LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
Data File : BP012463.D
Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

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: BP012463.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.028	3	7	17	rBV	3394526	5264013	9.20%	0.883%
2	3.975	155	168	185	rBV2	21331426	57242844	100.00%	9.601%
3	5.099	349	359	369	rBV	8245232	13879588	24.25%	2.328%
4	5.293	385	392	403	rVV	16128888	29625586	51.75%	4.969%
5	5.434	407	416	425	rVV	18024835	54443693	95.11%	9.131%
6	5.799	470	478	484	rBV	15263809	28221130	49.30%	4.733%
7	6.263	547	557	563	rBV	3349595	5646556	9.86%	0.947%
8	6.452	583	589	596	rBV2	955714	1503570	2.63%	0.252%
9	6.758	628	641	653	rBV	2101877	3583231	6.26%	0.601%
10	6.869	653	660	665	rBV	4638383	7276585	12.71%	1.220%
11	6.928	665	670	676	rVV2	2774076	4502978	7.87%	0.755%
12	6.999	676	682	687	rVV2	3008970	5329400	9.31%	0.894%
13	7.046	687	690	696	rVB2	864269	1254099	2.19%	0.210%
14	7.122	696	703	707	rBV2	1724898	3008470	5.26%	0.505%
15	7.169	707	711	715	rVV	2698389	4105364	7.17%	0.689%
16	7.216	715	719	726	rVB2	1667291	2576890	4.50%	0.432%
17	7.322	732	737	743	rVV	1346313	2662650	4.65%	0.447%
18	7.428	749	755	761	rVV	10134303	16583955	28.97%	2.781%
19	7.481	761	764	771	rVB	775150	1136244	1.98%	0.191%
20	7.787	803	816	819	rBV	1135848	2049341	3.58%	0.344%
21	7.816	819	821	827	rVV2	732290	1044167	1.82%	0.175%
22	7.893	828	834	843	rVB2	3580278	6425493	11.22%	1.078%
23	7.981	843	849	858	rVB	1532441	2339081	4.09%	0.392%
24	8.146	870	877	883	rBV2	1440736	3119885	5.45%	0.523%
25	8.234	883	892	896	rVV	12281379	24353115	42.54%	4.085%
26	8.269	896	898	903	rVB2	1382268	1634261	2.85%	0.274%
27	8.328	903	908	916	rBV2	863629	1432244	2.50%	0.240%
28	8.434	916	926	935	rBV	13955667	35293141	61.66%	5.919%
29	8.510	935	939	944	rVB	1185993	1844419	3.22%	0.309%
30	8.575	944	950	959	rBV3	956053	2358232	4.12%	0.396%
31	8.787	981	986	992	rVB	889629	1462118	2.55%	0.245%
32	8.887	992	1003	1009	rBV	1043620	1739346	3.04%	0.292%
33	8.957	1009	1015	1018	rBV	1776135	2999544	5.24%	0.503%
34	8.993	1018	1021	1025	rVB	1097174	1469008	2.57%	0.246%
35	9.105	1035	1040	1049	rVB	4608071	7201547	12.58%	1.208%
36	9.222	1050	1060	1072	rVB2	737851	1529214	2.67%	0.256%

LSC Area Percent Report

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Data File : BP012463.D
Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

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	9.410	1087	1092	1097	rBV	595646	937695	1.64%	0.157%
38	9.469	1097	1102	1107	rVB	778818	1109265	1.94%	0.186%
39	9.675	1131	1137	1144	rBV	1442131	2420933	4.23%	0.406%
40	9.904	1173	1176	1180	rVV	535730	921528	1.61%	0.155%
41	9.946	1180	1183	1186	rVV2	665173	1130605	1.98%	0.190%
42	10.004	1186	1193	1203	rVV4	1842800	4393934	7.68%	0.737%
43	10.093	1203	1208	1213	rVV	654804	980943	1.71%	0.165%
44	10.210	1217	1228	1236	rBV3	3443612	9923380	17.34%	1.664%
45	10.293	1236	1242	1250	rVV4	1193213	2648774	4.63%	0.444%
46	10.404	1257	1261	1267	rVB	793971	1225518	2.14%	0.206%
47	10.481	1268	1274	1279	rBV	1163286	2022801	3.53%	0.339%
48	10.587	1286	1292	1296	rVV	1509681	2490746	4.35%	0.418%
49	10.634	1296	1300	1304	rVV	1399338	2290518	4.00%	0.384%
50	10.699	1304	1311	1313	rVV	4545846	9334735	16.31%	1.566%
51	10.746	1313	1319	1332	rVB2	14707234	34843071	60.87%	5.844%
52	11.004	1349	1363	1368	rBV5	1033837	2904886	5.07%	0.487%
53	11.187	1389	1394	1403	rVB2	1053596	2049232	3.58%	0.344%
54	11.434	1427	1436	1439	rBV5	447268	962187	1.68%	0.161%
55	11.481	1439	1444	1451	rVB3	820766	1448131	2.53%	0.243%
56	11.610	1456	1466	1472	rVV3	1045436	2776738	4.85%	0.466%
57	11.687	1472	1479	1484	rVV	7466471	13045698	22.79%	2.188%
58	11.728	1484	1486	1493	rVB	905810	1331388	2.33%	0.223%
59	11.846	1501	1506	1513	rVB2	679127	1598028	2.79%	0.268%
60	11.940	1513	1522	1530	rBV2	1741002	4343422	7.59%	0.728%
61	12.351	1582	1592	1596	rBV4	349589	874173	1.53%	0.147%
62	12.404	1597	1601	1609	rVB6	410701	905835	1.58%	0.152%
63	12.922	1684	1689	1694	rBV2	1045267	1683068	2.94%	0.282%
64	13.028	1699	1707	1719	rVB2	998725	2690251	4.70%	0.451%
65	13.157	1722	1729	1734	rBV2	592190	1159849	2.03%	0.195%
66	13.381	1762	1767	1771	rBV	1045766	1299111	2.27%	0.218%
67	13.545	1789	1795	1797	rBV2	518541	916349	1.60%	0.154%
68	13.857	1840	1848	1856	rVB	4212682	6782310	11.85%	1.138%
69	13.945	1856	1863	1872	rBV	2022608	3882529	6.78%	0.651%
70	14.057	1877	1882	1885	rVV	919912	1363992	2.38%	0.229%
71	14.134	1889	1895	1904	rVB	4507774	7088673	12.38%	1.189%
72	14.440	1941	1947	1961	rVB2	2946683	5213962	9.11%	0.874%
73	14.687	1984	1989	1996	rVV3	525907	948161	1.66%	0.159%
74	14.987	2034	2040	2044	rVV2	1061480	1690156	2.95%	0.283%
75	15.445	2110	2118	2129	rBV2	13705383	27127088	47.39%	4.550%
76	15.575	2135	2140	2144	rBV	2014574	2708059	4.73%	0.454%

LSC Area Percent Report

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Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

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	15.957	2200	2205	2215	rVB3	1243107	2456461	4.29%	0.412%
78	16.087	2222	2227	2232	rBV2	625843	1013390	1.77%	0.170%
79	16.157	2234	2239	2243	rBV	870133	1297616	2.27%	0.218%
80	16.504	2295	2298	2307	rVB5	471155	989154	1.73%	0.166%
81	16.586	2307	2312	2322	rVB5	635937	1506230	2.63%	0.253%
82	17.198	2412	2416	2427	rVB	3387144	5342801	9.33%	0.896%
83	17.298	2427	2433	2440	rBV2	6391868	9675802	16.90%	1.623%
84	17.863	2525	2529	2533	rBV	817274	1085922	1.90%	0.182%
85	18.092	2564	2568	2578	rVB	1700744	2303389	4.02%	0.386%
86	18.310	2600	2605	2612	rVV4	1150573	2027830	3.54%	0.340%
87	18.657	2660	2664	2669	rBV2	656214	900851	1.57%	0.151%
88	18.757	2677	2681	2689	rVB2	751849	1141354	1.99%	0.191%
89	19.128	2740	2744	2748	rVB	1252646	1538078	2.69%	0.258%
90	19.328	2774	2778	2786	rBV	1808080	2293400	4.01%	0.385%
91	19.545	2809	2815	2820	rBV	7600664	10721506	18.73%	1.798%
92	19.598	2820	2824	2830	rVB	1365148	1781270	3.11%	0.299%
93	19.775	2849	2854	2859	rBV	2400659	2878480	5.03%	0.483%
94	20.039	2894	2899	2905	rBV2	859791	1282544	2.24%	0.215%
95	20.998	3058	3062	3073	rVB	1289517	1693748	2.96%	0.284%
96	21.292	3108	3112	3119	rVB	3797543	4948025	8.64%	0.830%
97	21.422	3129	3134	3139	rBV	3137582	4008400	7.00%	0.672%
98	22.521	3317	3321	3334	rVB2	655431	1201627	2.10%	0.202%
99	23.527	3485	3492	3505	rVB2	5015859	10085766	17.62%	1.692%
100	23.680	3512	3518	3528	rVB2	2175476	4441806	7.76%	0.745%

Sum of corrected areas: 596224174

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
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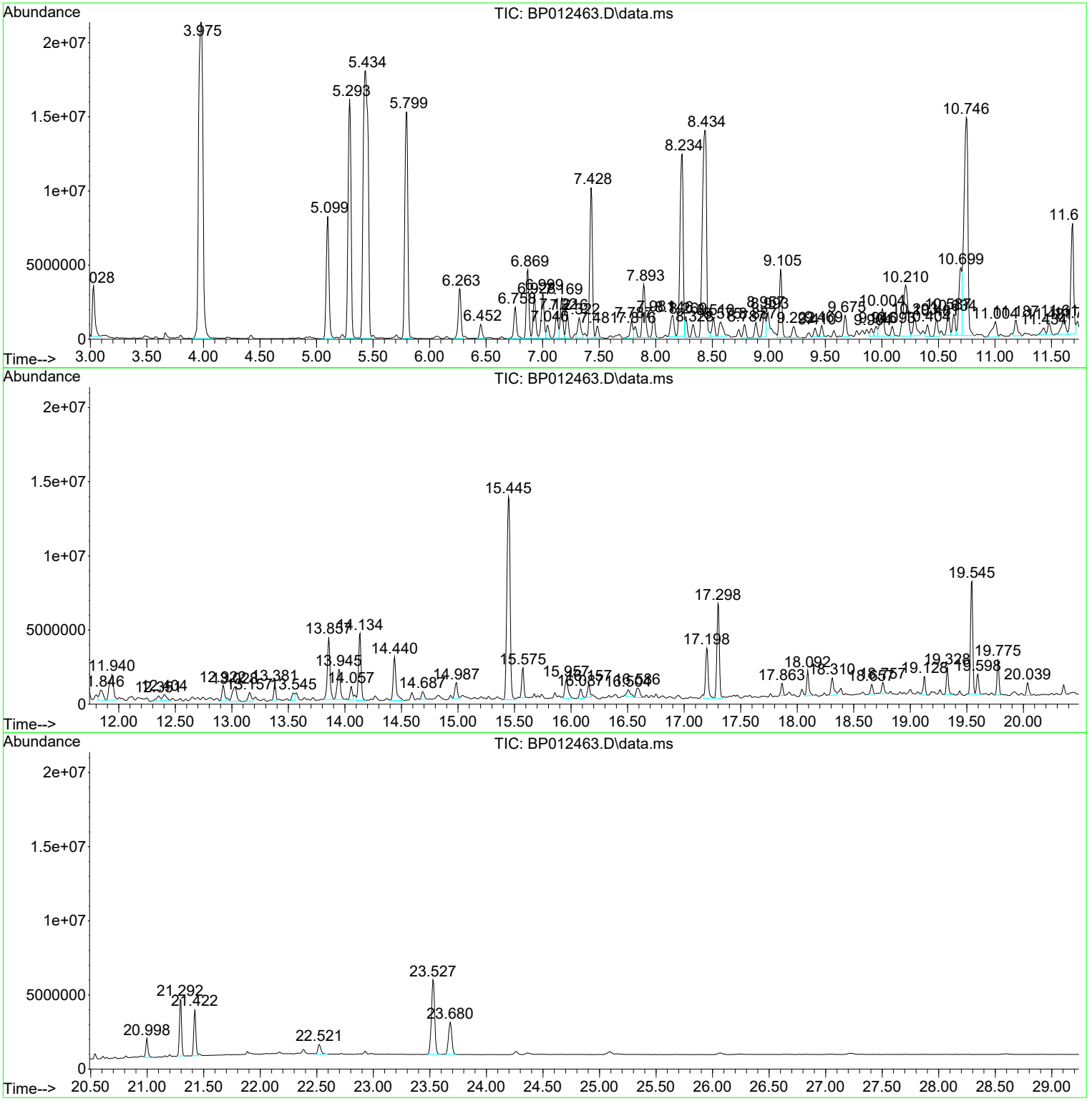
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

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\BP110222\
Data File : BP012463.D
Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
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Instrument :
BNA_P
ClientSampleId :
BHAC5

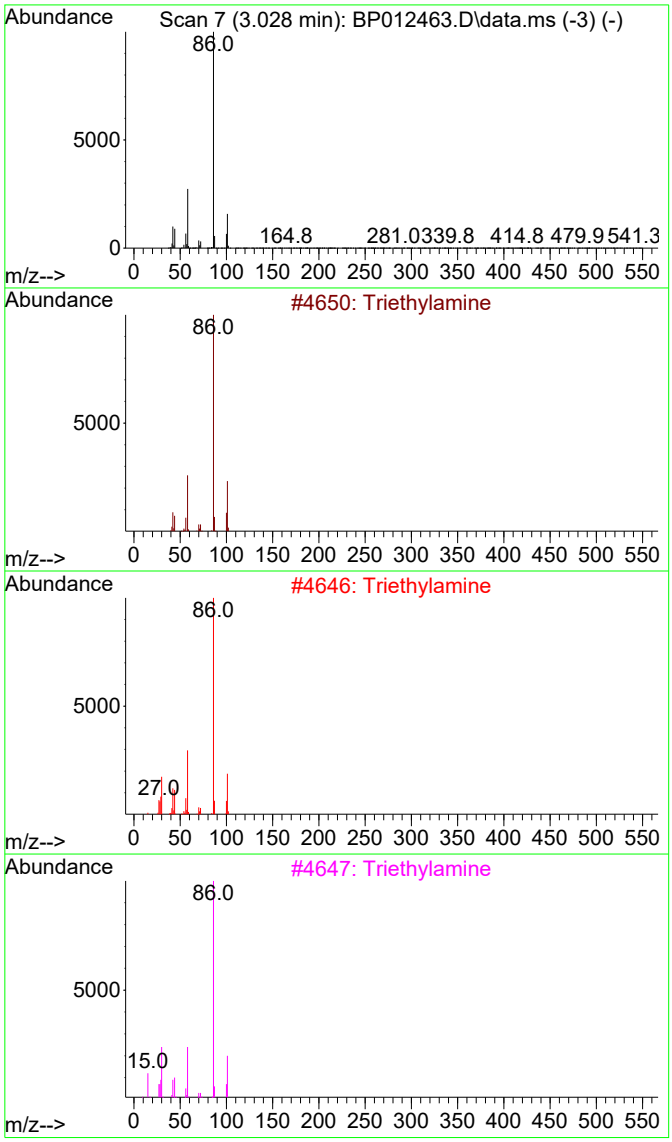
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

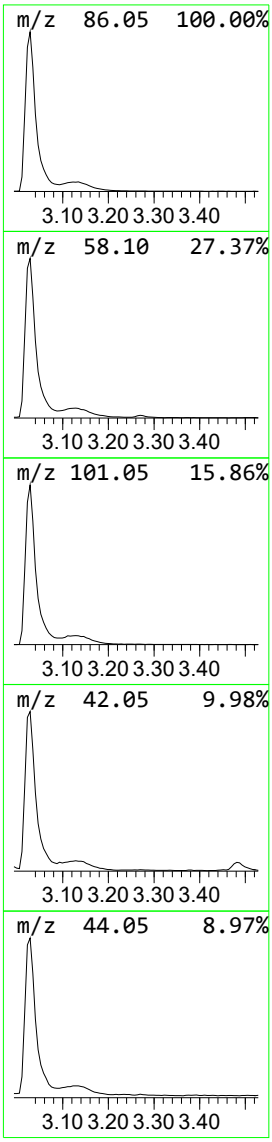
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Triethylamine Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.028	51.37 ng/ul	5264010	1,4-Dichlorobenzene-d4	7.787

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	91
5	Triethylamine	101	C6H15N	000121-44-8	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
Data File : BP012463.D
Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
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BHAC5

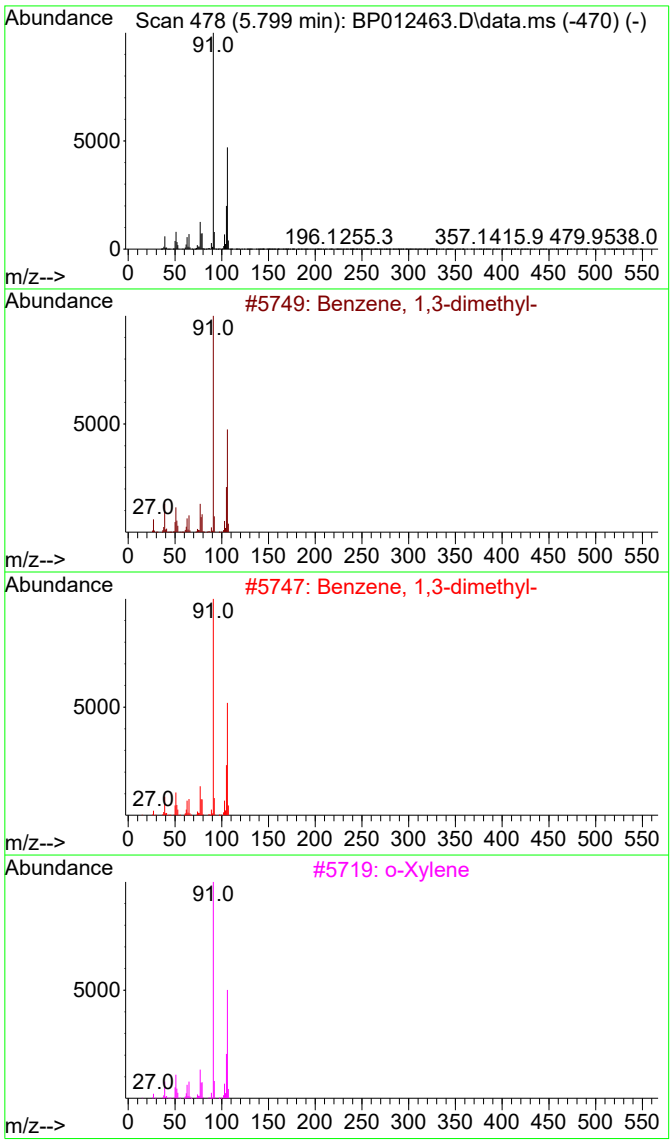
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

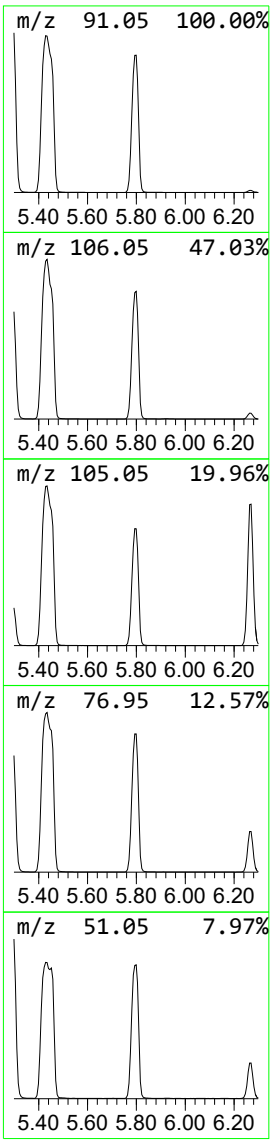
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.799	275.42 ng/ul	28221100	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
3	o-Xylene	106	C8H10	000095-47-6	97
4	p-Xylene	106	C8H10	000106-42-3	95
5	o-Xylene	106	C8H10	000095-47-6	95





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
Data File : BP012463.D
Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

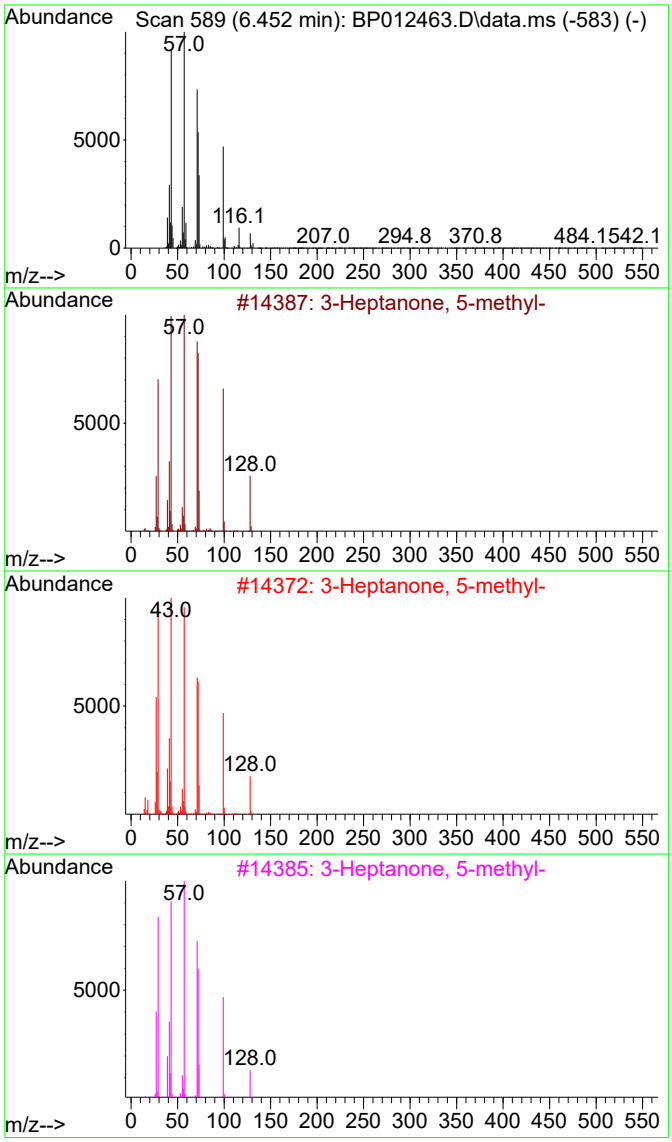
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

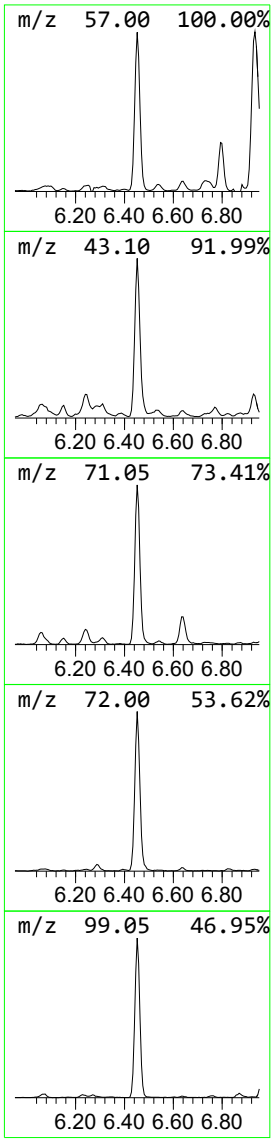
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 3-Heptanone, 5-methyl- Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.452	14.67 ng/ul	1503570	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Heptanone, 5-methyl-	128	C8H16O	000541-85-5	87
2	3-Heptanone, 5-methyl-	128	C8H16O	000541-85-5	72
3	3-Heptanone, 5-methyl-	128	C8H16O	000541-85-5	64
4	Ethyl amylketone	128	C8H16O	020616-93-7	64
5	3,6-Dimethyl-4-octanone	156	C10H20O	034645-03-9	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
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Acq On : 02 Nov 2022 22:19
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Misc :
ALS Vial : 17 Sample Multiplier: 1

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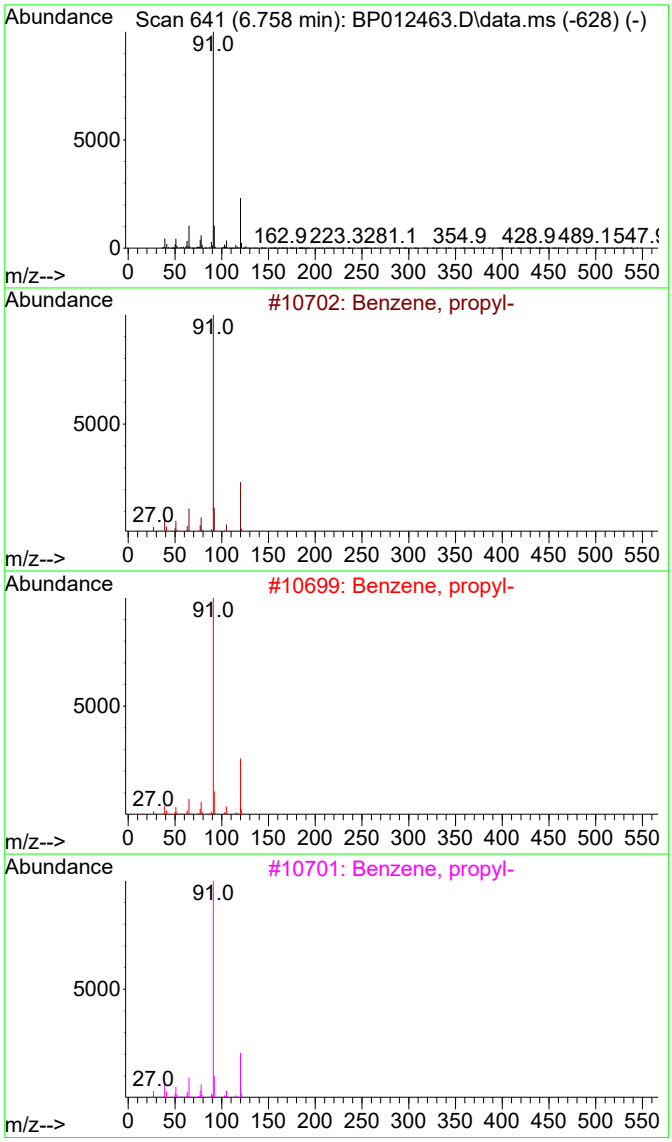
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

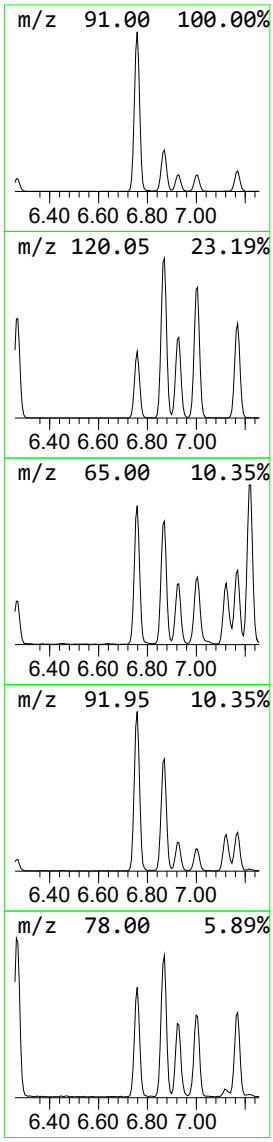
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, propyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.758	34.97 ng/ul	3583230	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	94
2	Benzene, propyl-	120	C9H12	000103-65-1	90
3	Benzene, propyl-	120	C9H12	000103-65-1	90
4	N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	78
5	1-(Benzylamino)propan-2-ol	165	C10H15NO	1000486-84-7	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
Data File : BP012463.D
Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

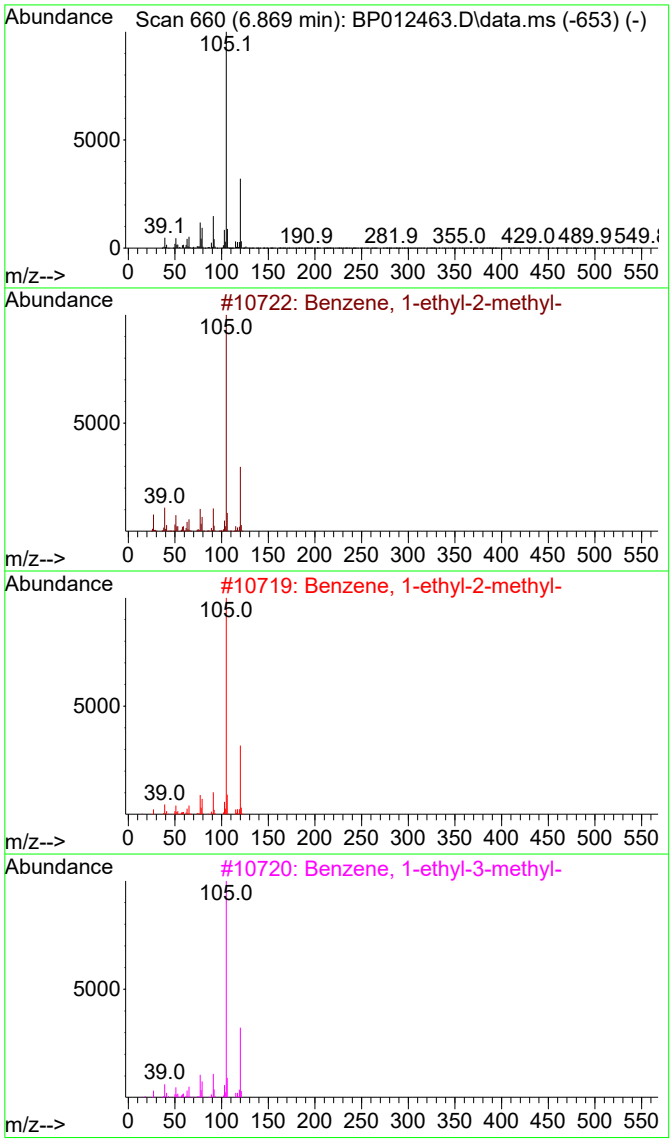
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

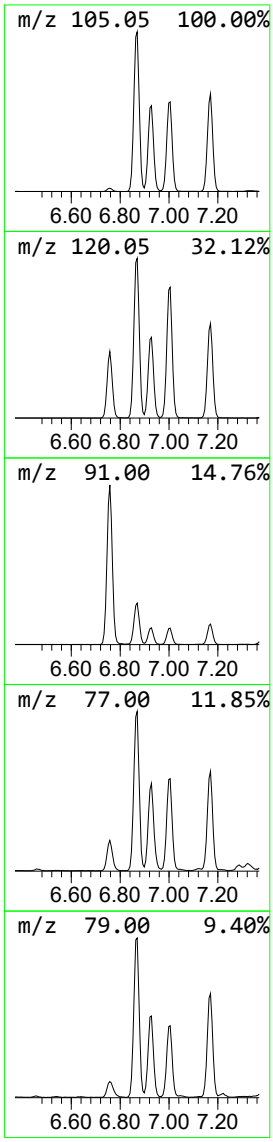
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1-ethyl-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.869	71.01 ng/ul	7276590	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





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Data File : BP012463.D
Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

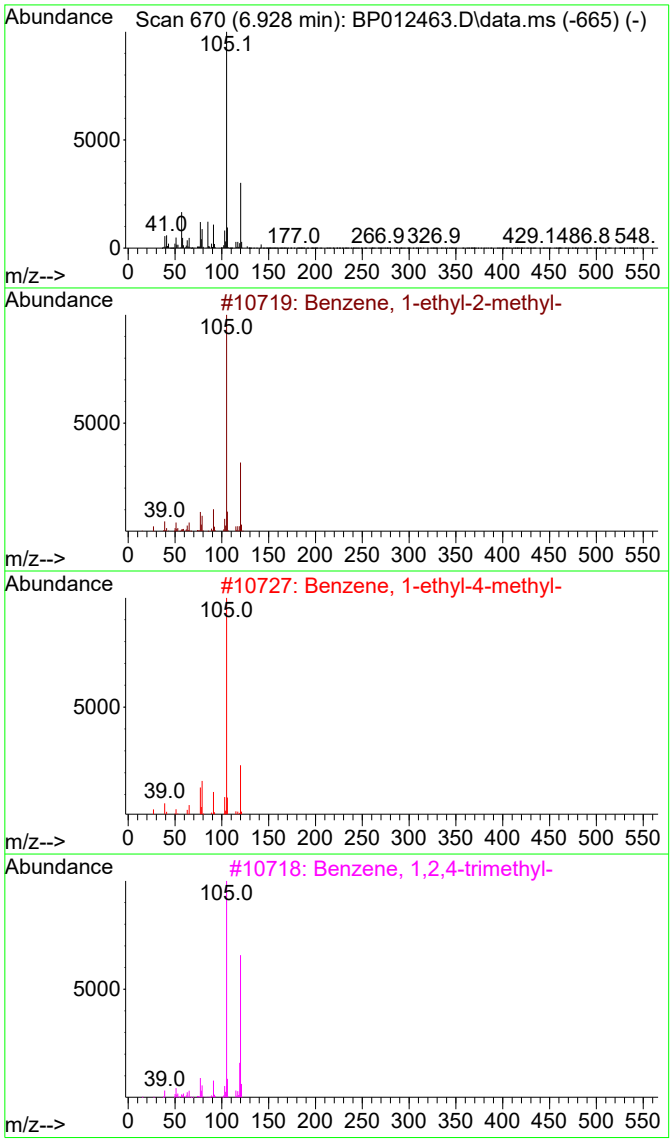
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

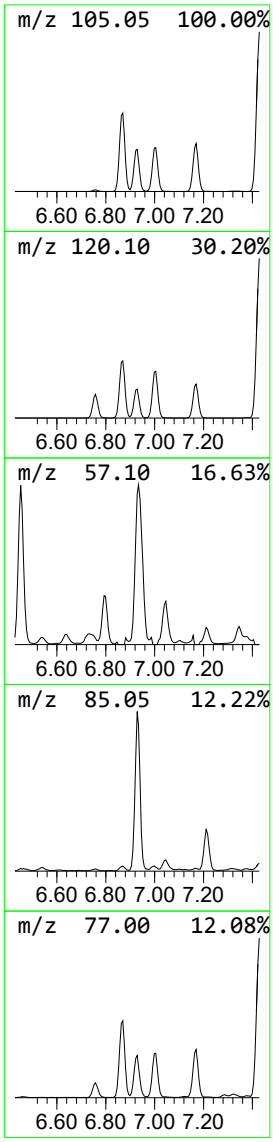
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzene, 1-ethyl-4-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.928	43.95 ng/ul	4502980	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	89
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	76
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	76
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	76





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Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

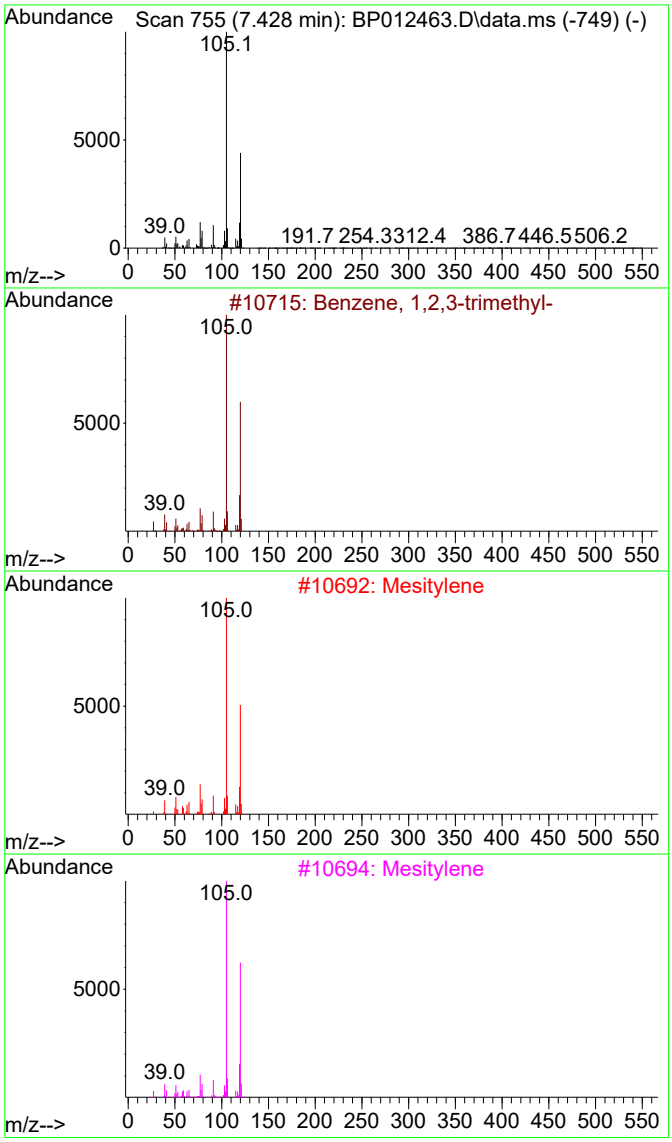
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

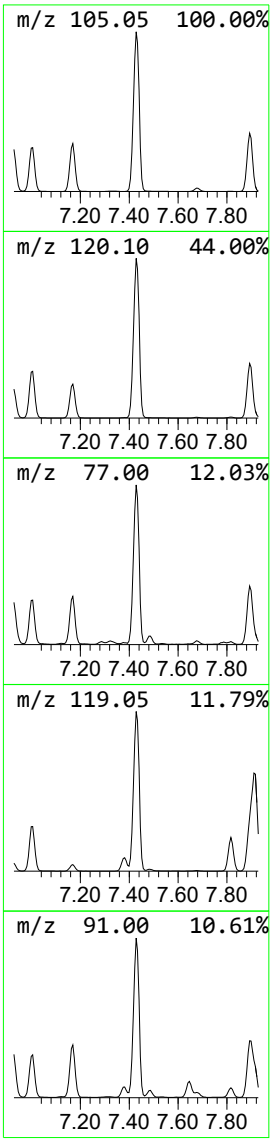
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Benzene, 1,2,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.428	161.85 ng/ul	16584000	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2	Mesitylene	120	C9H12	000108-67-8	97
3	Mesitylene	120	C9H12	000108-67-8	94
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94





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Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

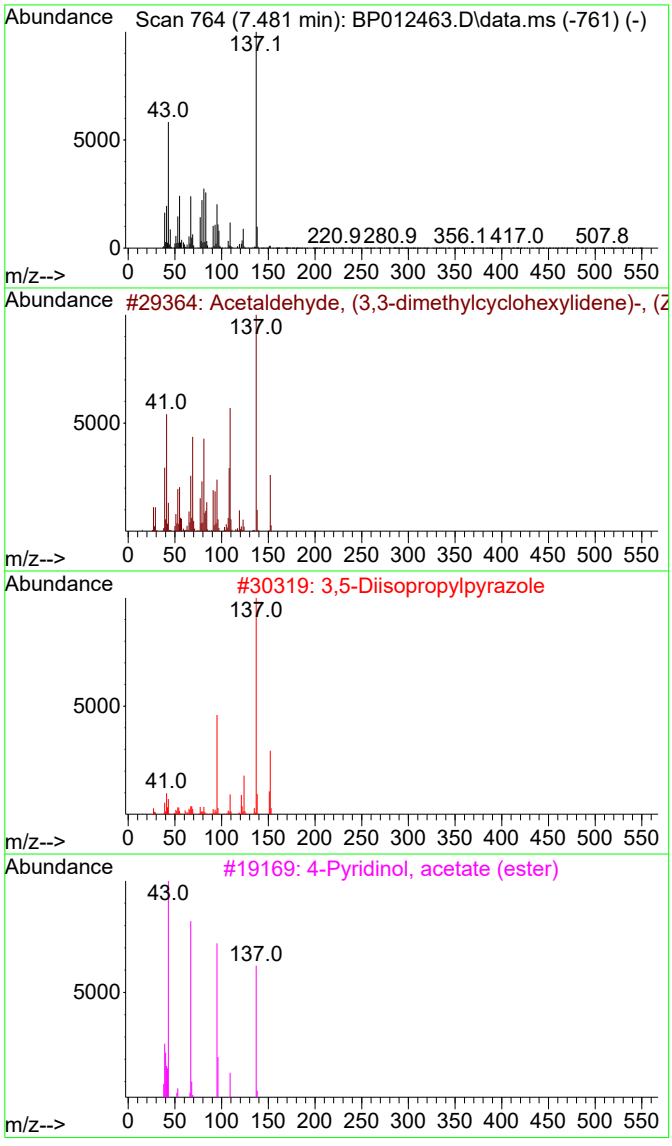
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

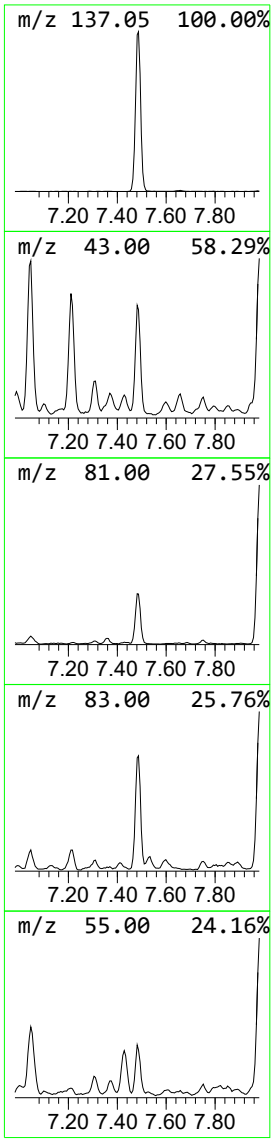
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Acetaldehyde, (3,3-dimethyl... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.481	11.09 ng/ul	1136240	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetaldehyde, (3,3-dimethylcyclo...	152	C10H16O	026532-24-1	50
2	3,5-Diisopropylpyrazole	152	C9H16N2	017536-00-4	47
3	4-Pyridinol, acetate (ester)	137	C7H7NO2	014210-20-9	42
4	3,6-Dimethyl-2,3,3a,4,5,7a-hexah...	152	C10H16O	070786-44-6	42
5	4,5-Heptadien-2-one, 3,3,6-trime...	152	C10H16O	081250-41-1	40





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Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

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

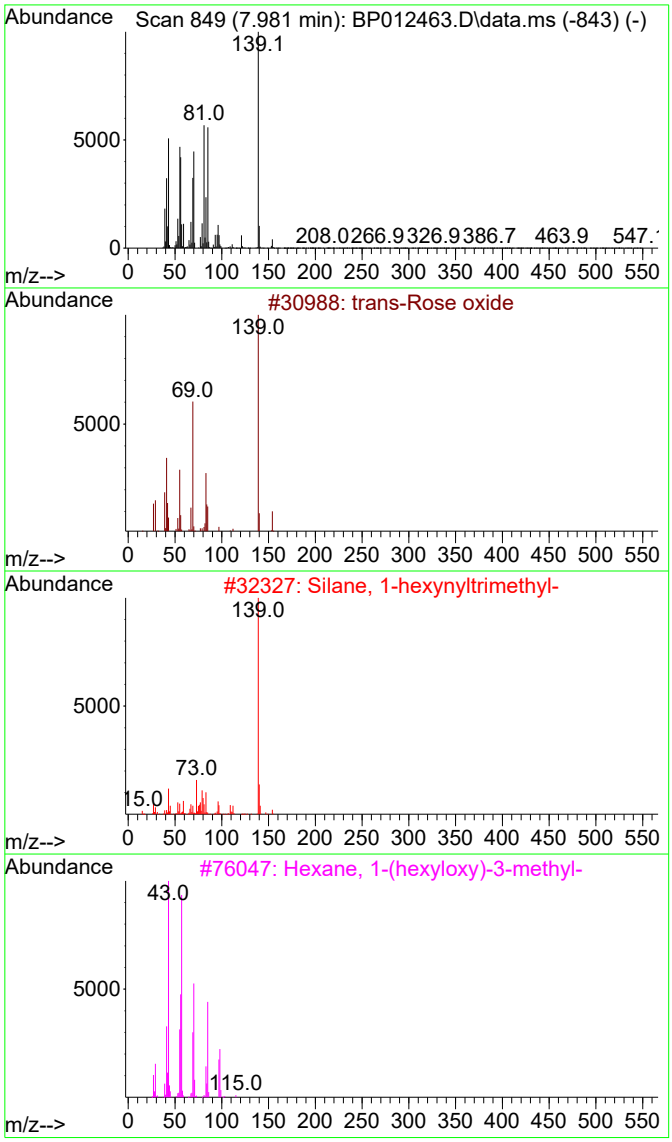
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

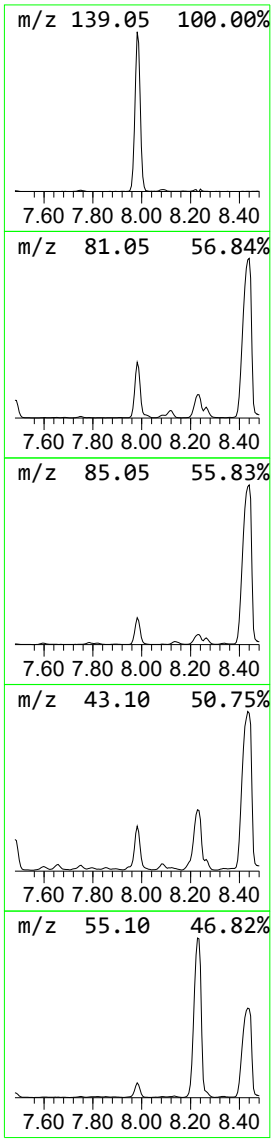
****

Peak Number 15 unknown-01 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.981	22.83 ng/ul	2339080	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	trans-Rose oxide	154	C10H18O	000876-18-6	35
2	Silane, 1-hexynyltrimethyl-	154	C9H18Si	003844-94-8	35
3	Hexane, 1-(hexyloxy)-3-methyl-	200	C13H28O	074421-18-4	32
4	2H-Pyran, tetrahydro-4-methyl-2-...	154	C10H18O	016409-43-1	25
5	(2S,4R)-4-Methyl-2-(2-methylprop...	154	C10H18O	003033-23-6	25





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Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

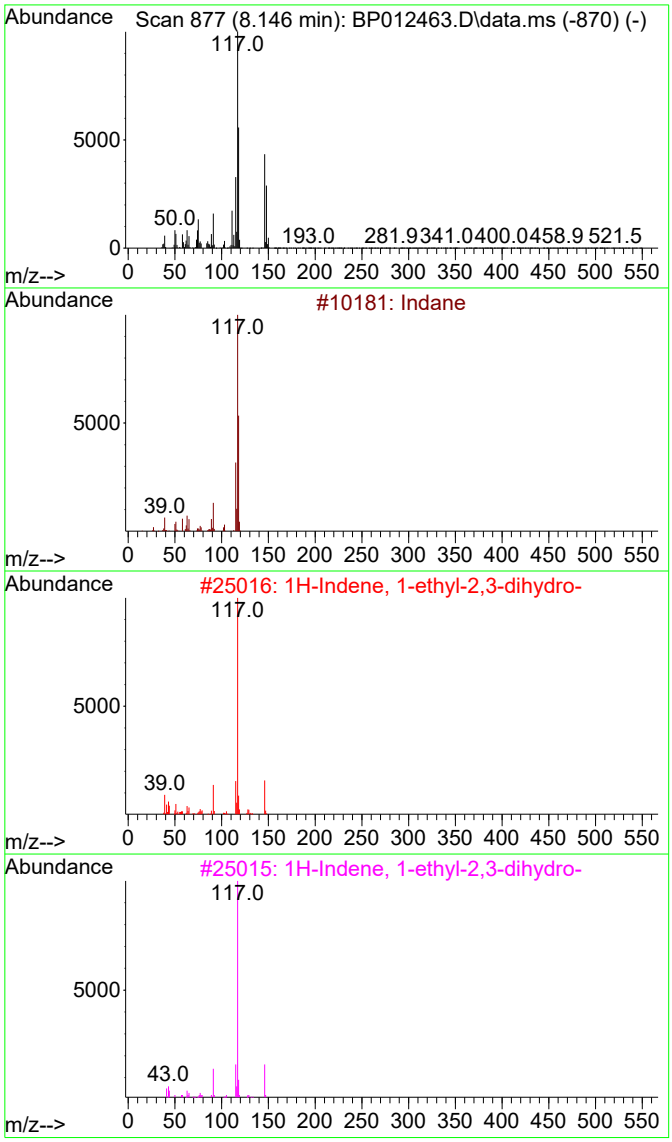
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

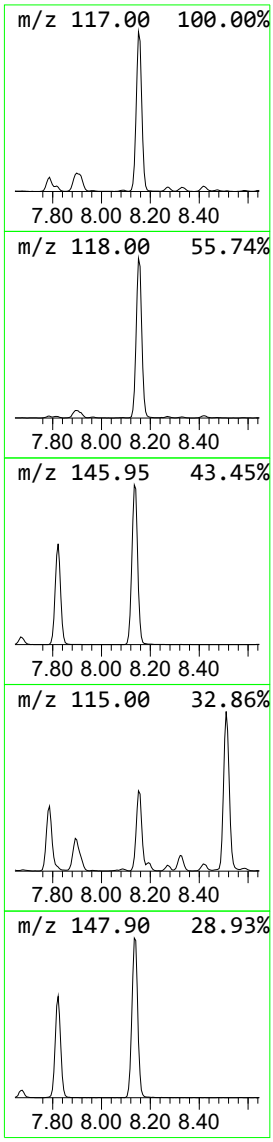
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Indane Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.146	30.45 ng/ul	3119890	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	87
2	1H-Indene, 1-ethyl-2,3-dihydro-	146	C11H14	004830-99-3	47
3	1H-Indene, 1-ethyl-2,3-dihydro-	146	C11H14	004830-99-3	47
4	Deltacyclene	118	C9H10	007785-10-6	46
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	46





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Acq On : 02 Nov 2022 22:19
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Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

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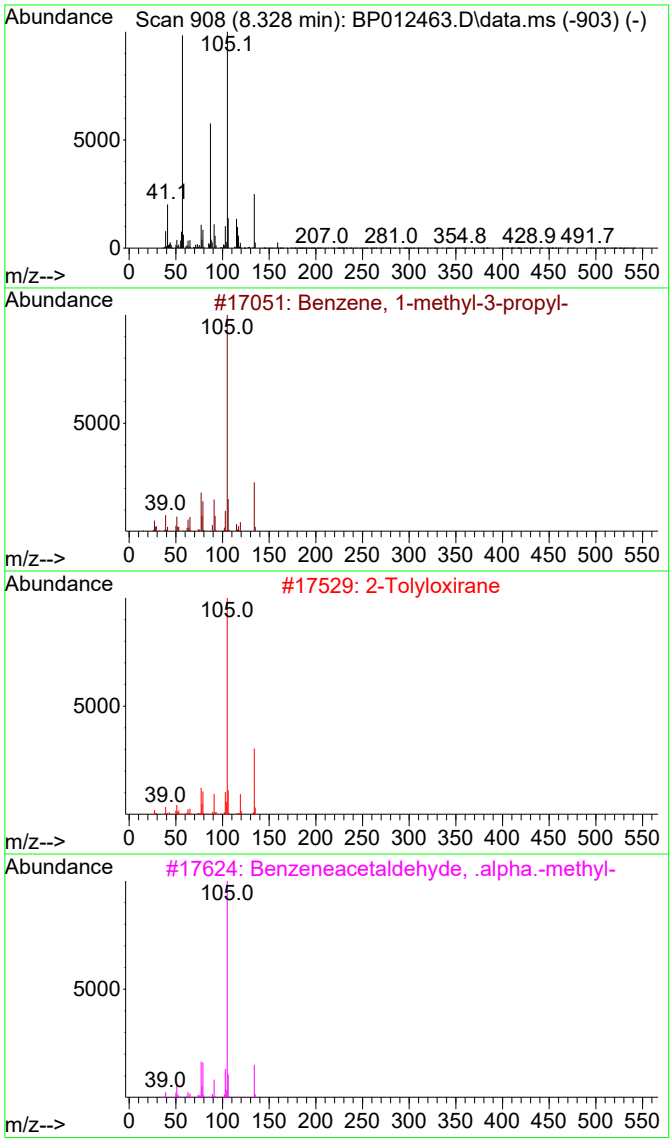
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

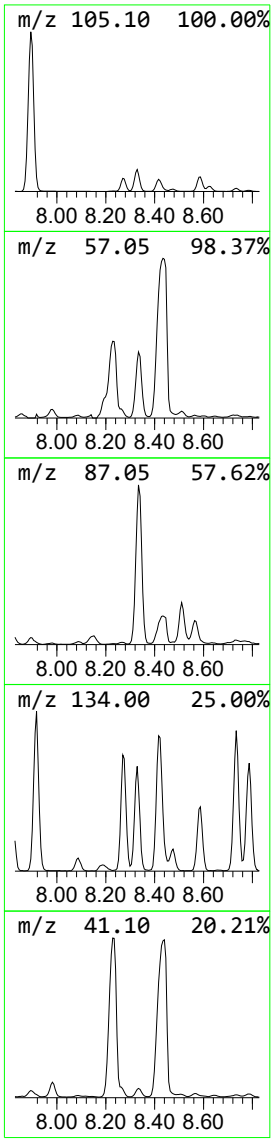
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 unknown-02 Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.328	13.98 ng/ul	1432240	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	46
2	2-Tolyloxirane	134	C9H10O	002783-26-8	42
3	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	42
4	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	42
5	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	38





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Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

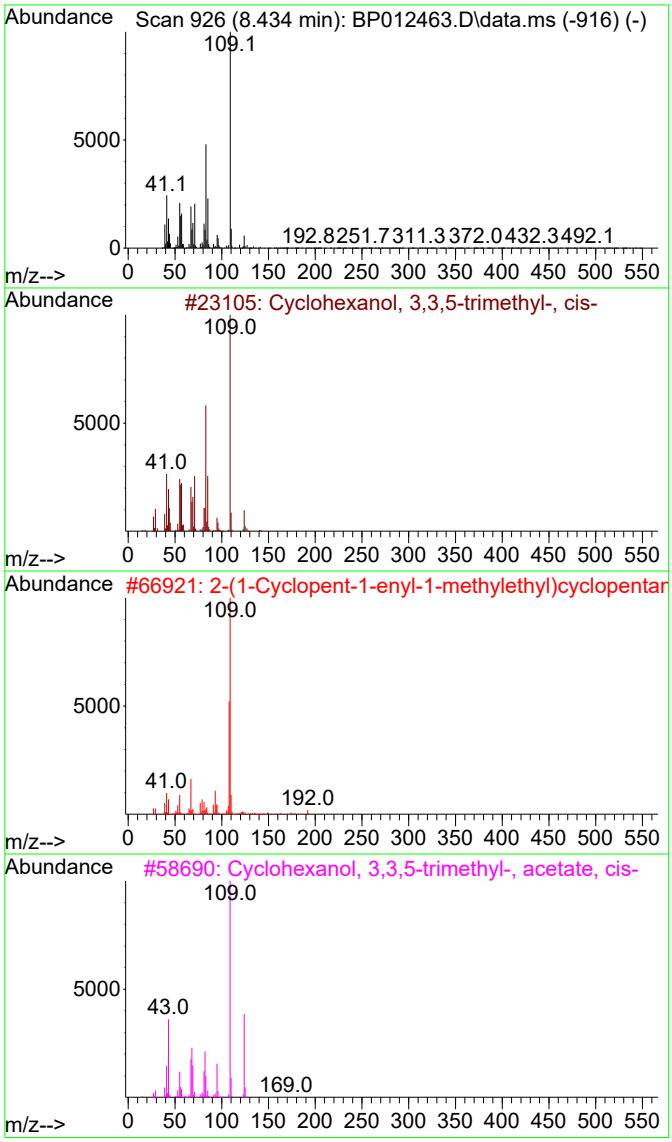
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

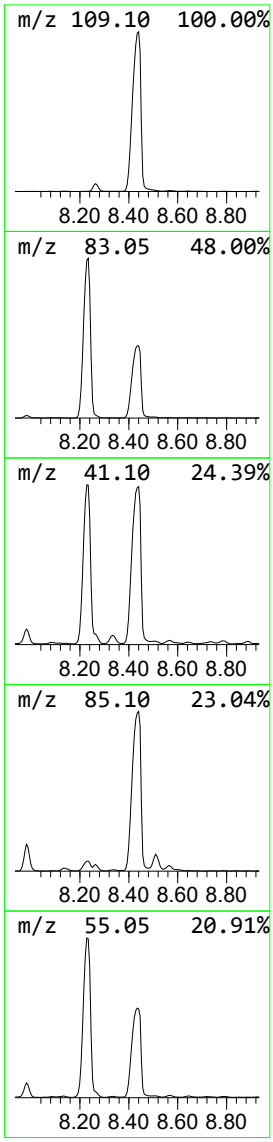
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Cyclohexanol, 3,3,5-trimeth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.434	344.43 ng/ul	35293100	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanol, 3,3,5-trimethyl-, ...	142	C9H18O	000933-48-2	64
2	2-(1-Cyclopent-1-enyl-1-methylet...	192	C13H20O	1000190-57-4	43
3	Cyclohexanol, 3,3,5-trimethyl-, ...	184	C11H20O2	024691-16-5	43
4	Cyclohexanol, 3,3,5-trimethyl-	142	C9H18O	000116-02-9	40
5	Allyltrivinylsilane	150	C9H14Si	115946-69-5	40





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Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

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

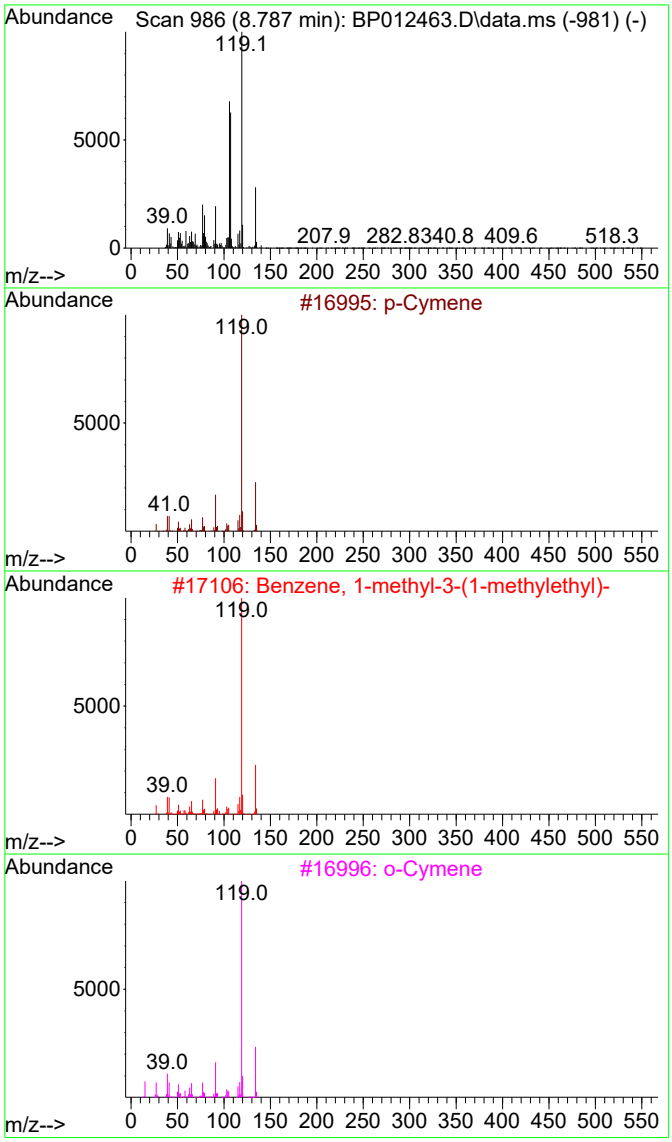
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

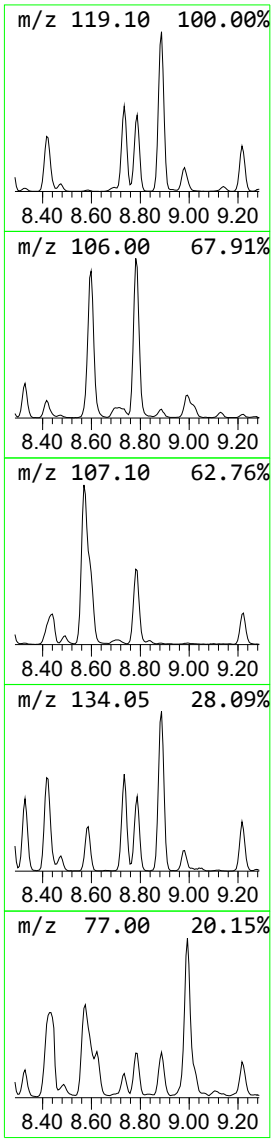
****

Peak Number 19 p-Cymene Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.787	14.27 ng/ul	1462120	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	90
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	90
3	o-Cymene	134	C10H14	000527-84-4	90
4	o-Cymene	134	C10H14	000527-84-4	87
5	o-Cymene	134	C10H14	000527-84-4	87





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Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

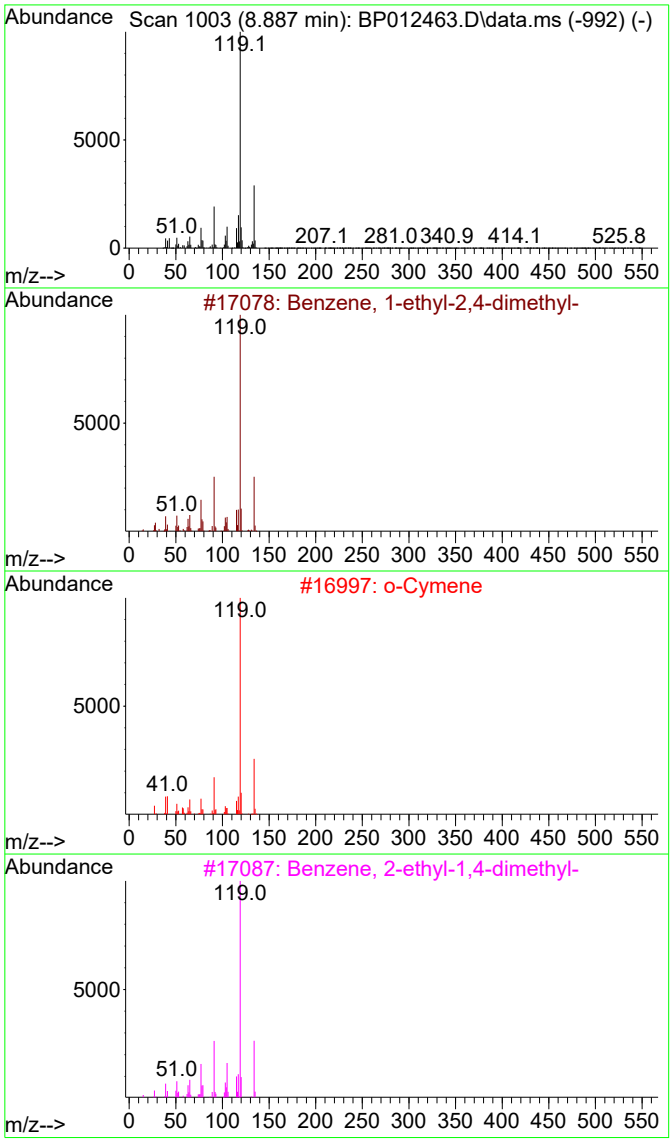
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

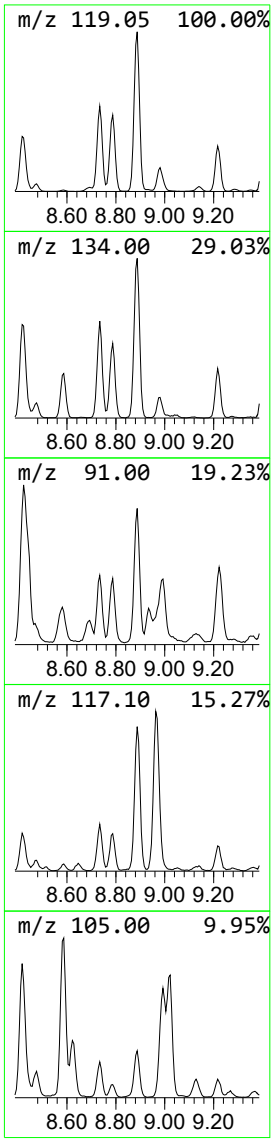
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.887	16.97 ng/ul	1739350	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
2	o-Cymene	134	C10H14	000527-84-4	95
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93





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Misc :
ALS Vial : 17 Sample Multiplier: 1

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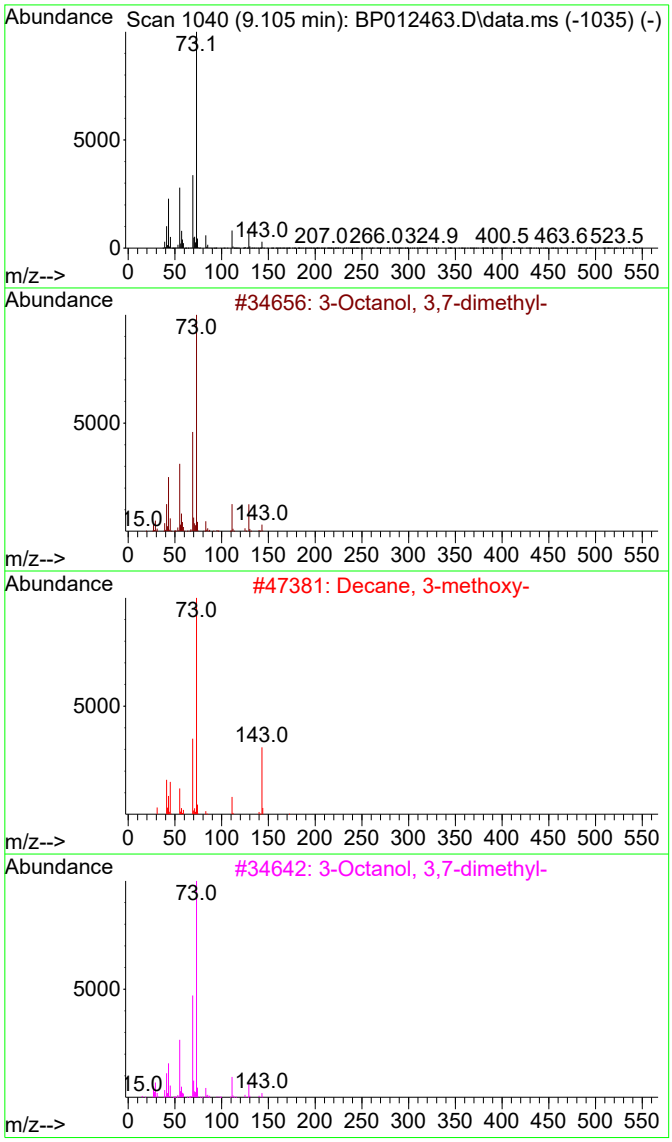
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

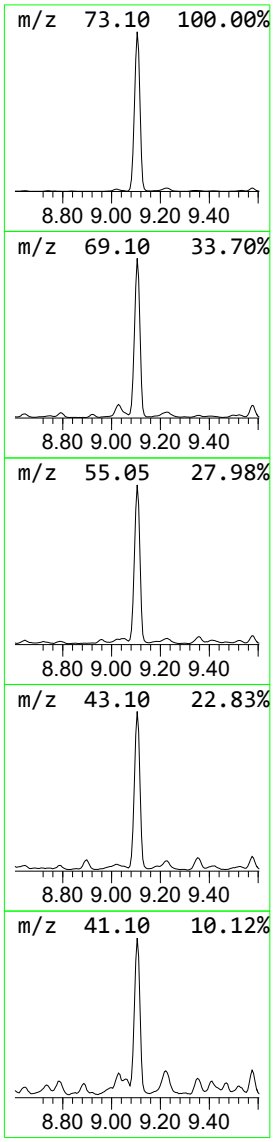
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 3-Octanol, 3,7-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.105	70.28 ng/ul	7201550	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	64
2	Decane, 3-methoxy-	172	C11H24O	055955-64-1	53
3	3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	45
4	3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	45
5	3-Octanol, 3,6-dimethyl-	158	C10H22O	000151-19-9	42





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ALS Vial : 17 Sample Multiplier: 1

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ClientSampleId :
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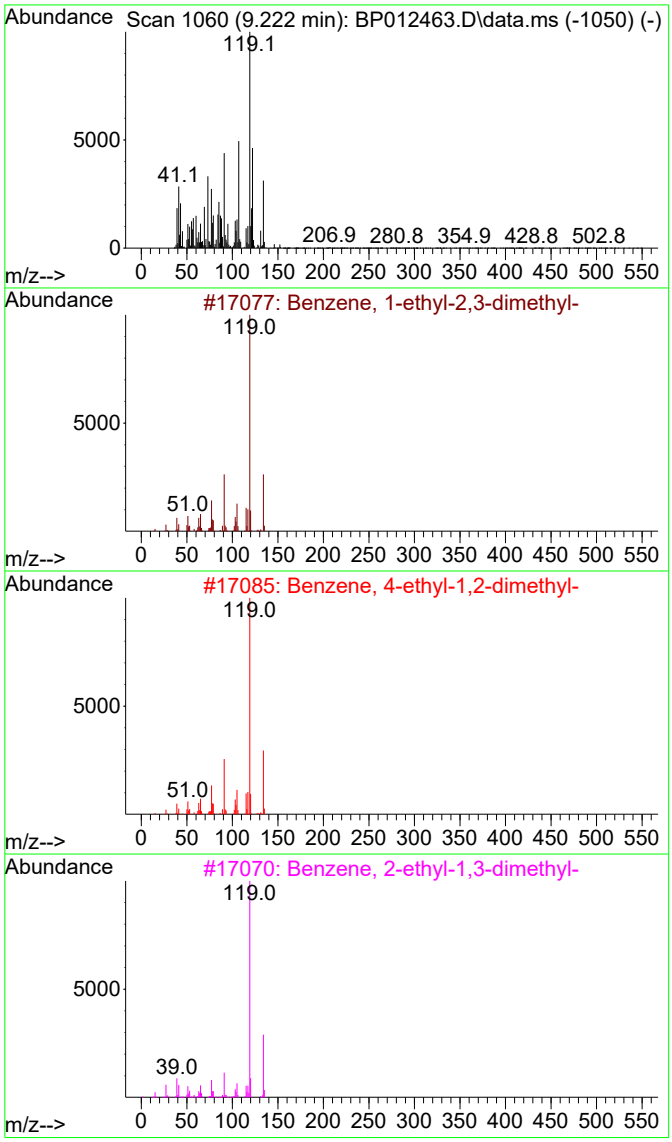
Quant Title : SVOA CALIBRATION

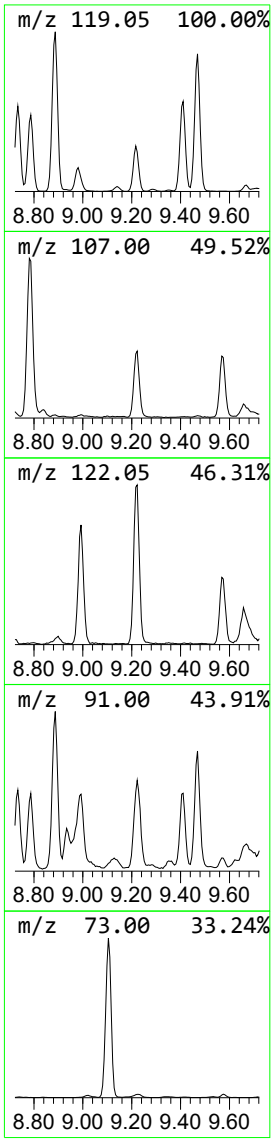
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.222	12.28 ng/ul	1529210	Naphthalene-d8	10.587

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	90
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	89
3	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	87
4	o-Cymene	134	C10H14	000527-84-4	84
5	o-Cymene	134	C10H14	000527-84-4	84





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Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

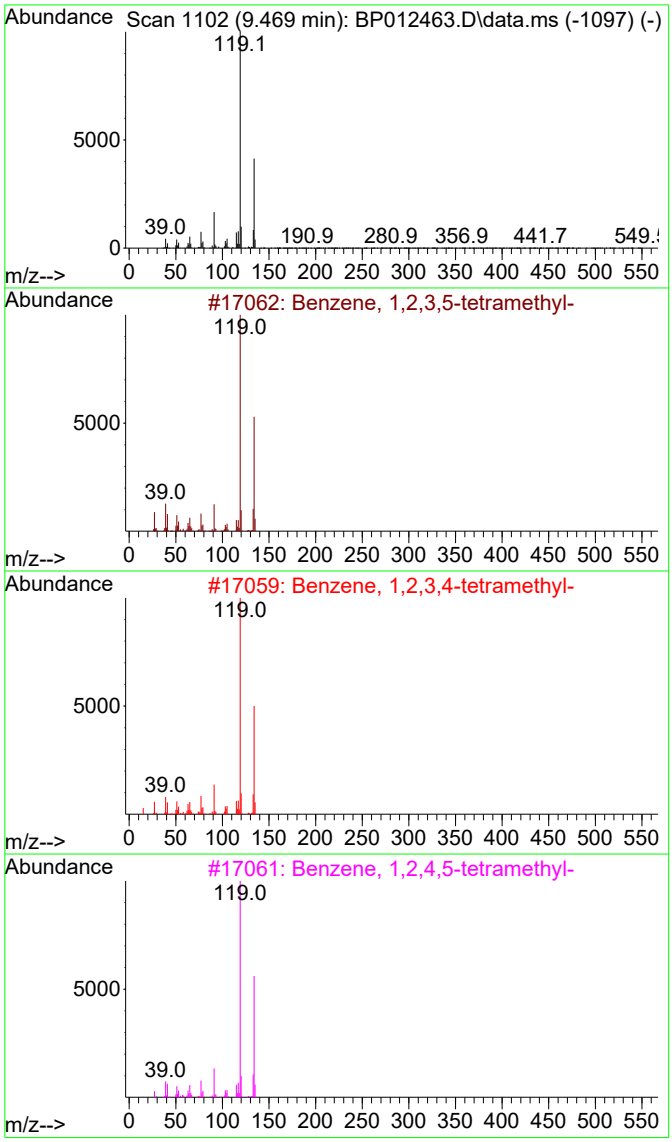
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

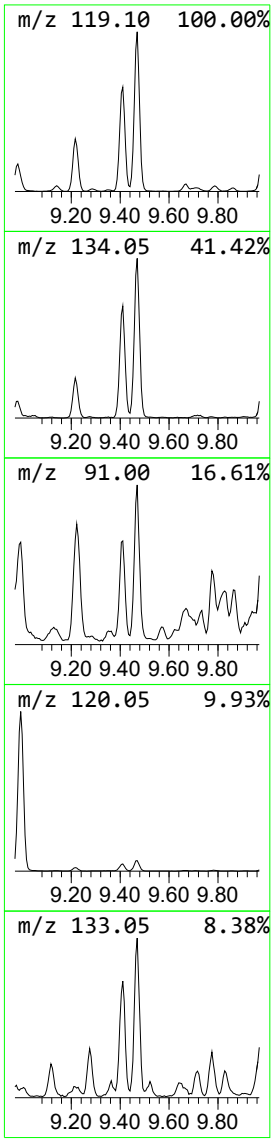
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.469	8.91 ng/ul	1109270	Naphthalene-d8	10.587

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95





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Data File : BP012463.D
Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

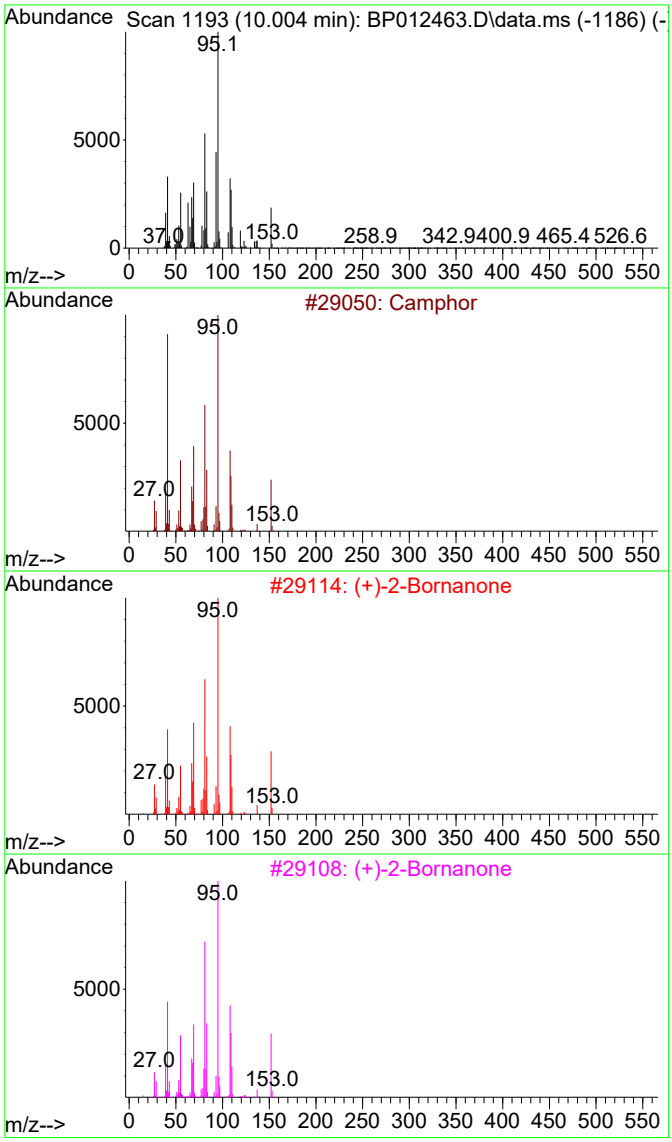
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

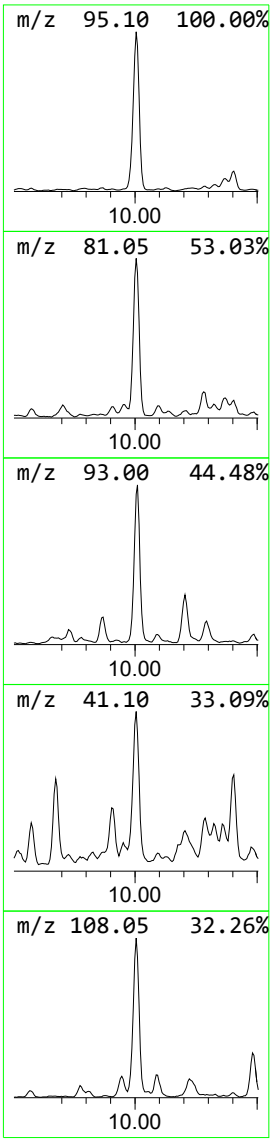
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 Camphor Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.005	35.28 ng/ul	4393930	Naphthalene-d8	10.587

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Camphor	152	C10H16O	000076-22-2	96
2	(+)-2-Bornanone	152	C10H16O	000464-49-3	93
3	(+)-2-Bornanone	152	C10H16O	000464-49-3	90
4	(+)-2-Bornanone	152	C10H16O	000464-49-3	90
5	Camphor	152	C10H16O	000076-22-2	90





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Data File : BP012463.D
Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

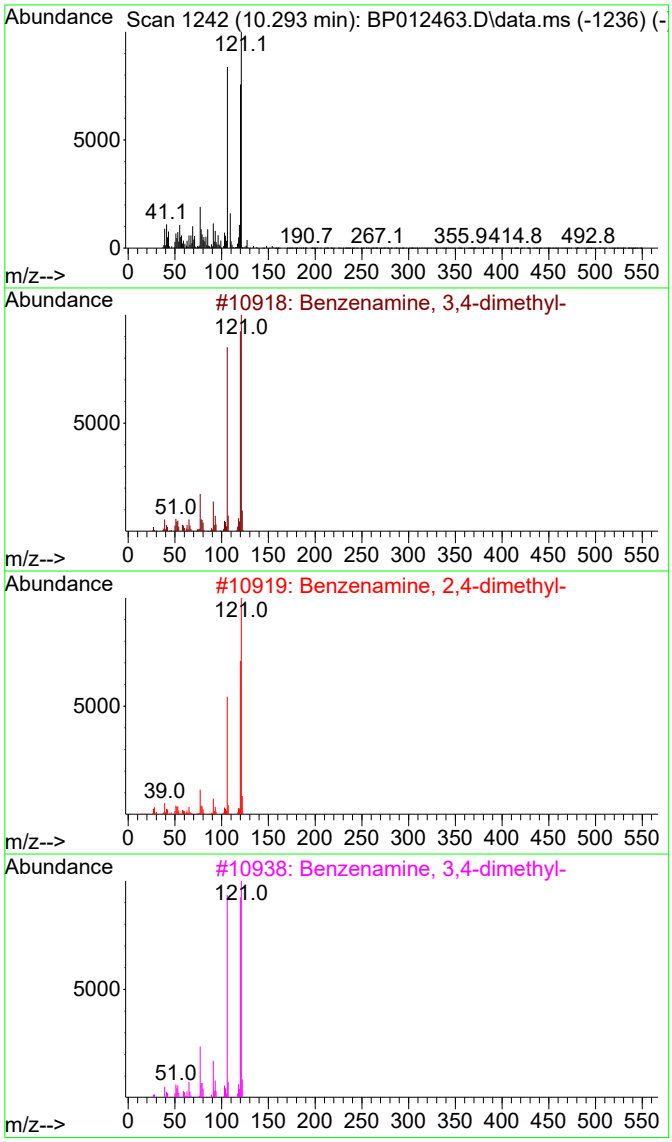
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

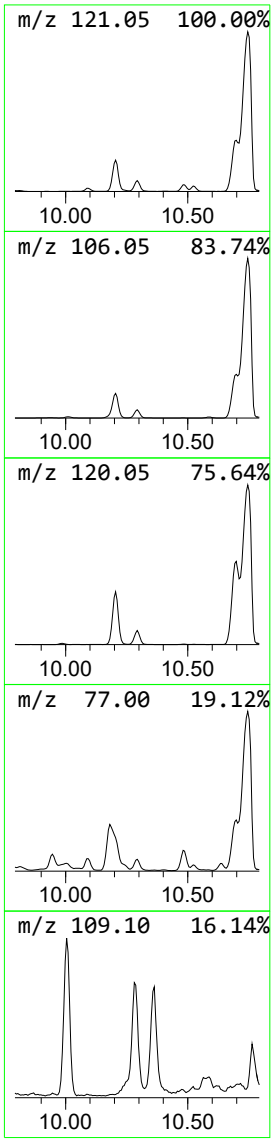
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 29 Benzenamine, 3,4-dimethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.293	21.27 ng/ul	2648770	Naphthalene-d8	10.587

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3,4-dimethyl-	121	C8H11N	000095-64-7	94
2	Benzenamine, 2,4-dimethyl-	121	C8H11N	000095-68-1	94
3	Benzenamine, 3,4-dimethyl-	121	C8H11N	000095-64-7	90
4	Benzenamine, 2,5-dimethyl-	121	C8H11N	000095-78-3	90
5	2,6-Xylidine	121	C8H11N	000087-62-7	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
Data File : BP012463.D
Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

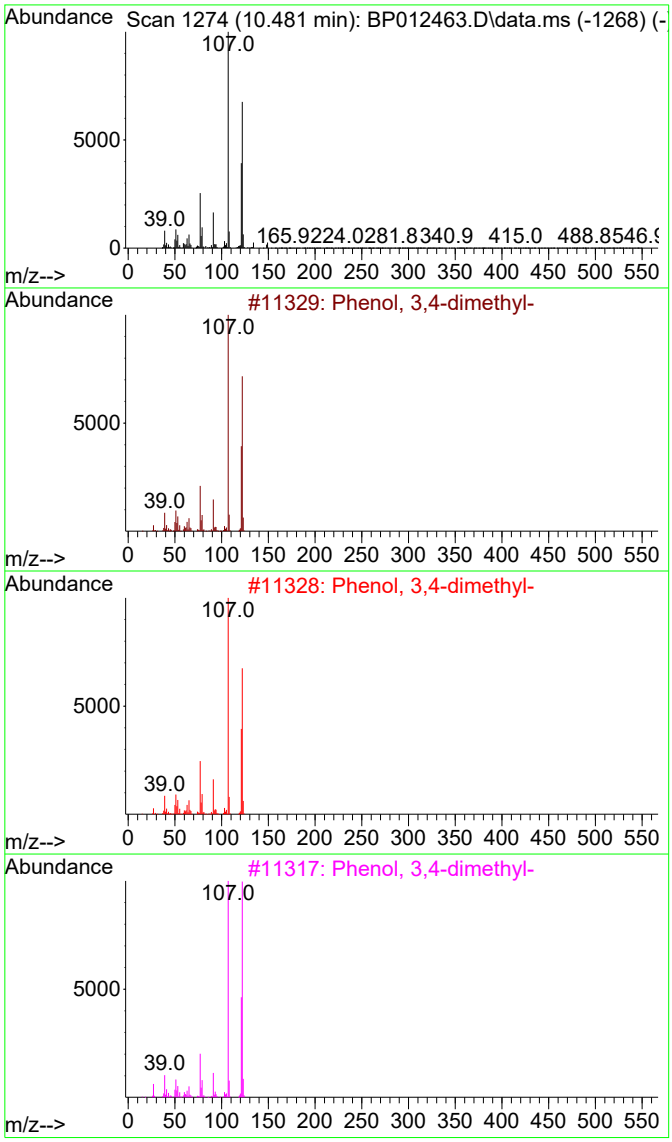
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

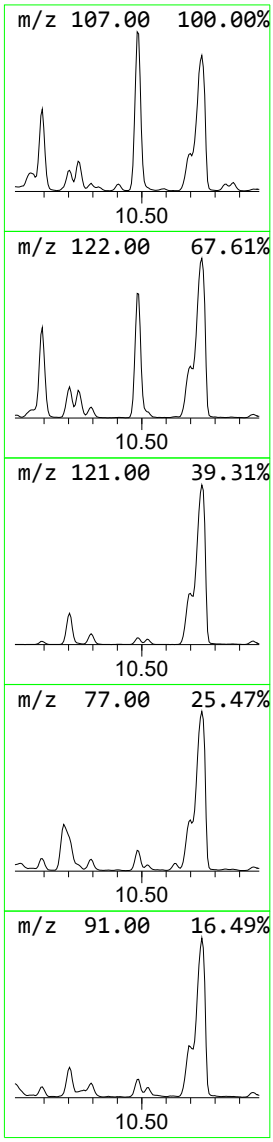
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 31 Phenol, 3,4-dimethyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.481	16.24 ng/ul	2022800	Naphthalene-d8	10.587

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	96
2	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	96
3	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	95
4	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	95
5	Phenol, 2-ethyl-	122	C8H10O	000090-00-6	93





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Data File : BP012463.D
Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

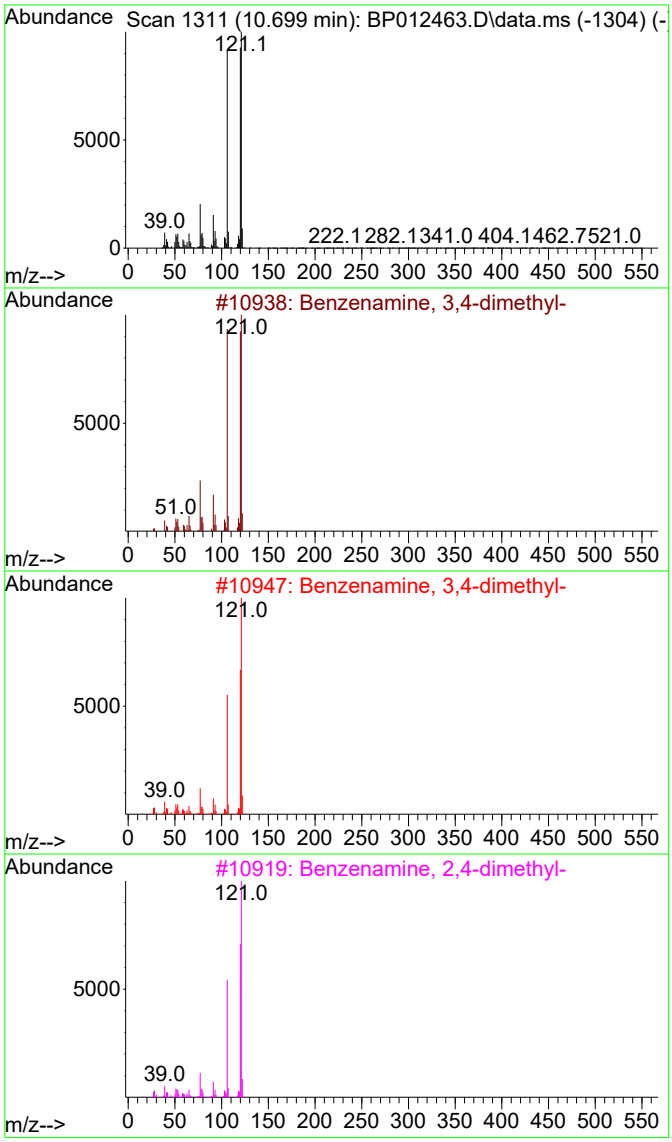
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

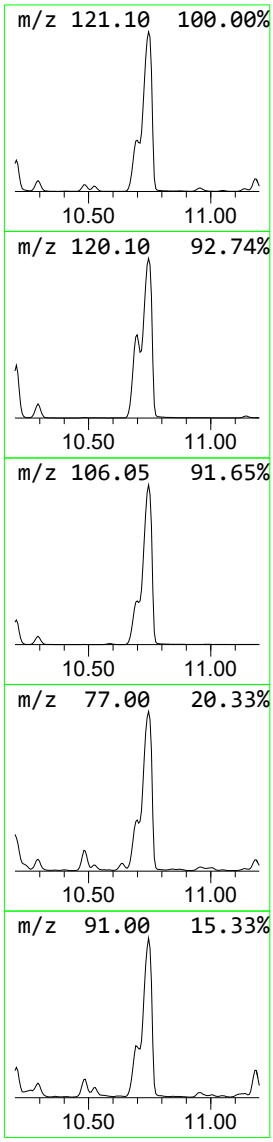
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 32 Benzenamine, 2,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.699	74.96 ng/ul	9334740	Naphthalene-d8	10.587

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3,4-dimethyl-	121	C8H11N	000095-64-7	96
2	Benzenamine, 3,4-dimethyl-	121	C8H11N	000095-64-7	96
3	Benzenamine, 2,4-dimethyl-	121	C8H11N	000095-68-1	96
4	Benzenamine, 3,4-dimethyl-	121	C8H11N	000095-64-7	96
5	Benzenamine, 2,4-dimethyl-	121	C8H11N	000095-68-1	96





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Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

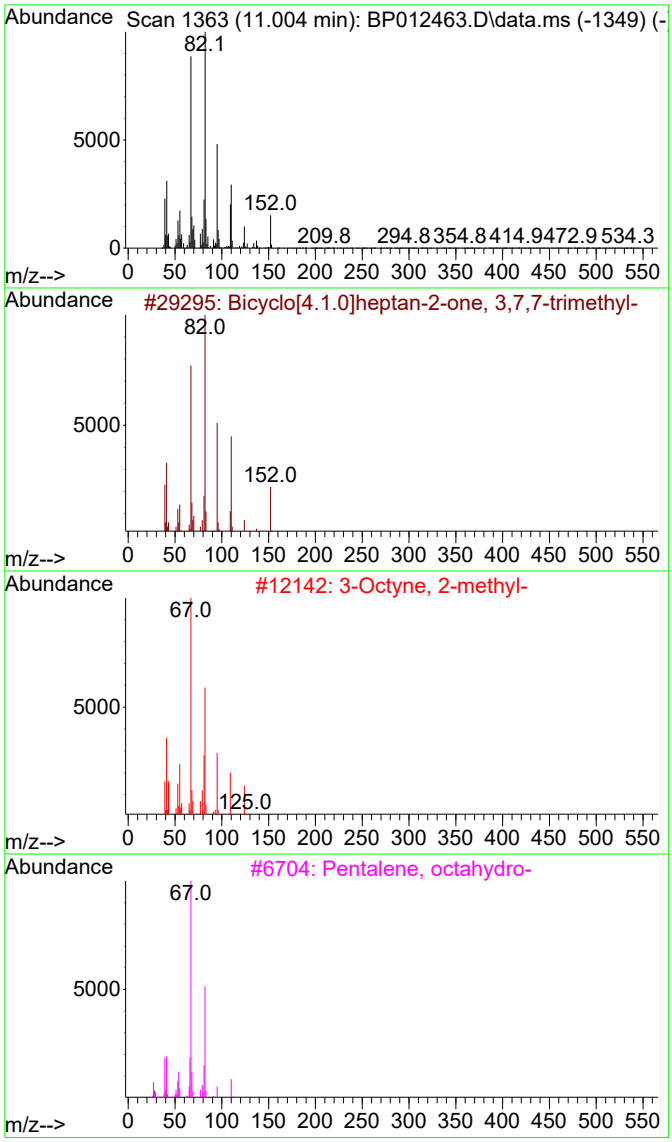
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

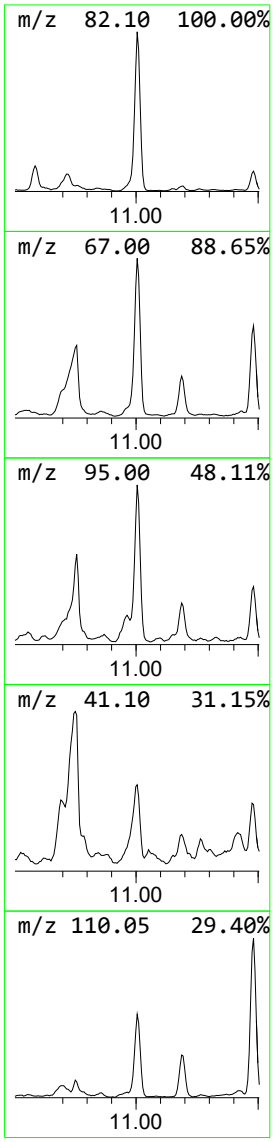
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 33 Bicyclo[4.1.0]heptan-2-one,... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.004	23.33 ng/ul	2904890	Naphthalene-d8	10.587

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[4.1.0]heptan-2-one, 3,7,...	152	C10H16O	000497-62-1	90
2	3-Octyne, 2-methyl-	124	C9H16	055402-15-8	60
3	Pentalene, octahydro-	110	C8H14	000694-72-4	58
4	1-Methylcycloheptene	110	C8H14	055308-20-8	49
5	cis-Hexenyl octyne carbonate	236	C15H24O2	068480-29-5	47





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Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

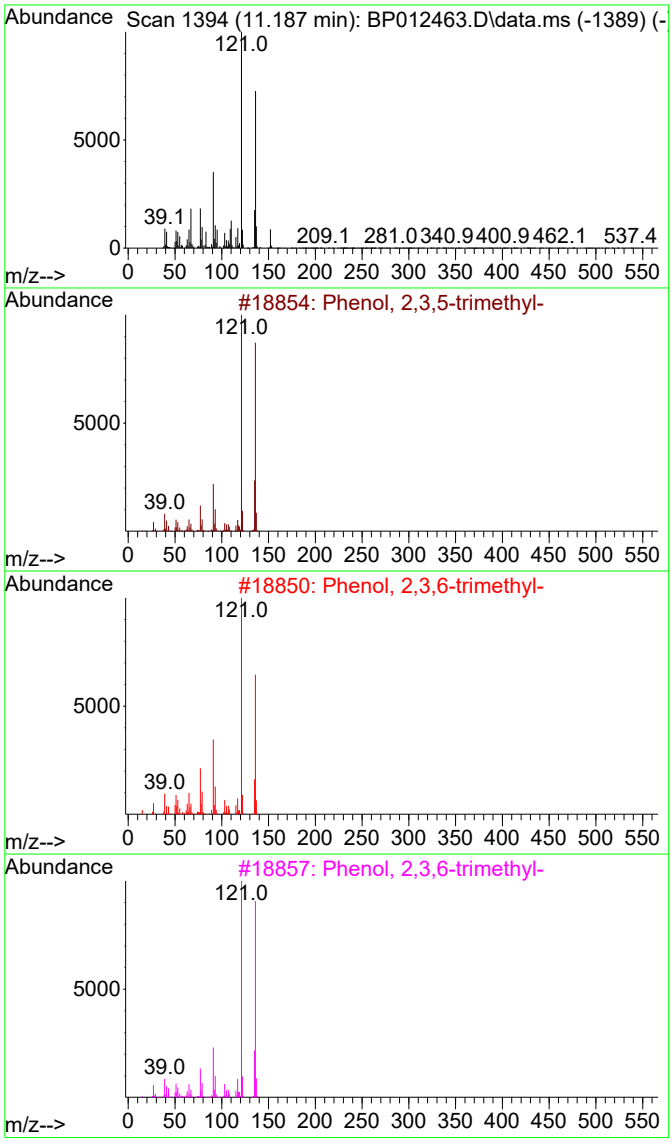
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

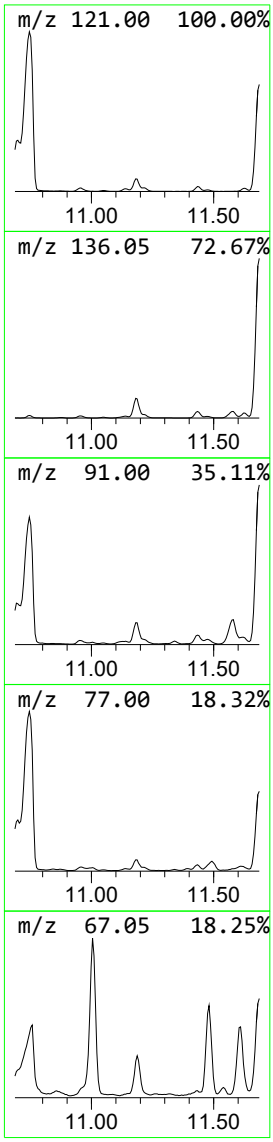
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 34 Phenol, 2,3,5-trimethyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.187	16.45 ng/ul	2049230	Naphthalene-d8	10.587

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,3,5-trimethyl-	136	C9H12O	000697-82-5	95
2	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	94
3	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	93
4	Phenol, 2,4,5-trimethyl-	136	C9H12O	000496-78-6	93
5	Phenol, 2,4,6-trimethyl-	136	C9H12O	000527-60-6	90





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Data File : BP012463.D
Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

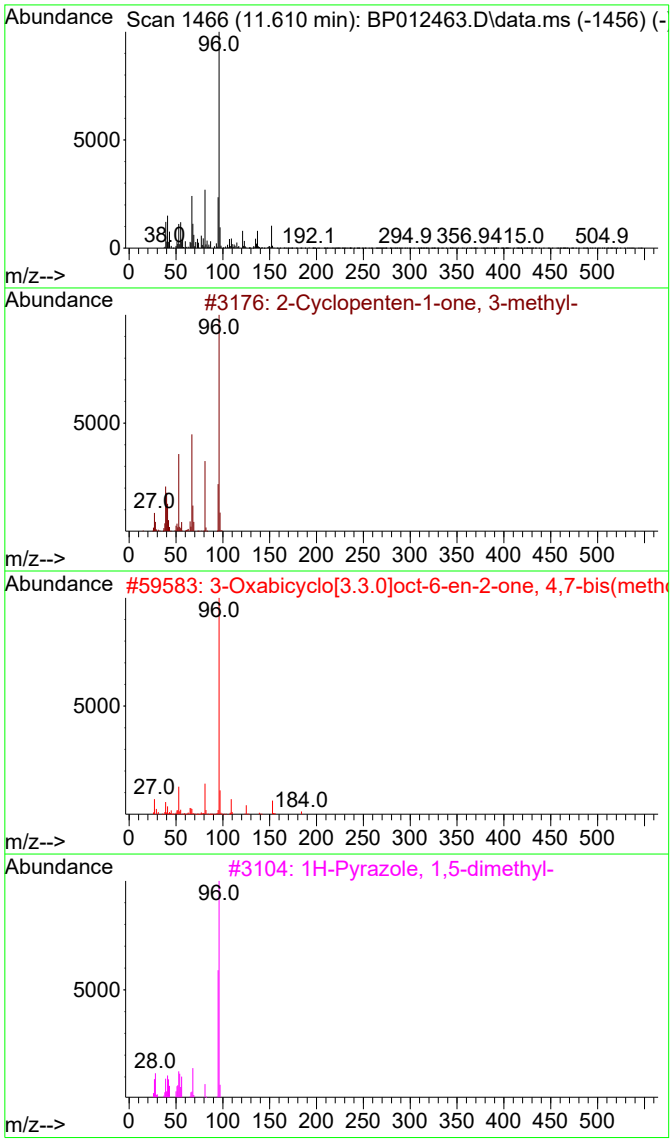
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

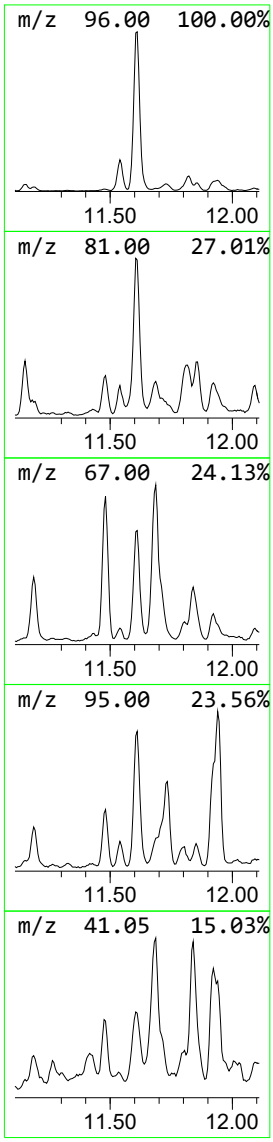
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 37 2-Cyclopenten-1-one, 3-methyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.610	22.30 ng/ul	2776740	Naphthalene-d8	10.587

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Cyclopenten-1-one, 3-methyl-	96	C6H8O	002758-18-1	59
2	3-Oxabicyclo[3.3.0]oct-6-en-2-on...	184	C9H12O4	1000155-61-6	47
3	1H-Pyrazole, 1,5-dimethyl-	96	C5H8N2	000694-31-5	43
4	3,4-dimethylfuran	96	C6H8O	1000458-50-4	40
5	2(1H)-Pyrimidinone	96	C4H4N2O	000557-01-7	40





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Data File : BP012463.D
Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

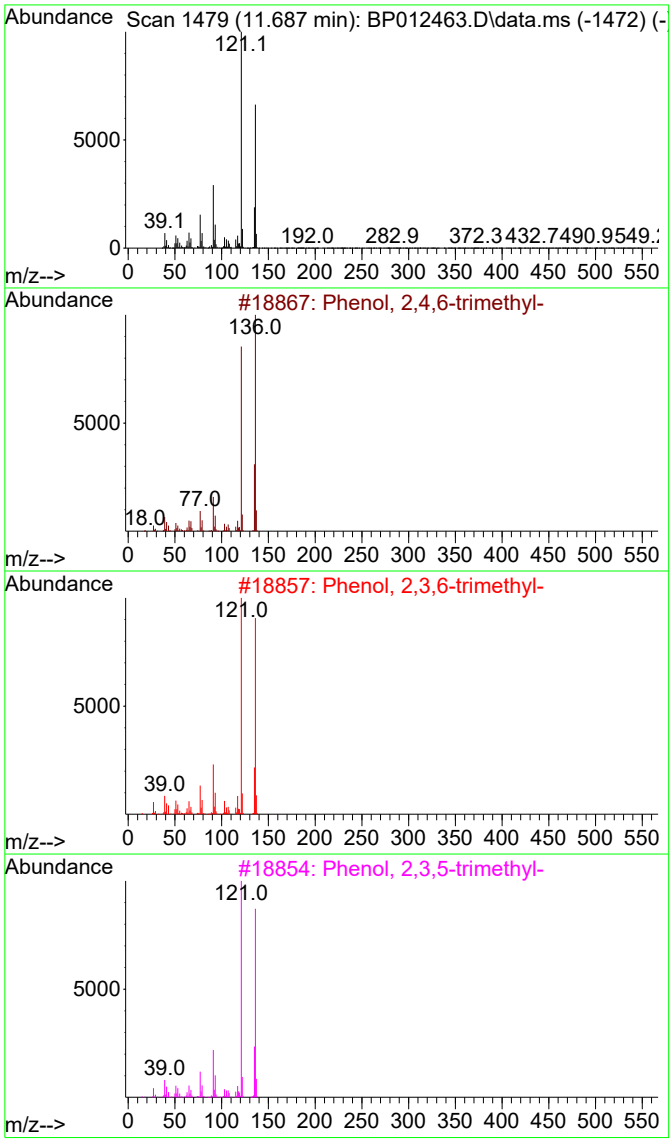
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

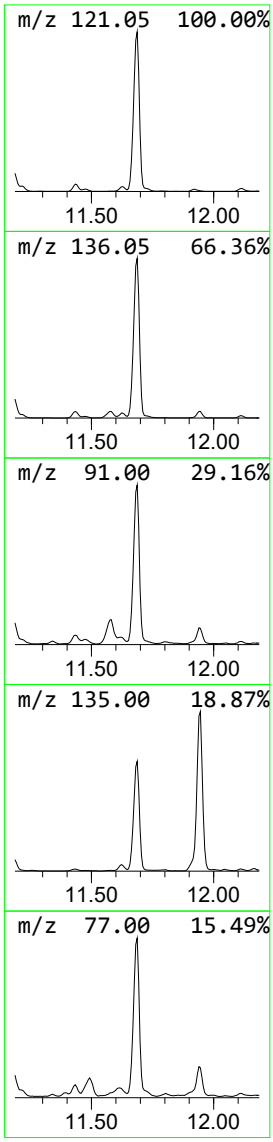
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 38 Phenol, 2,4,6-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.687	104.75 ng/ul	13045700	Naphthalene-d8	10.587

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,4,6-trimethyl-	136	C9H12O	000527-60-6	97
2	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	97
3	Phenol, 2,3,5-trimethyl-	136	C9H12O	000697-82-5	97
4	Phenol, 2,4,6-trimethyl-	136	C9H12O	000527-60-6	96
5	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	95





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Data File : BP012463.D
Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

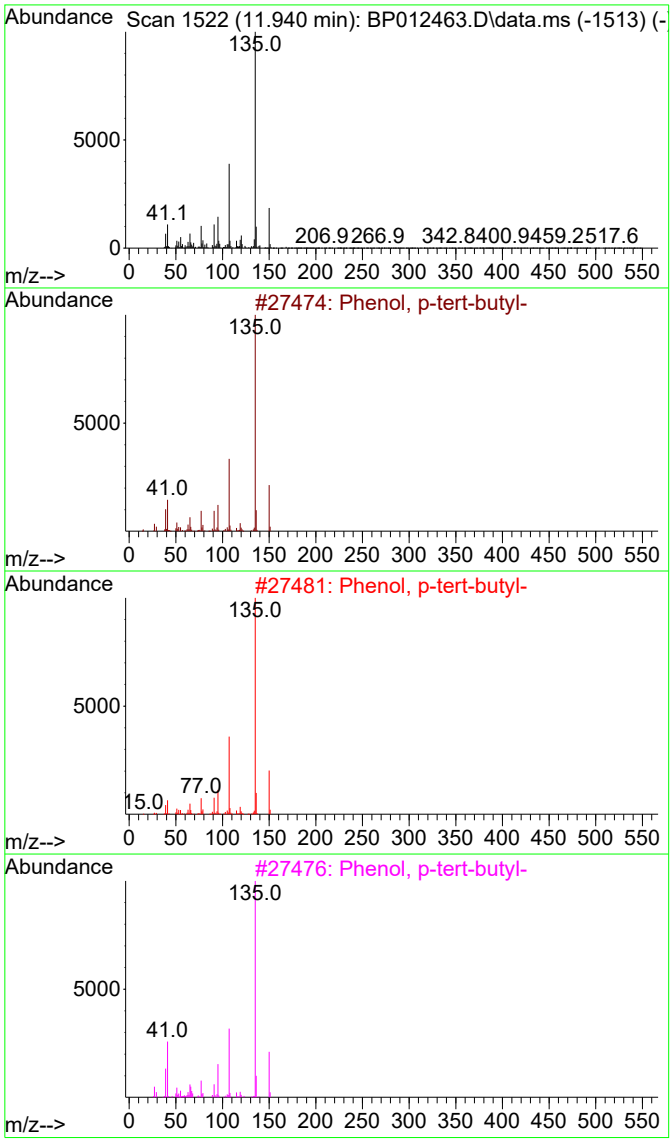
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

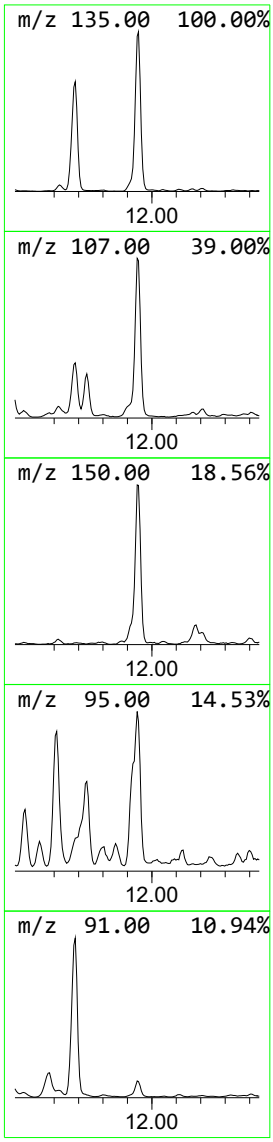
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 41 Phenol, p-tert-butyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.940	34.88 ng/ul	4343420	Naphthalene-d8	10.587

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





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
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Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

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

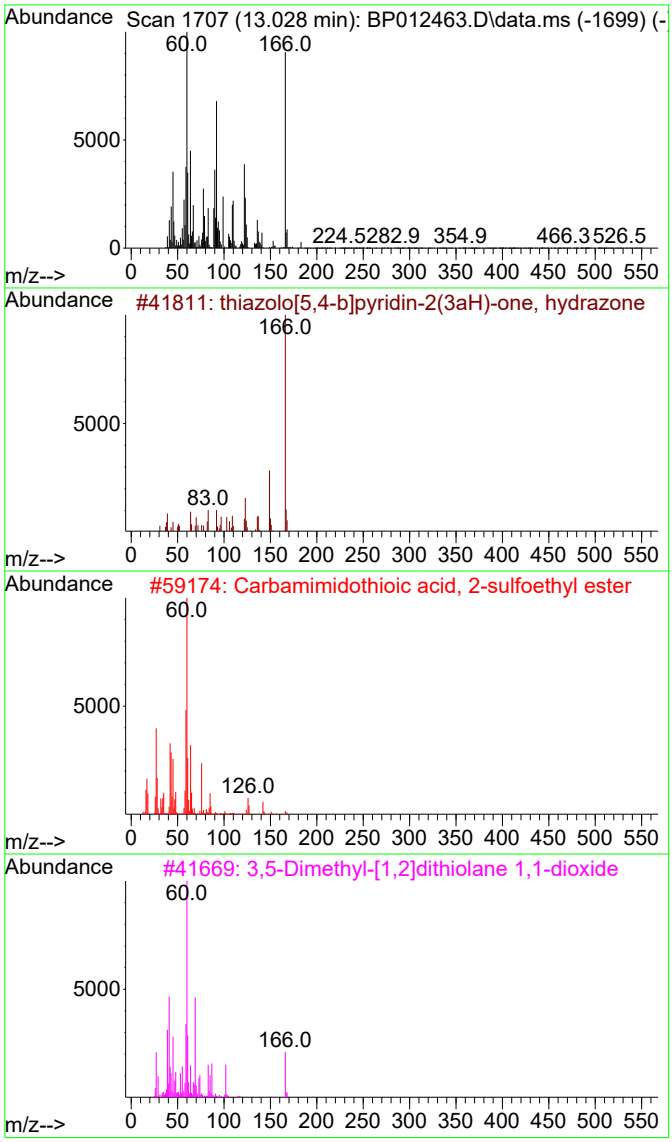
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

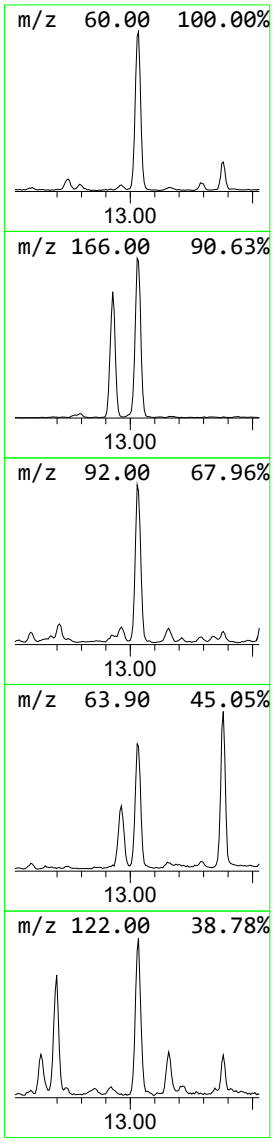
****

Peak Number 45 unknown-03 Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.028	10.32 ng/ul	2690250	Acenaphthene-d10	14.440

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	thiazolo[5,4-b]pyridin-2(3aH)-on...	166	C6H6N4S	1000404-51-7	43
2	Carbamimidothioic acid, 2-sulfoe...	184	C3H8N2O3S2	025985-57-3	17
3	3,5-Dimethyl-[1,2]dithiolane 1,1...	166	C5H10O2S2	1000185-07-5	14
4	Bicyclo[2.2.1]heptane-2,5-dione,...	166	C10H14O2	004230-32-4	10
5	Propanoic acid, 2-chloro-	108	C3H5ClO2	000598-78-7	10





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Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

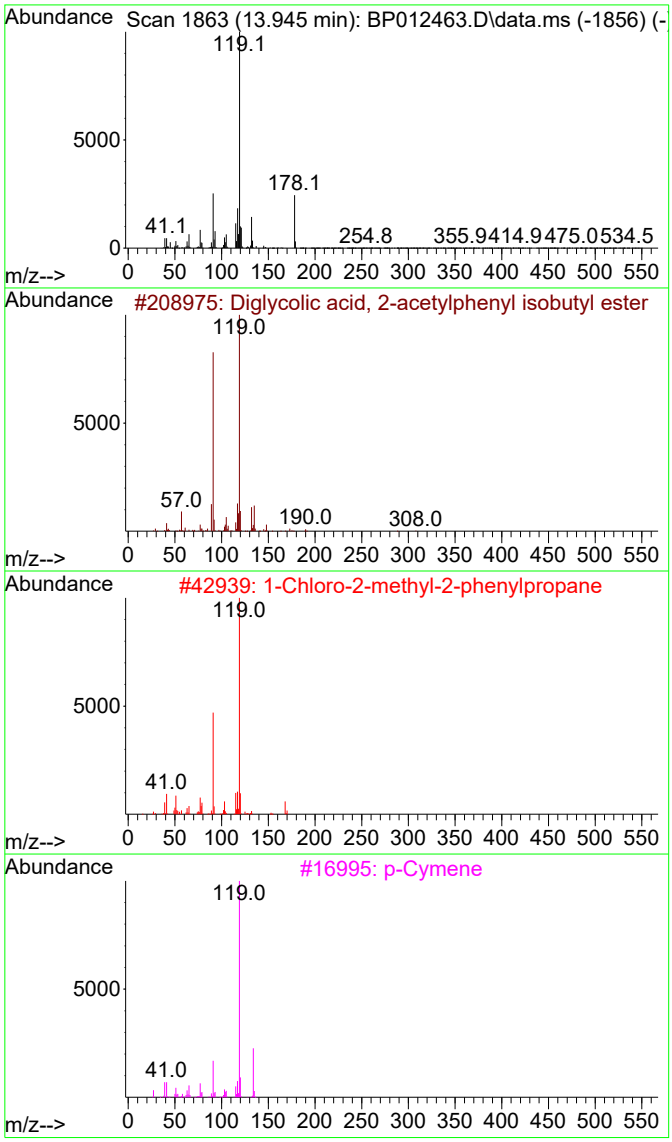
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

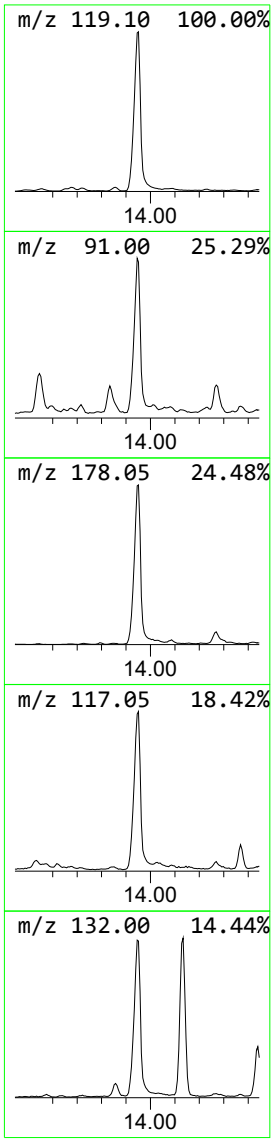
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 49 Diglycolic acid, 2-acetylph... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.945	14.89 ng/ul	3882530	Acenaphthene-d10	14.440

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diglycolic acid, 2-acetylphenyl ...	308	C16H20O6	1000382-71-8	53
2	1-Chloro-2-methyl-2-phenylpropane	168	C10H13Cl	000515-40-2	53
3	p-Cymene	134	C10H14	000099-87-6	52
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	52
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	50





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Data File : BP012463.D
Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

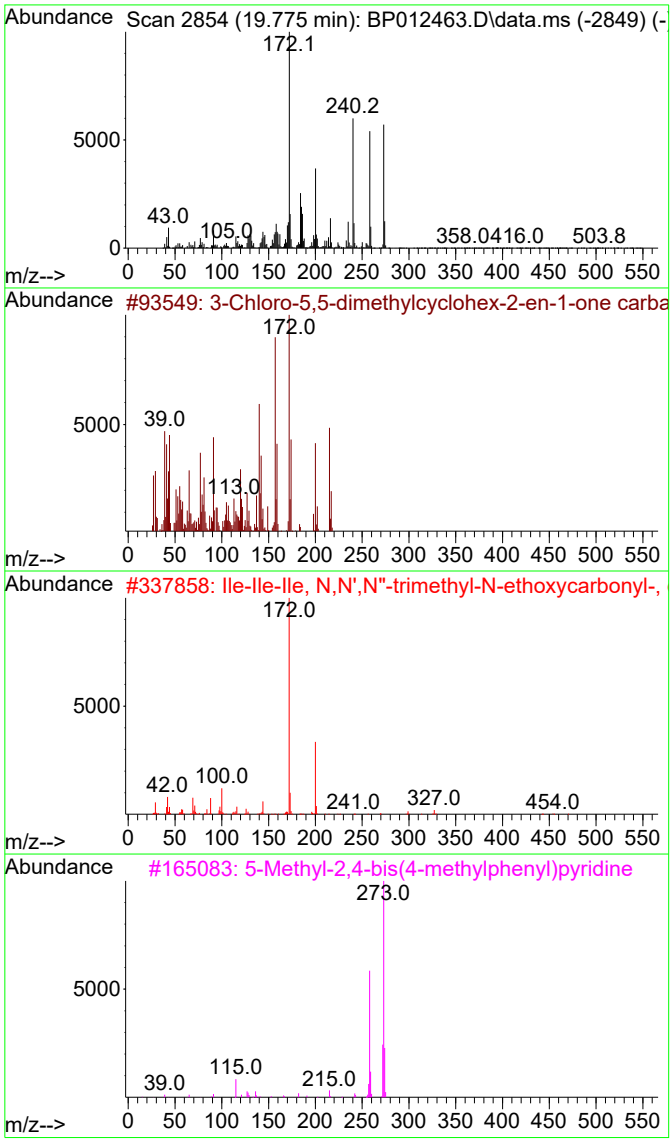
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

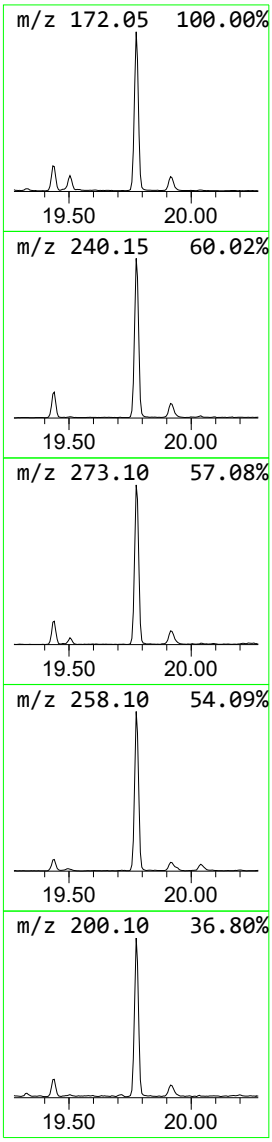
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 65 3-Chloro-5,5-dimethylcycloh... Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.775	11.63 ng/ul	2878480	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-5,5-dimethylcyclohex-2-...	215	C9H14ClN3O	1000185-25-7	93
2	Ile-Ile-Ile, N,N',N"-trimethyl-N...	499	C26H49N3O6	1000454-85-6	22
3	5-Methyl-2,4-bis(4-methylphenyl)...	273	C20H19N	1010445-08-8	20
4	3-Ethyl-4-phenylimidazo[1,5-a]qu...	273	C18H15N3	1360871-21-1	20
5	6-Chloro-N-sec-butyl-1,3,5-triaz...	201	C7H12ClN5	037019-18-4	18





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
Data File : BP012463.D
Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

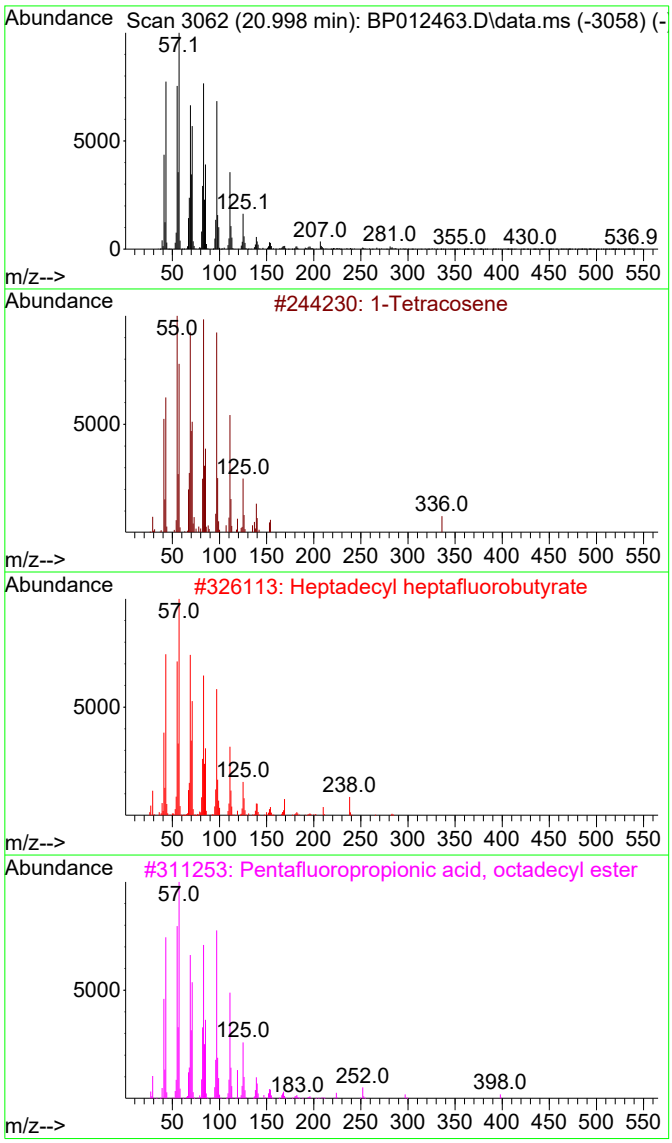
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

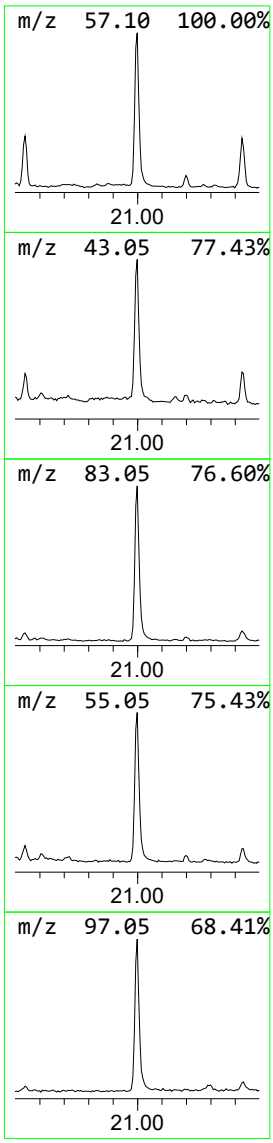
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 67 (DEL) Alkane: Straight-Chai... Concentration Rank 53

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.998	6.85 ng/ul	1693750	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetracosene	336	C24H48	010192-32-2	98
2	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
3	Pentafluoropropionic acid, octad...	416	C21H37F5O2	959261-25-7	94
4	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
5	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
Data File : BP012463.D
Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

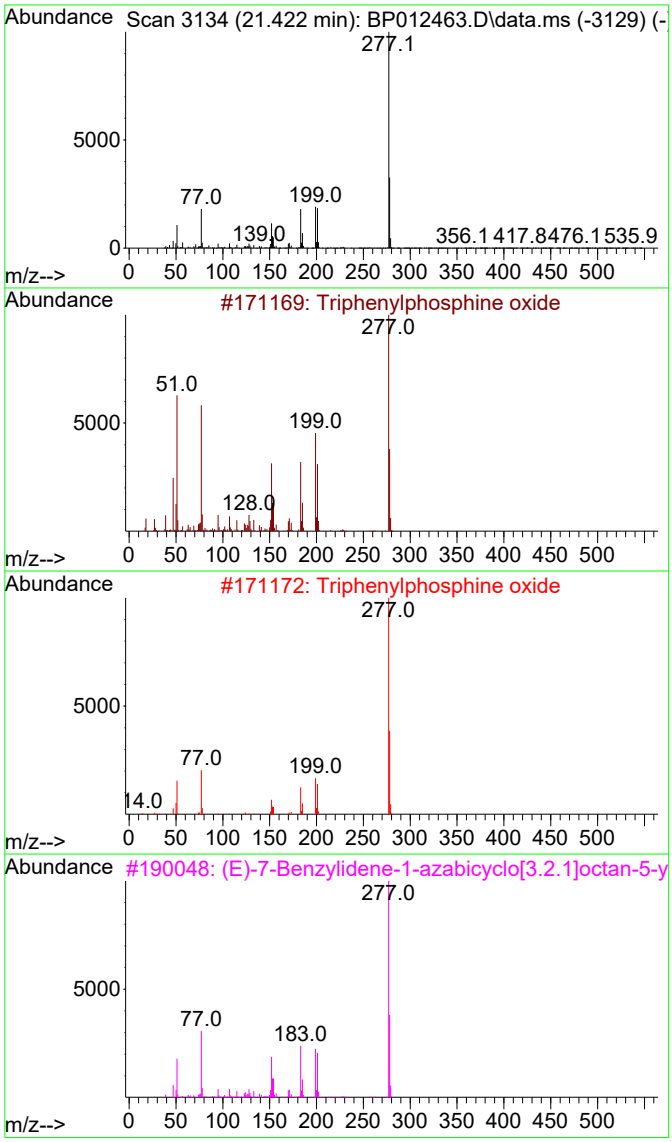
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

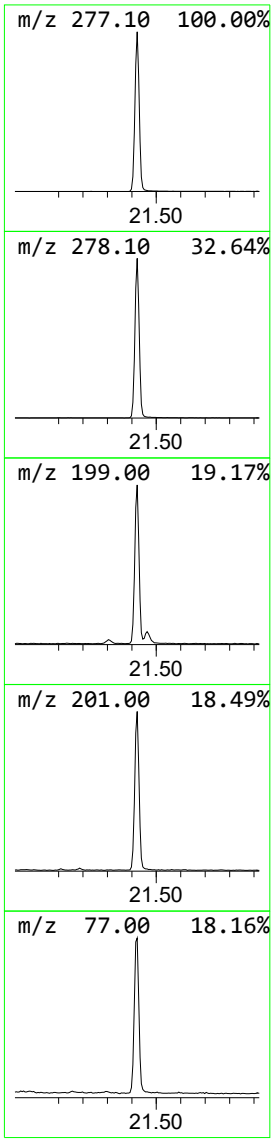
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 68 Triphenylphosphine oxide Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.422	16.20 ng/ul	4008400	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	95
2	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	93
3	(E)-7-Benzylidene-1-azabicyclo[3...	293	C15H19NO3S	1000483-83-4	91
4	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	52
5	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	50





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
Data File : BP012463.D
Acq On : 02 Nov 2022 22:19
Operator : CG/JU
Sample : N5300-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC5

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
Triethylamine	3.028	51.4	ng/ul	5264010	1	7.787	2049340	20.0
Benzene, 1,3-di...	5.799	275.4	ng/ul	28221100	1	7.787	2049340	20.0
3-Heptanone, 5-...	6.452	14.7	ng/ul	1503570	1	7.787	2049340	20.0
Benzene, propyl-	6.758	35.0	ng/ul	3583230	1	7.787	2049340	20.0
Benzene, 1-ethy...	6.869	71.0	ng/ul	7276590	1	7.787	2049340	20.0
Benzene, 1-ethy...	6.928	44.0	ng/ul	4502980	1	7.787	2049340	20.0
Benzene, 1,2,3-...	7.428	161.8	ng/ul	16584000	1	7.787	2049340	20.0
Acetaldehyde, (...	7.481	11.1	ng/ul	1136240	1	7.787	2049340	20.0
unknown-01	7.981	22.8	ng/ul	2339080	1	7.787	2049340	20.0
Indane	8.146	30.4	ng/ul	3119890	1	7.787	2049340	20.0
unknown-02	8.328	14.0	ng/ul	1432240	1	7.787	2049340	20.0
Cyclohexanol, 3...	8.434	344.4	ng/ul	35293100	1	7.787	2049340	20.0
p-Cymene	8.787	14.3	ng/ul	1462120	1	7.787	2049340	20.0
Benzene, 1-ethy...	8.887	17.0	ng/ul	1739350	1	7.787	2049340	20.0
3-Octanol, 3,7-...	9.105	70.3	ng/ul	7201550	1	7.787	2049340	20.0
Benzene, 1-ethy...	9.222	12.3	ng/ul	1529210	2	10.587	2490750	20.0
Benzene, 1,2,3,...	9.469	8.9	ng/ul	1109270	2	10.587	2490750	20.0
Camphor	10.005	35.3	ng/ul	4393930	2	10.587	2490750	20.0
Benzenamine, 3,...	10.293	21.3	ng/ul	2648770	2	10.587	2490750	20.0
Phenol, 3,4-dim...	10.481	16.2	ng/ul	2022800	2	10.587	2490750	20.0
Benzenamine, 2,...	10.699	75.0	ng/ul	9334740	2	10.587	2490750	20.0
Bicyclo[4.1.0]h...	11.004	23.3	ng/ul	2904890	2	10.587	2490750	20.0
Phenol, 2,3,5-t...	11.187	16.4	ng/ul	2049230	2	10.587	2490750	20.0
2-Cyclopenten-1...	11.610	22.3	ng/ul	2776740	2	10.587	2490750	20.0
Phenol, 2,4,6-t...	11.687	104.8	ng/ul	13045700	2	10.587	2490750	20.0
Phenol, p-tert-...	11.940	34.9	ng/ul	4343420	2	10.587	2490750	20.0
unknown-03	13.028	10.3	ng/ul	2690250	3	14.440	5213960	20.0
Diglycolic acid...	13.945	14.9	ng/ul	3882530	3	14.440	5213960	20.0
3-Chloro-5,5-di...	19.775	11.6	ng/ul	2878480	5	21.298	4948030	20.0
(DEL) Alkane: S...	20.998	6.8	ng/ul	1693750	5	21.298	4948030	20.0
Triphenylphosph...	21.422	16.2	ng/ul	4008400	5	21.298	4948030	20.0