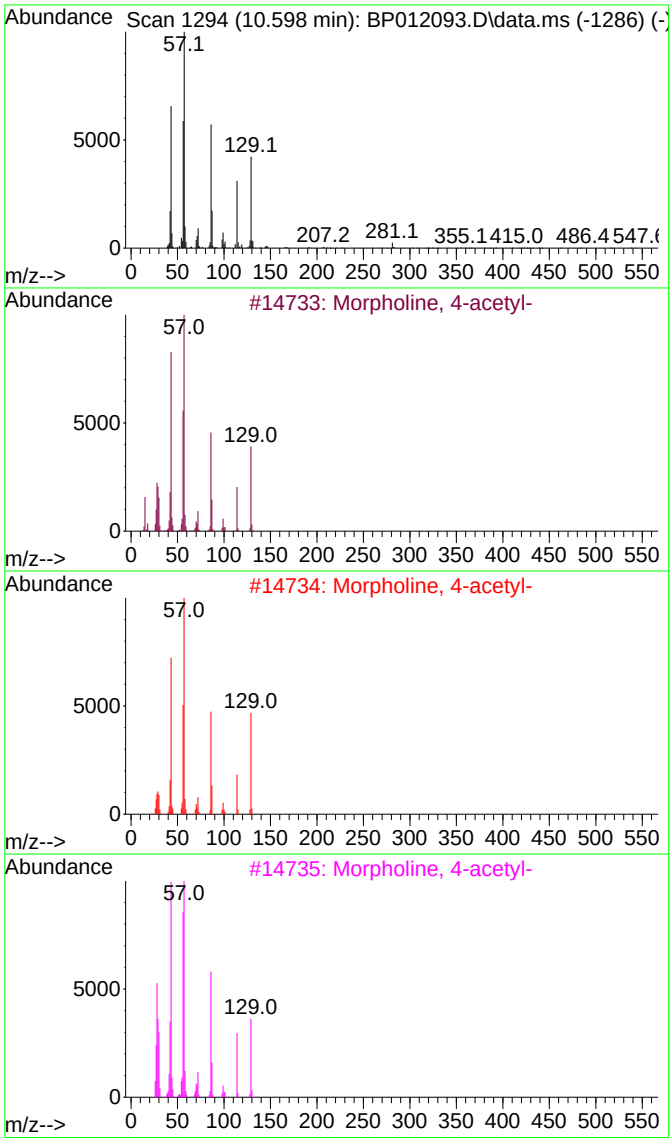
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

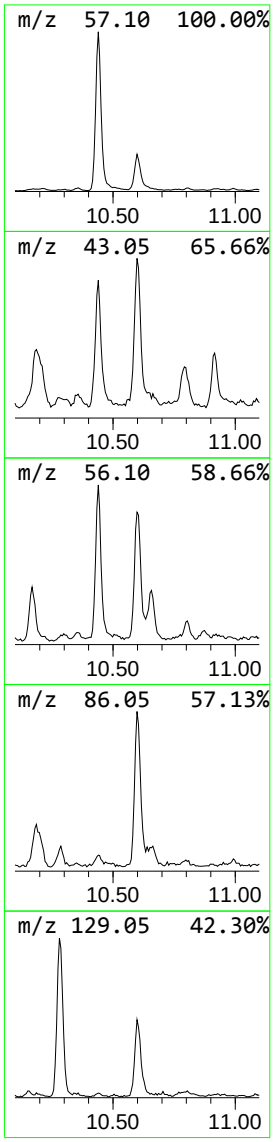
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Morpholine, 4-acetyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.598	2.20 ng/ul	229722	Naphthalene-d8	10.657

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Morpholine, 4-acetyl-	129	C6H11NO2	001696-20-4	91
2	Morpholine, 4-acetyl-	129	C6H11NO2	001696-20-4	90
3	Morpholine, 4-acetyl-	129	C6H11NO2	001696-20-4	53
4	2-Ethyl-oxetane	86	C5H10O	1010386-40-2	37
5	2-Methylpropionic acid, morpholide	157	C8H15NO2	1000307-31-7	37





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Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

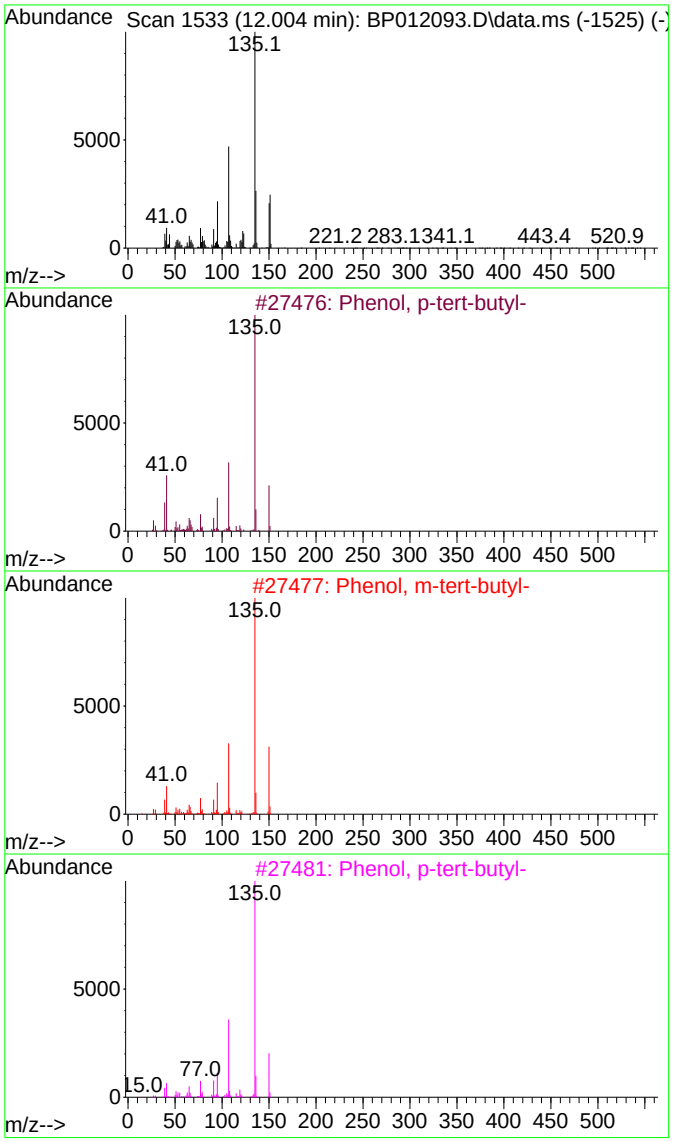
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

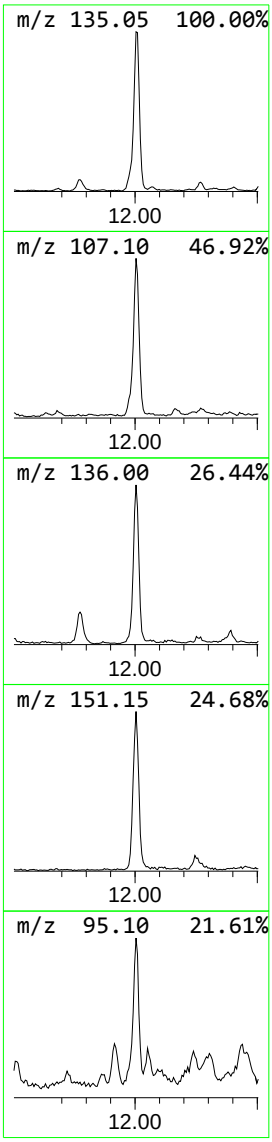
****

Peak Number 9 Phenol, p-tert-butyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.004	4.60 ng/ul	479580	Naphthalene-d8	10.657

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	74
2	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	68
3	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	68
4	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	68
5	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	53





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Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y7

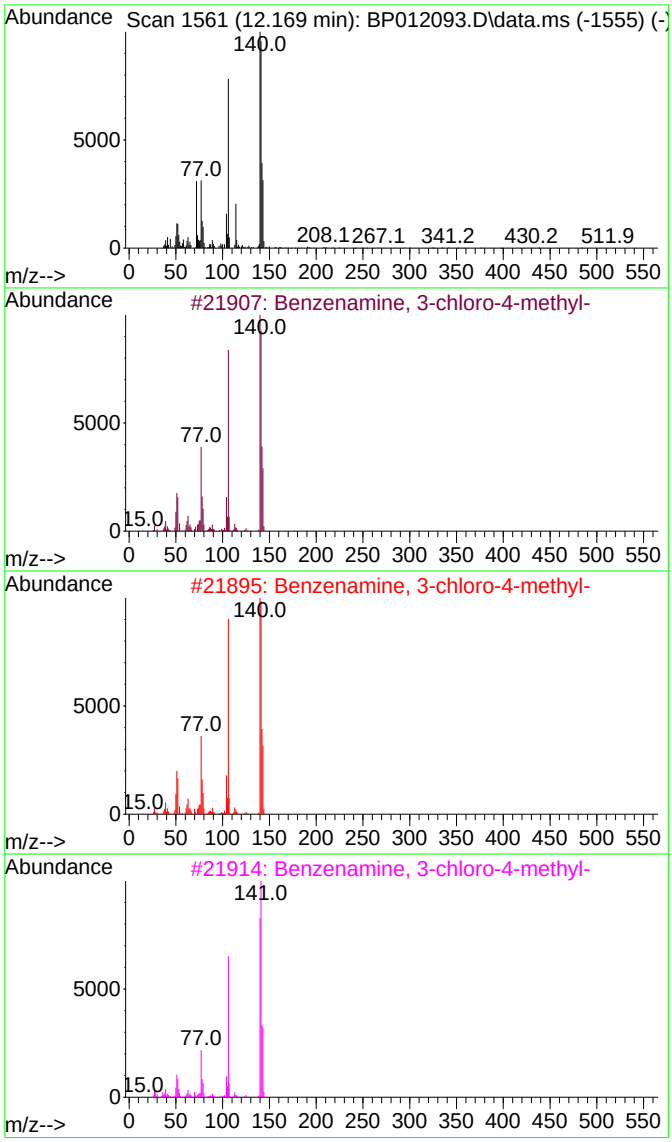
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

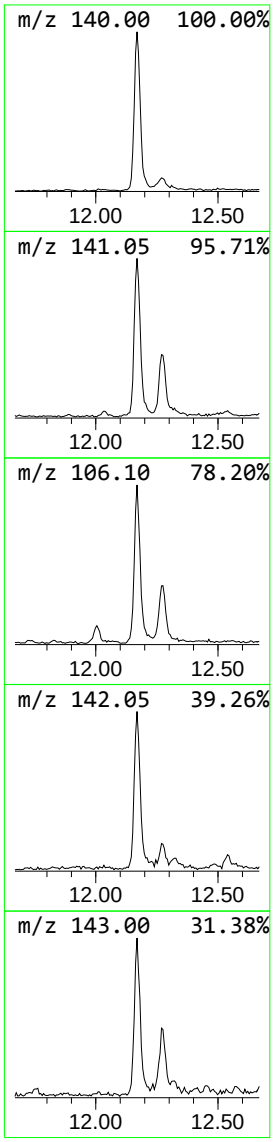
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzenamine, 3-chloro-4-met... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.169	2.23 ng/ul	232958	Naphthalene-d8	10.657

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	95
2	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	94
3	Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	87
4	Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	87
5	Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	76





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Acq On : 12 Oct 2022 22:41
Operator : CG/JU
Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y7

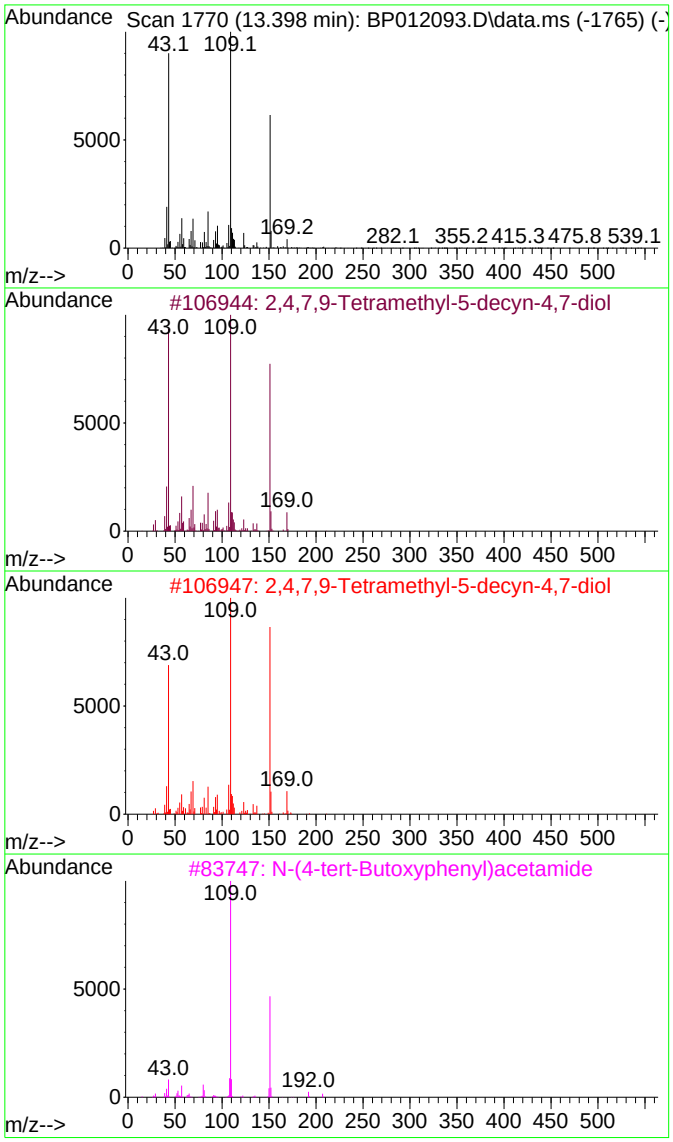
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

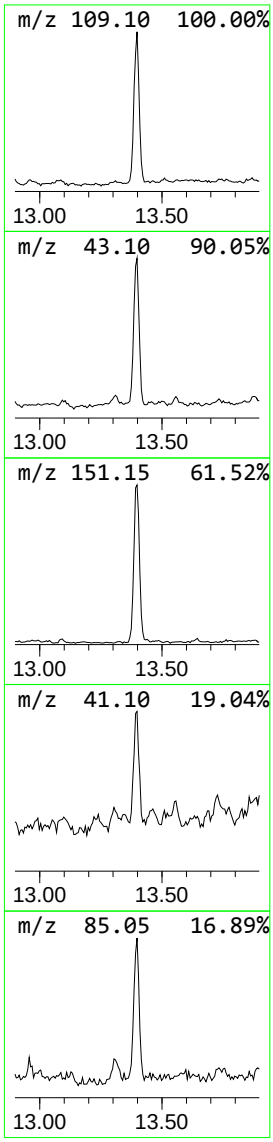
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.398	2.28 ng/ul	313584	Acenaphthene-d10	14.498

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91
2	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	90
3	N-(4-tert-Butoxyphenyl)acetamide	207	C12H17NO2	002109-73-1	59
4	Metacetamol	151	C8H9NO2	000621-42-1	59
5	Metacetamol	151	C8H9NO2	000621-42-1	59





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Operator : CG/JU
Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

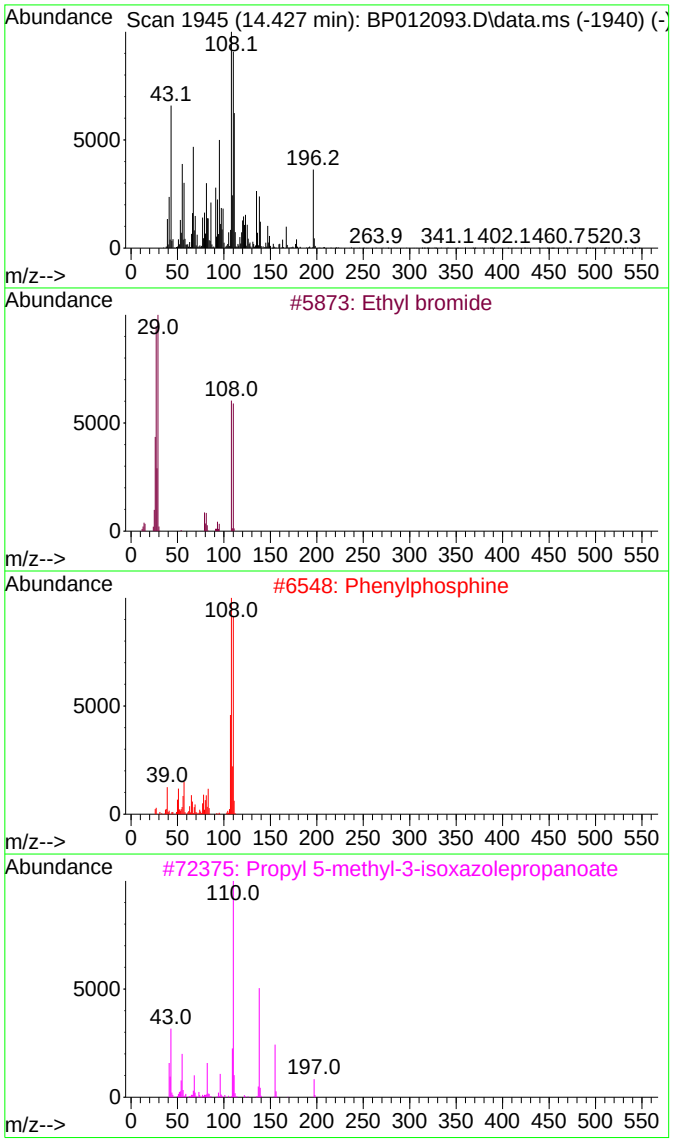
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

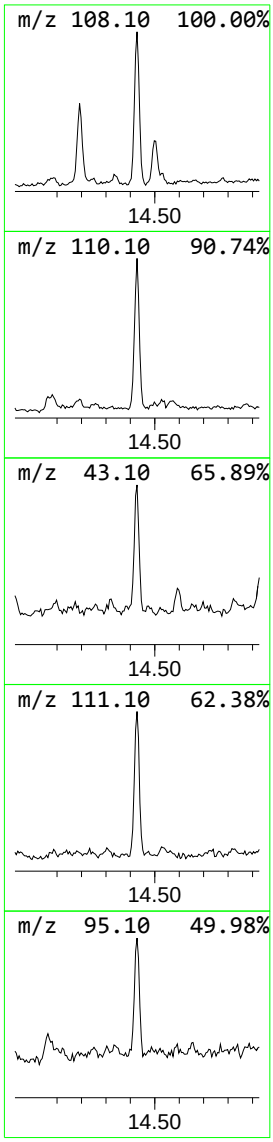
****

Peak Number 12 unknown-02 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.427	2.88 ng/ul	395641	Acenaphthene-d10	14.498

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethyl bromide	108	C2H5Br	000074-96-4	38
2	Phenylphosphine	110	C6H7P	000638-21-1	38
3	Propyl 5-methyl-3-isoxazolepropa...	197	C10H15NO3	1010099-85-3	14
4	Imidazole, 1,4,5-trimethyl-	110	C6H10N2	020185-22-2	14
5	Ethanone, 2-cyclopentyl-1-(1H-im...	178	C10H14N2O	069393-25-5	14





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Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CB8Y7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

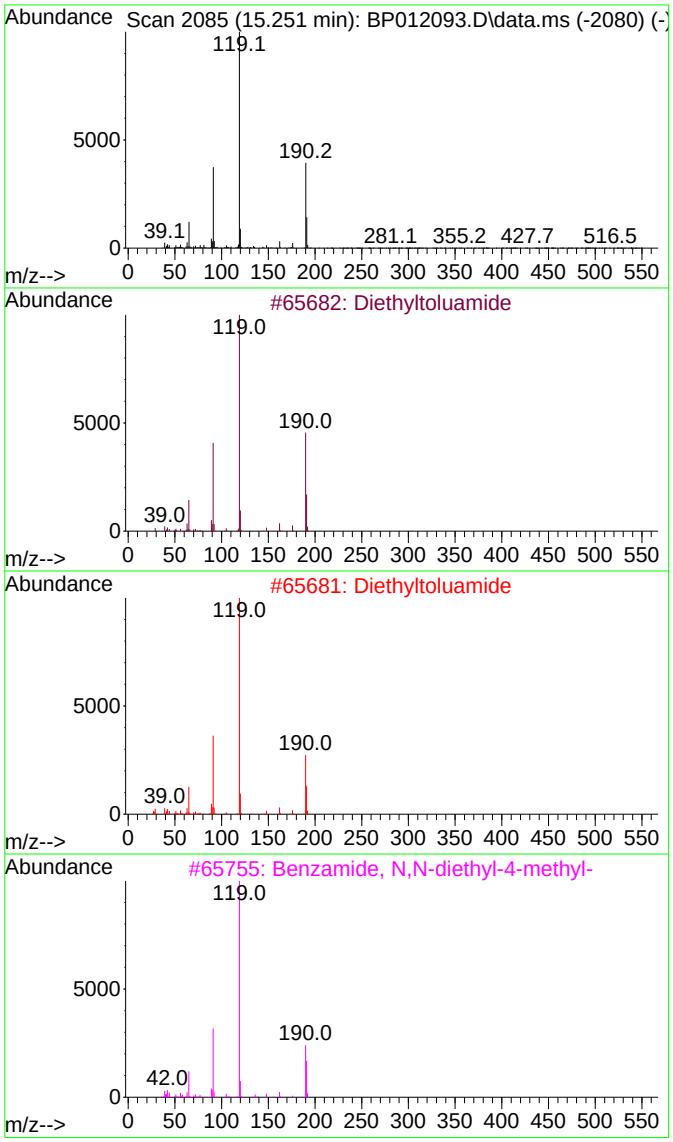
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

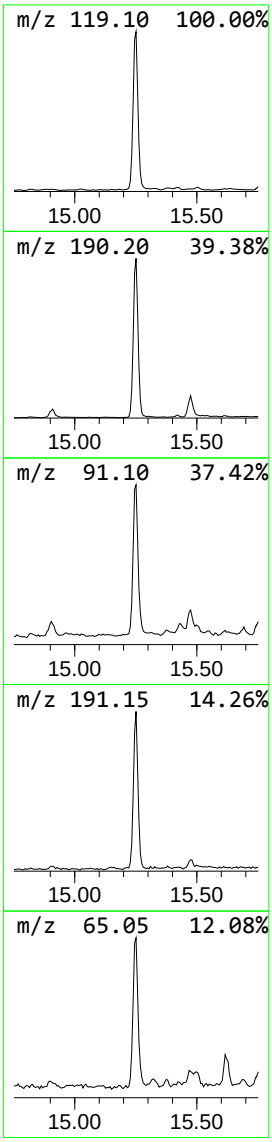
****

Peak Number 13 Diethyltoluamide Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.251	6.06 ng/ul	832961	Acenaphthene-d10	14.498

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethyltoluamide	191	C12H17NO	000134-62-3	97
2	Diethyltoluamide	191	C12H17NO	000134-62-3	95
3	Benzamide, N,N-diethyl-4-methyl-	191	C12H17NO	002728-05-4	93
4	Benzamide, 4-methyl-N-butyl-	191	C12H17NO	1000407-46-6	76
5	Benzamide, 3-methyl-N-methyl-N-p...	191	C12H17NO	1000421-46-2	76





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ALS Vial : 63 Sample Multiplier: 1

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ClientSampleId :
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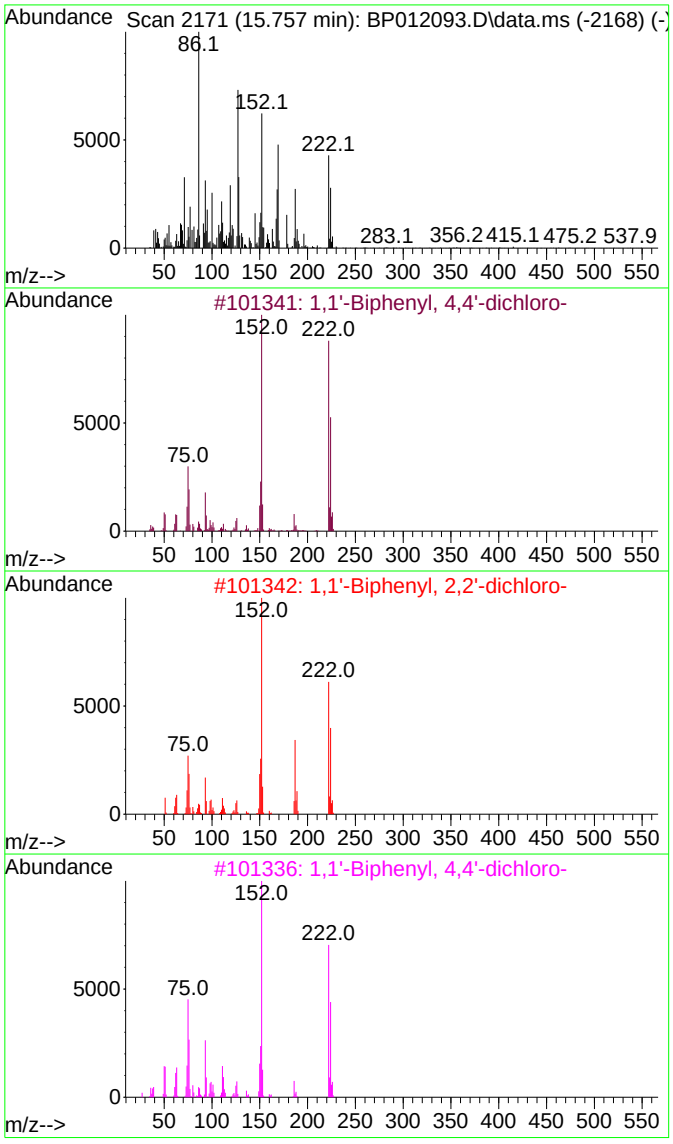
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

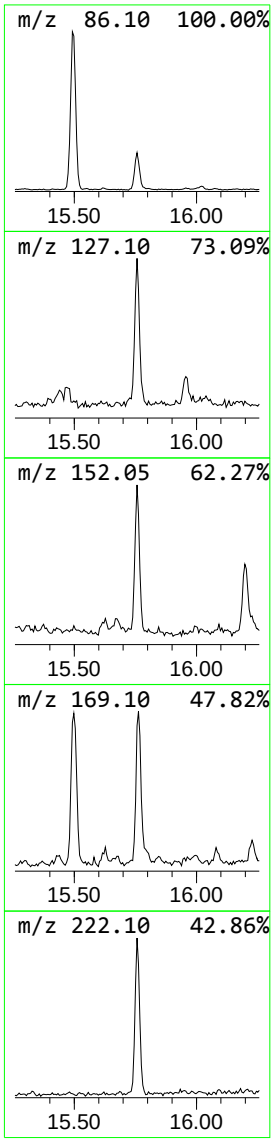
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 1,1'-Biphenyl, 4,4'-dichloro- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.757	2.46 ng/ul	337513	Acenaphthene-d10	14.498

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 4,4'-dichloro-	222	C12H8Cl2	002050-68-2	68
2	1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	50
3	1,1'-Biphenyl, 4,4'-dichloro-	222	C12H8Cl2	002050-68-2	45
4	1,1'-Biphenyl, 2,6-dichloro-	222	C12H8Cl2	033146-45-1	45
5	1,1'-Biphenyl, 3,3'-dichloro-	222	C12H8Cl2	002050-67-1	45





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ClientSampleId :
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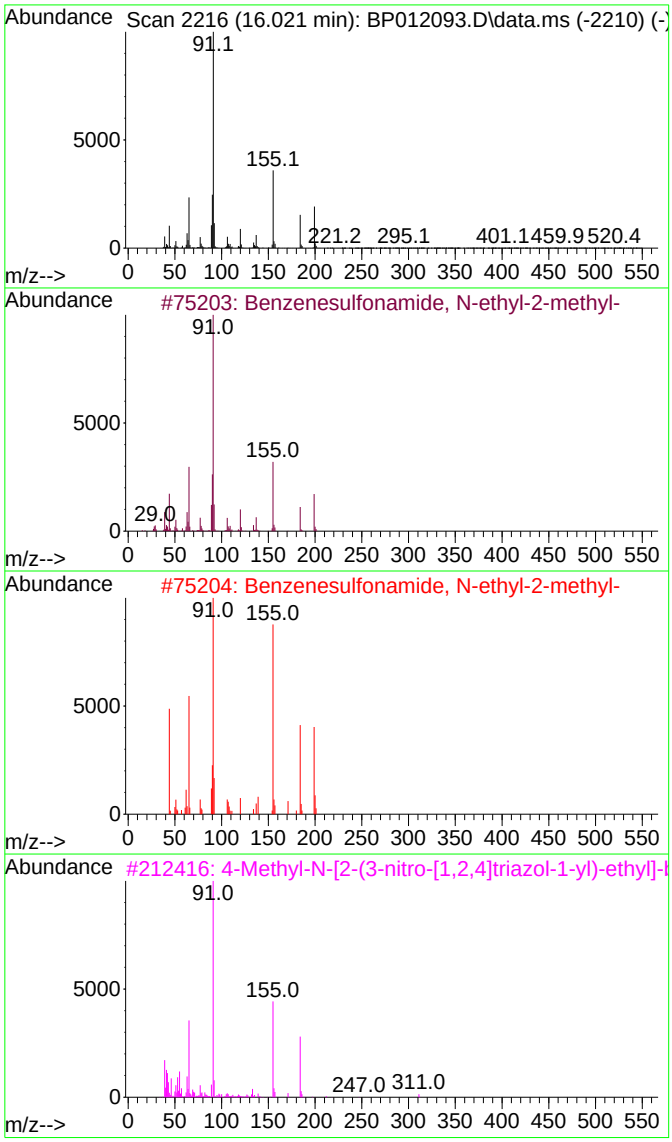
Quant Title : SVOA CALIBRATION

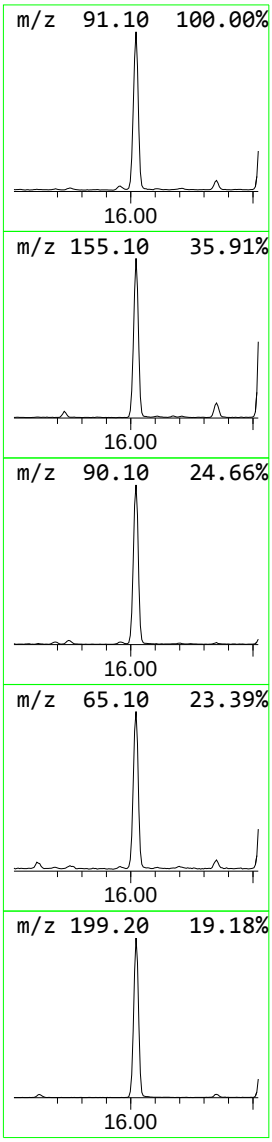
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Benzenesulfonamide, N-ethyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.021	15.58 ng/ul	2505900	Phenanthrene-d10	17.257

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	96
2	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	62
3	4-Methyl-N-[2-(3-nitro-[1,2,4]tr...	311	C11H13N5O4S	1000301-03-5	50
4	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	49
5	benzenesulfonamide, 4-methyl-N-[...	369	C16H19NO5S2	1000402-61-7	47





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Data File : BP012093.D
Acq On : 12 Oct 2022 22:41
Operator : CG/JU
Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y7

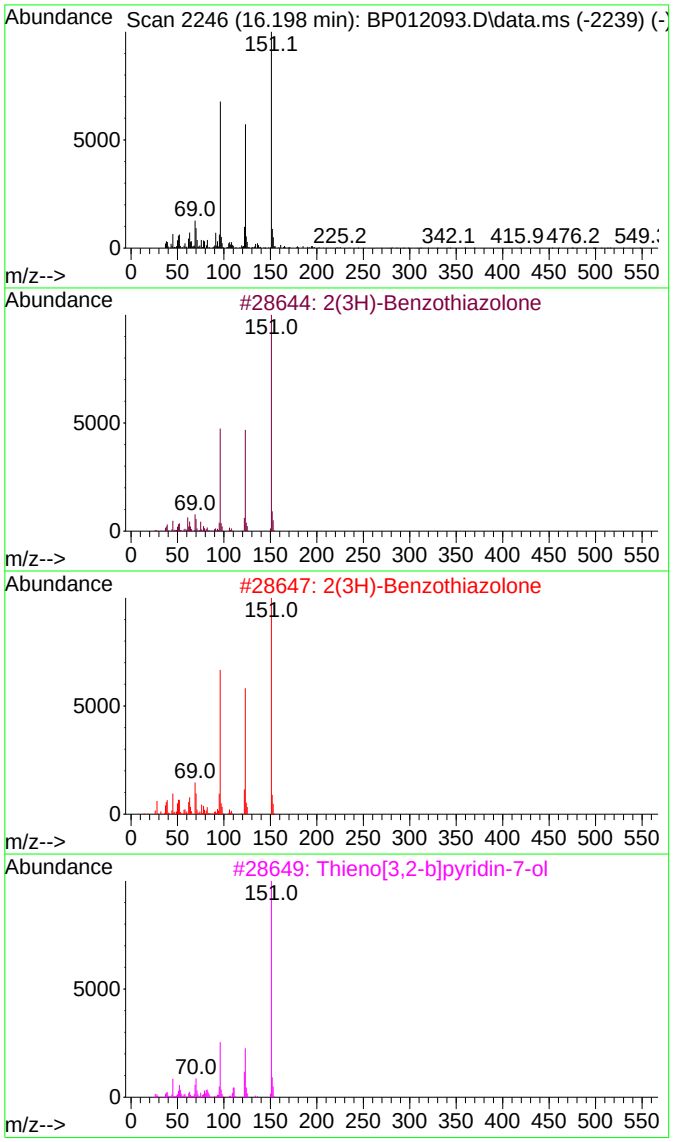
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

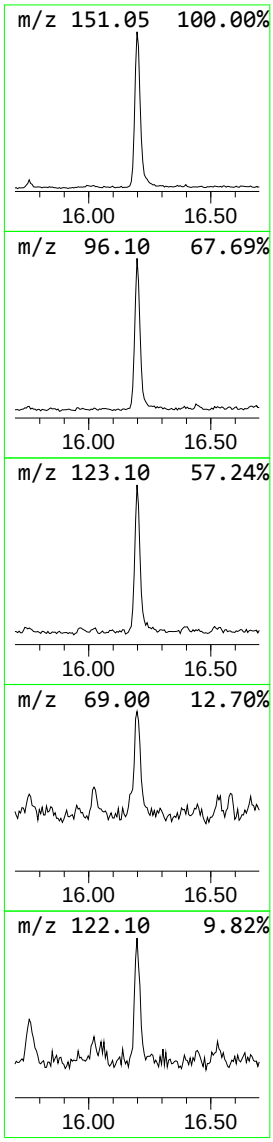
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 2(3H)-Benzothiazolone Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.198	2.61 ng/ul	419177	Phenanthrene-d10	17.257

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	94
2	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	94
3	Thieno[3,2-b]pyridin-7-ol	151	C7H5NOS	107818-20-2	58
4	1,2-Benzisothiazol-3(2H)-one	151	C7H5NOS	002634-33-5	47
5	6-Methyl-7-oxo-4,7-dihydro-triaz...	151	C5H5N5O	057250-39-2	47





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Acq On : 12 Oct 2022 22:41
Operator : CG/JU
Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y7

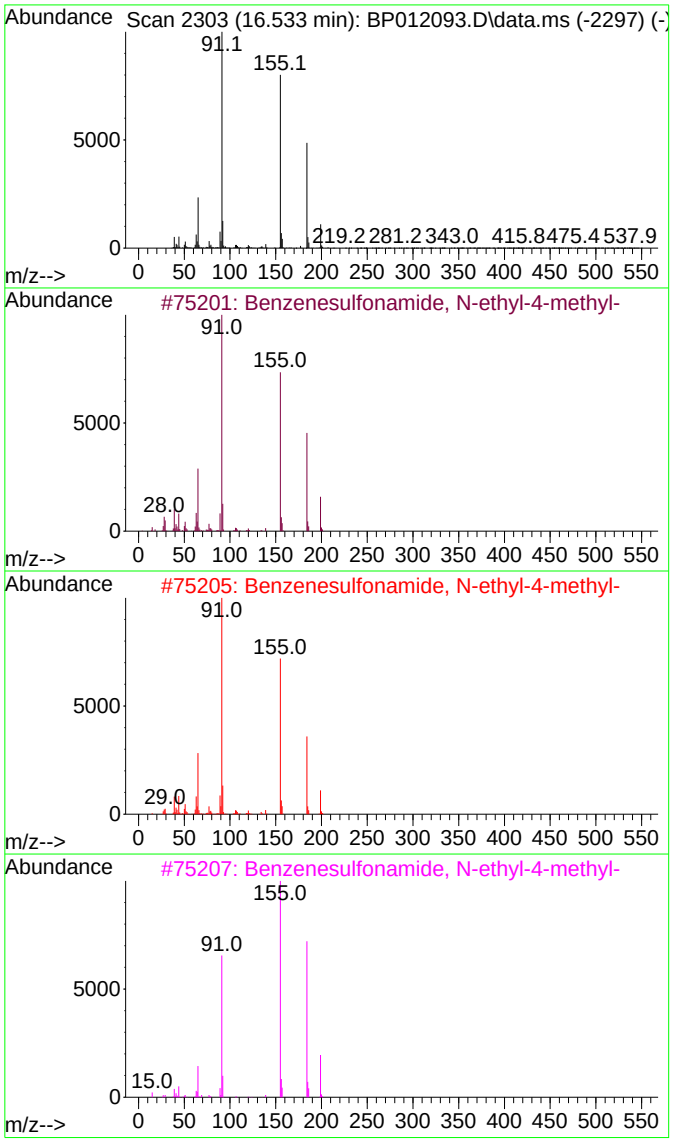
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

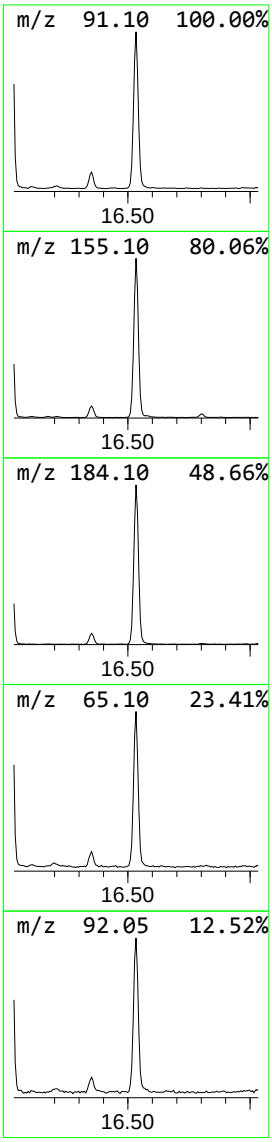
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Benzenesulfonamide, N-ethyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.533	8.94 ng/ul	1438380	Phenanthrene-d10	17.257

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	95
3	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	91
4	N-Hexyl-4-methylbenzenesulfonamide	255	C13H21NO2S	001143-01-7	80
5	N-isobutyl-4-methylbenzenesulfon...	227	C11H17NO2S	023705-38-6	80





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ALS Vial : 63 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

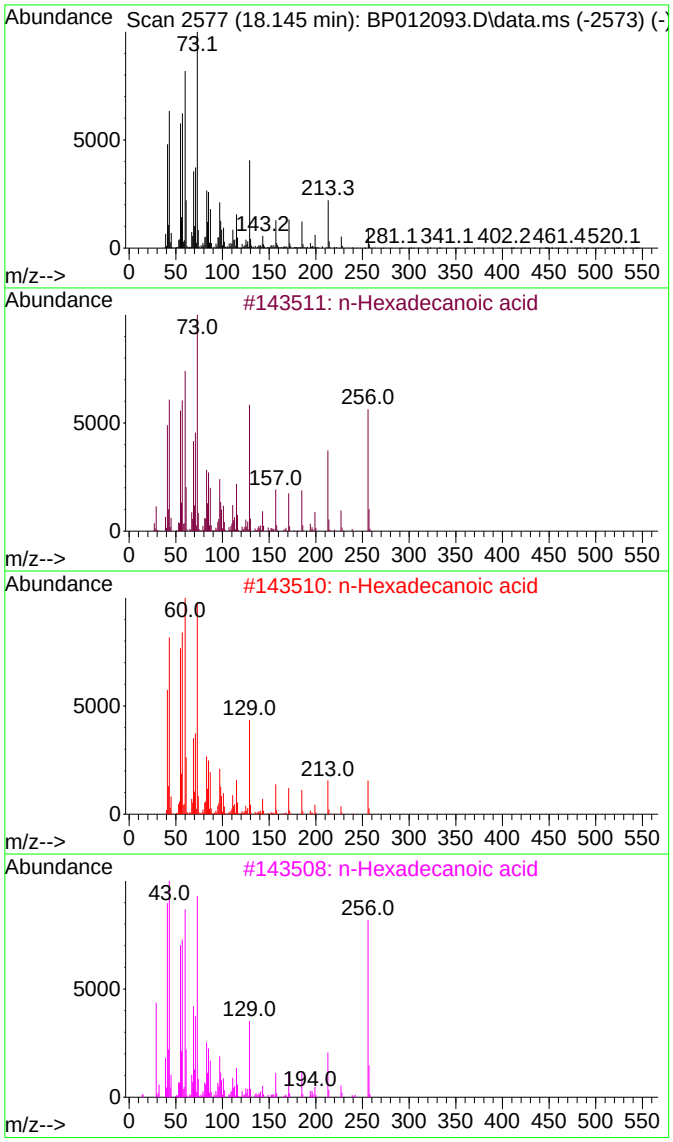
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

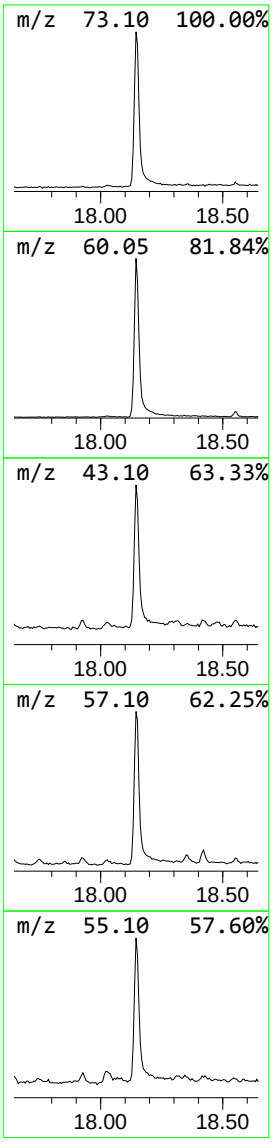
****

Peak Number 18 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.145	9.79 ng/ul	1574750	Phenanthrene-d10	17.257

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	94





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Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y7

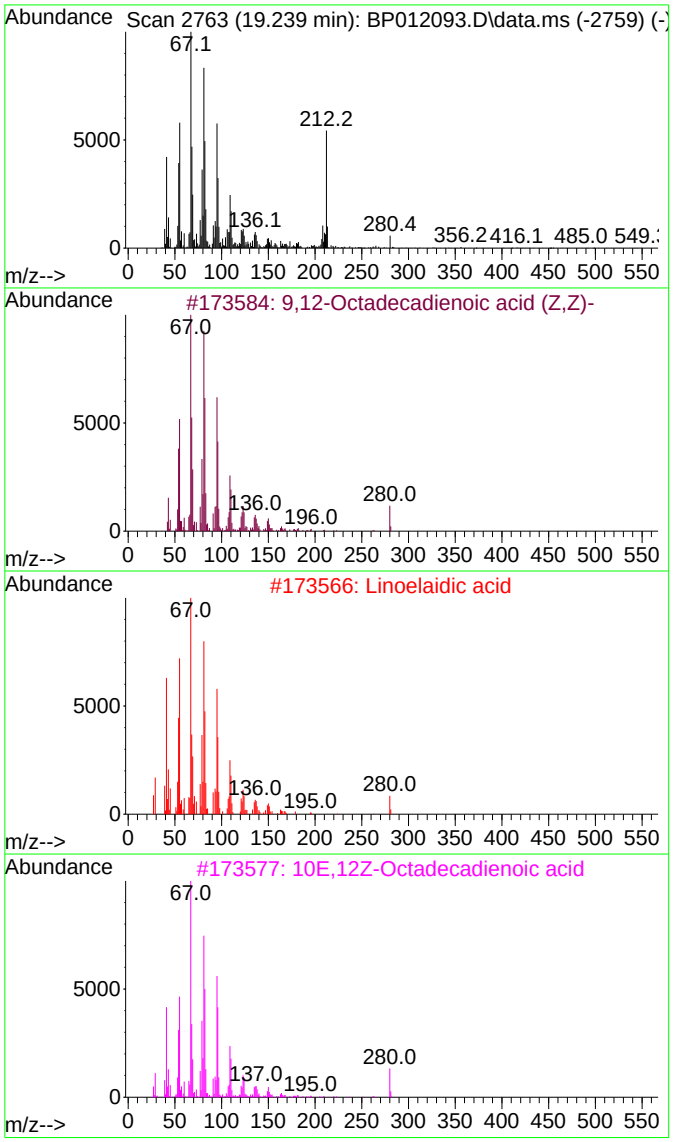
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

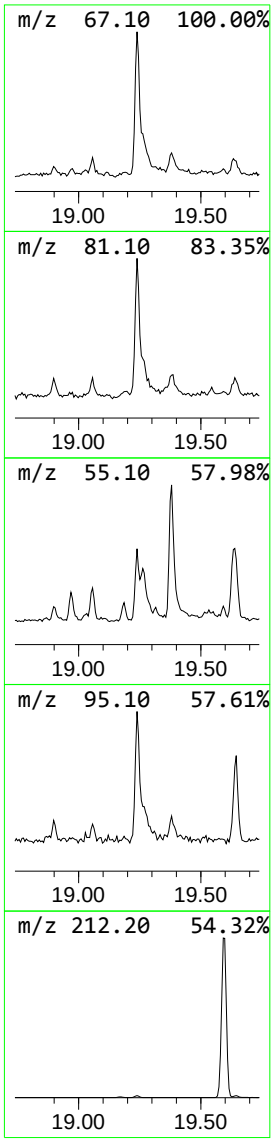
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 9,12-Octadecadienoic acid (... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.239	3.51 ng/ul	565201	Phenanthrene-d10	17.257

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid (Z,Z)-	280	C18H32O2	000060-33-3	99
2	Linoelaidic acid	280	C18H32O2	000506-21-8	98
3	10E,12Z-Octadecadienoic acid	280	C18H32O2	002420-56-6	98
4	(9E,11E)-Octadecadienoic acid	280	C18H32O2	000544-71-8	95
5	9,12-Octadecadienoic acid (Z,Z)-	280	C18H32O2	000060-33-3	91





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Operator : CG/JU
Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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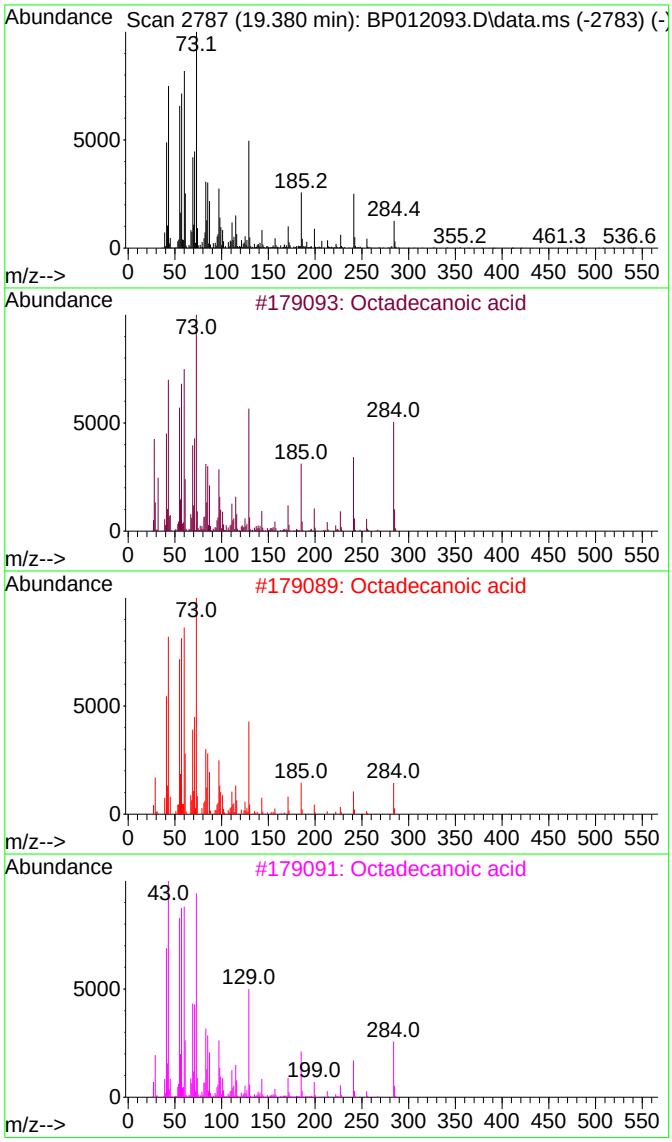
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

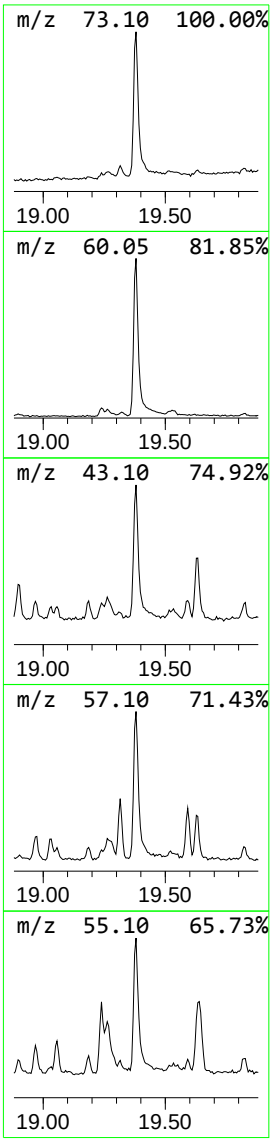
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.380	5.28 ng/ul	1008540	Chrysene-d12	21.351

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	99
4	Octadecanoic acid	284	C18H36O2	000057-11-4	98
5	Octadecanoic acid	284	C18H36O2	000057-11-4	93





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Acq On : 12 Oct 2022 22:41
Operator : CG/JU
Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y7

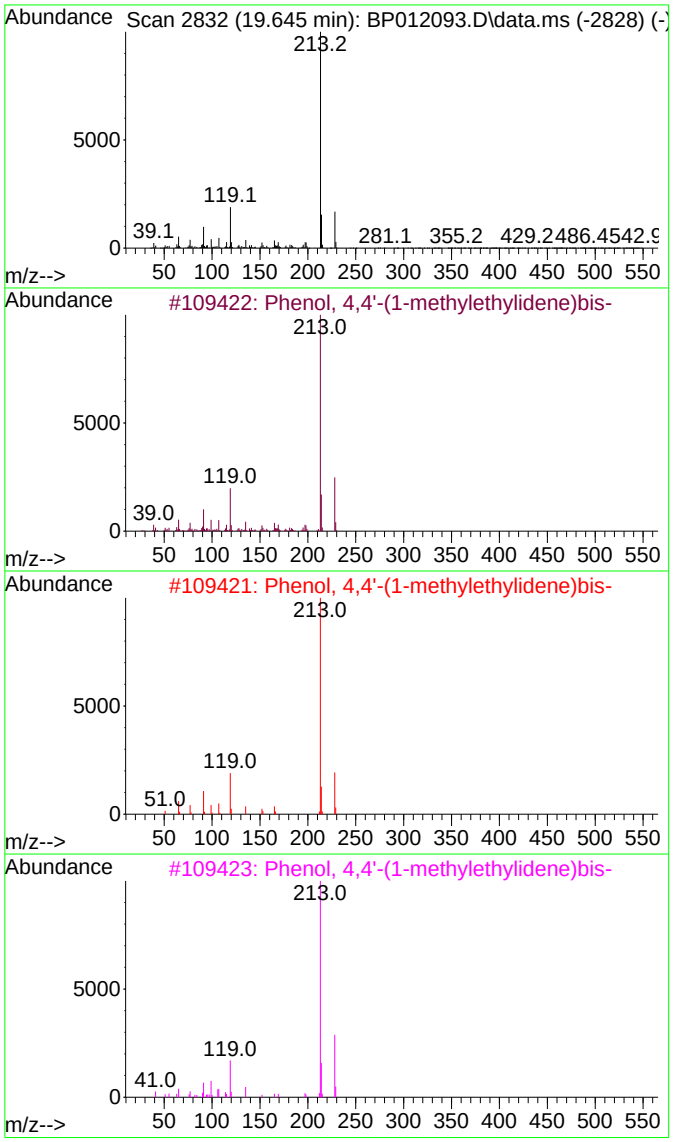
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

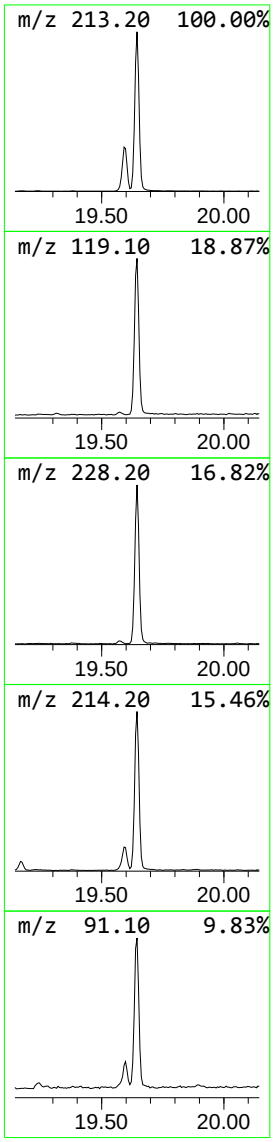
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 Phenol, 4,4'-(1-methylethyl... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.645	19.14 ng/ul	3658630	Chrysene-d12	21.351

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	96
2	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	96
3	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	90
4	4,4'-(1,3-Dimethylbutylidene)bis...	270	C18H22O2	006807-17-6	74
5	Diphenolic acid	286	C17H18O4	000126-00-1	64





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Acq On : 12 Oct 2022 22:41
Operator : CG/JU
Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Y7

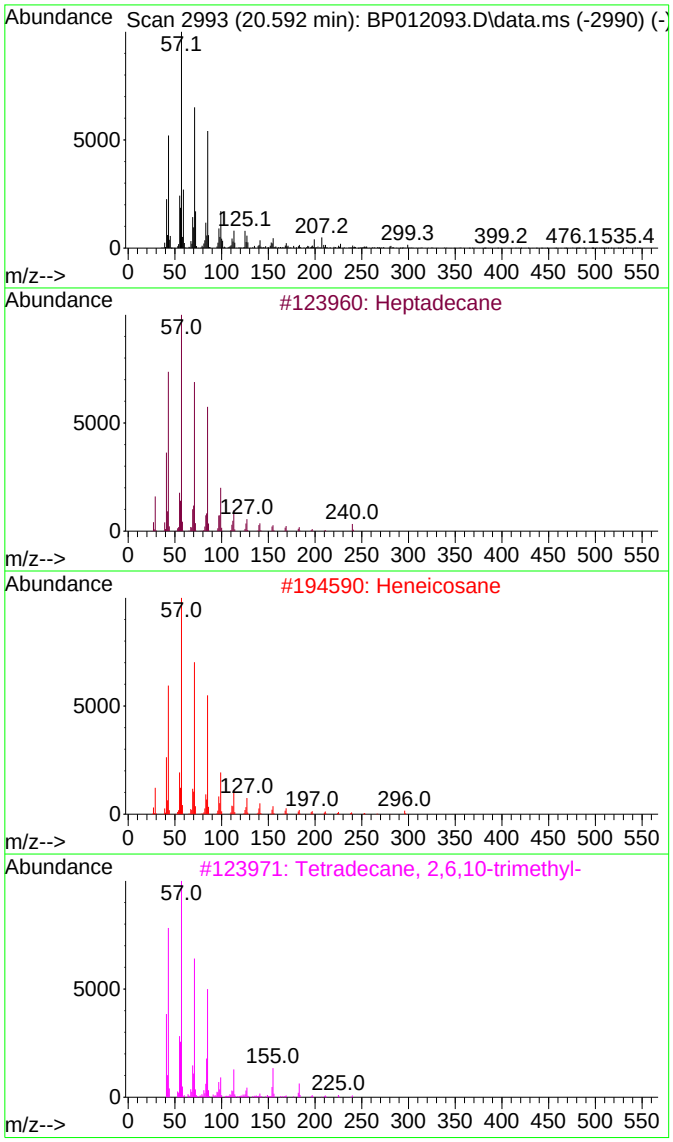
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

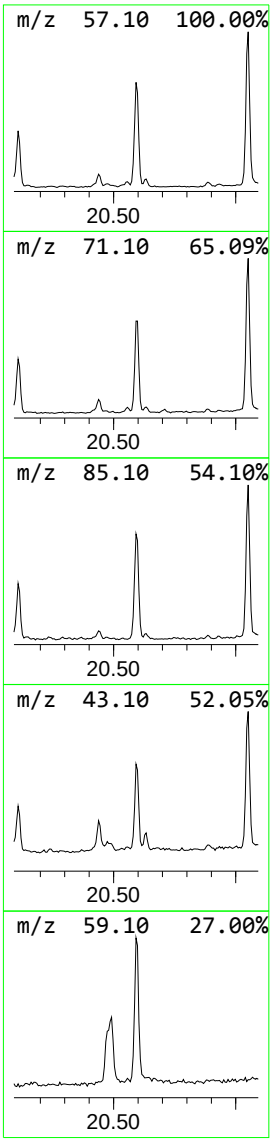
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.592	2.96 ng/ul	565440	Chrysene-d12	21.351

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	95
2	Heneicosane	296	C21H44	000629-94-7	76
3	Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	76
4	Tricosane	324	C23H48	000638-67-5	70
5	Octadecane	254	C18H38	000593-45-3	68





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Operator : CG/JU
Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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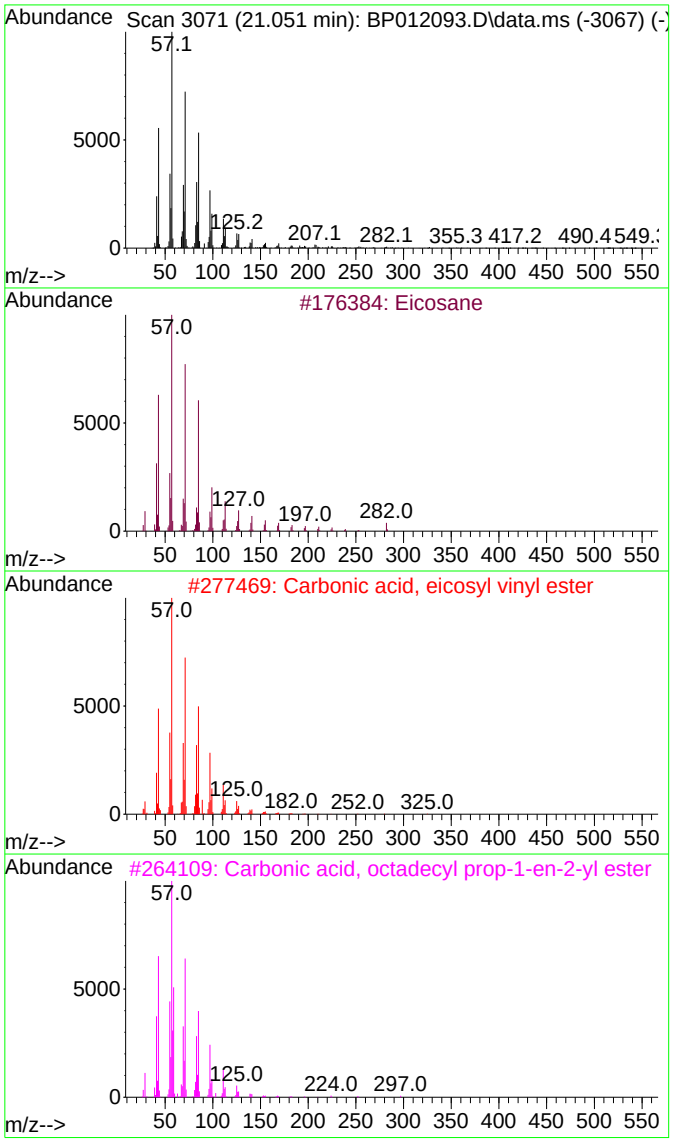
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

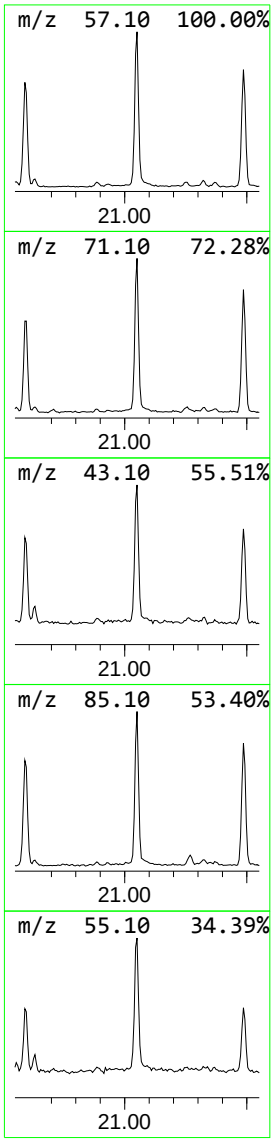
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.051	4.64 ng/ul	887305	Chrysene-d12	21.351

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	96
2	Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	95
3	Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	91
4	Carbonic acid, hexadecyl prop-1-...	326	C20H38O3	1000382-90-3	91
5	Carbonic acid, tetradecyl vinyl ...	284	C17H32O3	1000382-54-5	91





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Misc :
ALS Vial : 63 Sample Multiplier: 1

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ClientSampleId :
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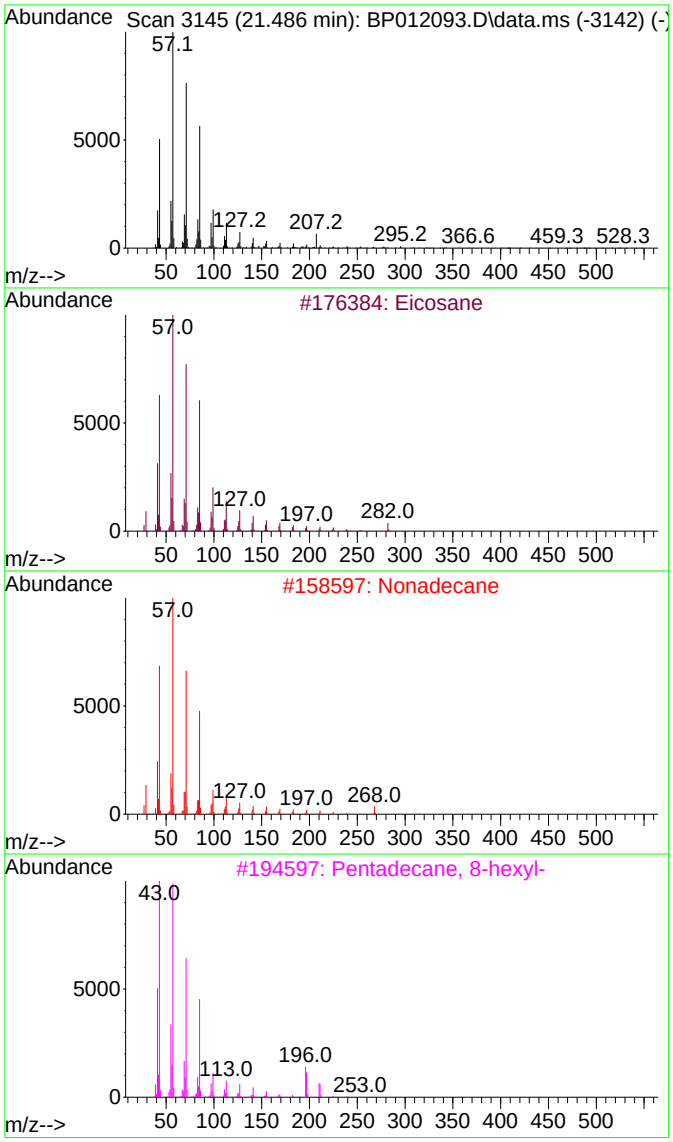
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

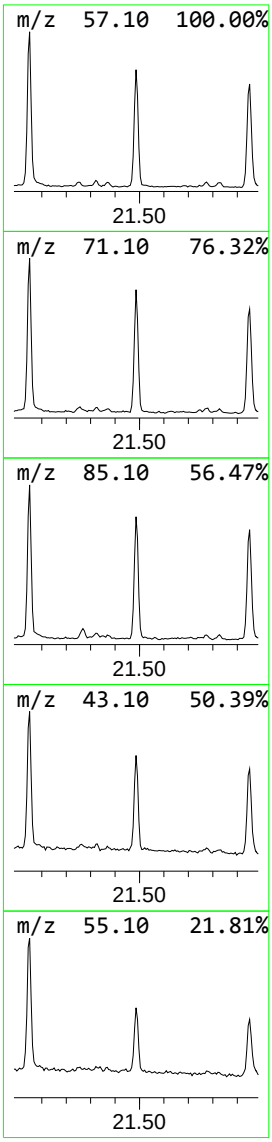
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.486	2.85 ng/ul	544309	Chrysene-d12	21.351

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	98
2	Nonadecane	268	C19H40	000629-92-5	95
3	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
4	Heneicosane	296	C21H44	000629-94-7	91
5	Hexacosane	366	C26H54	000630-01-3	91





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Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

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ClientSampleId :
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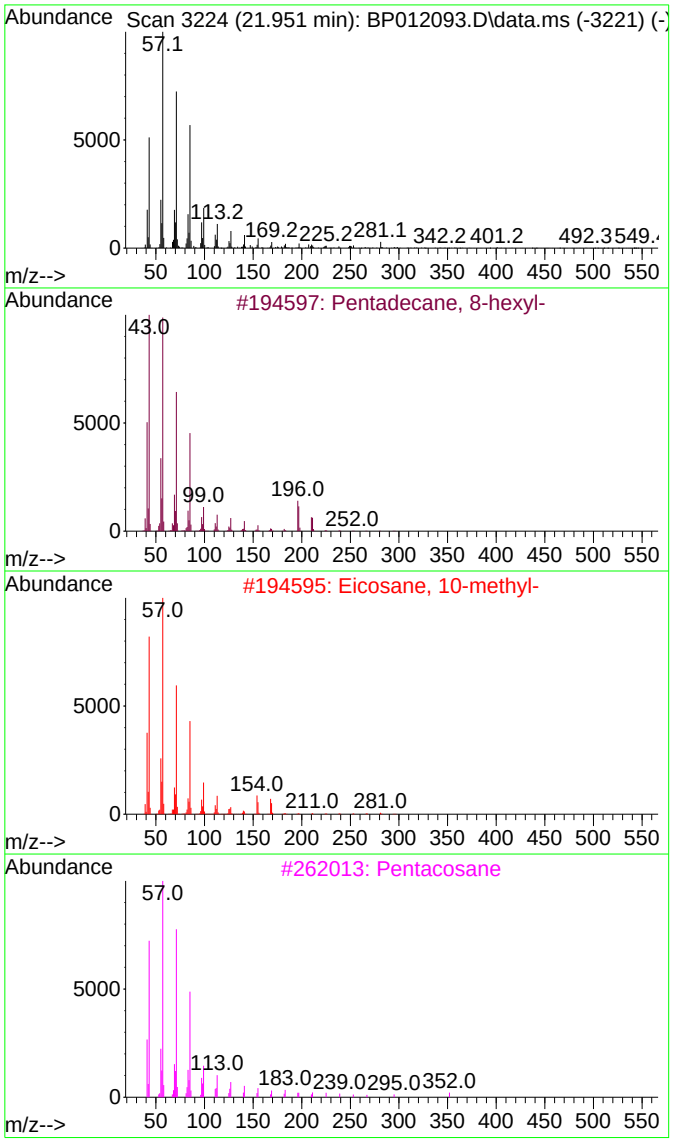
Quant Title : SVOA CALIBRATION

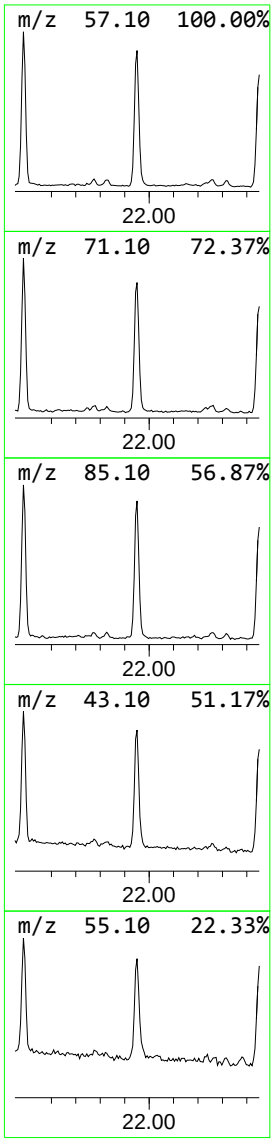
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.951	2.98 ng/ul	569794	Chrysene-d12	21.351

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93
2	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
3	Pentacosane	352	C25H52	000629-99-2	91
4	Eicosane	282	C20H42	000112-95-8	90
5	Heneicosane	296	C21H44	000629-94-7	90





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Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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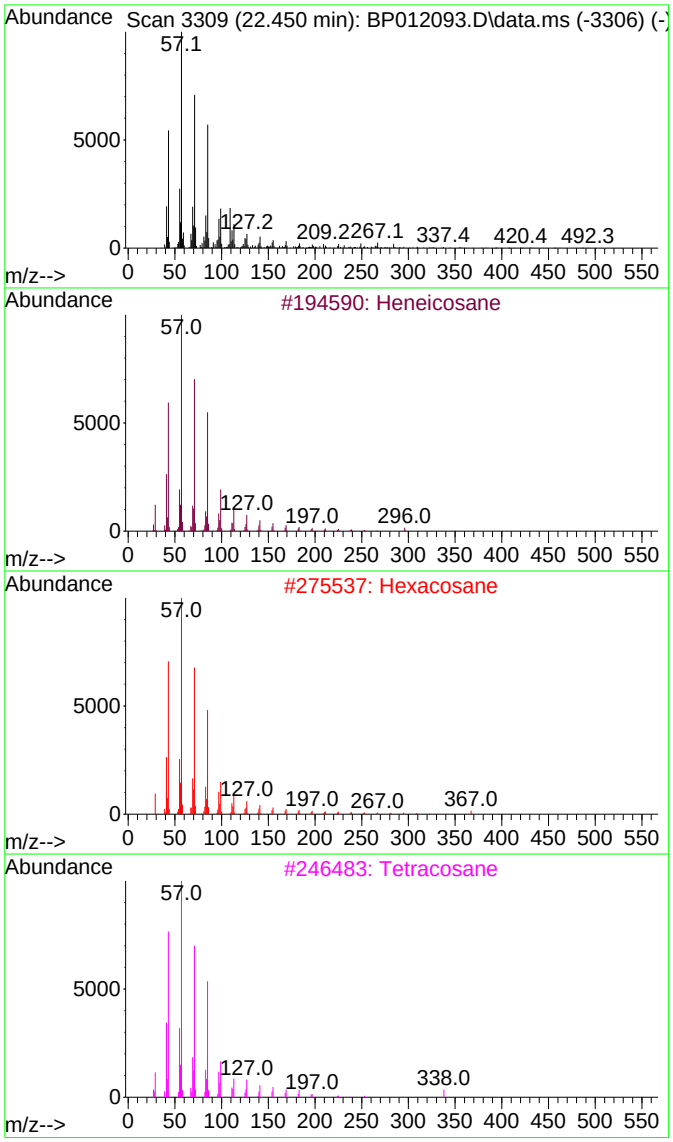
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

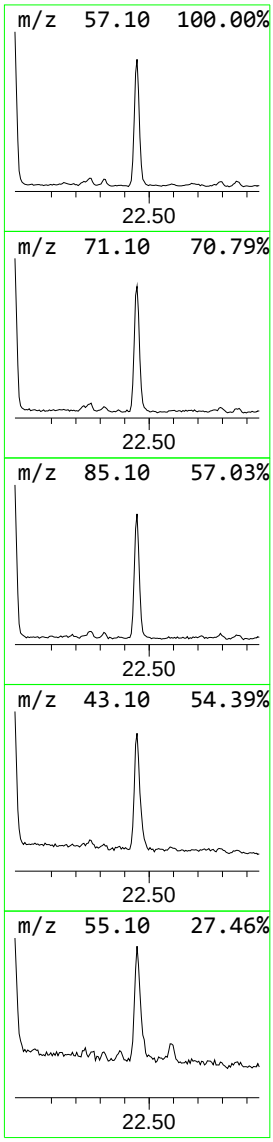
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.451	4.51 ng/ul	862175	Chrysene-d12	21.351

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	87
2	Hexacosane	366	C26H54	000630-01-3	87
3	Tetracosane	338	C24H50	000646-31-1	87
4	Pentacosane	352	C25H52	000629-99-2	83
5	Octacosane, 2-methyl-	408	C29H60	001560-98-1	81





Tentatively Identified Compound (LSC) summary

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Data File : BP012093.D
Acq On : 12 Oct 2022 22:41
Operator : CG/JU
Sample : N4976-11
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CB8Y7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.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
Triethylamine	3.075	3.9	ng/ul	321460	1	7.851	1631220	20.0
1,3-Oxathiolane	4.469	5.8	ng/ul	473608	1	7.851	1631220	20.0
Benzene, chloro-	5.157	5.6	ng/ul	454856	1	7.851	1631220	20.0
Indane	8.222	4.6	ng/ul	377843	1	7.851	1631220	20.0
Benzenamine, 3-...	8.845	6.5	ng/ul	527429	1	7.851	1631220	20.0
unknown-01	10.181	3.4	ng/ul	352703	2	10.657	2085510	20.0
Ethanol, 2-(2-b...	10.439	8.7	ng/ul	903003	2	10.657	2085510	20.0
Morpholine, 4-a...	10.598	2.2	ng/ul	229722	2	10.657	2085510	20.0
Phenol, p-tert-...	12.004	4.6	ng/ul	479580	2	10.657	2085510	20.0
Benzenamine, 3-...	12.169	2.2	ng/ul	232958	2	10.657	2085510	20.0
2,4,7,9-Tetrame...	13.398	2.3	ng/ul	313584	3	14.498	2747050	20.0
unknown-02	14.427	2.9	ng/ul	395641	3	14.498	2747050	20.0
Diethyltoluamide	15.251	6.1	ng/ul	832961	3	14.498	2747050	20.0
1,1'-Biphenyl, ...	15.757	2.5	ng/ul	337513	3	14.498	2747050	20.0
Benzenesulfonam...	16.021	15.6	ng/ul	2505900	4	17.257	3217520	20.0
2(3H)-Benzothia...	16.198	2.6	ng/ul	419177	4	17.257	3217520	20.0
Benzenesulfonam...	16.533	8.9	ng/ul	1438380	4	17.257	3217520	20.0
n-Hexadecanoic ...	18.145	9.8	ng/ul	1574750	4	17.257	3217520	20.0
9,12-Octadecadi...	19.239	3.5	ng/ul	565201	4	17.257	3217520	20.0
Octadecanoic acid	19.380	5.3	ng/ul	1008540	5	21.351	3823000	20.0
Phenol, 4,4'-(1...	19.645	19.1	ng/ul	3658630	5	21.351	3823000	20.0
(DEL) Alkane: S...	20.592	3.0	ng/ul	565440	5	21.351	3823000	20.0
(DEL) Alkane: S...	21.051	4.6	ng/ul	887305	5	21.351	3823000	20.0
(DEL) Alkane: S...	21.486	2.9	ng/ul	544309	5	21.351	3823000	20.0
(DEL) Alkane: S...	21.951	3.0	ng/ul	569794	5	21.351	3823000	20.0
(DEL) Alkane: S...	22.451	4.5	ng/ul	862175	5	21.351	3823000	20.0