LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110822\
Data File : BP012548.D
Acq On : 08 Nov 2022 12:21
Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

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-BP102522.M
Title : SVOA CALIBRATION

Signal : TIC: BP012548.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.034	5	8	13	rVB2	755610	1331891	1.63%	0.312%
2	3.146	20	27	38	rVB3	264011	584144	0.71%	0.137%
3	3.564	92	98	106	rBV3	1374145	2691951	3.29%	0.631%
4	3.728	114	126	128	rBV2	842519	2389867	2.92%	0.560%
5	3.811	132	140	148	rBV	8384993	19390129	23.72%	4.547%
6	3.881	148	152	159	rVB	1383134	2375545	2.91%	0.557%
7	3.964	160	166	170	rBV3	259350	635596	0.78%	0.149%
8	5.099	284	359	360	rBV4	2642852	32623663	39.91%	7.651%
9	5.287	370	391	393	rVB	2118641	10050305	12.29%	2.357%
10	5.517	418	430	433	rBV	518126	1365276	1.67%	0.320%
11	6.375	566	576	579	rVV	339335	843617	1.03%	0.198%
12	6.440	579	587	591	rVV	467054	1098190	1.34%	0.258%
13	7.011	651	684	695	rBV	10516532	29178864	35.70%	6.843%
14	7.122	695	703	708	rVV	1168137	2017859	2.47%	0.473%
15	7.311	728	735	745	rVV2	1144598	2628432	3.22%	0.616%
16	7.769	804	813	822	rVB4	1389513	2922166	3.57%	0.685%
17	7.864	822	829	843	rBV	3390387	8439549	10.32%	1.979%
18	8.093	855	868	872	rBV3	583156	1734988	2.12%	0.407%
19	8.499	919	937	941	rBV3	1638435	4552724	5.57%	1.068%
20	8.569	941	949	955	rVB2	1791986	3517626	4.30%	0.825%
21	8.811	977	990	997	rBV2	872243	2417658	2.96%	0.567%
22	8.881	997	1002	1007	rVB2	404312	683303	0.84%	0.160%
23	8.946	1007	1013	1020	rVB	1458658	2393501	2.93%	0.561%
24	9.022	1020	1026	1033	rBV3	268559	676642	0.83%	0.159%
25	9.258	1040	1066	1070	rBV	1978982	9977327	12.21%	2.340%
26	9.346	1075	1081	1084	rBV4	366651	563357	0.69%	0.132%
27	9.387	1084	1088	1094	rVB4	768393	1435606	1.76%	0.337%
28	9.446	1094	1098	1103	rBV2	443405	617521	0.76%	0.145%
29	9.516	1103	1110	1117	rBV3	1070694	2278560	2.79%	0.534%
30	9.605	1120	1125	1129	rVB	376754	576209	0.70%	0.135%
31	9.663	1129	1135	1143	rVB	949804	1588903	1.94%	0.373%
32	9.763	1147	1152	1155	rBV	586238	983836	1.20%	0.231%
33	9.799	1155	1158	1163	rVB	654610	898272	1.10%	0.211%
34	10.040	1186	1199	1202	rBV2	1988229	6364785	7.79%	1.493%
35	10.081	1202	1206	1209	rBV3	702149	1202641	1.47%	0.282%
36	10.210	1224	1228	1236	rVB3	5023018	9419727	11.52%	2.209%

LSC Area Percent Report

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Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

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-BP102522.M
Title : SVOA CALIBRATION

37	10.299	1237	1243	1246	rBV2	529006	908063	1.11%	0.213%
38	10.369	1251	1255	1264	rVB4	724355	1538174	1.88%	0.361%
39	10.510	1274	1279	1285	rVB	3886607	6030800	7.38%	1.414%
40	10.575	1285	1290	1296	rBV	1700160	2799963	3.43%	0.657%
41	10.640	1296	1301	1305	rVV	1019355	1715310	2.10%	0.402%
42	10.687	1305	1309	1313	rVV2	470075	828469	1.01%	0.194%
43	10.769	1320	1323	1331	rVV2	486937	823018	1.01%	0.193%
44	10.852	1332	1337	1344	rVV4	566238	1057823	1.29%	0.248%
45	10.946	1345	1353	1368	rVB	1751519	4358673	5.33%	1.022%
46	11.105	1373	1380	1390	rVB8	426495	1156619	1.41%	0.271%
47	11.522	1396	1451	1454	rBV	7821204	81744015	100.00%	19.170%
48	11.628	1462	1469	1472	rVV2	1343717	2775599	3.40%	0.651%
49	11.663	1472	1475	1479	rVB2	1426869	1766372	2.16%	0.414%
50	11.746	1484	1489	1495	rVV5	969861	1933649	2.37%	0.453%
51	11.834	1501	1504	1508	rVB4	393574	532327	0.65%	0.125%
52	11.940	1517	1522	1533	rVB	2457262	4028007	4.93%	0.945%
53	12.093	1543	1548	1555	rVB	2184168	3730453	4.56%	0.875%
54	12.163	1556	1560	1568	rVB5	374176	722170	0.88%	0.169%
55	12.275	1568	1579	1584	rBV7	575134	1588440	1.94%	0.373%
56	12.510	1613	1619	1623	rBV	1264958	2058266	2.52%	0.483%
57	12.610	1631	1636	1642	rBV4	444496	1136168	1.39%	0.266%
58	12.675	1642	1647	1650	rVV3	492542	783336	0.96%	0.184%
59	12.828	1661	1673	1676	rBV	2492414	6315026	7.73%	1.481%
60	12.857	1676	1678	1683	rVB	618922	774316	0.95%	0.182%
61	12.910	1683	1687	1696	rBV7	390334	990862	1.21%	0.232%
62	13.016	1701	1705	1710	rBV2	439414	730446	0.89%	0.171%
63	13.269	1743	1748	1754	rBV3	1816257	2944688	3.60%	0.691%
64	13.334	1755	1759	1765	rBV5	384090	786260	0.96%	0.184%
65	13.493	1782	1786	1797	rVB5	654007	1441218	1.76%	0.338%
66	13.640	1807	1811	1816	rVB	865553	1194285	1.46%	0.280%
67	13.722	1821	1825	1828	rBV4	699739	999164	1.22%	0.234%
68	13.851	1840	1847	1851	rBV	3589759	5866535	7.18%	1.376%
69	13.969	1862	1867	1872	rBV2	1212614	2077203	2.54%	0.487%
70	14.122	1888	1893	1900	rVB	3954561	6281935	7.68%	1.473%
71	14.316	1920	1926	1935	rVB2	2785733	7002983	8.57%	1.642%
72	14.428	1940	1945	1950	rVV	2711210	3860150	4.72%	0.905%
73	14.534	1960	1963	1969	rVB3	755467	1179133	1.44%	0.277%
74	14.669	1982	1986	1988	rBV2	408315	615550	0.75%	0.144%
75	14.710	1988	1993	2003	rVB4	1367666	2919240	3.57%	0.685%
76	14.910	2020	2027	2031	rBV	1922642	3230839	3.95%	0.758%

LSC Area Percent Report

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Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

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-BP102522.M
Title : SVOA CALIBRATION

77	15.010	2039	2044	2048	rVB	1047271	1502771	1.84%	0.352%
78	15.428	2108	2115	2121	rBV2	5154504	8226498	10.06%	1.929%
79	15.510	2125	2129	2134	rVV4	693580	1306427	1.60%	0.306%
80	15.563	2134	2138	2143	rVB	1384060	1949867	2.39%	0.457%
81	15.793	2174	2177	2182	rBV2	415259	620231	0.76%	0.145%
82	15.869	2186	2190	2194	rVB2	1124351	1586790	1.94%	0.372%
83	16.128	2230	2234	2237	rVB	1262791	1399538	1.71%	0.328%
84	16.328	2265	2268	2275	rVB3	755679	1388993	1.70%	0.326%
85	16.410	2276	2282	2290	rVV2	3287345	4855784	5.94%	1.139%
86	17.192	2410	2415	2420	rVV	3083724	4350846	5.32%	1.020%
87	17.292	2427	2432	2443	rVB2	5452843	8939867	10.94%	2.097%
88	17.575	2477	2480	2487	rVB3	584142	830515	1.02%	0.195%
89	17.704	2498	2502	2511	rBV2	1865490	3427602	4.19%	0.804%
90	17.939	2539	2542	2546	rBV3	866915	1137591	1.39%	0.267%
91	18.104	2564	2570	2576	rBV	5597039	9042197	11.06%	2.121%
92	18.222	2588	2590	2595	rVB	545715	609443	0.75%	0.143%
93	18.445	2625	2628	2632	rBV	462647	632734	0.77%	0.148%
94	18.863	2696	2699	2703	rBV	765037	945830	1.16%	0.222%
95	19.322	2774	2777	2784	rBV	916856	1280628	1.57%	0.300%
96	19.533	2809	2813	2821	rBV	5847010	8431833	10.31%	1.977%
97	20.986	3056	3060	3071	rVB	1652552	2640839	3.23%	0.619%
98	21.292	3107	3112	3117	rVB	3208069	4248517	5.20%	0.996%
99	23.522	3484	3491	3503	rVB2	3547368	7361901	9.01%	1.726%
100	23.674	3511	3517	3527	rVB2	1988485	4024099	4.92%	0.944%

Sum of corrected areas: 426417118

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110822\
Data File : BP012548.D
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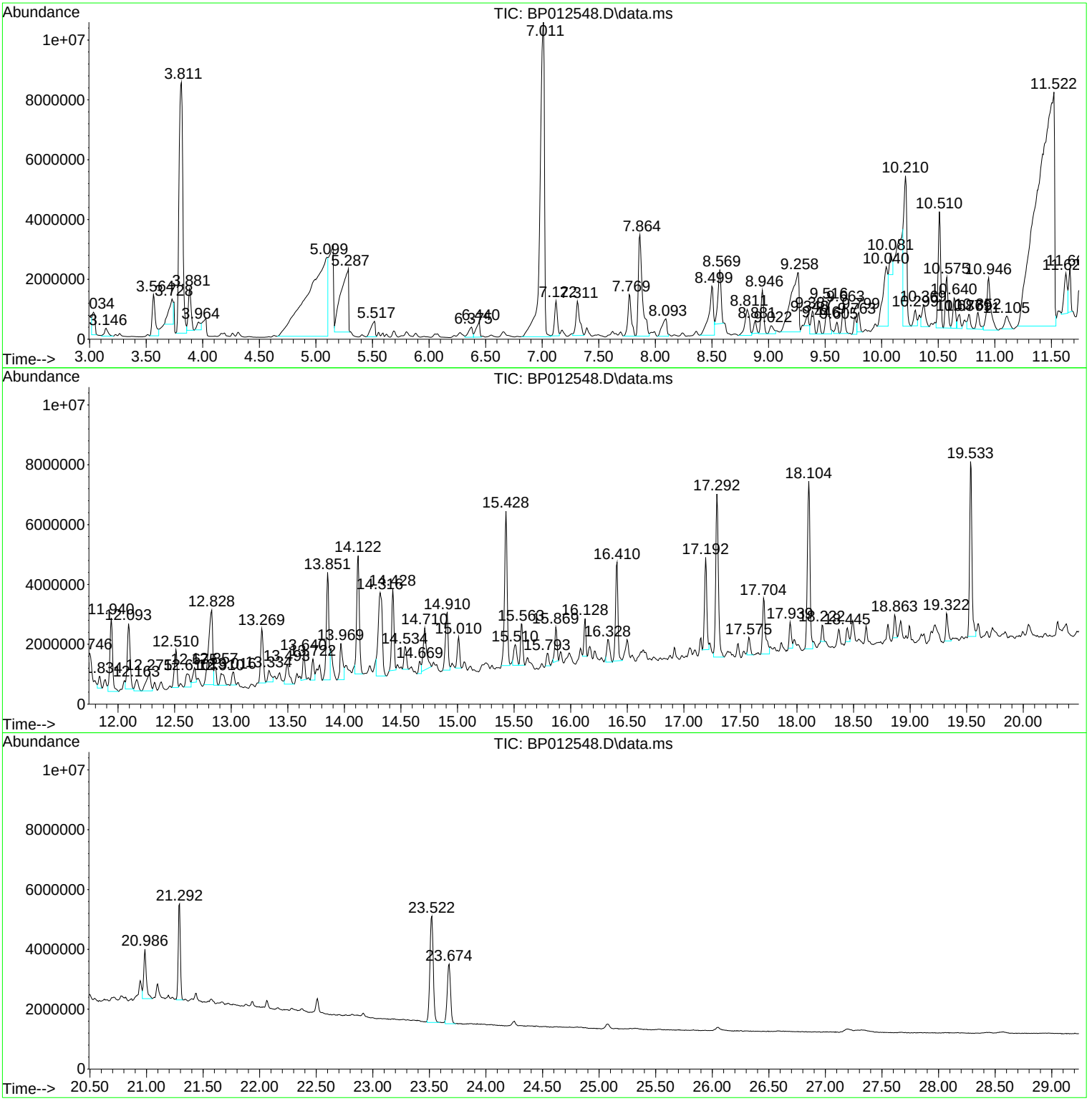
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110822\
Data File : BP012548.D
Acq On : 08 Nov 2022 12:21
Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

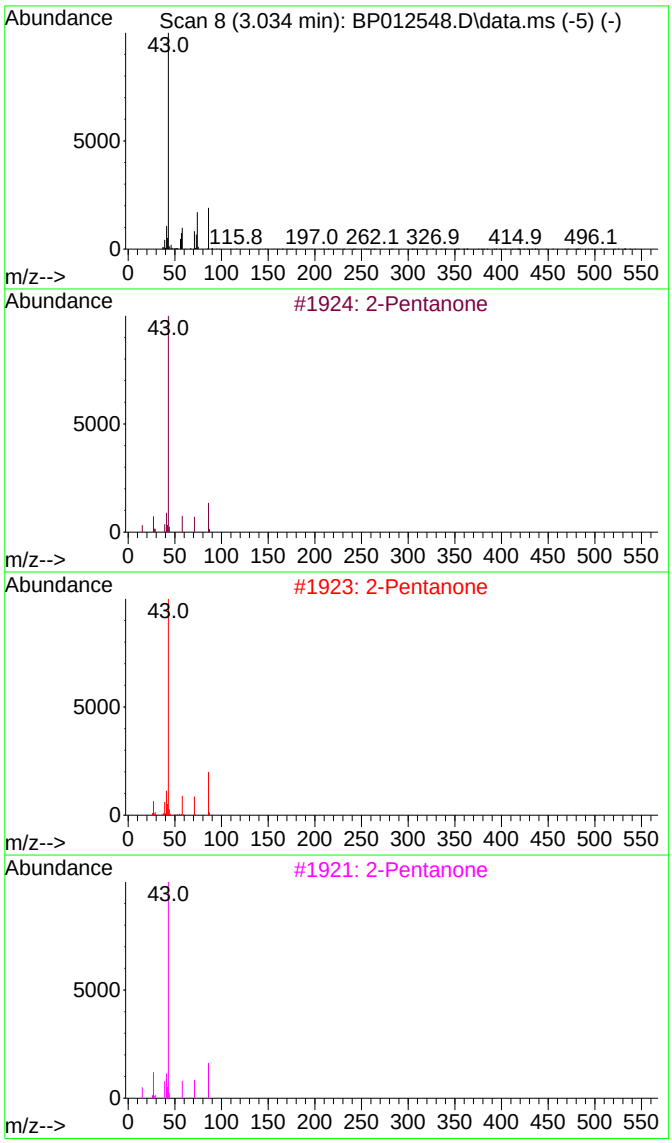
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

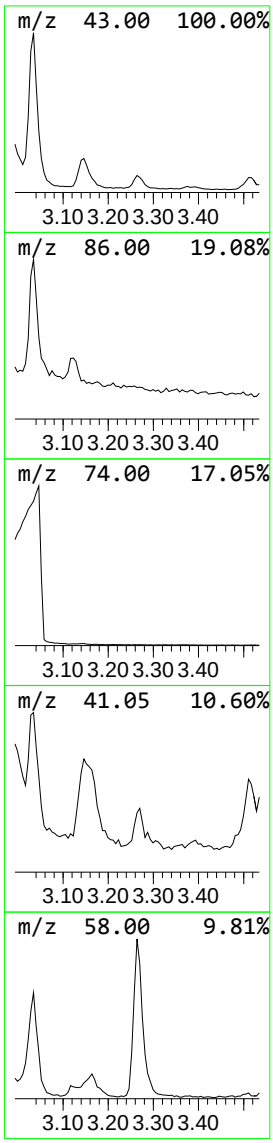
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 2-Pentanone Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.034	9.12 ng/ul	1331890	1,4-Dichlorobenzene-d4	7.775

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone	86	C5H10O	000107-87-9	52
2	2-Pentanone	86	C5H10O	000107-87-9	52
3	2-Pentanone	86	C5H10O	000107-87-9	40
4	2-Pentanone	86	C5H10O	000107-87-9	38
5	2-Pentanone	86	C5H10O	000107-87-9	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110822\
Data File : BP012548.D
Acq On : 08 Nov 2022 12:21
Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

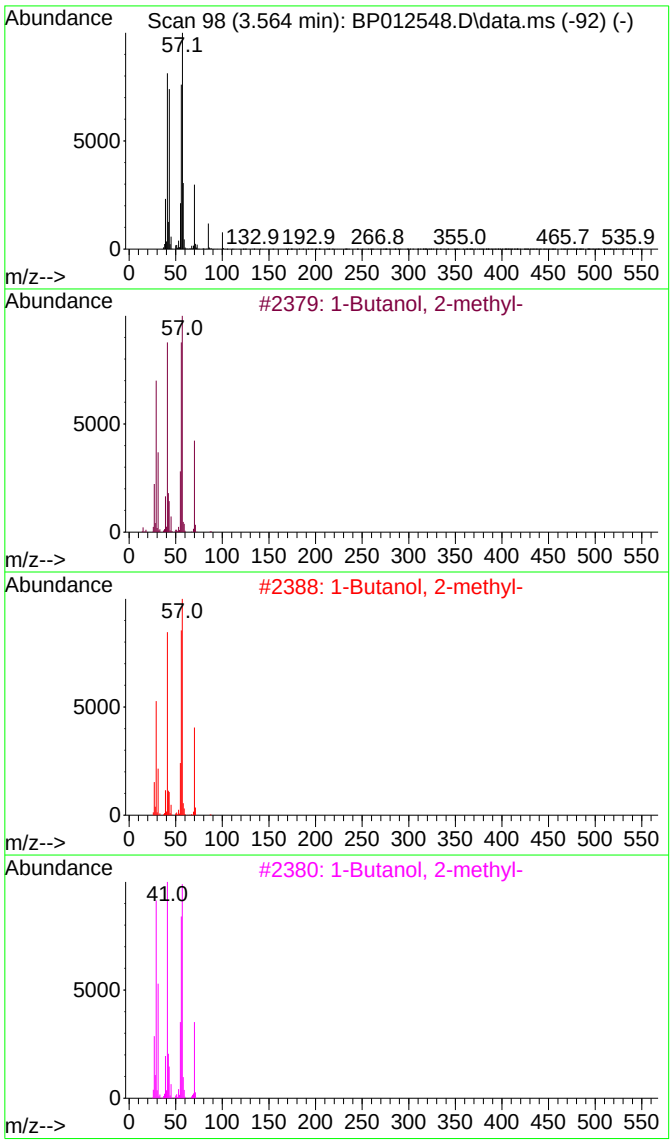
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

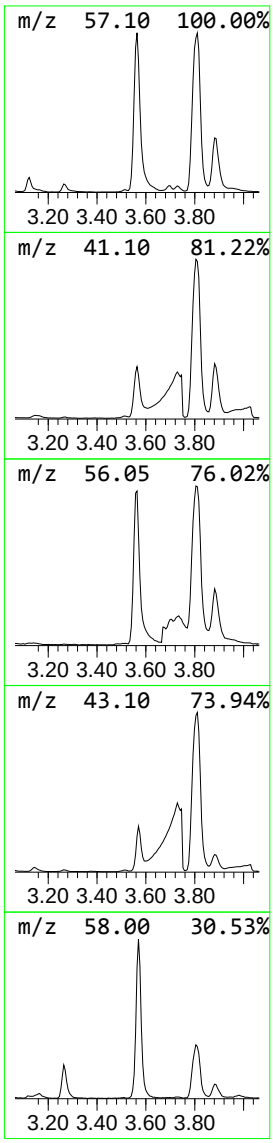
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 1-Butanol, 2-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.564	18.42 ng/ul	2691950	1,4-Dichlorobenzene-d4	7.775

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butanol, 2-methyl-	88	C5H12O	000137-32-6	68
2	1-Butanol, 2-methyl-	88	C5H12O	000137-32-6	59
3	1-Butanol, 2-methyl-	88	C5H12O	000137-32-6	53
4	Propane, 1-(ethenyloxy)-2-methyl-	100	C6H12O	000109-53-5	50
5	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	45





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110822\
Data File : BP012548.D
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ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

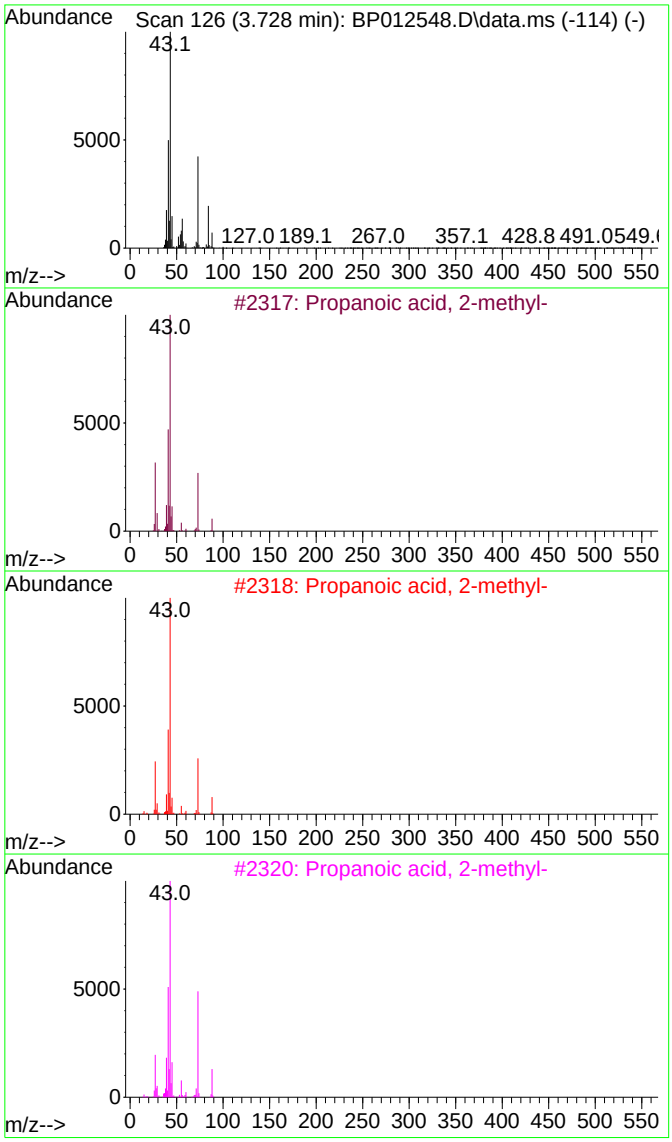
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

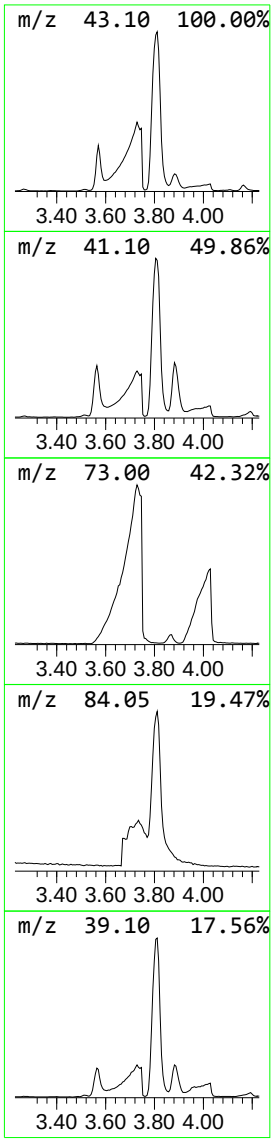
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Propanoic acid, 2-methyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.728	16.36 ng/ul	2389870	1,4-Dichlorobenzene-d4	7.775

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	58
2	Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	58
3	Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	49
4	Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	49
5	Ethanone, 1-cyclopropyl-	84	C5H8O	000765-43-5	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110822\
Data File : BP012548.D
Acq On : 08 Nov 2022 12:21
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Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

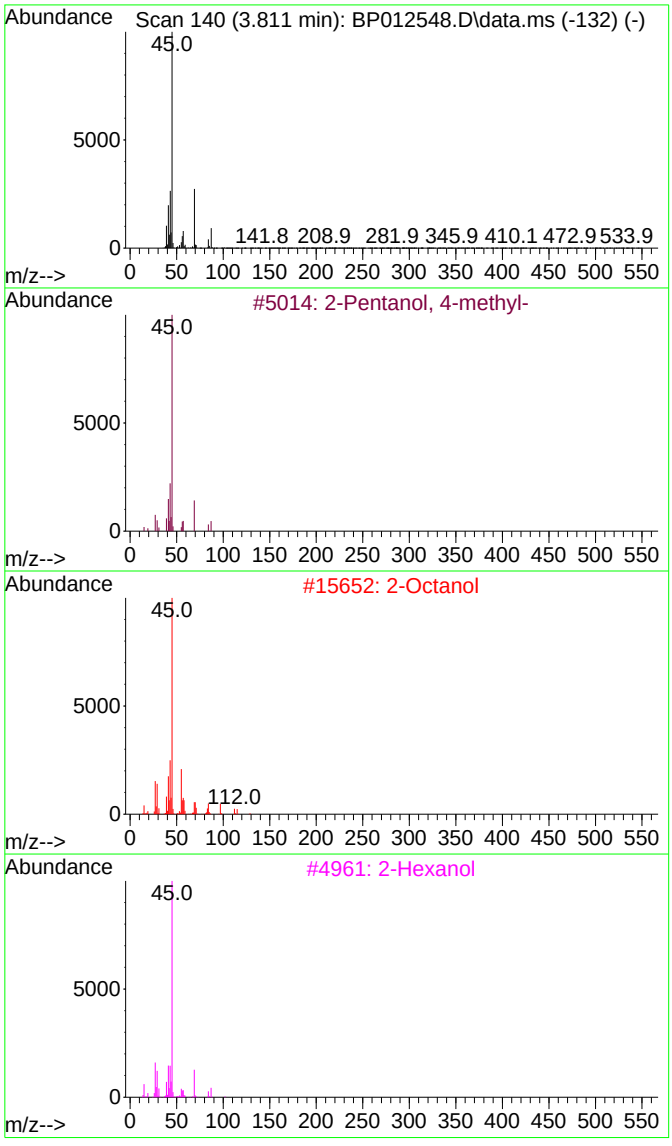
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

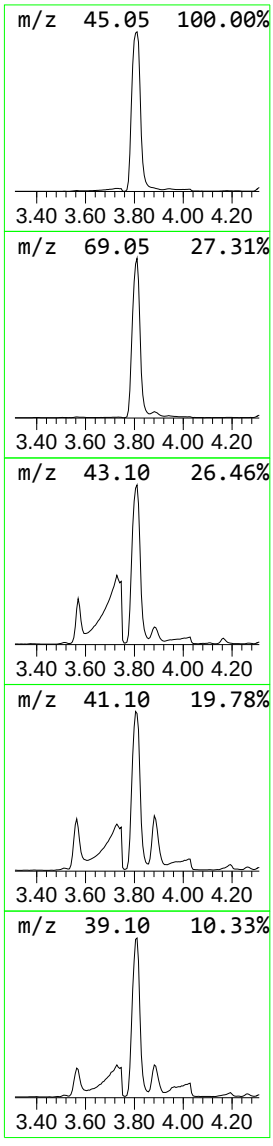
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 2-Pentanol, 4-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.811	132.71 ng/ul	19390100	1,4-Dichlorobenzene-d4	7.775

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	83
2	2-Octanol	130	C8H18O	000123-96-6	78
3	2-Hexanol	102	C6H14O	000626-93-7	74
4	2-Octanol, (S)-	130	C8H18O	006169-06-8	72
5	2-Octanol, (R)-	130	C8H18O	005978-70-1	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110822\
Data File : BP012548.D
Acq On : 08 Nov 2022 12:21
Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

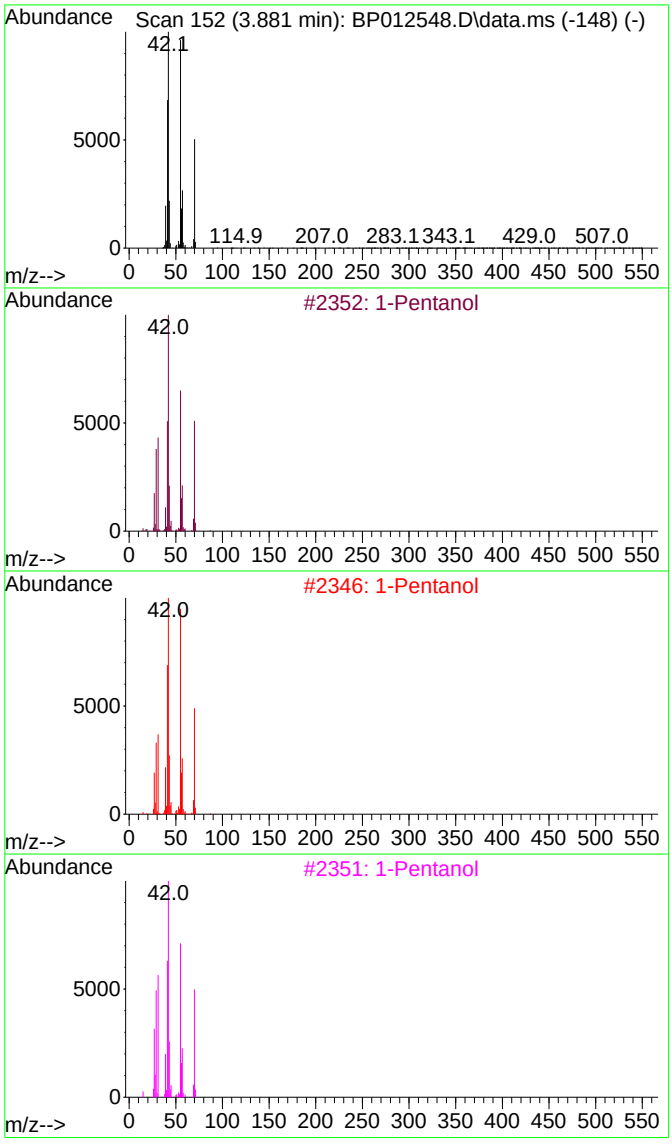
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

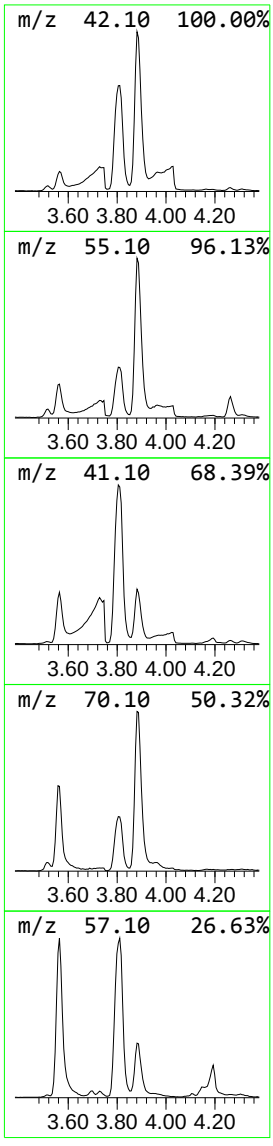
****

Peak Number 6 1-Pentanol Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.881	16.26 ng/ul	2375550	1,4-Dichlorobenzene-d4	7.775

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentanol	88	C5H12O	000071-41-0	83
2	1-Pentanol	88	C5H12O	000071-41-0	74
3	1-Pentanol	88	C5H12O	000071-41-0	64
4	1-Pentene	70	C5H10	000109-67-1	56
5	1-Pentene	70	C5H10	000109-67-1	56





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Acq On : 08 Nov 2022 12:21
Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

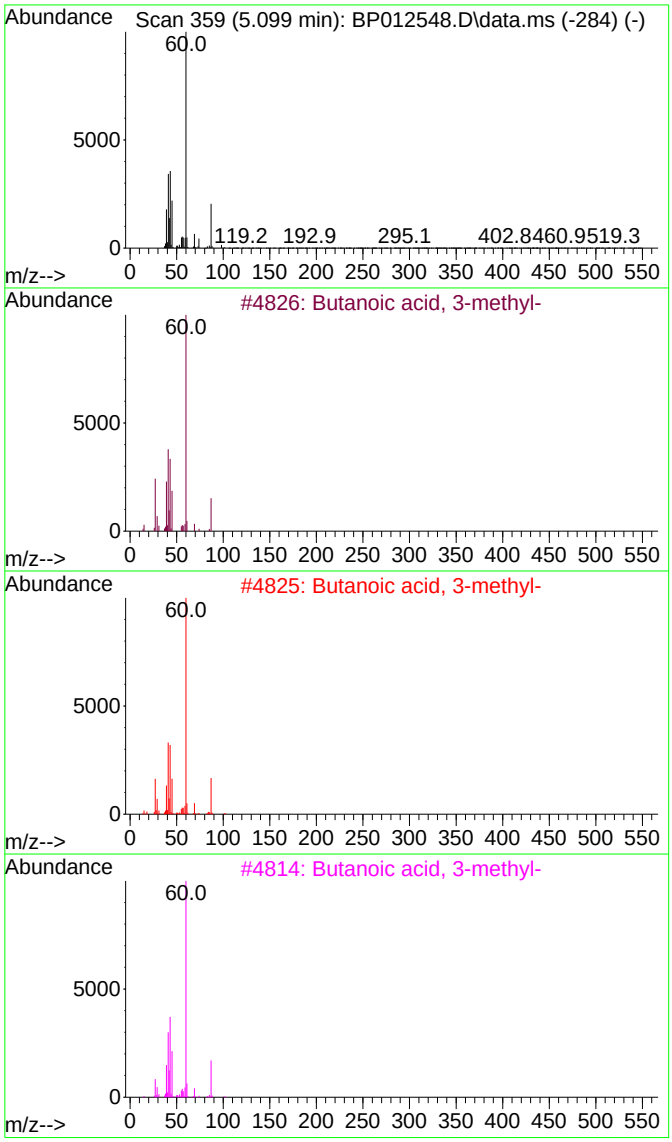
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

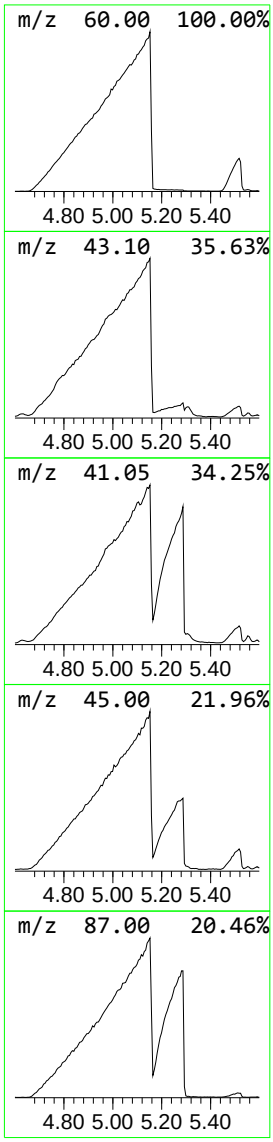
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Butanoic acid, 3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.099	223.28 ng/ul	32623700	1,4-Dichlorobenzene-d4	7.775

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	83
2	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	78
3	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	78
4	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	50
5	Pentanoic acid	102	C5H10O2	000109-52-4	42





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Acq On : 08 Nov 2022 12:21
Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

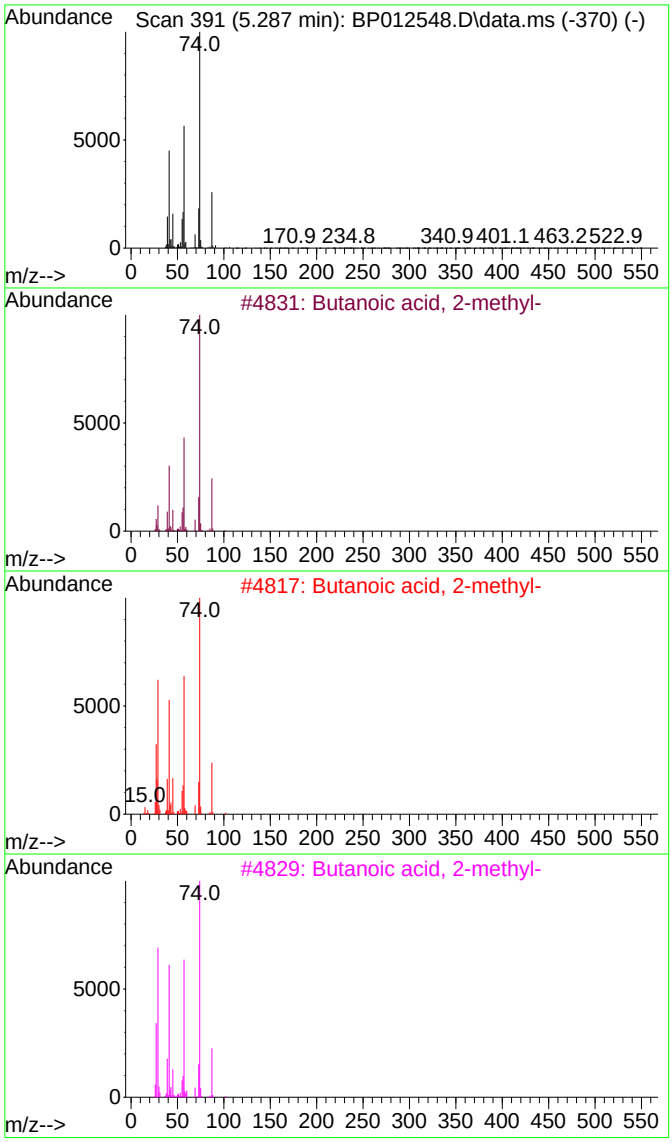
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

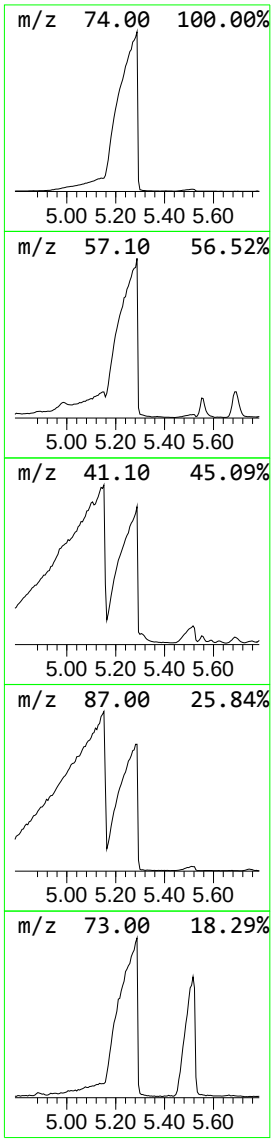
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Butanoic acid, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.287	68.79 ng/ul	10050300	1,4-Dichlorobenzene-d4	7.775

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, 2-methyl-	102	C5H10O2	000116-53-0	90
2	Butanoic acid, 2-methyl-	102	C5H10O2	000116-53-0	83
3	Butanoic acid, 2-methyl-	102	C5H10O2	000116-53-0	74
4	Pentanoic acid, 2-methyl-	116	C6H12O2	000097-61-0	64
5	Butanoic acid, 2-methyl-	102	C5H10O2	000116-53-0	56





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Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

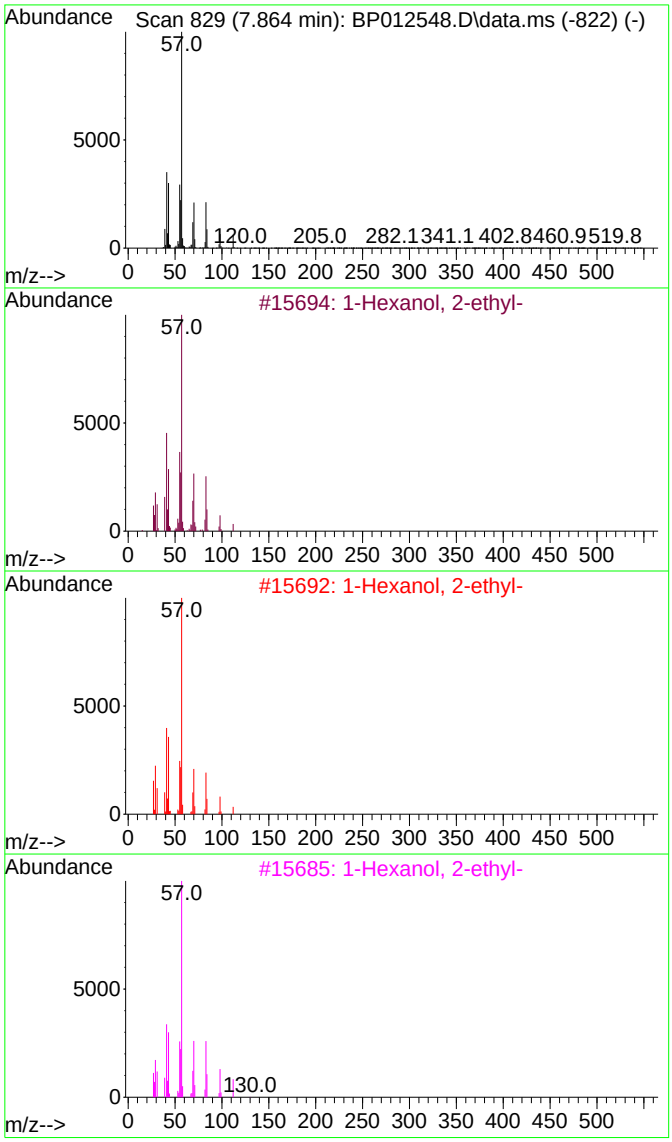
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

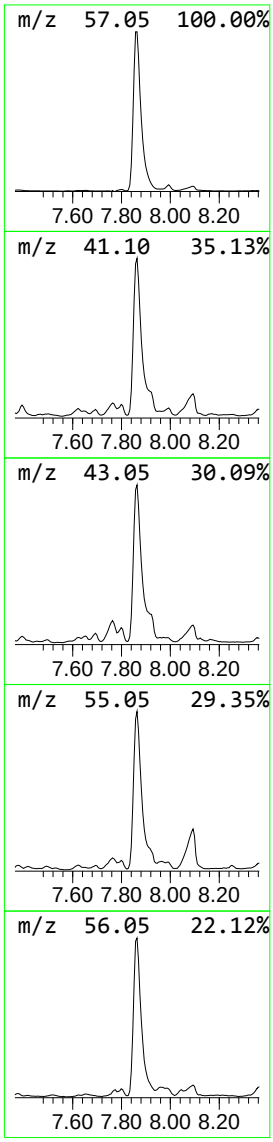
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 1-Hexanol, 2-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.864	57.76 ng/ul	8439550	1,4-Dichlorobenzene-d4	7.775

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
3	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	74
4	1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	56
5	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	56





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Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

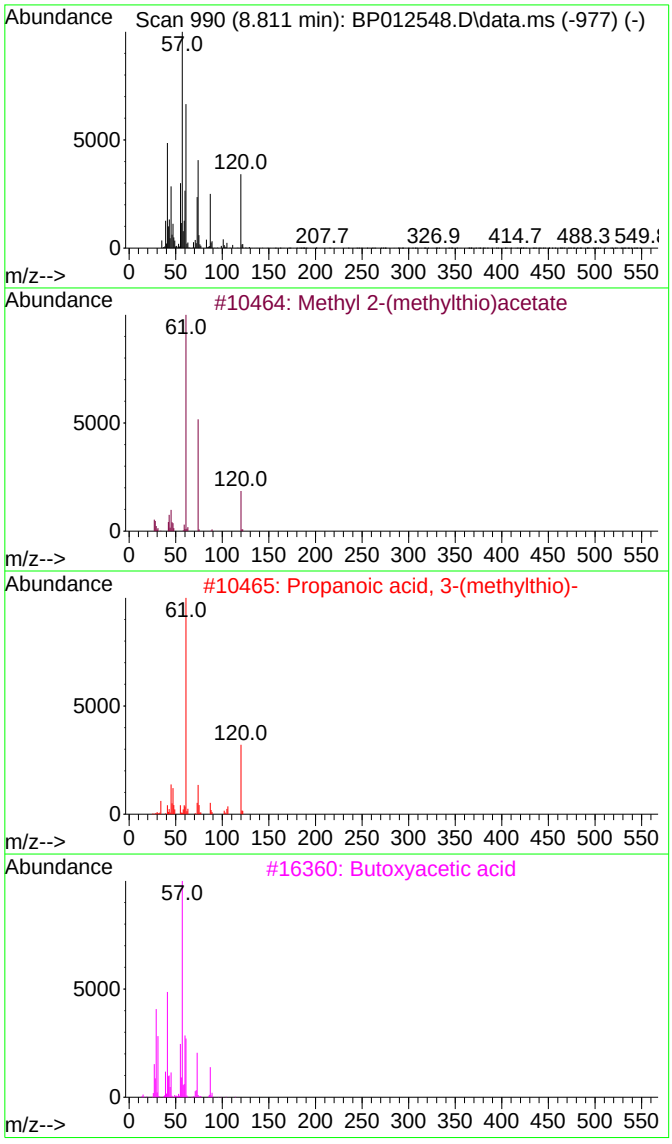
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

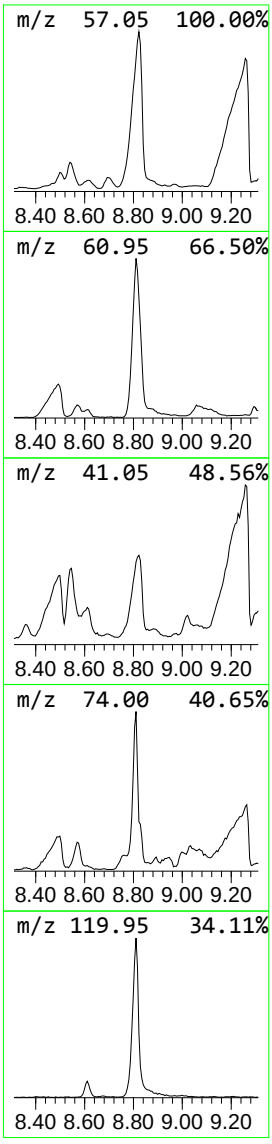
****

Peak Number 15 unknown-01 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.811	16.55 ng/ul	2417660	1,4-Dichlorobenzene-d4	7.775

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl 2-(methylthio)acetate	120	C4H8O2S	016630-66-3	46
2	Propanoic acid, 3-(methylthio)-	120	C4H8O2S	000646-01-5	38
3	Butoxyacetic acid	132	C6H12O3	002516-93-0	35
4	Xylitol	152	C5H12O5	000087-99-0	25
5	Butoxyacetic acid	132	C6H12O3	002516-93-0	16





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Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

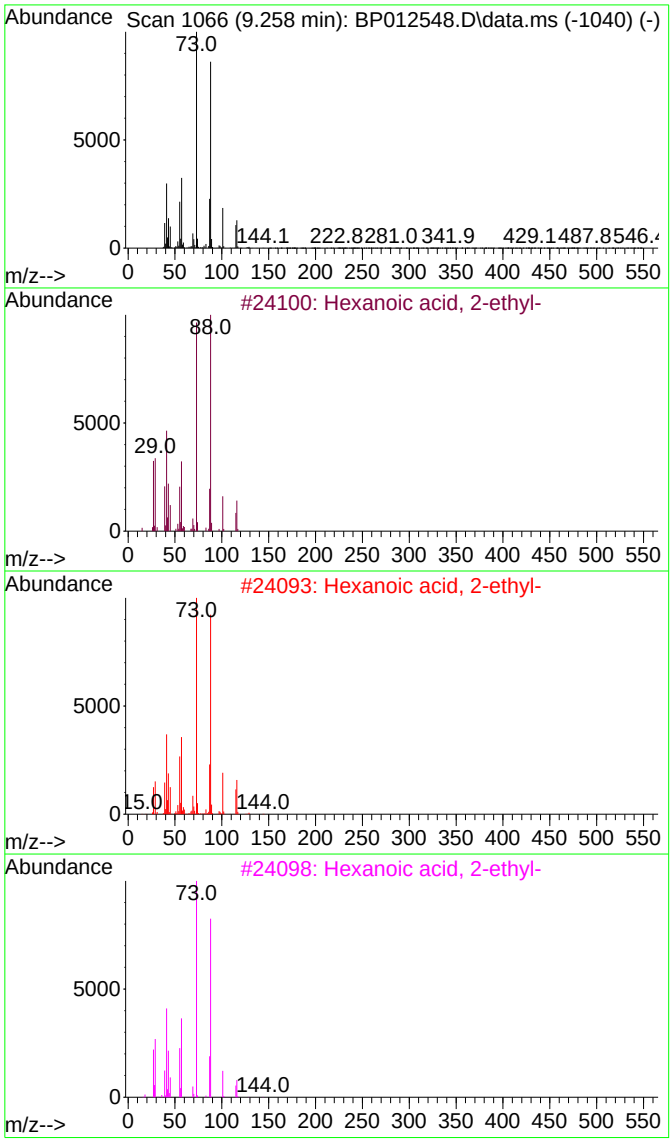
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

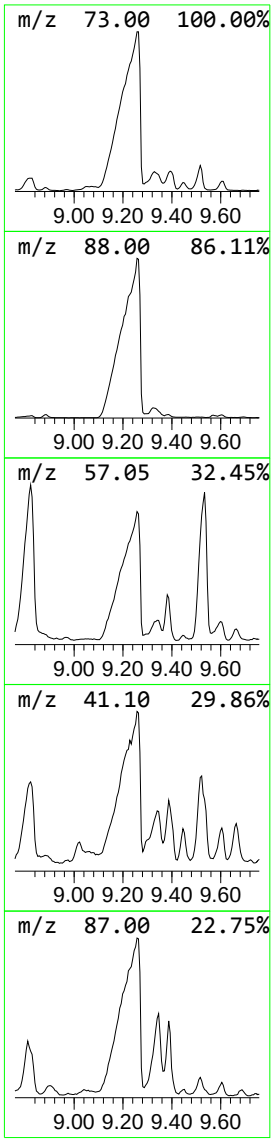
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Hexanoic acid, 2-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.258	71.27 ng/ul	9977330	Naphthalene-d8	10.575

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
2	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
3	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	78
4	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	78
5	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	78





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Acq On : 08 Nov 2022 12:21
Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

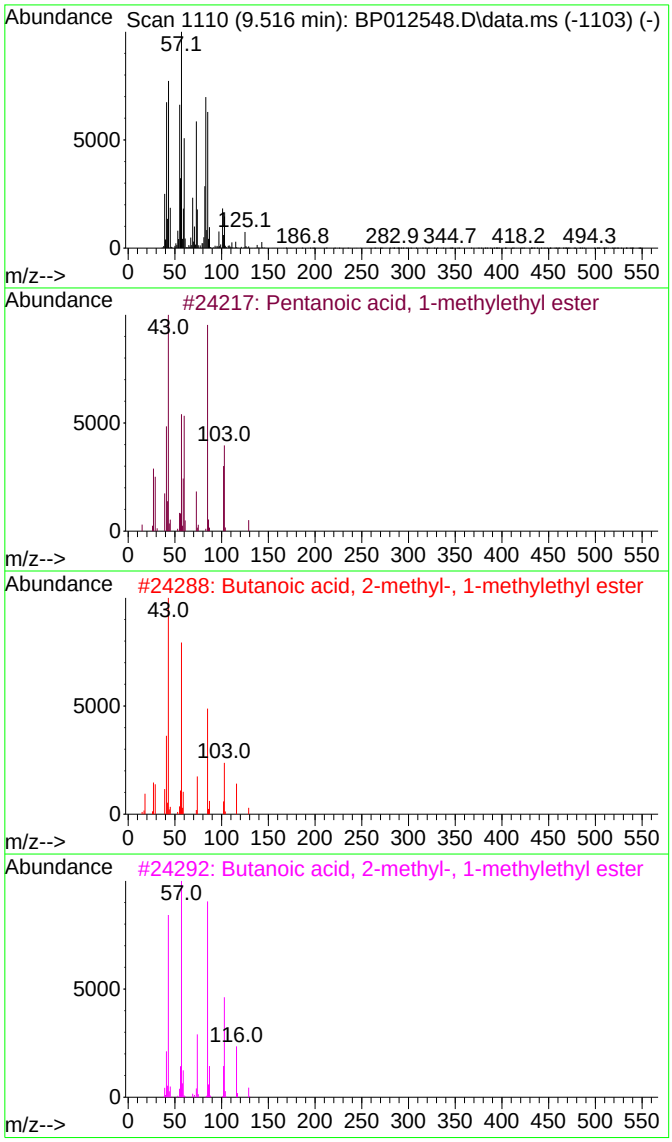
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

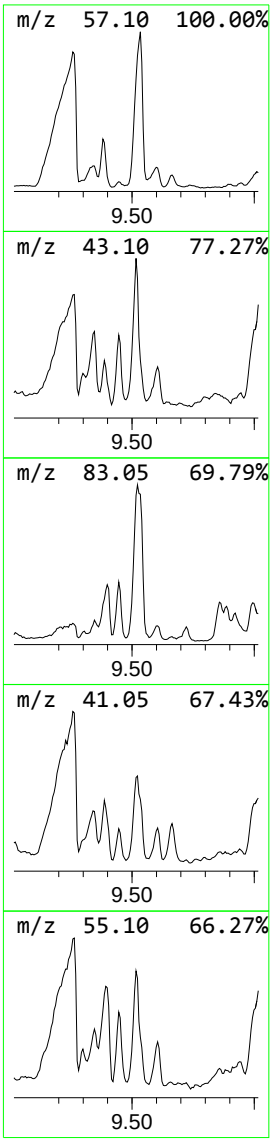
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 unknown-02 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.516	16.28 ng/ul	2278560	Naphthalene-d8	10.575

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentanoic acid, 1-methylethyl ester	144	C8H16O2	018362-97-5	38
2	Butanoic acid, 2-methyl-, 1-meth...	144	C8H16O2	066576-71-4	35
3	Butanoic acid, 2-methyl-, 1-meth...	144	C8H16O2	066576-71-4	35
4	Isobutyl isovalerate	158	C9H18O2	000589-59-3	27
5	Butanoic acid, 2-methyl-, 2-meth...	158	C9H18O2	002445-67-2	22





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Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

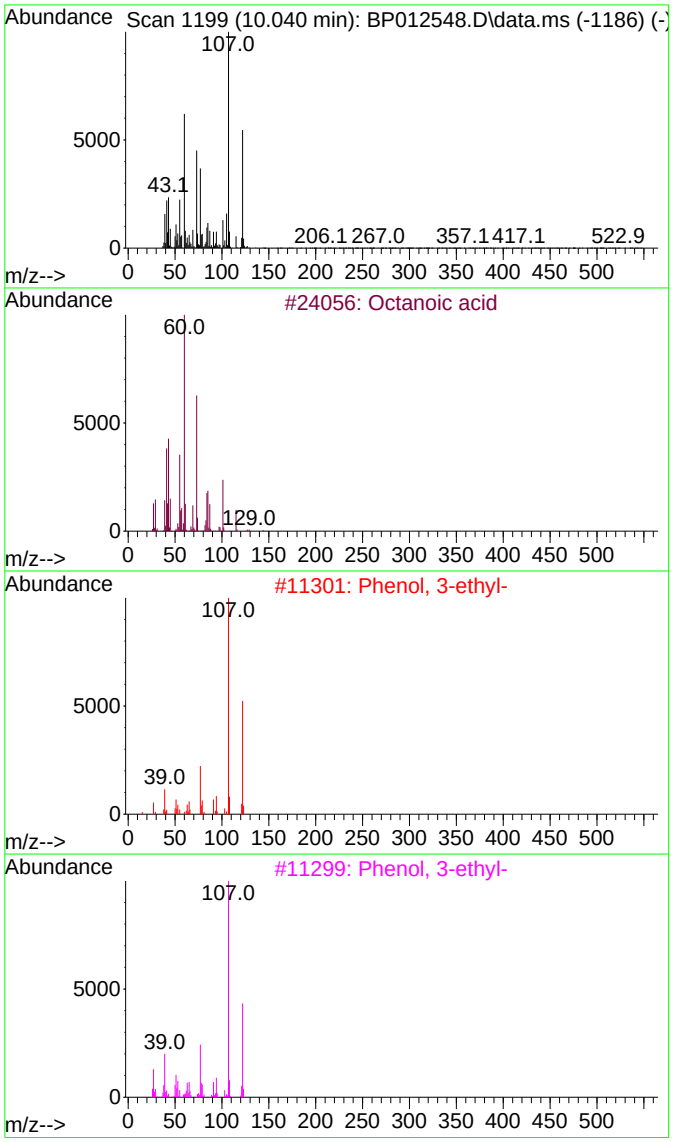
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

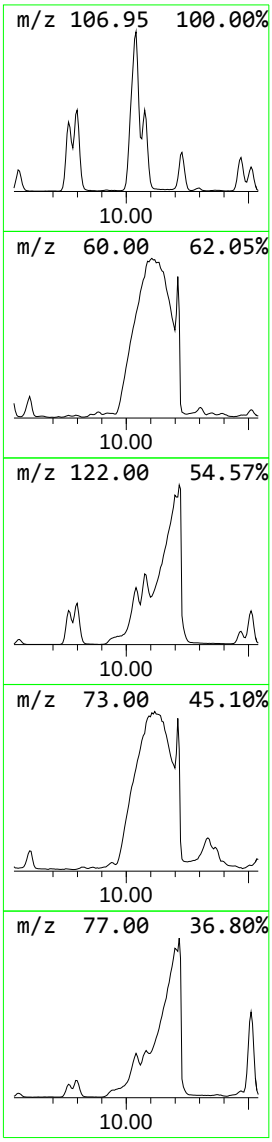
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Octanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.040	45.46 ng/ul	6364790	Naphthalene-d8	10.575

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octanoic acid	144	C8H16O2	000124-07-2	92
2	Phenol, 3-ethyl-	122	C8H10O	000620-17-7	91
3	Phenol, 3-ethyl-	122	C8H10O	000620-17-7	64
4	Phenol, 2-ethyl-	122	C8H10O	000090-00-6	60
5	Phenol, 3-ethyl-	122	C8H10O	000620-17-7	60





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Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

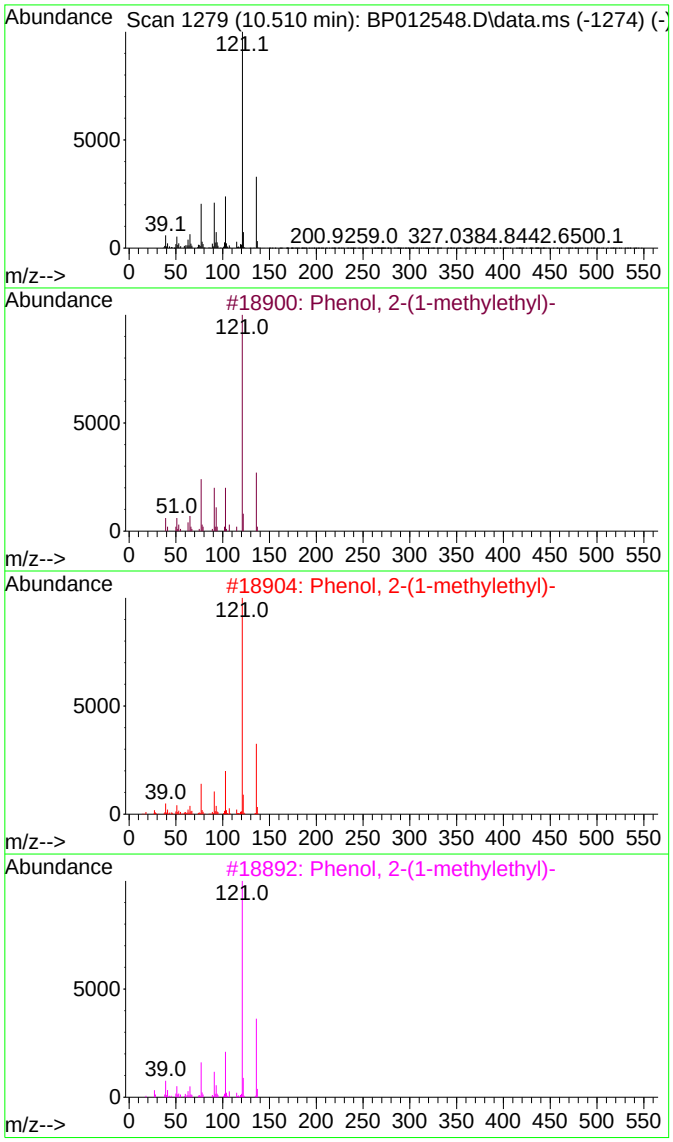
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

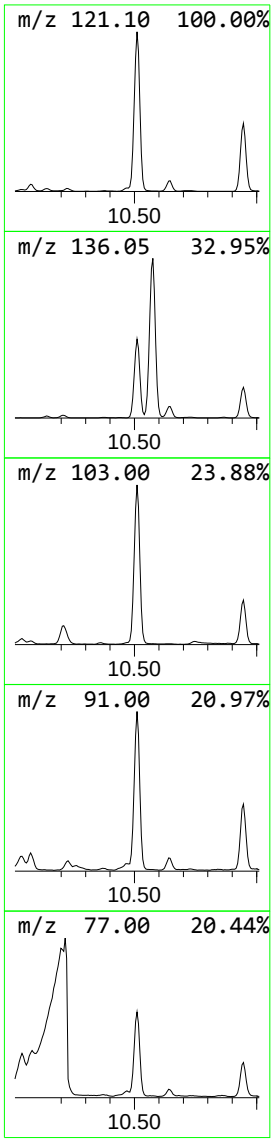
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 28 Phenol, 2-(1-methylethyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.510	43.08 ng/ul	6030800	Naphthalene-d8	10.575

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	93
2	Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	91
3	Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	90
4	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	86
5	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	86





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Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

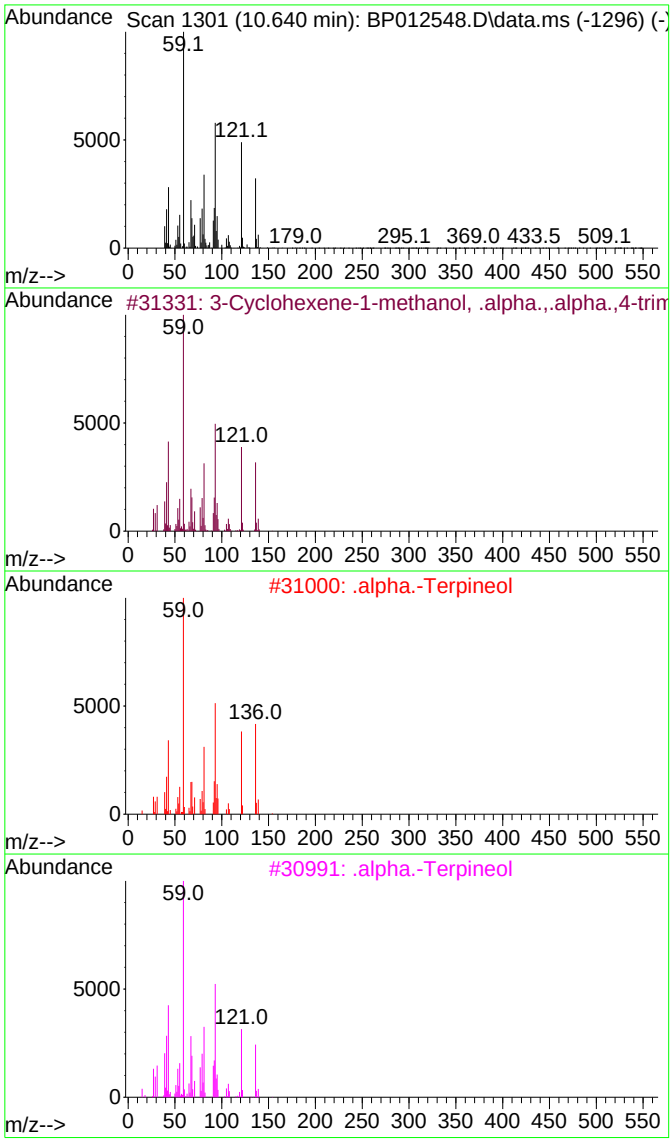
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

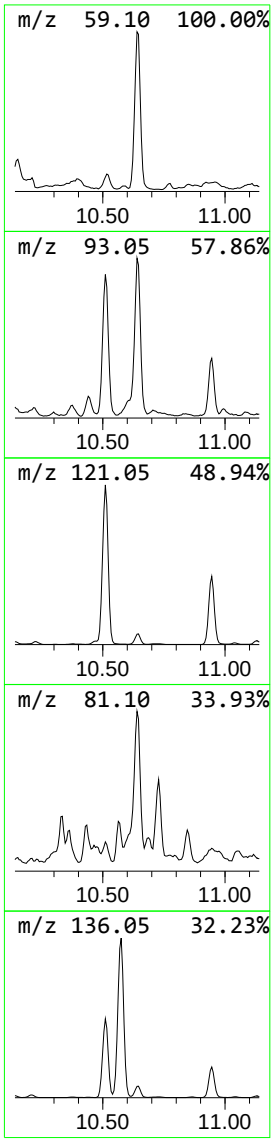
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 29 3-Cyclohexene-1-methanol, ... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.640	12.25 ng/ul	1715310	Naphthalene-d8	10.575

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Cyclohexene-1-methanol, .alpha...	154	C10H18O	007785-53-7	91
2	.alpha.-Terpineol	154	C10H18O	000098-55-5	90
3	.alpha.-Terpineol	154	C10H18O	000098-55-5	86
4	.alpha.-Terpineol	154	C10H18O	000098-55-5	74
5	.alpha.-Terpineol	154	C10H18O	000098-55-5	52





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Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

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Quant Title : SVOA CALIBRATION

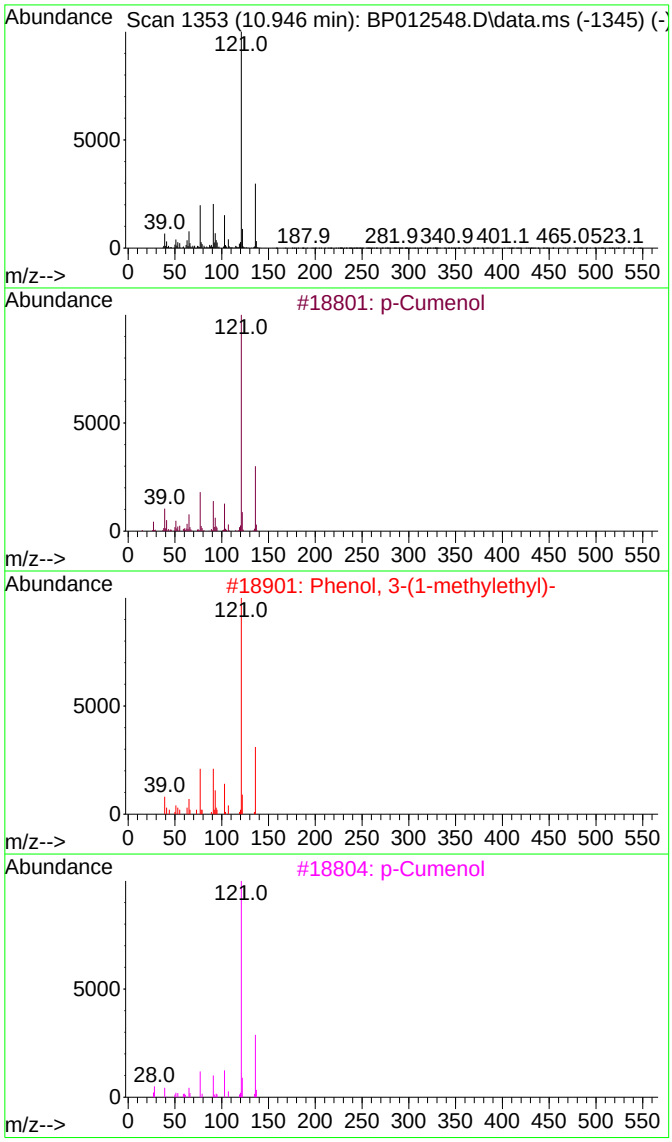
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

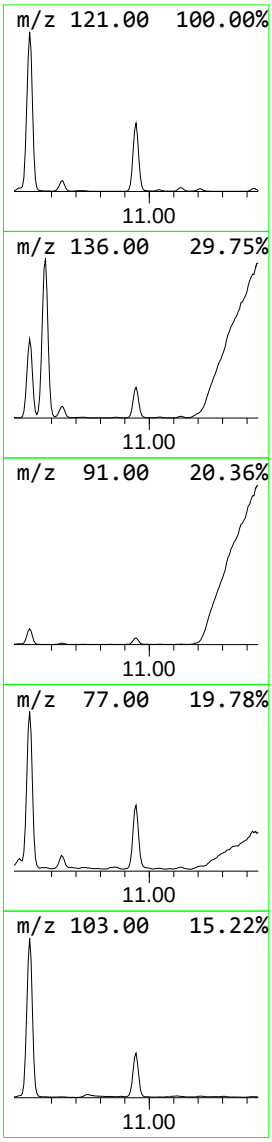
****

Peak Number 32 p-Cumenol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.946	31.13 ng/ul	4358670	Naphthalene-d8	10.575

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cumenol	136	C9H12O	000099-89-8	95
2	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	94
3	p-Cumenol	136	C9H12O	000099-89-8	91
4	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	91
5	Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	91





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ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EW4Y2

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Quant Title : SVOA CALIBRATION

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

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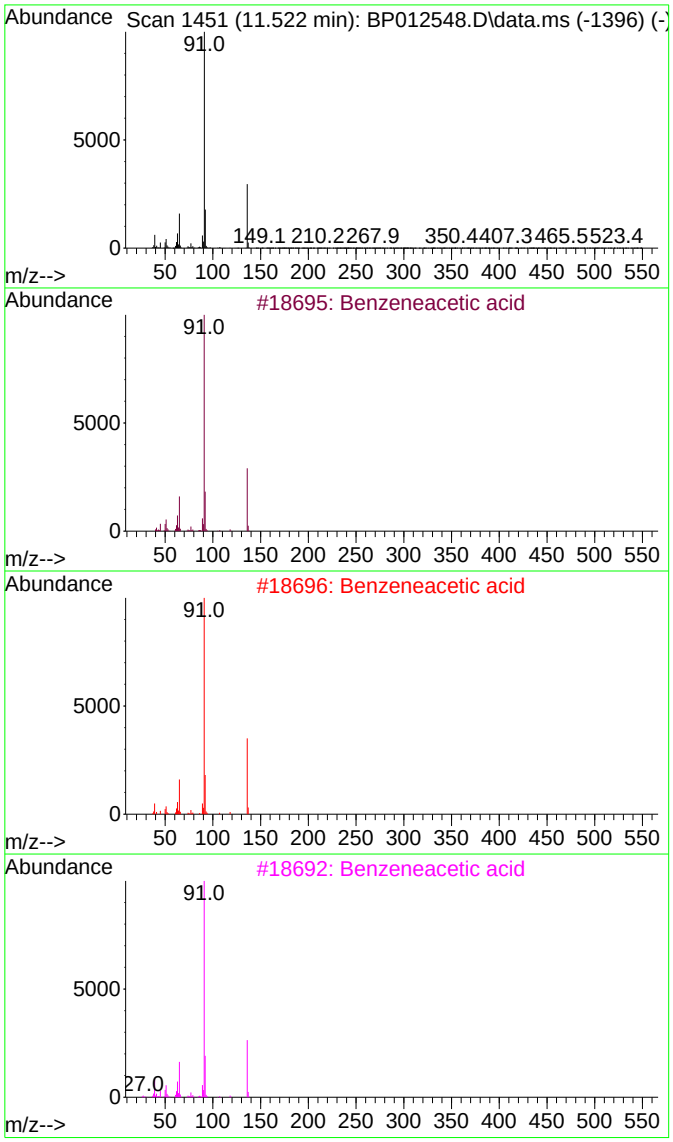
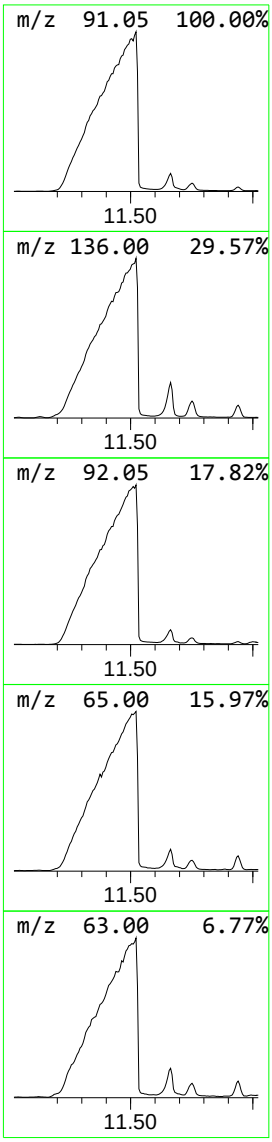
Peak Number 34 Benzeneacetic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.522	583.89 ng/ul	81744000	Naphthalene-d8	10.575

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzeneacetic acid	136	C8H8O2	000103-82-2	94
2	Benzeneacetic acid	136	C8H8O2	000103-82-2	94
3	Benzeneacetic acid	136	C8H8O2	000103-82-2	94
4	Benzeneacetic acid	136	C8H8O2	000103-82-2	90
5	Benzeneacetic acid	136	C8H8O2	000103-82-2	64





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Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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

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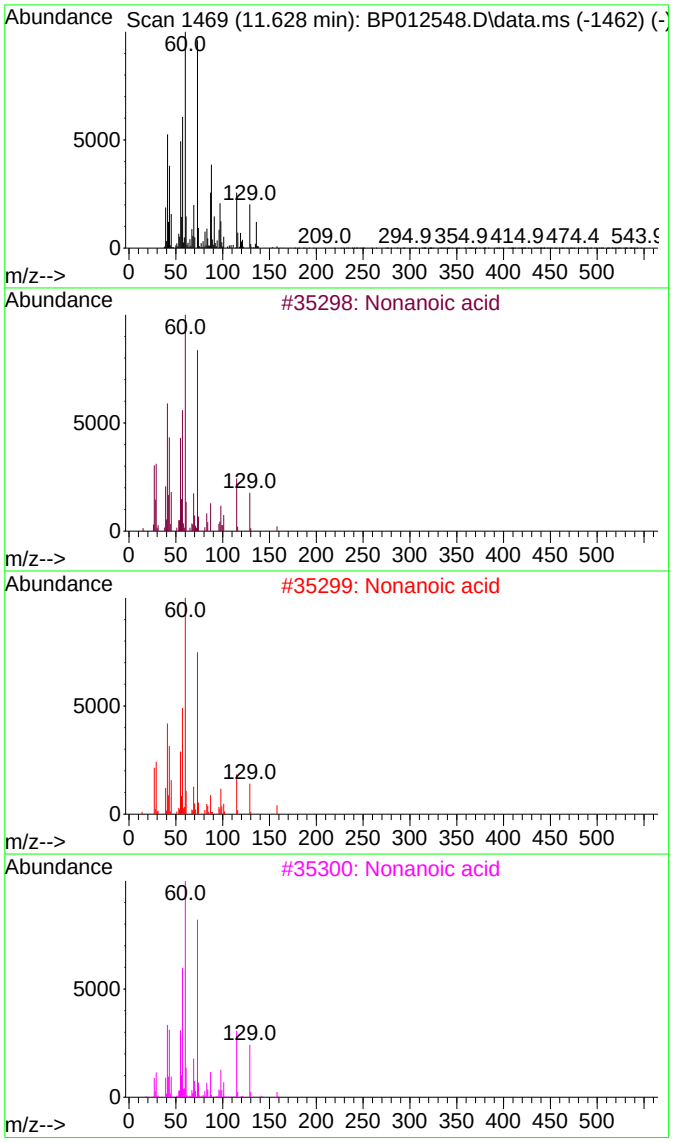
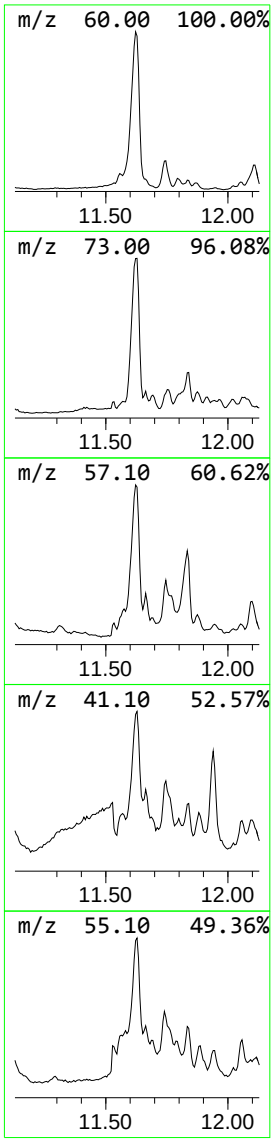
Peak Number 35 Nonanoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.628	19.83 ng/ul	2775600	Naphthalene-d8	10.575

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonanoic acid	158	C9H18O2	000112-05-0	53
2	Nonanoic acid	158	C9H18O2	000112-05-0	50
3	Nonanoic acid	158	C9H18O2	000112-05-0	49
4	Nonanoic acid	158	C9H18O2	000112-05-0	38
5	Nonanoic acid	158	C9H18O2	000112-05-0	38





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Acq On : 08 Nov 2022 12:21
Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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

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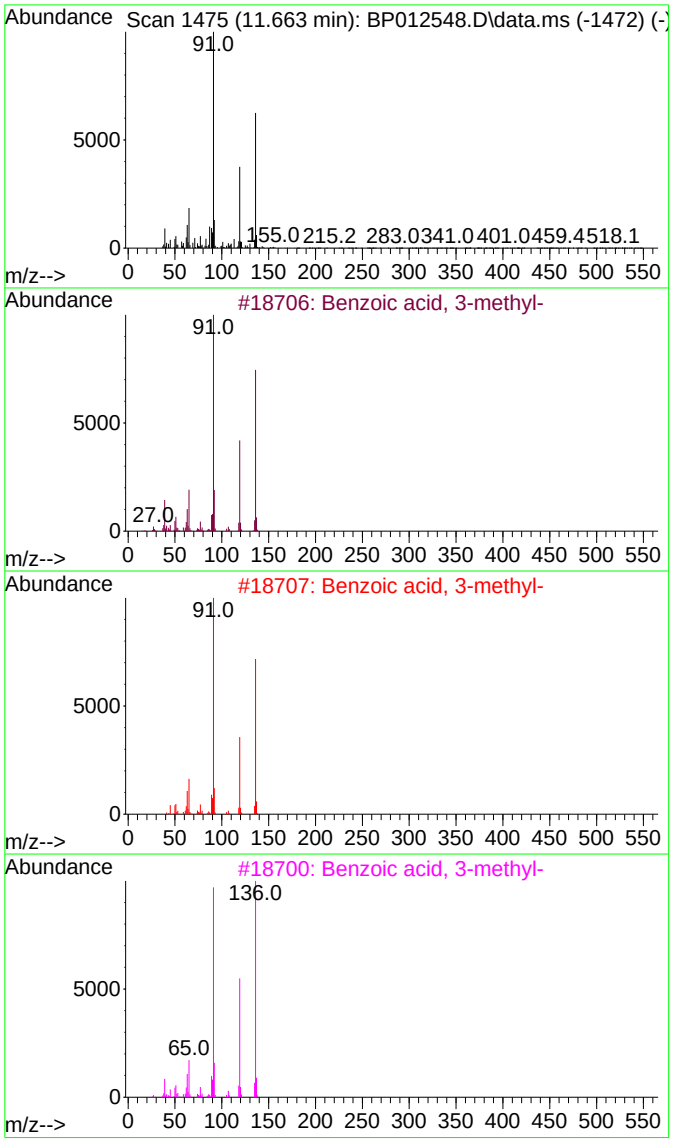
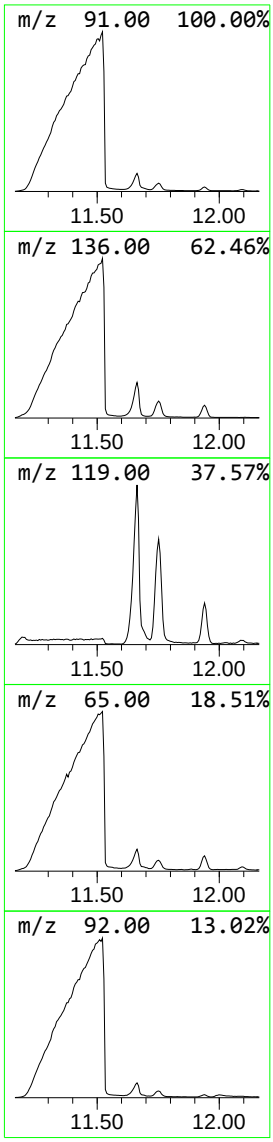
Peak Number 36 Benzoic acid, 3-methyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.663	12.62 ng/ul	1766370	Naphthalene-d8	10.575

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 3-methyl-	136	C8H8O2	000099-04-7	96
2	Benzoic acid, 3-methyl-	136	C8H8O2	000099-04-7	96
3	Benzoic acid, 3-methyl-	136	C8H8O2	000099-04-7	95
4	Benzoic acid, 4-methyl-	136	C8H8O2	000099-94-5	87
5	Benzoic acid, 4-methyl-	136	C8H8O2	000099-94-5	87





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Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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

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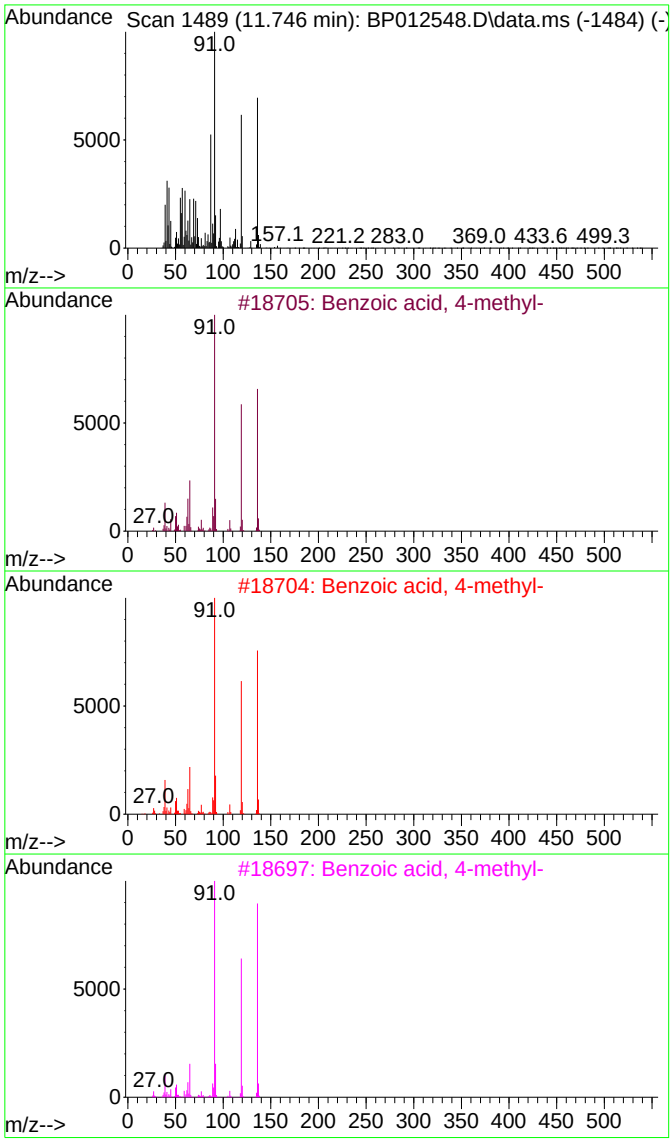
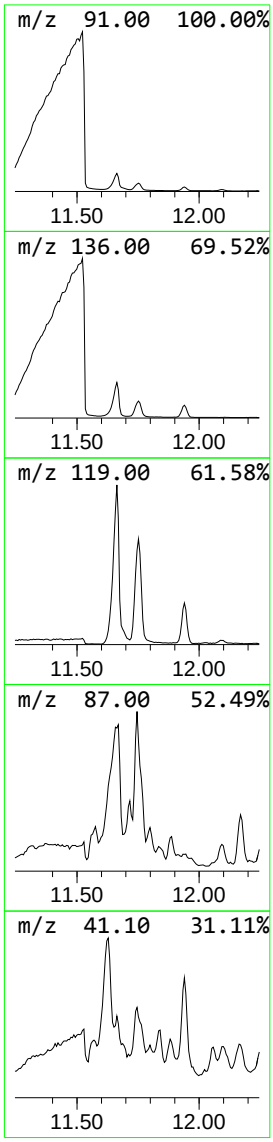
Peak Number 37 Benzoic acid, 4-methyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.746	13.81 ng/ul	1933650	Naphthalene-d8	10.575

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 4-methyl-	136	C8H8O2	000099-94-5	95
2	Benzoic acid, 4-methyl-	136	C8H8O2	000099-94-5	95
3	Benzoic acid, 4-methyl-	136	C8H8O2	000099-94-5	89
4	Benzoic acid, 4-methyl-	136	C8H8O2	000099-94-5	89
5	Benzoic acid, 4-methyl-	136	C8H8O2	000099-94-5	70





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Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

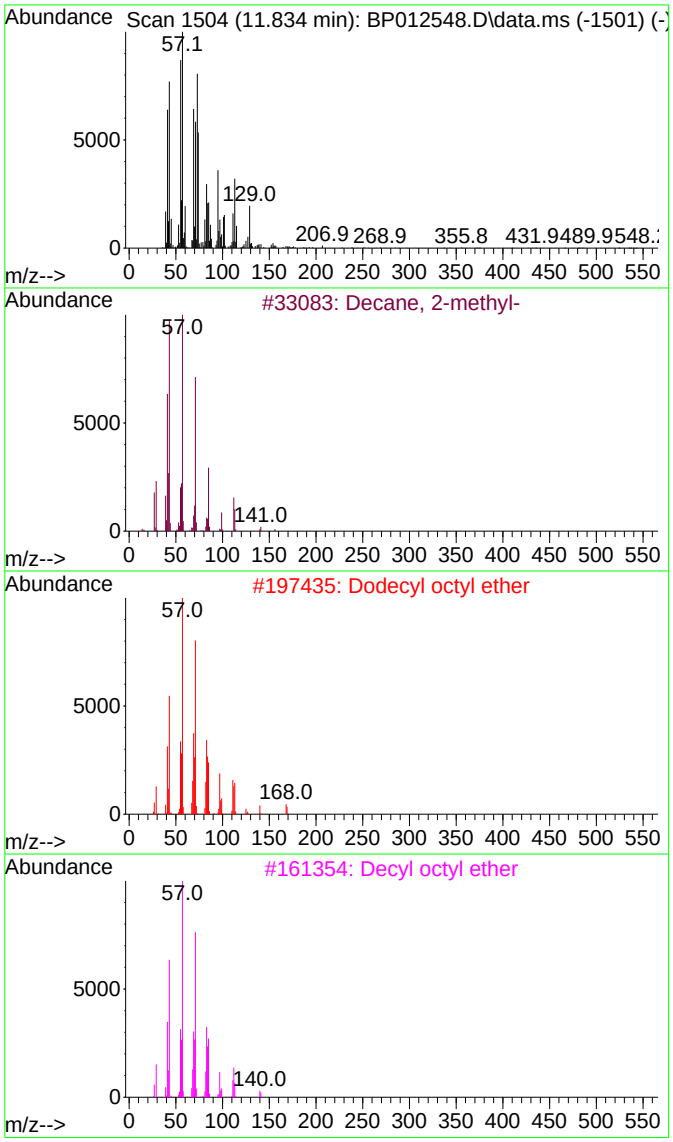
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

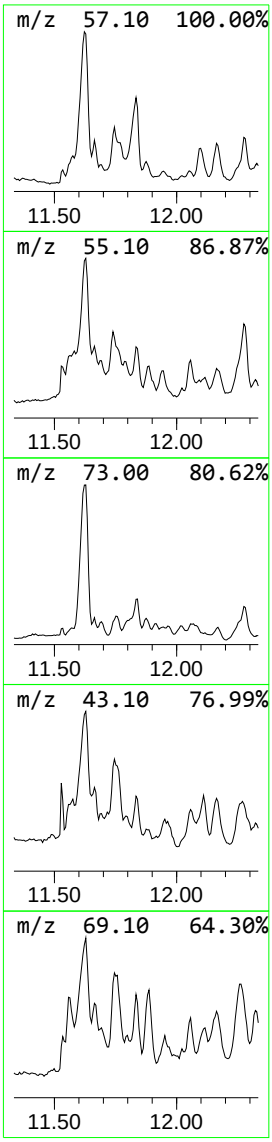
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 69

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.834	3.80 ng/ul	532327	Naphthalene-d8	10.575

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 2-methyl-	156	C11H24	006975-98-0	35
2	Dodecyl octyl ether	298	C20H42O	1000406-38-4	30
3	Decyl octyl ether	270	C18H38O	1000406-38-3	30
4	Carbonic acid, dodecyl 2-ethylhe...	342	C21H42O3	1000383-13-9	30
5	Hexadecyl octyl ether	354	C24H50O	1000406-38-6	30





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Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

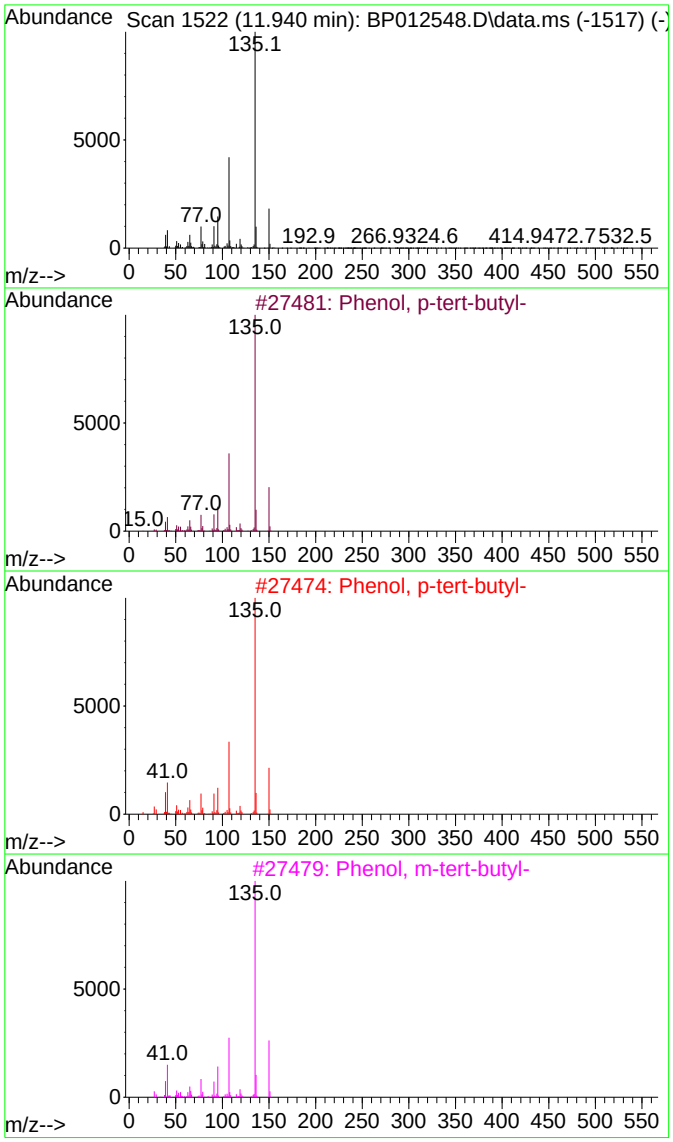
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

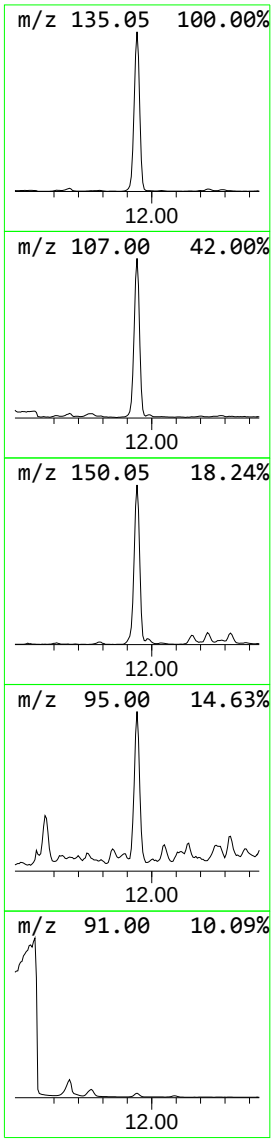
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 39 Phenol, p-tert-butyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.940	28.77 ng/ul	4028010	Naphthalene-d8	10.575

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	97
2	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	96
3	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	93
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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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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

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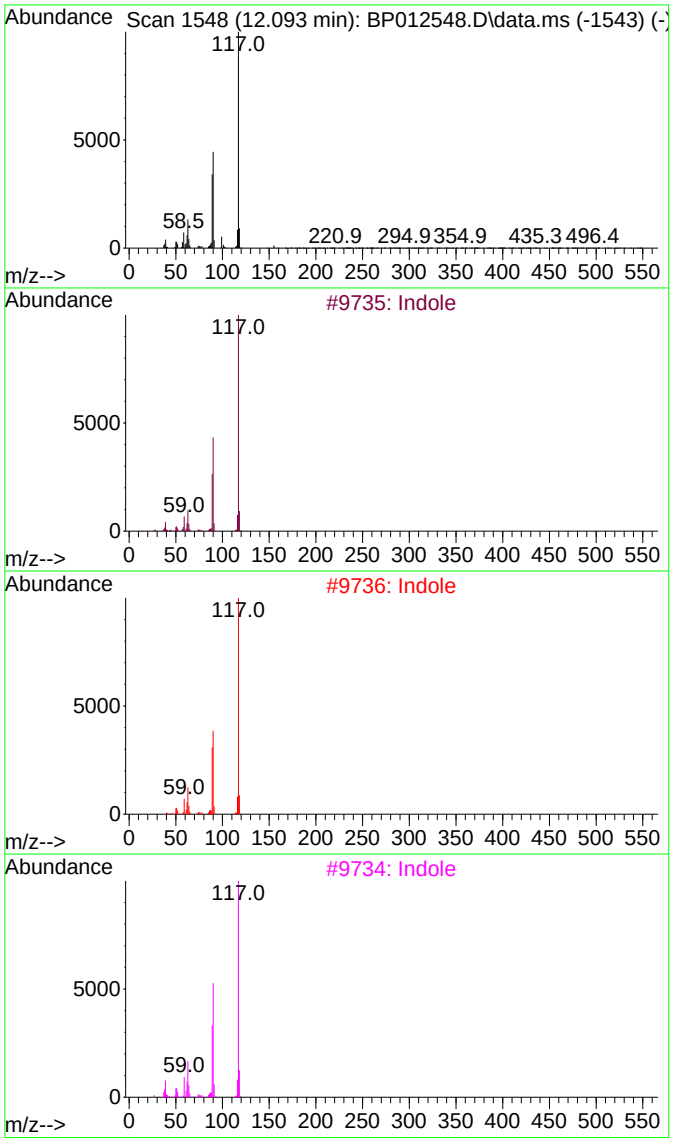
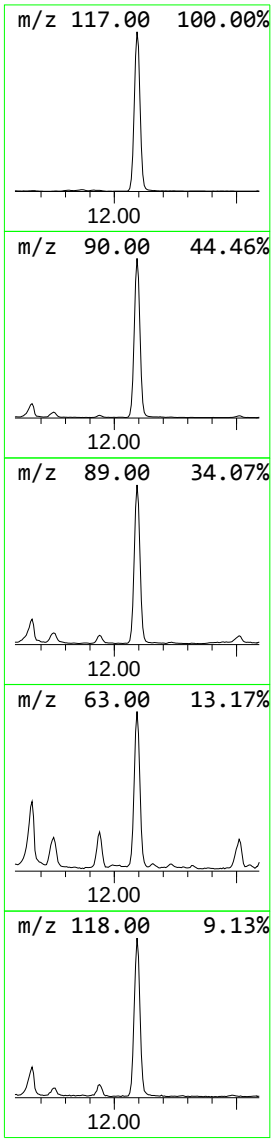
Peak Number 40 Indole Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.093	26.65 ng/ul	3730450	Naphthalene-d8	10.575

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indole	117	C8H7N	000120-72-9	95
2	Indole	117	C8H7N	000120-72-9	91
3	Indole	117	C8H7N	000120-72-9	91
4	Indole	117	C8H7N	000120-72-9	91
5	Benzyl nitrile	117	C8H7N	000140-29-4	90





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Acq On : 08 Nov 2022 12:21
Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

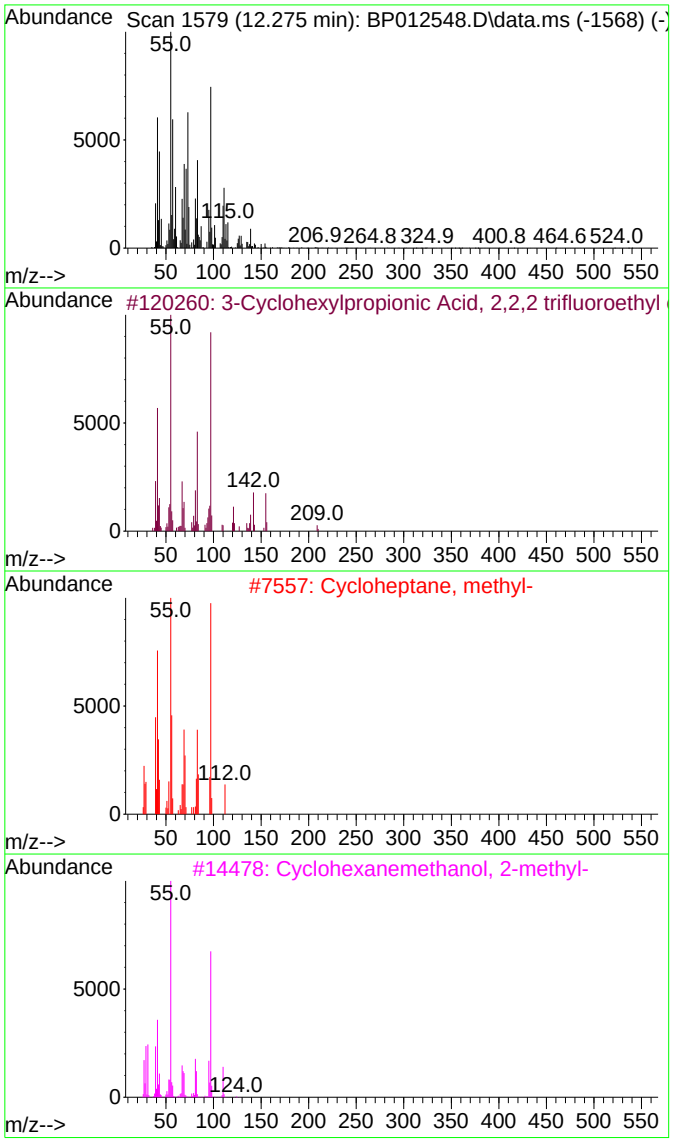
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

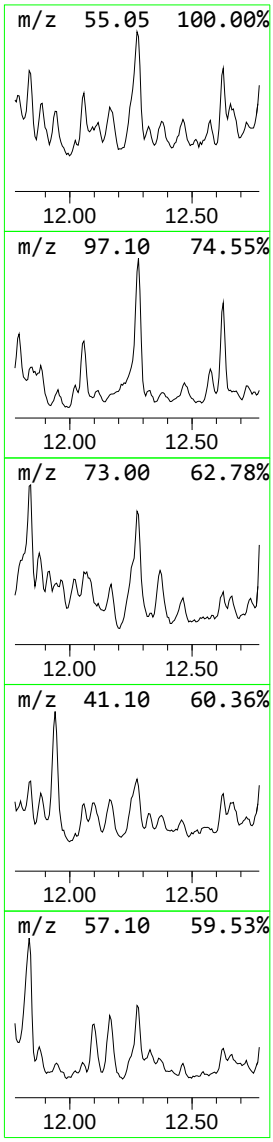
****

Peak Number 42 unknown-03 Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.275	11.35 ng/ul	1588440	Naphthalene-d8	10.575

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Cyclohexylpropionic Acid, 2,2,...	238	C11H17F3O2	1000376-54-1	47
2	Cycloheptane, methyl-	112	C8H16	004126-78-7	46
3	Cyclohexanemethanol, 2-methyl-	128	C8H16O	002105-40-0	43
4	4,4-Dimethyl-oct-5-enal	154	C10H18O	1010143-73-6	38
5	Sulfurous acid, cyclohexylmethyl...	402	C23H46O3S	1000309-22-4	35





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Acq On : 08 Nov 2022 12:21
Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

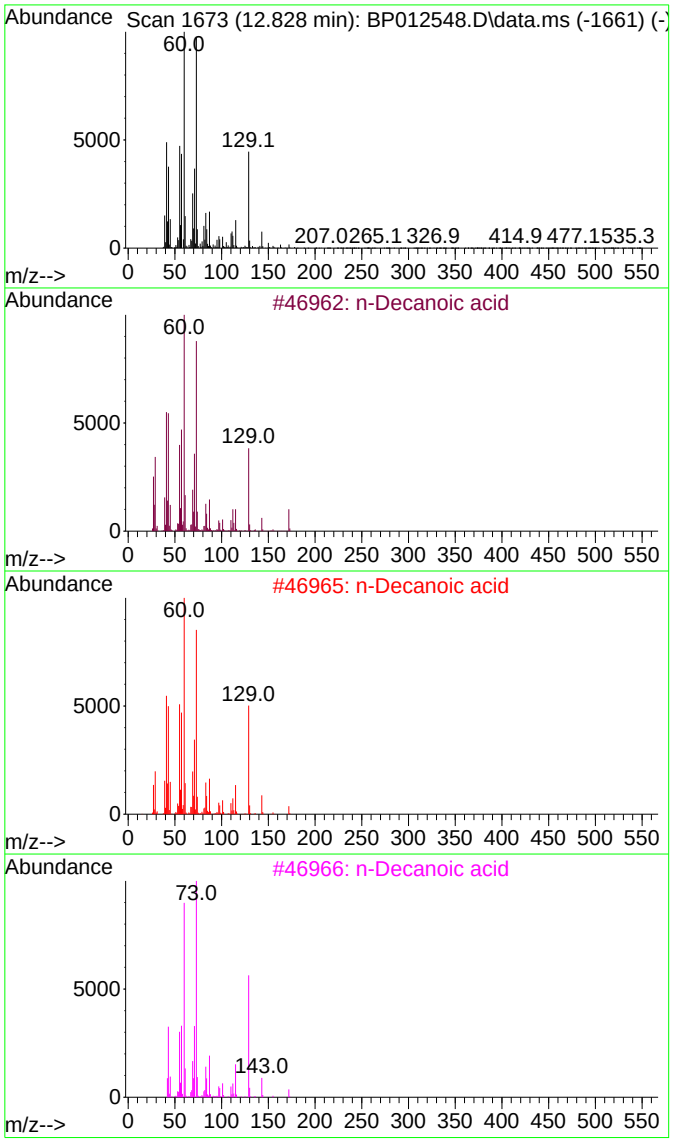
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

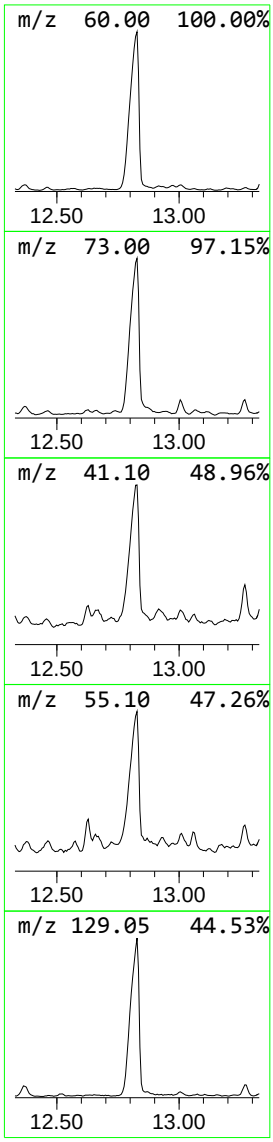
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 46 n-Decanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.828	32.72 ng/ul	6315030	Acenaphthene-d10	14.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Decanoic acid	172	C10H20O2	000334-48-5	94
2	n-Decanoic acid	172	C10H20O2	000334-48-5	91
3	n-Decanoic acid	172	C10H20O2	000334-48-5	91
4	n-Decanoic acid	172	C10H20O2	000334-48-5	90
5	n-Decanoic acid	172	C10H20O2	000334-48-5	72





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Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

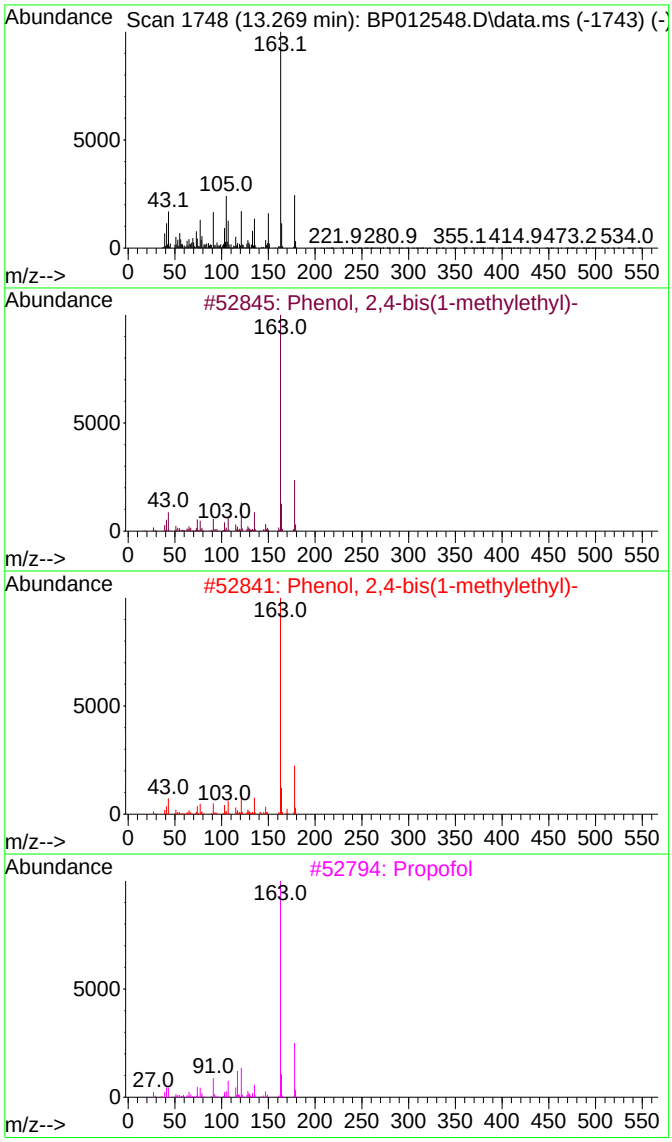
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

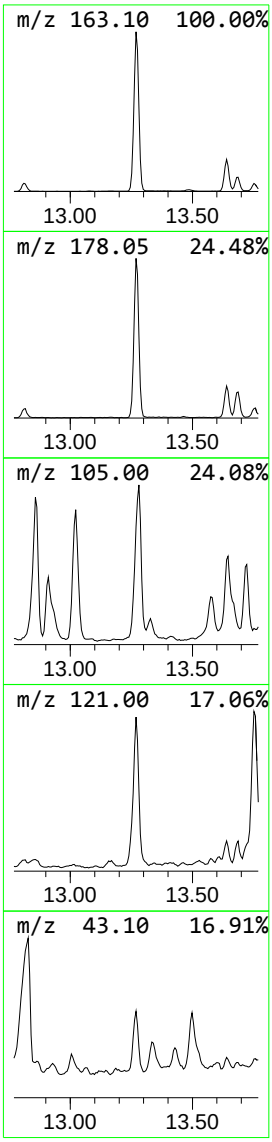
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 49 Phenol, 2,4-bis(1-methyleth... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.269	15.26 ng/ul	2944690	Acenaphthene-d10	14.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,4-bis(1-methylethyl)-	178	C12H18O	002934-05-6	94
2	Phenol, 2,4-bis(1-methylethyl)-	178	C12H18O	002934-05-6	90
3	Propofol	178	C12H18O	002078-54-8	76
4	Propofol	178	C12H18O	002078-54-8	72
5	Propofol	178	C12H18O	002078-54-8	70





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Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

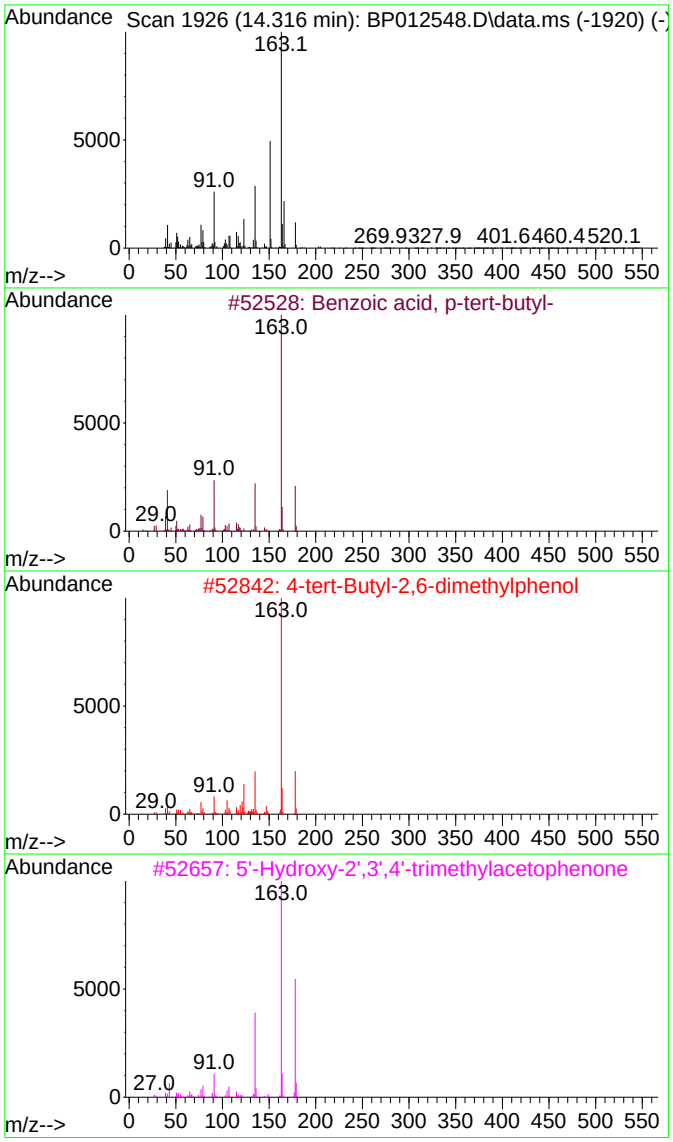
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

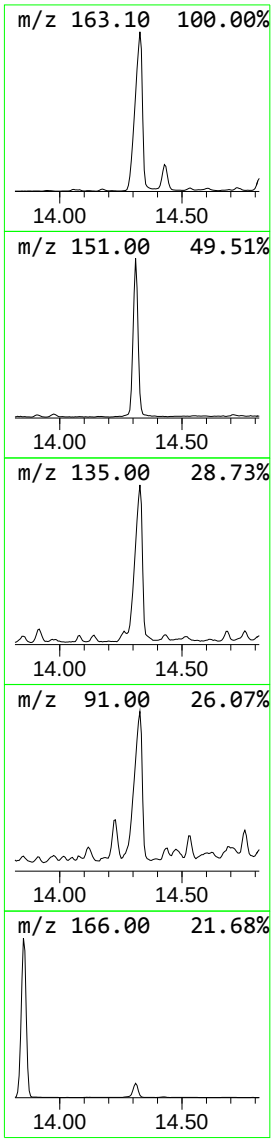
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 55 Benzoic acid, p-tert-butyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.316	36.28 ng/ul	7002980	Acenaphthene-d10	14.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, p-tert-butyl-	178	C11H14O2	000098-73-7	91
2	4-tert-Butyl-2,6-dimethylphenol	178	C12H18O	000879-97-0	52
3	5'-Hydroxy-2',3',4'-trimethylace...	178	C11H14O2	013667-28-2	43
4	Benzoic acid, 4-acetyl-, methyl ...	178	C10H10O3	003609-53-8	43
5	1,3,4-Trimethyladamantane	178	C13H22	1000214-98-3	43





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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

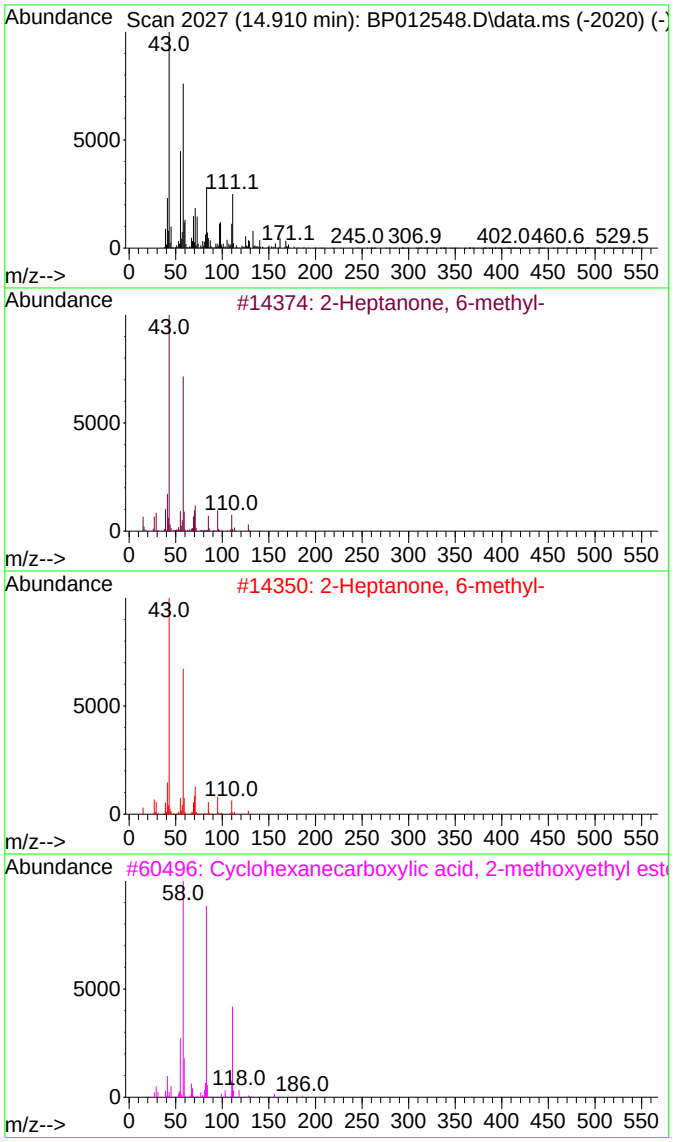
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

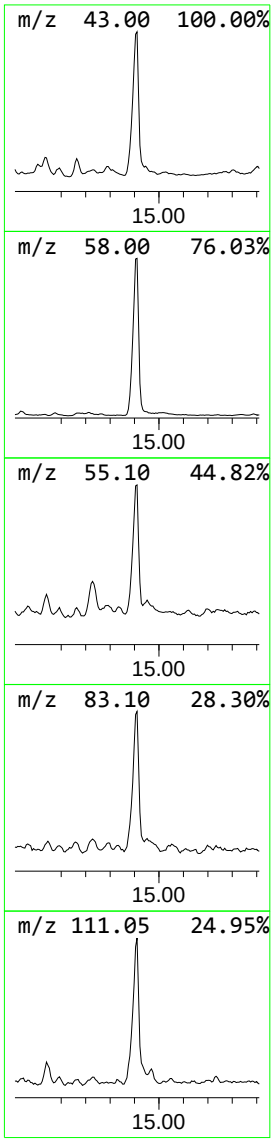
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 57 unknown-04 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.910	16.74 ng/ul	3230840	Acenaphthene-d10	14.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone, 6-methyl-	128	C8H16O	000928-68-7	43
2	2-Heptanone, 6-methyl-	128	C8H16O	000928-68-7	43
3	Cyclohexanecarboxylic acid, 2-me...	186	C10H18O3	1000330-87-5	40
4	2-Octanone	128	C8H16O	000111-13-7	38
5	Propanenitrile, 3-(ethylamino)-	98	C5H10N2	021539-47-9	38





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Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

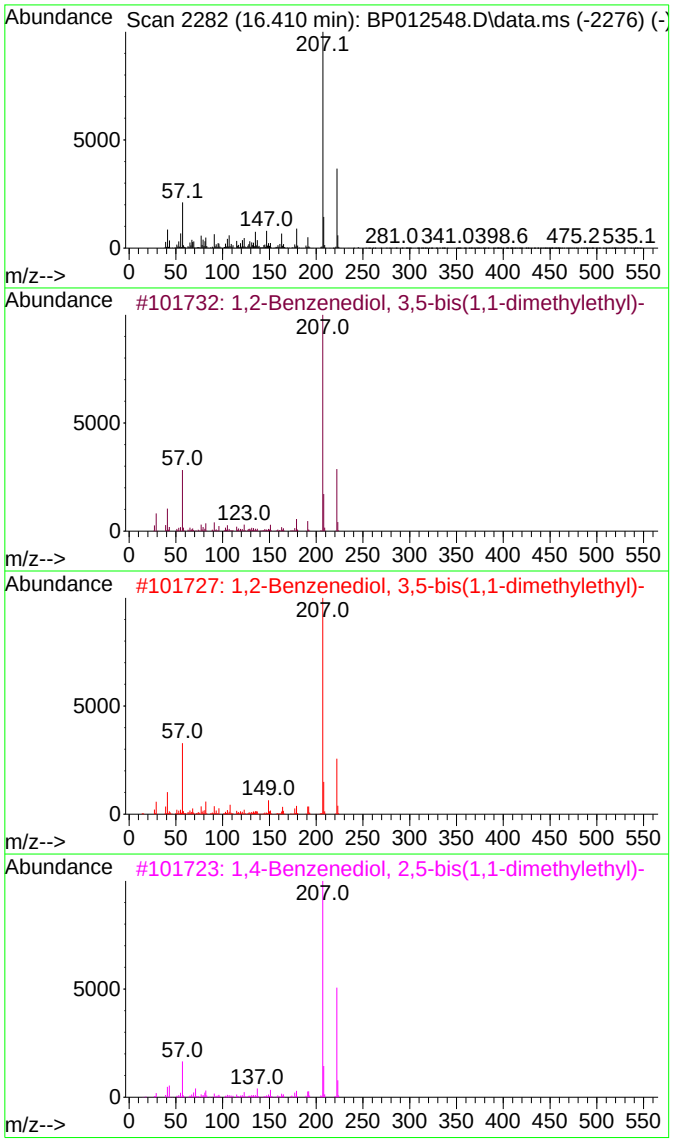
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

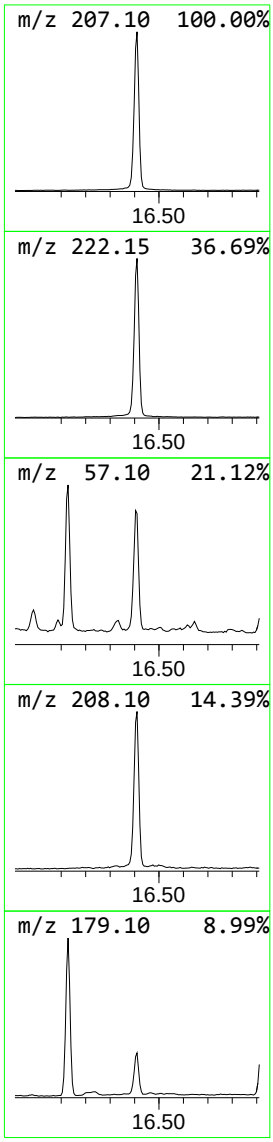
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 64 1,2-Benzenediol, 3,5-bis(1,... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.410	22.32 ng/ul	4855780	Phenanthrene-d10	17.192

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Benzenediol, 3,5-bis(1,1-dim...	222	C14H22O2	001020-31-1	87
2	1,2-Benzenediol, 3,5-bis(1,1-dim...	222	C14H22O2	001020-31-1	86
3	1,4-Benzenediol, 2,5-bis(1,1-dim...	222	C14H22O2	000088-58-4	83
4	1,4-Benzenediol, 2,5-bis(1,1-dim...	222	C14H22O2	000088-58-4	80
5	1,2-Benzenediol, 3,5-bis(1,1-dim...	222	C14H22O2	001020-31-1	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110822\
Data File : BP012548.D
Acq On : 08 Nov 2022 12:21
Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

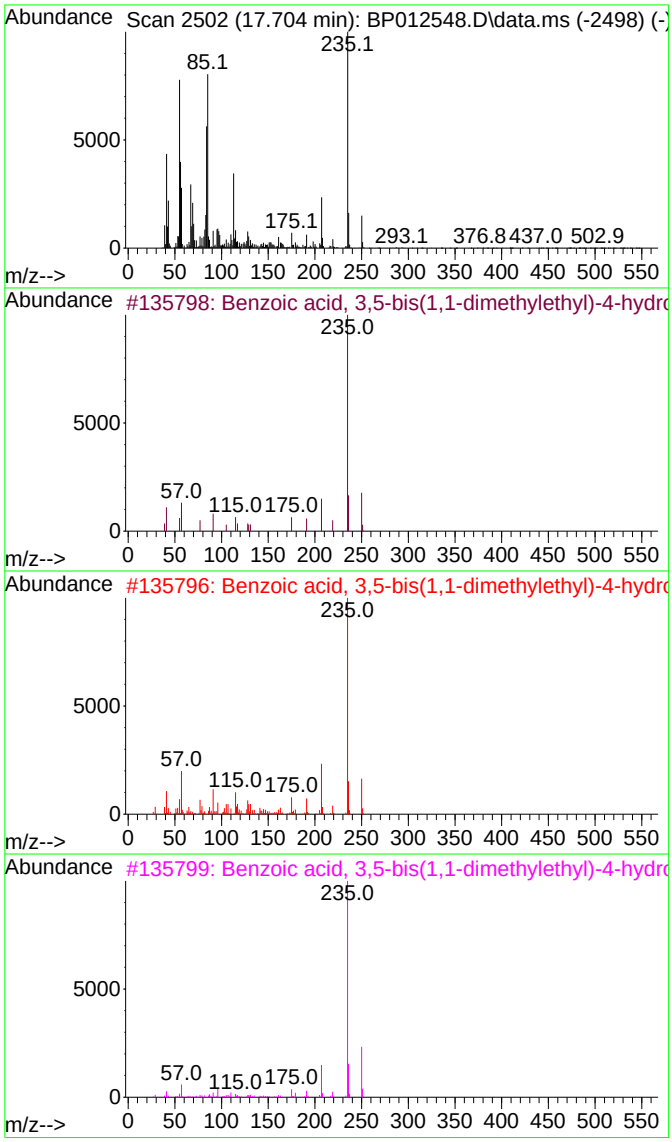
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

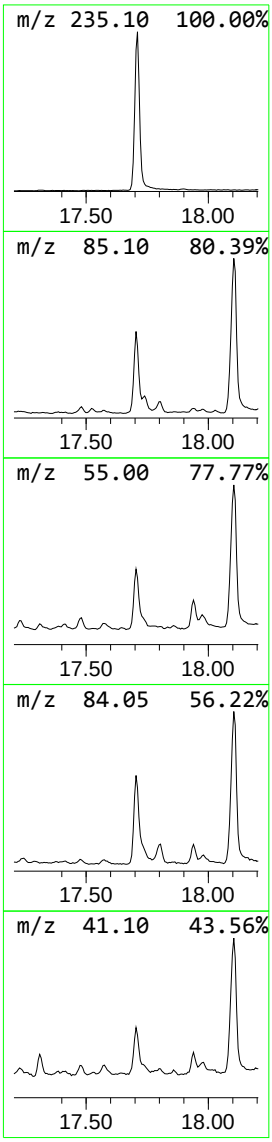
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 66 Benzoic acid, 3,5-bis(1,1-d... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.704	15.76 ng/ul	3427600	Phenanthrene-d10	17.192

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 3,5-bis(1,1-dimeth...	250	C15H22O3	001421-49-4	74
2	Benzoic acid, 3,5-bis(1,1-dimeth...	250	C15H22O3	001421-49-4	56
3	Benzoic acid, 3,5-bis(1,1-dimeth...	250	C15H22O3	001421-49-4	56
4	2-Oxepanone, 7-hexyl-	198	C12H22O2	016429-21-3	53
5	4-tert-Butyl-2,6-dimethylphenol,...	250	C15H26OSi	1000495-37-2	35





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110822\
Data File : BP012548.D
Acq On : 08 Nov 2022 12:21
Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

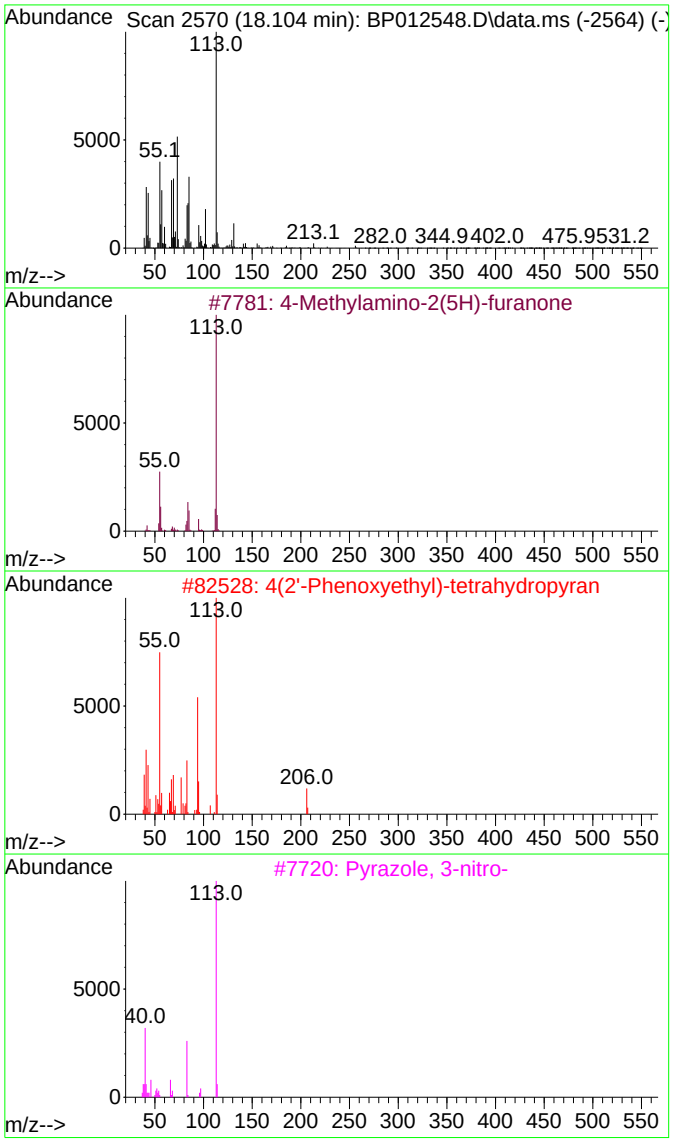
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

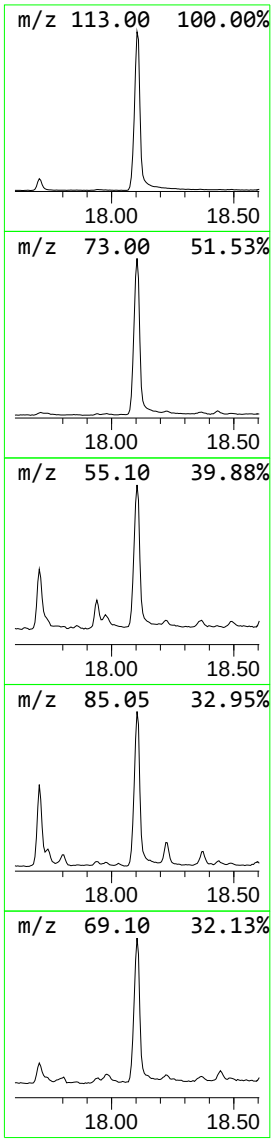
****

Peak Number 68 unknown-05 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.104	41.57 ng/ul	9042200	Phenanthrene-d10	17.192

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Methylamino-2(5H)-furanone	113	C5H7NO2	1000427-45-5	47
2	4(2'-Phenoxyethyl)-tetrahydropyran	206	C13H18O2	062635-45-4	43
3	Pyrazole, 3-nitro-	113	C3H3N3O2	026621-44-3	43
4	2-Thiazoleacetamide, 4-methyl-, ...	228	C9H16N2OSSi	1000479-77-7	43
5	.beta.-Guanidinopropionic acid	131	C4H9N3O2	000353-09-3	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110822\
Data File : BP012548.D
Acq On : 08 Nov 2022 12:21
Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

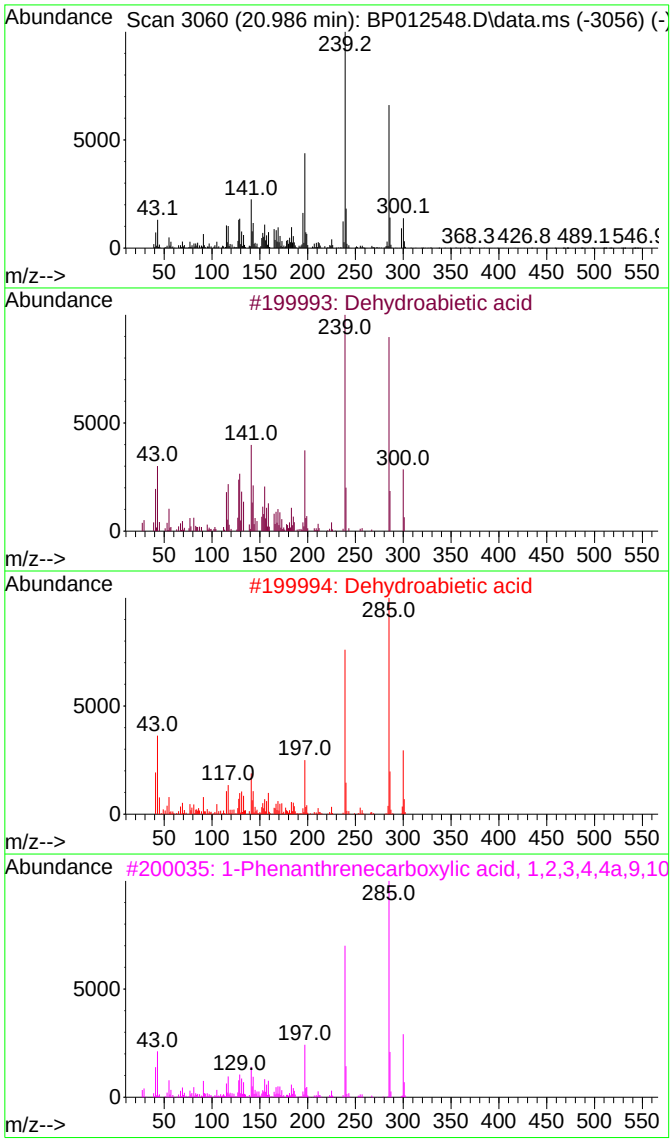
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

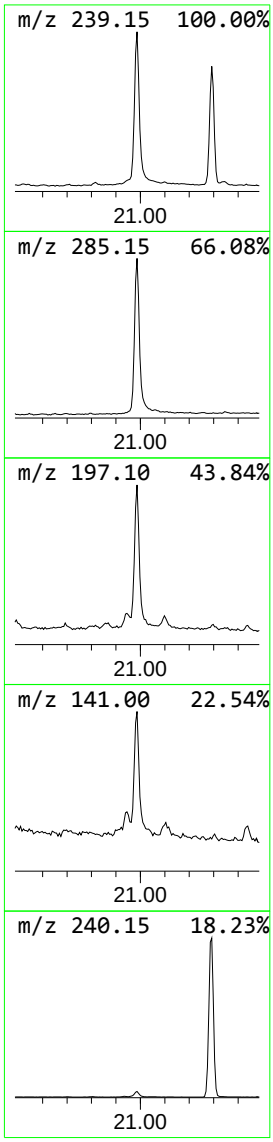
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 73 Dehydroabietic acid Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.986	12.43 ng/ul	2640840	Chrysene-d12	21.292

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dehydroabietic acid	300	C20H28O2	001740-19-8	96
2	Dehydroabietic acid	300	C20H28O2	001740-19-8	93
3	1-Phenanthrenecarboxylic acid, 1...	300	C20H28O2	005155-70-4	93
4	Dehydroabietic acid	300	C20H28O2	001740-19-8	90
5	Dehydroabietic acid	300	C20H28O2	001740-19-8	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110822\
Data File : BP012548.D
Acq On : 08 Nov 2022 12:21
Operator : CG/JU
Sample : N5407-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4Y2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.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
2-Pentanone	3.034	9.1	ng/ul	1331890	1	7.775	2922170	20.0
1-Butanol, 2-me...	3.564	18.4	ng/ul	2691950	1	7.775	2922170	20.0
Propanoic acid,...	3.728	16.4	ng/ul	2389870	1	7.775	2922170	20.0
2-Pentanol, 4-m...	3.811	132.7	ng/ul	19390100	1	7.775	2922170	20.0
1-Pentanol	3.881	16.3	ng/ul	2375550	1	7.775	2922170	20.0
Butanoic acid, ...	5.099	223.3	ng/ul	32623700	1	7.775	2922170	20.0
Butanoic acid, ...	5.287	68.8	ng/ul	10050300	1	7.775	2922170	20.0
1-Hexanol, 2-et...	7.864	57.8	ng/ul	8439550	1	7.775	2922170	20.0
unknown-01	8.811	16.6	ng/ul	2417660	1	7.775	2922170	20.0
Hexanoic acid, ...	9.258	71.3	ng/ul	9977330	2	10.575	2799960	20.0
unknown-02	9.516	16.3	ng/ul	2278560	2	10.575	2799960	20.0
Octanoic acid	10.040	45.5	ng/ul	6364790	2	10.575	2799960	20.0
Phenol, 2-(1-me...	10.510	43.1	ng/ul	6030800	2	10.575	2799960	20.0
3-Cyclohexene-1...	10.640	12.3	ng/ul	1715310	2	10.575	2799960	20.0
p-Cumenol	10.946	31.1	ng/ul	4358670	2	10.575	2799960	20.0
Benzeneacetic acid	11.522	583.9	ng/ul	81744000	2	10.575	2799960	20.0
Nonanoic acid	11.628	19.8	ng/ul	2775600	2	10.575	2799960	20.0
Benzoic acid, 3...	11.663	12.6	ng/ul	1766370	2	10.575	2799960	20.0
Benzoic acid, 4...	11.746	13.8	ng/ul	1933650	2	10.575	2799960	20.0
(DEL) Alkane: S...	11.834	3.8	ng/ul	532327	2	10.575	2799960	20.0
Phenol, p-tert-...	11.940	28.8	ng/ul	4028010	2	10.575	2799960	20.0
Indole	12.093	26.6	ng/ul	3730450	2	10.575	2799960	20.0
unknown-03	12.275	11.4	ng/ul	1588440	2	10.575	2799960	20.0
n-Decanoic acid	12.828	32.7	ng/ul	6315030	3	14.428	3860150	20.0
Phenol, 2,4-bis...	13.269	15.3	ng/ul	2944690	3	14.428	3860150	20.0
Benzoic acid, p...	14.316	36.3	ng/ul	7002980	3	14.428	3860150	20.0
unknown-04	14.910	16.7	ng/ul	3230840	3	14.428	3860150	20.0
1,2-Benzenediol...	16.410	22.3	ng/ul	4855780	4	17.192	4350850	20.0
Benzoic acid, 3...	17.704	15.8	ng/ul	3427600	4	17.192	4350850	20.0
unknown-05	18.104	41.6	ng/ul	9042200	4	17.192	4350850	20.0
Dehydroabietic ...	20.986	12.4	ng/ul	2640840	5	21.292	4248520	20.0