LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
Data File : BP012484.D
Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BHA73DL2

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: BP012484.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.422	234	244	252	rVB	76654	142101	2.30%	0.209%
2	4.940	329	332	346	rVB2	44145	103319	1.67%	0.152%
3	5.093	352	358	367	rBV	279553	444713	7.18%	0.655%
4	5.222	375	380	385	rBV	95880	139361	2.25%	0.205%
5	5.287	385	391	400	rVV	471525	717537	11.59%	1.058%
6	5.416	408	413	421	rBV	120660	203118	3.28%	0.299%
7	5.499	421	427	437	rVV	551856	799168	12.91%	1.178%
8	5.705	455	462	464	rBV	69149	118550	1.91%	0.175%
9	5.728	464	466	472	rVV2	68259	83117	1.34%	0.123%
10	5.793	472	477	488	rVB2	52214	87205	1.41%	0.129%
11	6.058	511	522	532	rBV3	53226	149156	2.41%	0.220%
12	6.152	532	538	544	rVB	44302	77172	1.25%	0.114%
13	6.263	547	557	562	rBV2	230109	455560	7.36%	0.671%
14	6.452	580	589	598	rBV3	137880	282725	4.57%	0.417%
15	6.758	635	641	651	rBV	119493	203484	3.29%	0.300%
16	6.993	669	681	687	rBV5	48063	125242	2.02%	0.185%
17	7.128	696	704	708	rBV	170081	305859	4.94%	0.451%
18	7.322	731	737	748	rBV2	97991	211304	3.41%	0.311%
19	7.428	748	755	761	rBV	940429	1458084	23.55%	2.149%
20	7.787	805	816	827	rBV	874427	1578525	25.49%	2.327%
21	7.893	827	834	845	rVB	354666	610627	9.86%	0.900%
22	7.981	845	849	855	rVB2	55771	80631	1.30%	0.119%
23	8.152	870	878	885	rBV	1336651	2257862	36.47%	3.328%
24	8.222	885	890	895	rVB	447244	678375	10.96%	1.000%
25	8.416	916	923	935	rBV	419545	781709	12.63%	1.152%
26	8.516	935	940	946	rVB	72057	127371	2.06%	0.188%
27	8.693	965	970	974	rBV2	101021	166282	2.69%	0.245%
28	8.787	981	986	995	rVB3	85741	146309	2.36%	0.216%
29	8.887	997	1003	1010	rBV3	147571	270127	4.36%	0.398%
30	8.963	1010	1016	1021	rBV3	152677	301049	4.86%	0.444%
31	9.016	1021	1025	1041	rVB4	218814	479086	7.74%	0.706%
32	9.222	1054	1060	1066	rBV2	46490	93733	1.51%	0.138%
33	9.410	1087	1092	1097	rBV	77609	121081	1.96%	0.178%
34	9.469	1097	1102	1108	rBV	86028	142495	2.30%	0.210%
35	9.575	1115	1120	1125	rBV	261886	377981	6.10%	0.557%
36	9.681	1133	1138	1142	rBV	88841	175052	2.83%	0.258%

LSC Area Percent Report

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Data File : BP012484.D
Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

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	9.828	1157	1163	1174	rVB	197714	394653	6.37%	0.582%
38	9.957	1179	1185	1186	rBV	160603	239107	3.86%	0.352%
39	9.987	1186	1190	1201	rVB4	261315	773261	12.49%	1.140%
40	10.087	1201	1207	1211	rBV2	148082	316004	5.10%	0.466%
41	10.128	1211	1214	1221	rVB4	146914	235728	3.81%	0.347%
42	10.216	1221	1229	1237	rBV2	215389	479630	7.75%	0.707%
43	10.322	1237	1247	1251	rBV6	116324	349282	5.64%	0.515%
44	10.404	1257	1261	1269	rVB2	102530	182267	2.94%	0.269%
45	10.587	1285	1292	1296	rVV	1317547	2300740	37.16%	3.391%
46	10.640	1296	1301	1311	rVV	3498057	6191720	100.00%	9.126%
47	10.728	1311	1316	1333	rVB4	514024	1546127	24.97%	2.279%
48	11.004	1358	1363	1369	rVB	162900	262855	4.25%	0.387%
49	11.110	1375	1381	1390	rBV	169049	345845	5.59%	0.510%
50	11.787	1490	1496	1503	rBV8	56244	104409	1.69%	0.154%
51	11.945	1512	1523	1528	rBV2	179475	433977	7.01%	0.640%
52	12.010	1528	1534	1542	rVB8	51901	132445	2.14%	0.195%
53	12.116	1545	1552	1555	rBV8	35267	78240	1.26%	0.115%
54	12.257	1569	1576	1589	rVB	673147	1124444	18.16%	1.657%
55	12.475	1607	1613	1623	rVB	880815	1432873	23.14%	2.112%
56	12.587	1626	1632	1639	rVV2	149892	284842	4.60%	0.420%
57	12.751	1655	1660	1669	rVB6	48098	85714	1.38%	0.126%
58	13.075	1710	1715	1727	rVB	312091	469267	7.58%	0.692%
59	13.575	1793	1800	1805	rBV2	117441	229761	3.71%	0.339%
60	13.722	1821	1825	1829	rVV3	82716	146482	2.37%	0.216%
61	13.757	1829	1831	1836	rVV3	54689	82849	1.34%	0.122%
62	13.857	1843	1848	1856	rVV2	422111	897170	14.49%	1.322%
63	13.928	1856	1860	1867	rVV	149376	280908	4.54%	0.414%
64	13.992	1868	1871	1876	rVV4	52257	93014	1.50%	0.137%
65	14.057	1877	1882	1885	rVV	137991	236947	3.83%	0.349%
66	14.092	1885	1888	1890	rVV2	113482	159579	2.58%	0.235%
67	14.134	1890	1895	1903	rVV	483420	801972	12.95%	1.182%
68	14.234	1905	1912	1924	rVB3	97795	234878	3.79%	0.346%
69	14.439	1941	1947	1955	rVV	2300991	3708951	59.90%	5.467%
70	14.504	1955	1958	1964	rVV	293076	426995	6.90%	0.629%
71	14.839	2010	2015	2028	rVB6	75419	207834	3.36%	0.306%
72	15.192	2071	2075	2080	rBV	187782	269431	4.35%	0.397%
73	15.445	2112	2118	2124	rBV2	1673650	2509590	40.53%	3.699%
74	15.581	2137	2141	2149	rVB2	107148	190443	3.08%	0.281%
75	15.739	2164	2168	2178	rVB3	200097	367381	5.93%	0.542%
76	15.963	2200	2206	2220	rVB2	332061	613416	9.91%	0.904%

LSC Area Percent Report

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Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

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	16.151	2233	2238	2247	rBV2	292989	589131	9.51%	0.868%
78	16.339	2266	2270	2274	rBV3	61577	79262	1.28%	0.117%
79	16.486	2291	2295	2296	rBV	105262	131630	2.13%	0.194%
80	16.610	2309	2316	2321	rVB7	58011	139514	2.25%	0.206%
81	16.957	2371	2375	2378	rVB	83139	111097	1.79%	0.164%
82	17.204	2411	2417	2428	rBV	2999676	4400757	71.07%	6.487%
83	17.304	2429	2434	2442	rVB	666801	933823	15.08%	1.376%
84	17.475	2460	2463	2468	rVB3	82015	110504	1.78%	0.163%
85	17.733	2502	2507	2524	rBV	1116551	1808579	29.21%	2.666%
86	17.880	2528	2532	2541	rBV5	121458	293788	4.74%	0.433%
87	18.075	2561	2565	2576	rVB2	233038	399906	6.46%	0.589%
88	18.245	2590	2594	2598	rBV	128084	176370	2.85%	0.260%
89	18.345	2607	2611	2614	rBV	73791	113256	1.83%	0.167%
90	18.392	2614	2619	2628	rVB	244759	391413	6.32%	0.577%
91	19.269	2764	2768	2772	rBV	208823	249372	4.03%	0.368%
92	19.545	2810	2815	2820	rBV	909278	1158486	18.71%	1.708%
93	19.598	2820	2824	2834	rVV	668482	946555	15.29%	1.395%
94	20.039	2895	2899	2906	rVB	1037711	1269177	20.50%	1.871%
95	20.398	2957	2960	2964	rVB4	109198	123990	2.00%	0.183%
96	20.539	2982	2984	2989	rVB4	79943	82538	1.33%	0.122%
97	21.298	3109	3113	3122	rBV	3681439	4723339	76.28%	6.962%
98	21.421	3131	3134	3139	rBV	433999	527266	8.52%	0.777%
99	23.533	3489	3493	3506	rVB2	475241	1060054	17.12%	1.562%
100	23.686	3513	3519	3535	rVB2	2075218	4307726	69.57%	6.349%

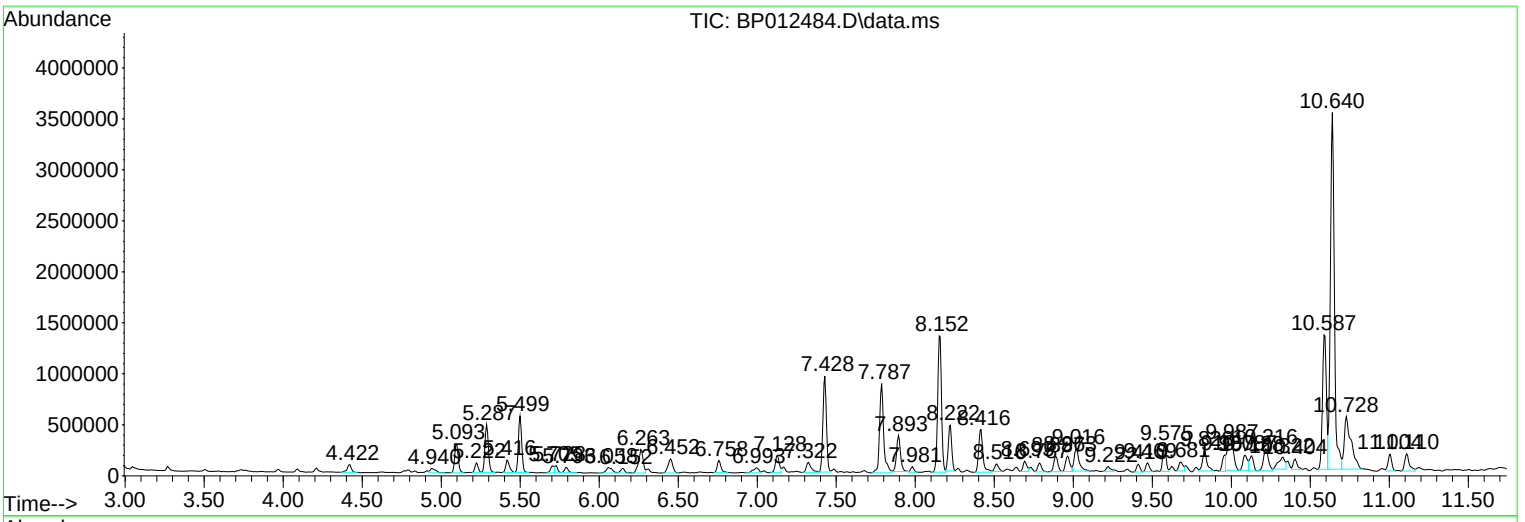
Sum of corrected areas: 67843634

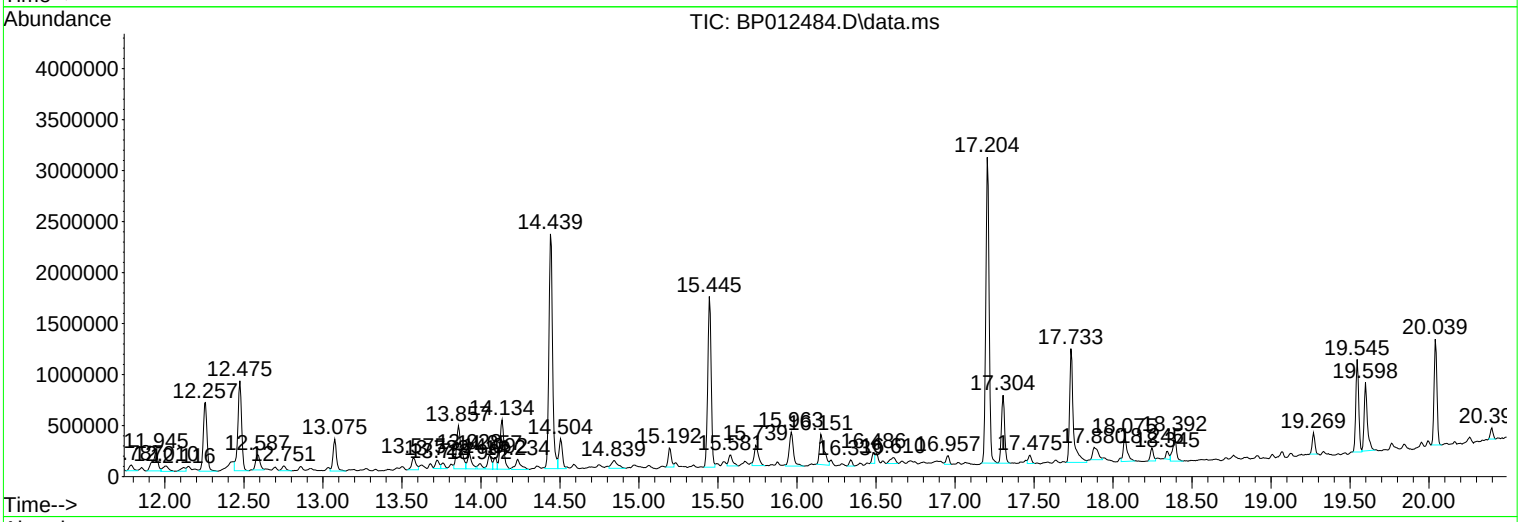
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
Data File : BP012484.D
Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

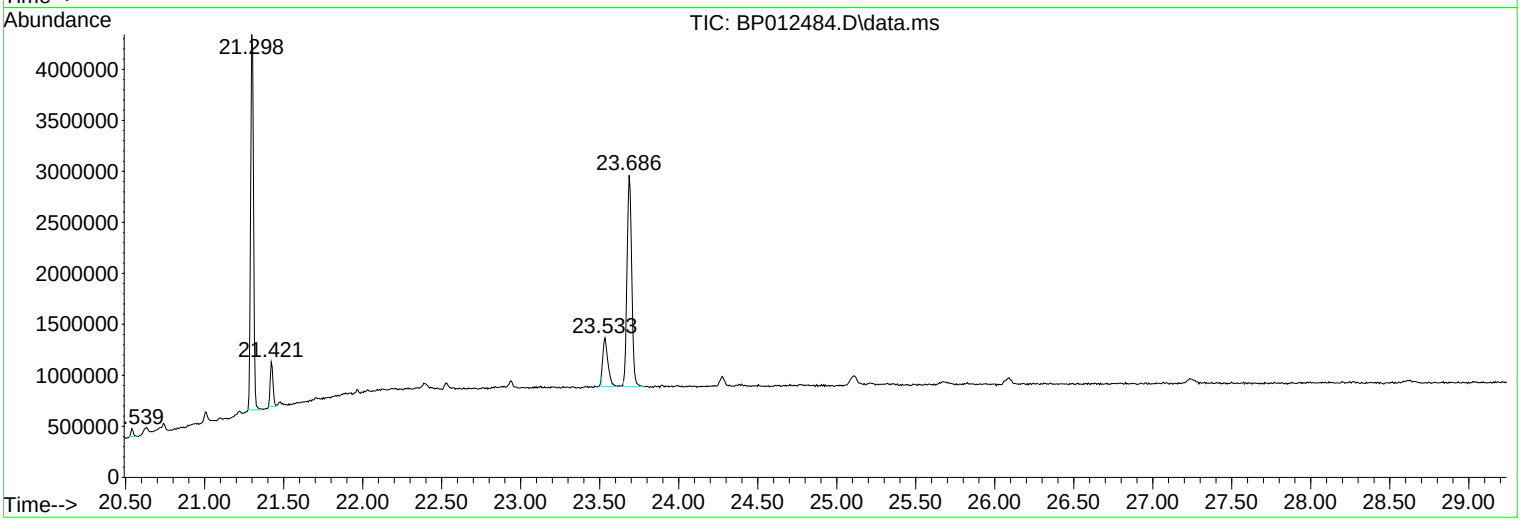
Instrument :
BNA_P
ClientSampleId :
BHA73DL2

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 : BP012484.D
Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

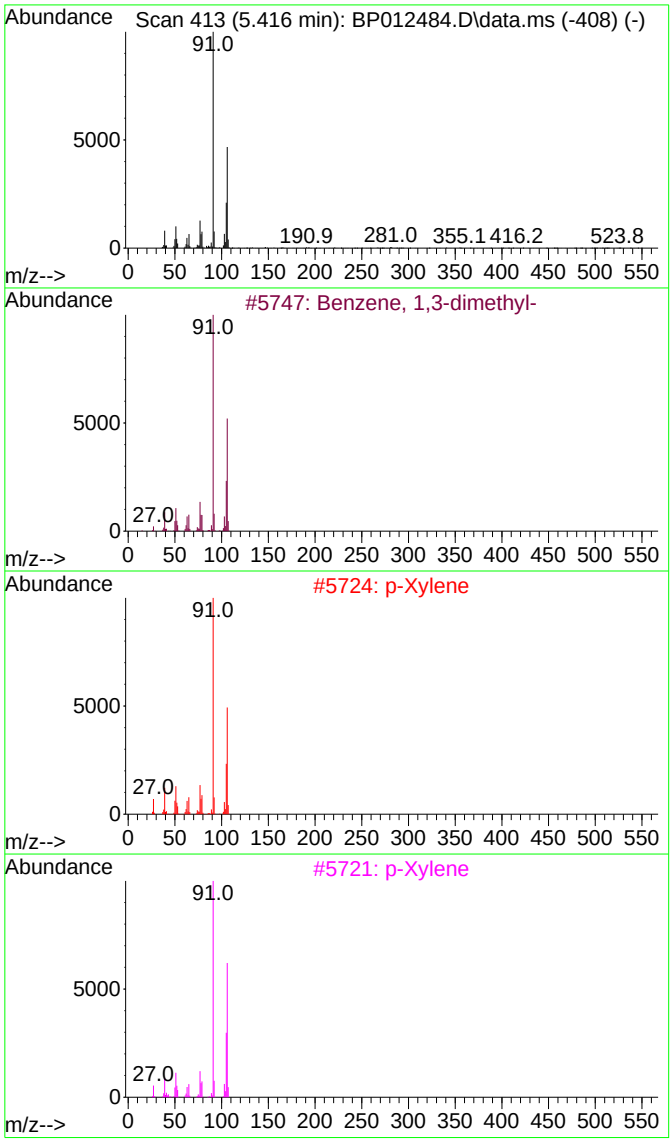
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

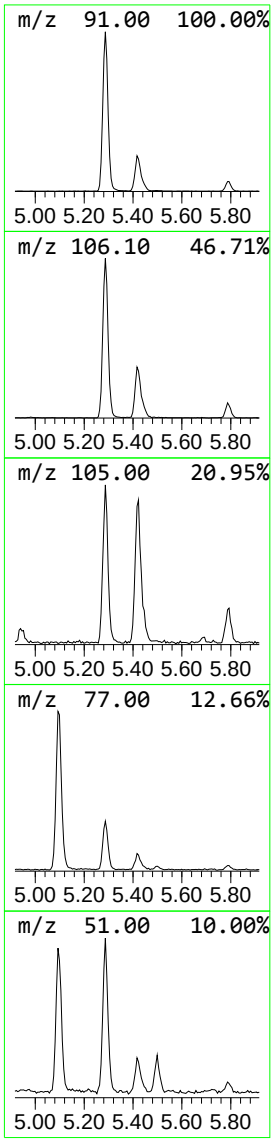
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1,3-dimethyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.416	2.57 ng/ul	203118	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	p-Xylene	106	C8H10	000106-42-3	97
3	p-Xylene	106	C8H10	000106-42-3	97
4	p-Xylene	106	C8H10	000106-42-3	97
5	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
Data File : BP012484.D
Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

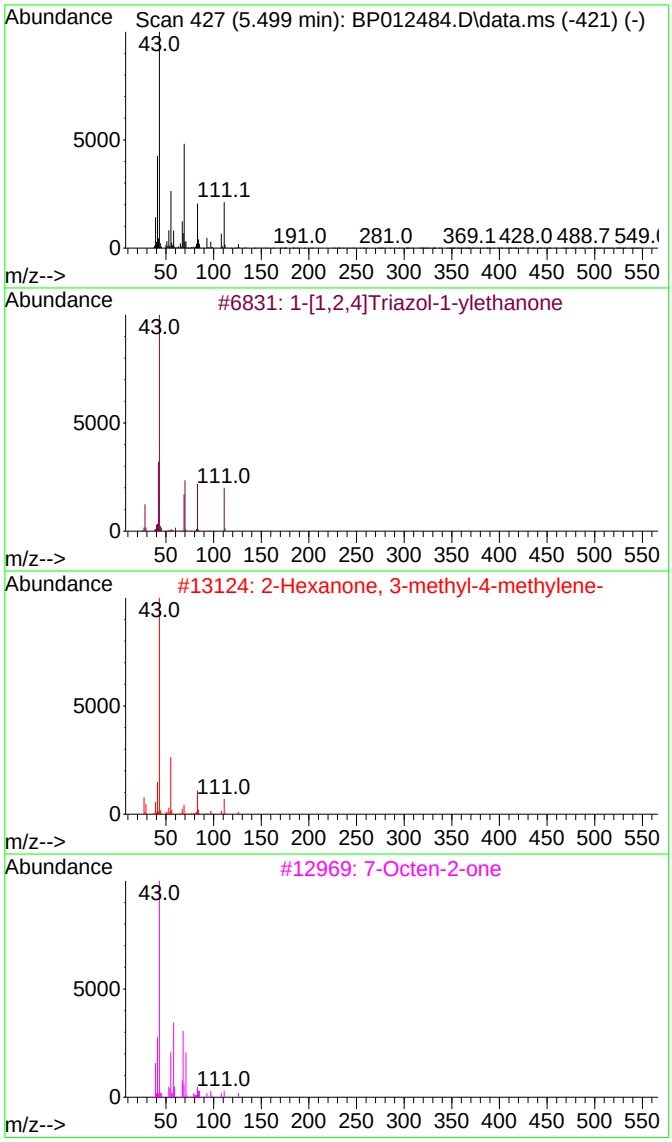
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

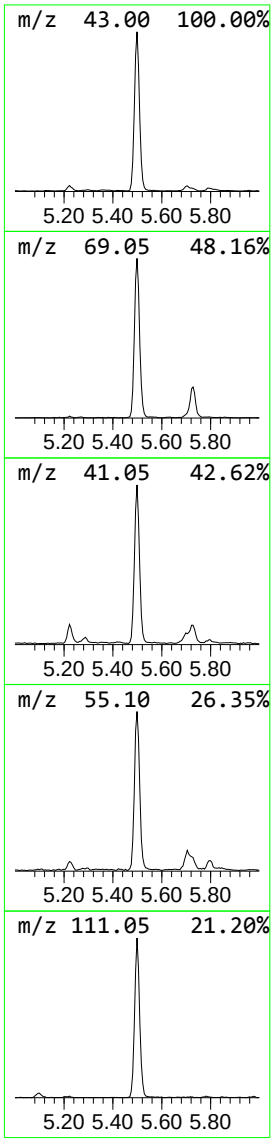
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 1-[1,2,4]Triazol-1-ylethanone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.499	10.13 ng/ul	799168	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-[1,2,4]Triazol-1-ylethanone	111	C4H5N3O	015625-88-4	64
2	2-Hexanone, 3-methyl-4-methylene-	126	C8H14O	020690-71-5	43
3	7-Octen-2-one	126	C8H14O	003664-60-6	37
4	3-Hexen-2-one, 3,4-dimethyl-, (E)-	126	C8H14O	020685-46-5	25
5	1-(1,3-Oxazol-2-yl)ethan-1-one	111	C5H5NO2	077311-07-0	12





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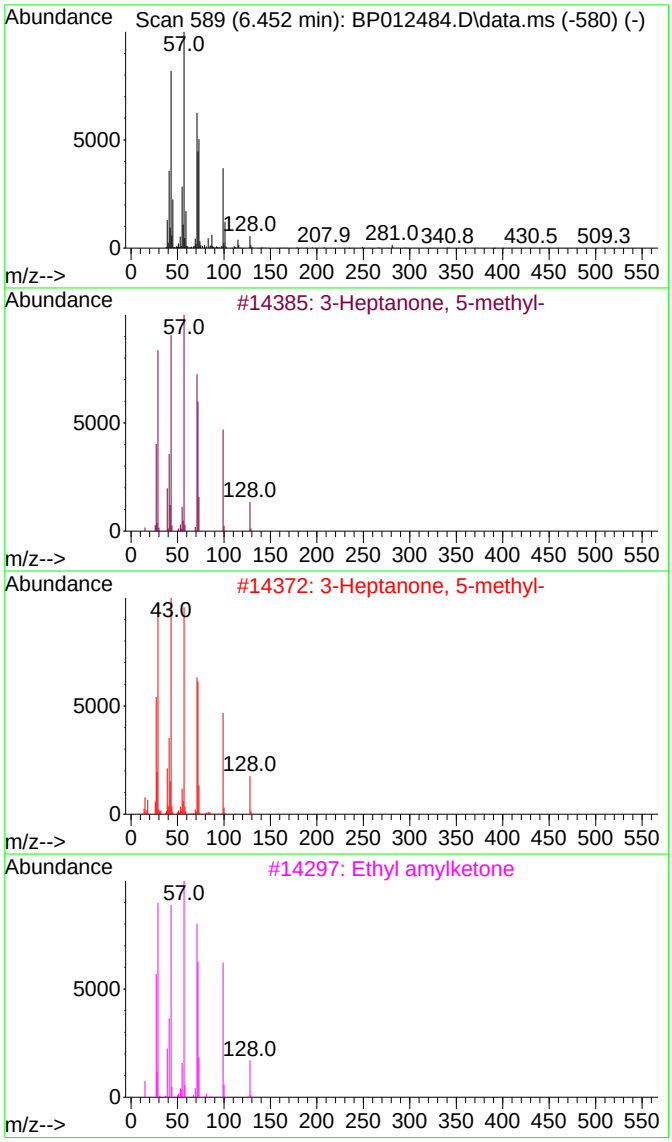
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

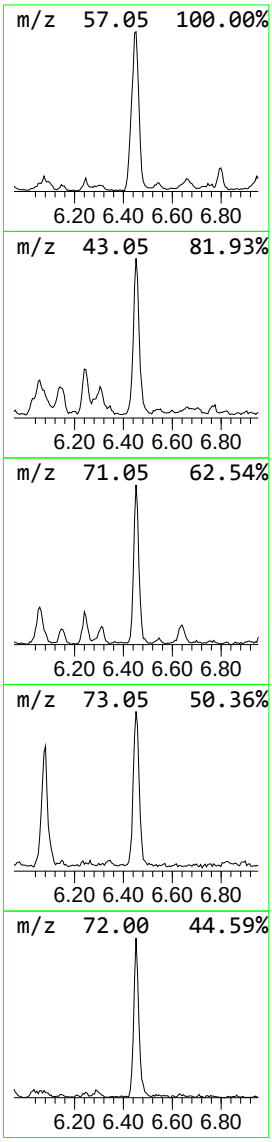
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 3-Heptanone, 5-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.452	3.58 ng/ul	282725	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	53
2	3-Heptanone, 5-methyl-	128	C8H16O	000541-85-5	50
3	Ethyl amylketone	128	C8H16O	020616-93-7	50
4	3-Octanone	128	C8H16O	000106-68-3	43
5	Hexaborane	76	B6H10	023777-80-2	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
Data File : BP012484.D
Acq On : 03 Nov 2022 11:25
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ALS Vial : 39 Sample Multiplier: 1

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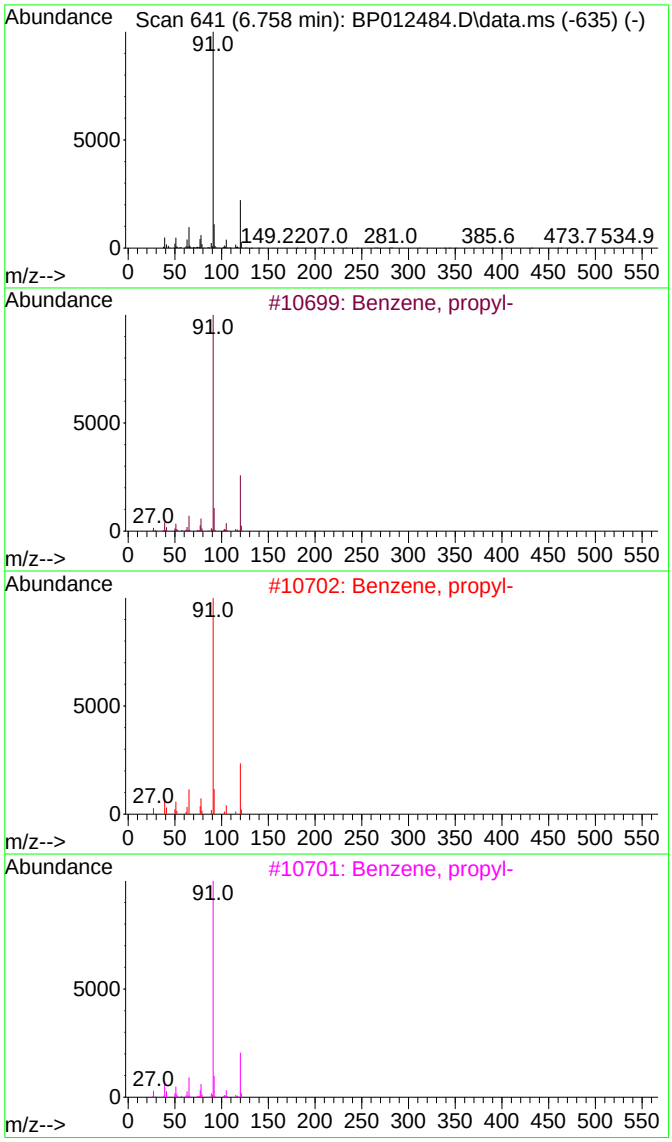
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

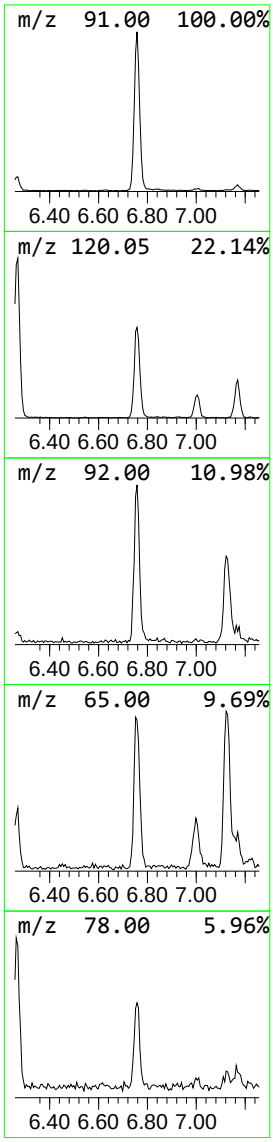
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, propyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.758	2.58 ng/ul	203484	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	91
2	Benzene, propyl-	120	C9H12	000103-65-1	91
3	Benzene, propyl-	120	C9H12	000103-65-1	91
4	Benzene, propyl-	120	C9H12	000103-65-1	87
5	N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	78





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
Data File : BP012484.D
Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

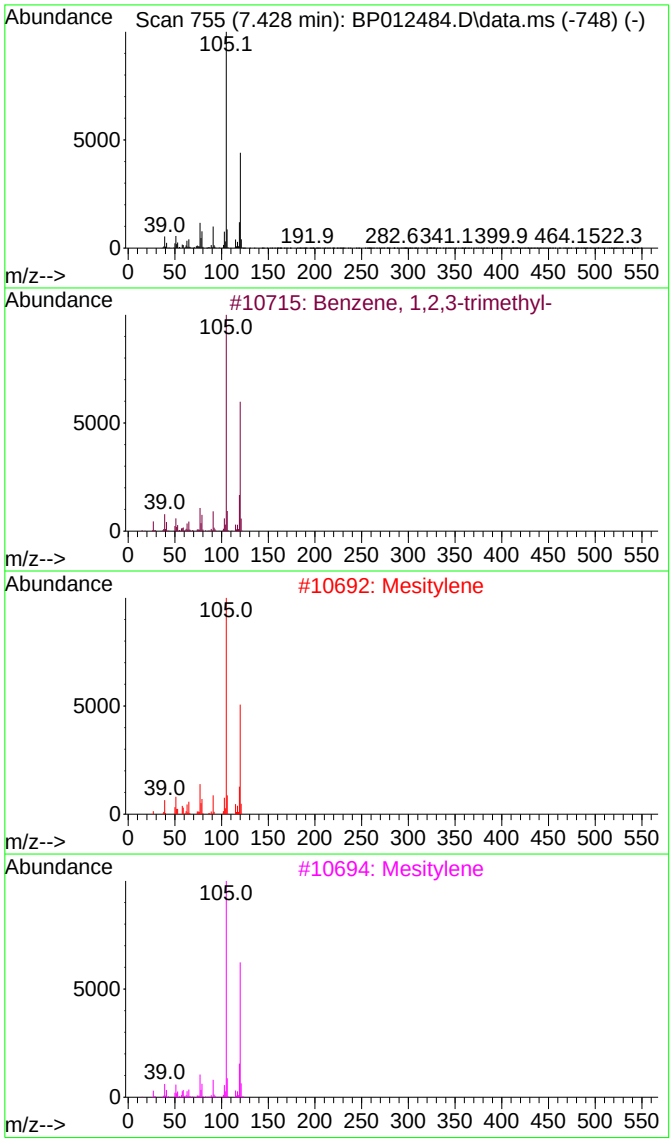
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

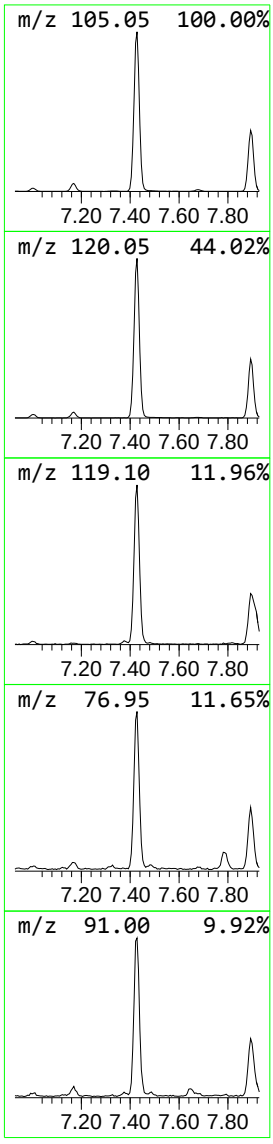
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.428	18.47 ng/ul	1458080	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, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94





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Data File : BP012484.D
Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

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

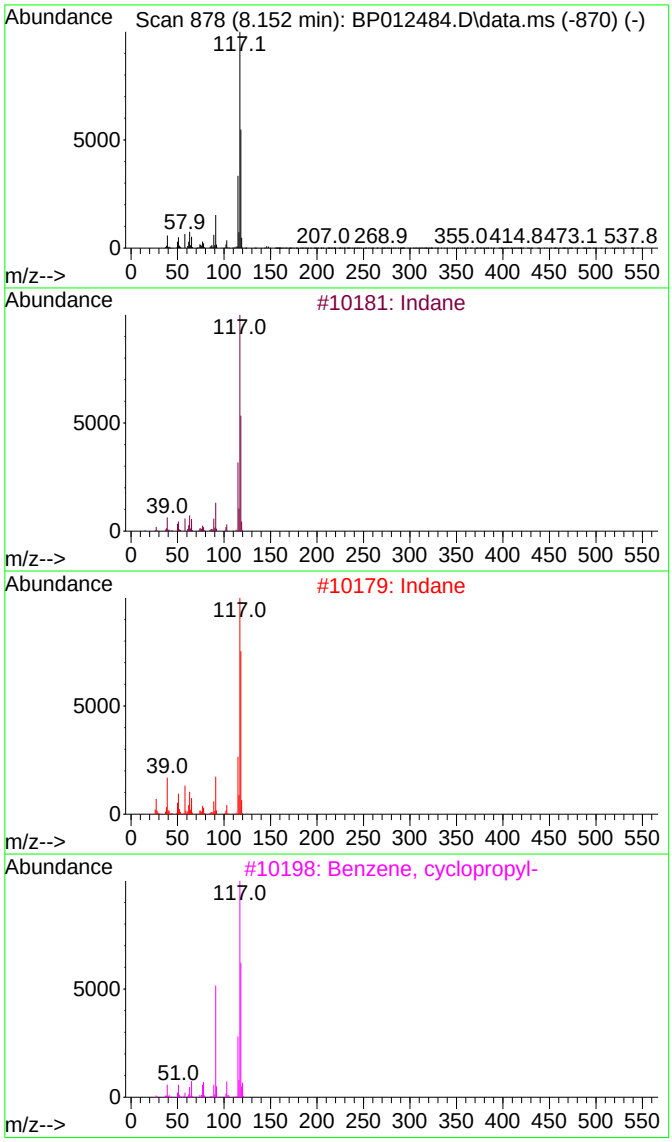
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

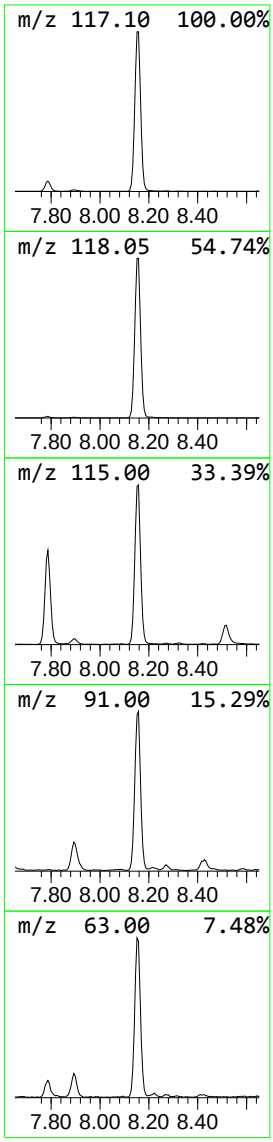
****

Peak Number 10 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.152	28.61 ng/ul	2257860	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	93
2	Indane	118	C9H10	000496-11-7	91
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	83
4	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	80
5	Deltacyclene	118	C9H10	007785-10-6	72





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Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

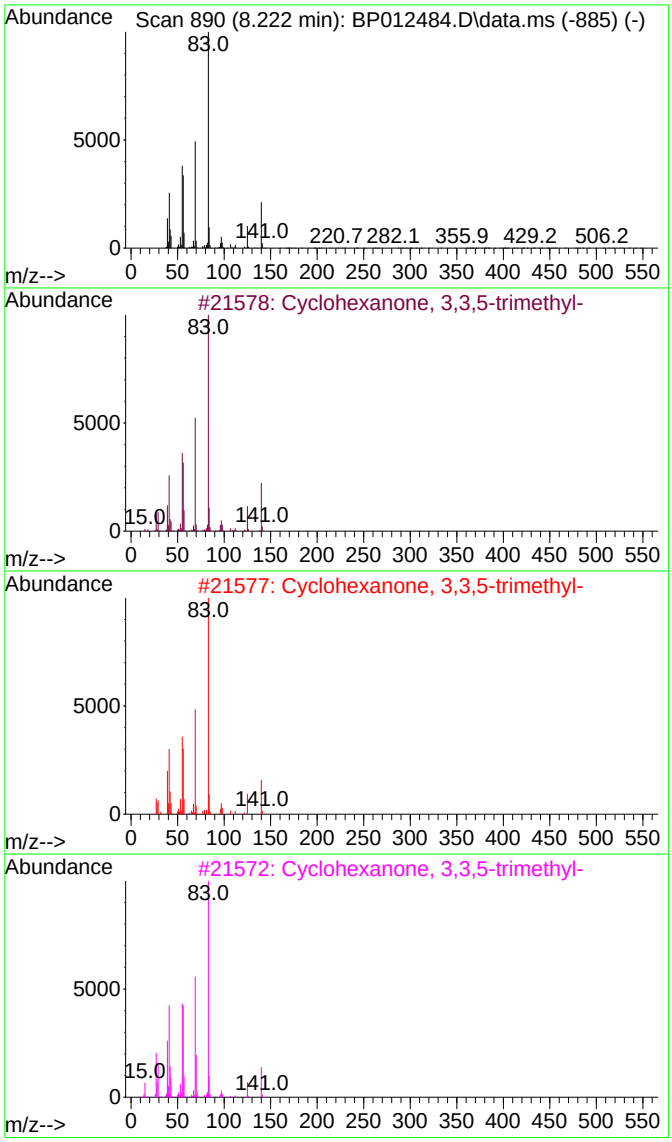
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

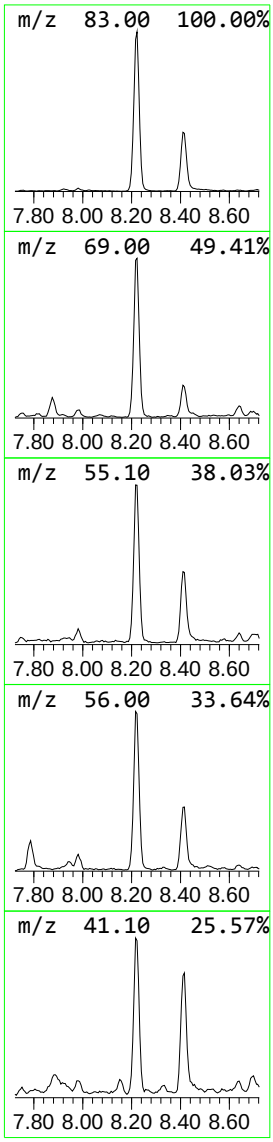
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Cyclohexanone, 3,3,5-trimet... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.222	8.60 ng/ul	678375	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	96
2	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	94
3	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	91
4	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	91
5	Undecenyl tiglate, 2E-	252	C16H28O2	1000383-56-5	50





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Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

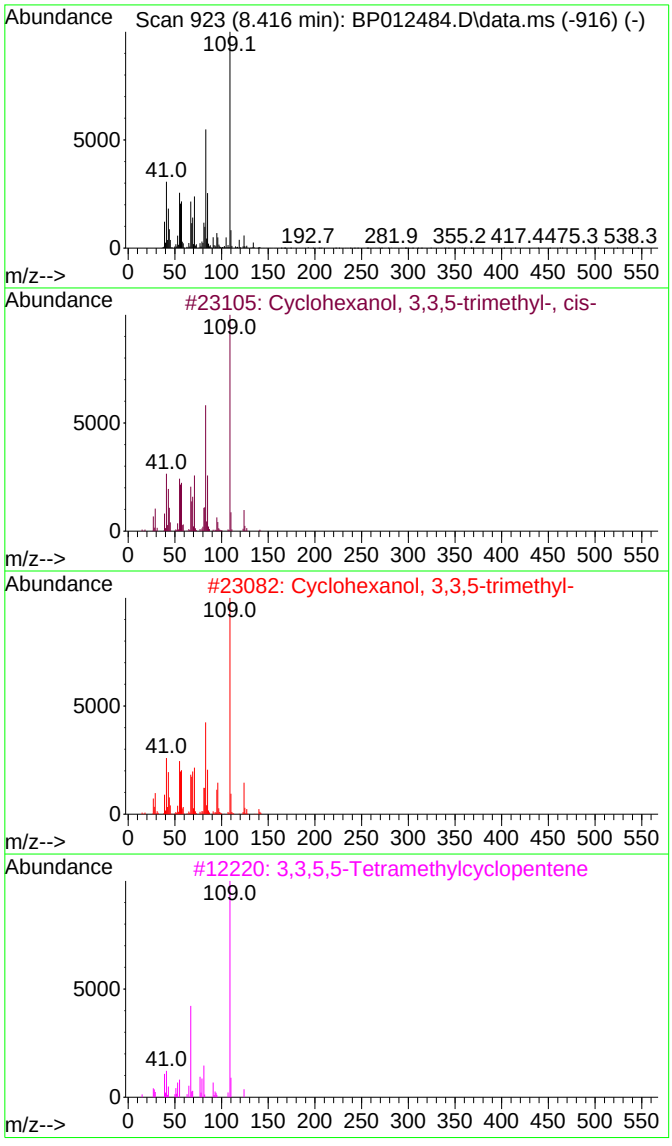
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

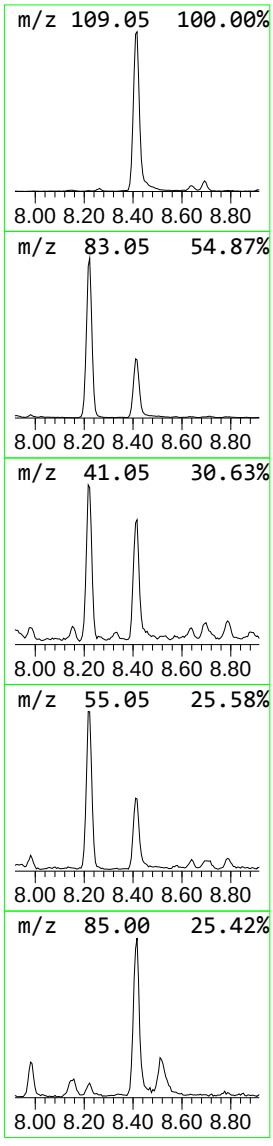
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Cyclohexanol, 3,3,5-trimeth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.416	9.90 ng/ul	781709	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	83
2	Cyclohexanol, 3,3,5-trimethyl-	142	C9H18O	000116-02-9	38
3	3,3,5,5-Tetramethylcyclopentene	124	C9H16	038667-10-6	38
4	2,4,4,7-Tetramethyl-octa-5,7-die...	180	C12H20O	1000190-70-6	37
5	1-(1,2,3-Trimethyl-cyclopent-2-e...	152	C10H16O	070987-81-4	35





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Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

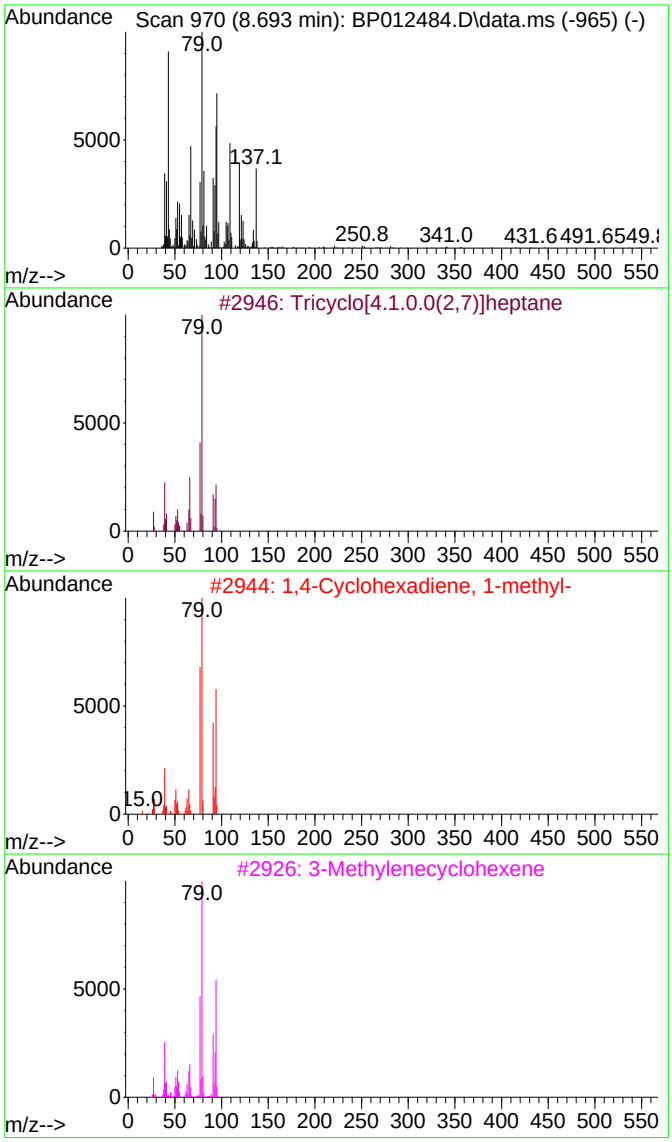
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

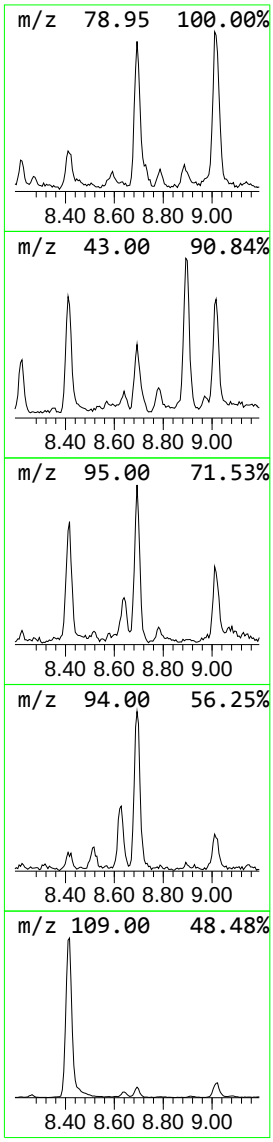
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 unknown-01 Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.693	2.11 ng/ul	166282	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tricyclo[4.1.0.0(2,7)]heptane	94	C7H10	000287-13-8	38
2	1,4-Cyclohexadiene, 1-methyl-	94	C7H10	004313-57-9	25
3	3-Methylenecyclohexene	94	C7H10	001888-90-0	25
4	Cyclopropane, (1-methyl-1,2-prop...	94	C7H10	051549-86-1	22
5	4,5,6,7-Tetrahydro-1H-1,2,3-benz...	123	C6H9N3	006789-99-7	22





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Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

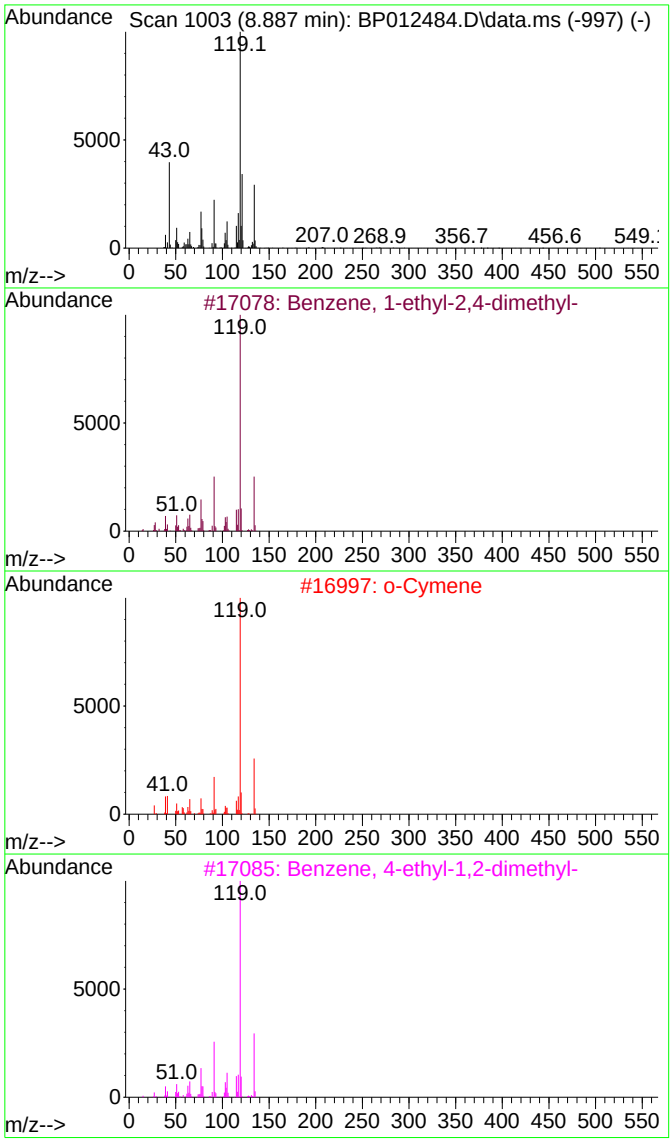
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

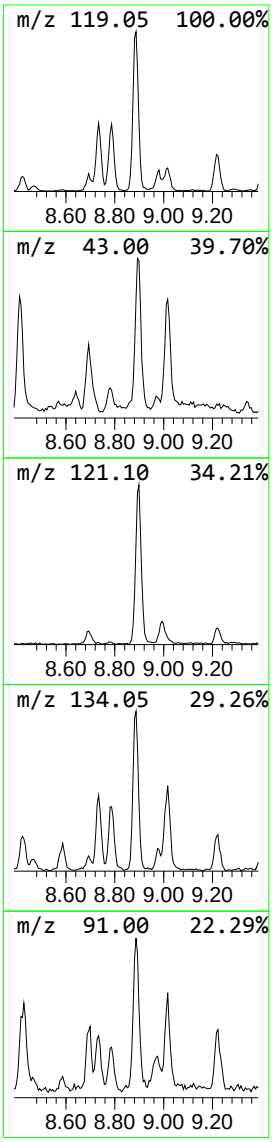
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.887	3.42 ng/ul	270127	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	96
2	o-Cymene	134	C10H14	000527-84-4	94
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
4	p-Cymene	134	C10H14	000099-87-6	94
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93





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Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

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

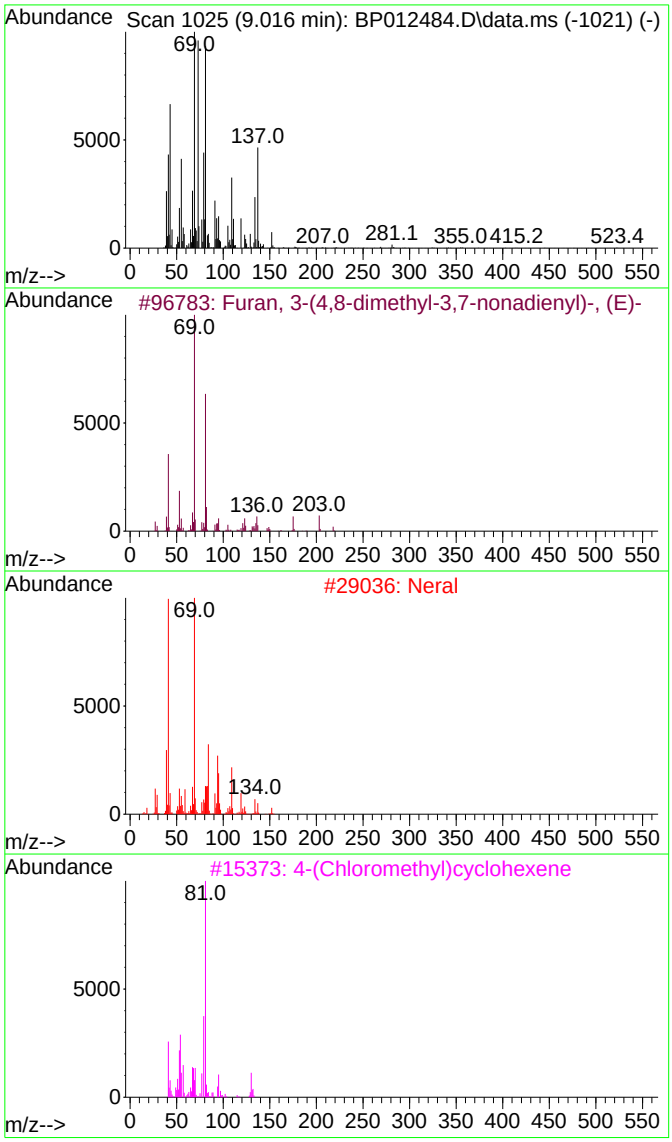
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

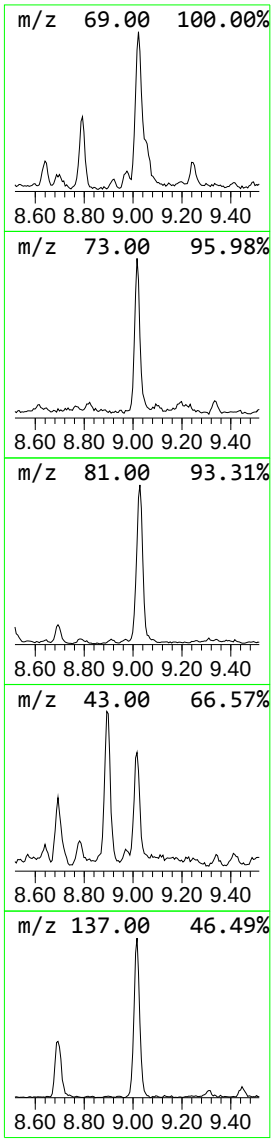
****

Peak Number 15 unknown-02 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.016	6.07 ng/ul	479086	1,4-Dichlorobenzene-d4	7.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Furan, 3-(4,8-dimethyl-3,7-nonad...	218	C15H22O	023262-34-2	38
2	Neral	152	C10H16O	000106-26-3	27
3	4-(Chloromethyl)cyclohexene	130	C7H11Cl	002555-08-0	27
4	Cyclopropanemethanol, 2-methyl-2...	168	C11H20O	098678-70-7	22
5	3,7-Octadien-2-one, (E)-	124	C8H12O	025172-06-9	22





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Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

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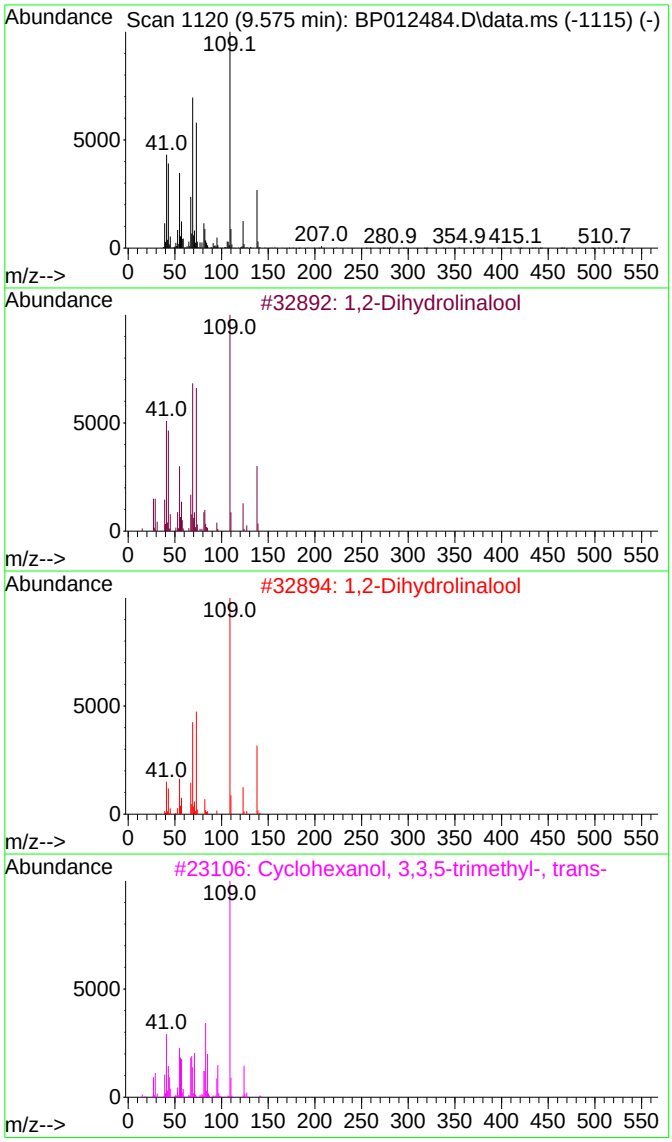
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

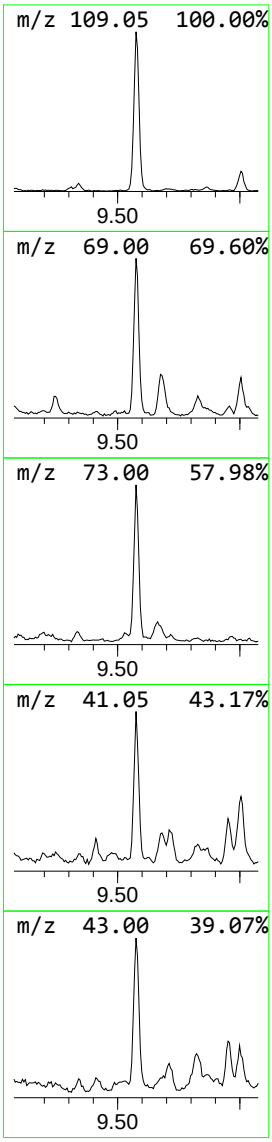
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 1,2-Dihydrolinalool Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.575	3.29 ng/ul	377981	Naphthalene-d8	10.593

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Dihydrolinalool	156	C10H20O	018479-51-1	80
2	1,2-Dihydrolinalool	156	C10H20O	018479-51-1	64
3	Cyclohexanol, 3,3,5-trimethyl-, ...	142	C9H18O	000767-54-4	59
4	2H-Pyrazolo[3,4-c]pyridin-3-amin...	138	C6H10N4	1000362-58-3	49
5	1-Propanone, 1-(5-methyl-2-furan...	138	C8H10O2	010599-69-6	47





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Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

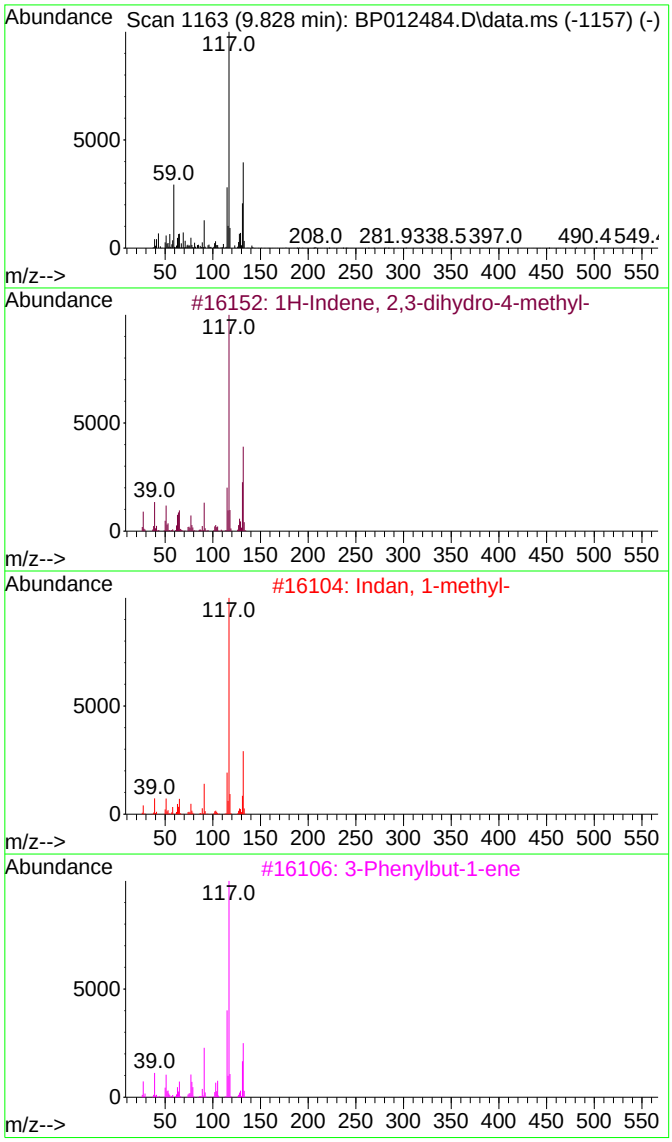
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

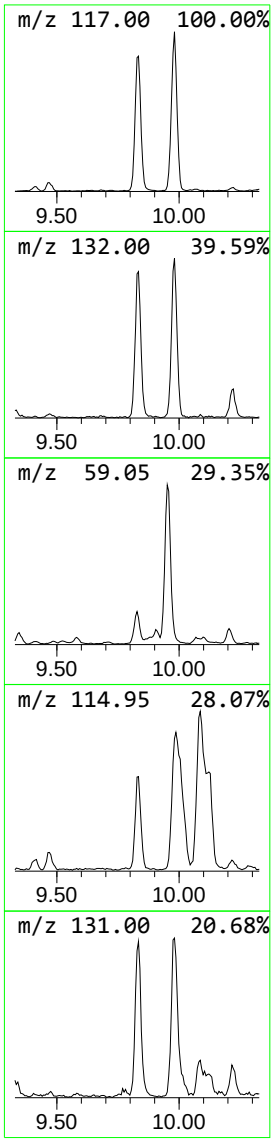
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.828	3.43 ng/ul	394653	Naphthalene-d8	10.593

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
2	Indan, 1-methyl-	132	C10H12	000767-58-8	76
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	74
4	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	72
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	72





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Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

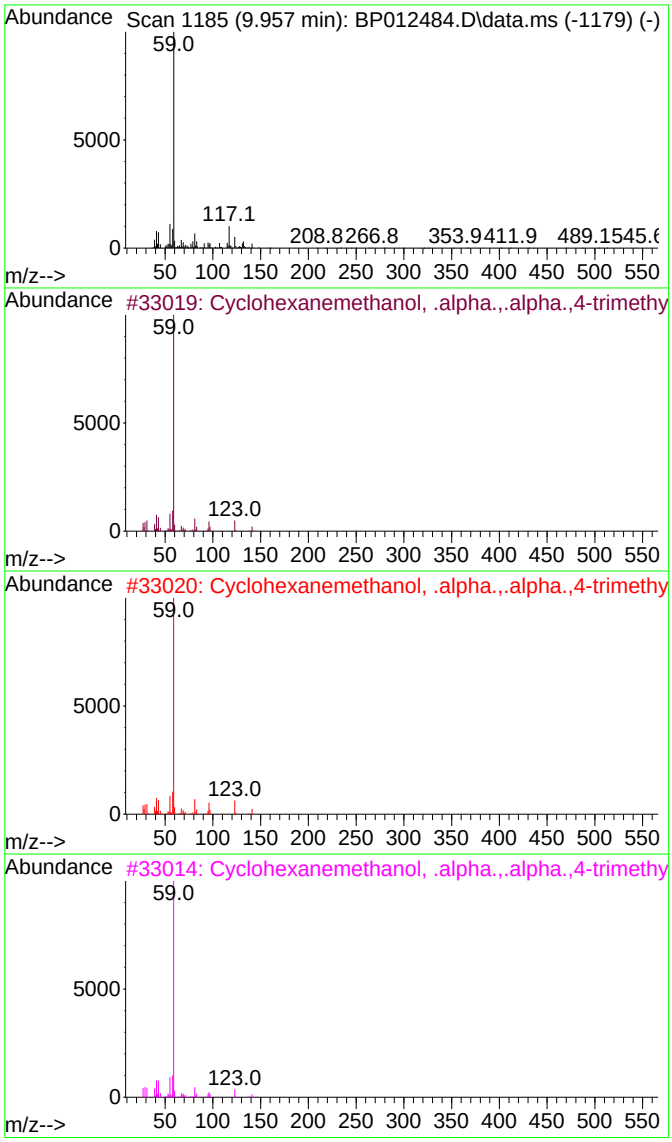
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

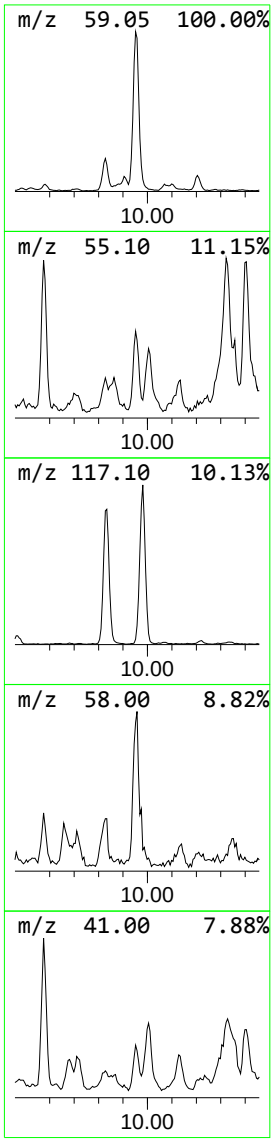
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Cyclohexanemethanol, .alpha... Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.957	2.08 ng/ul	239107	Naphthalene-d8	10.593

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	007322-63-6	64
2	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	005114-00-1	64
3	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	64
4	2,3-Dimethyl-4-penten-2-ol	114	C7H14O	019781-52-3	53
5	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	45





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Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

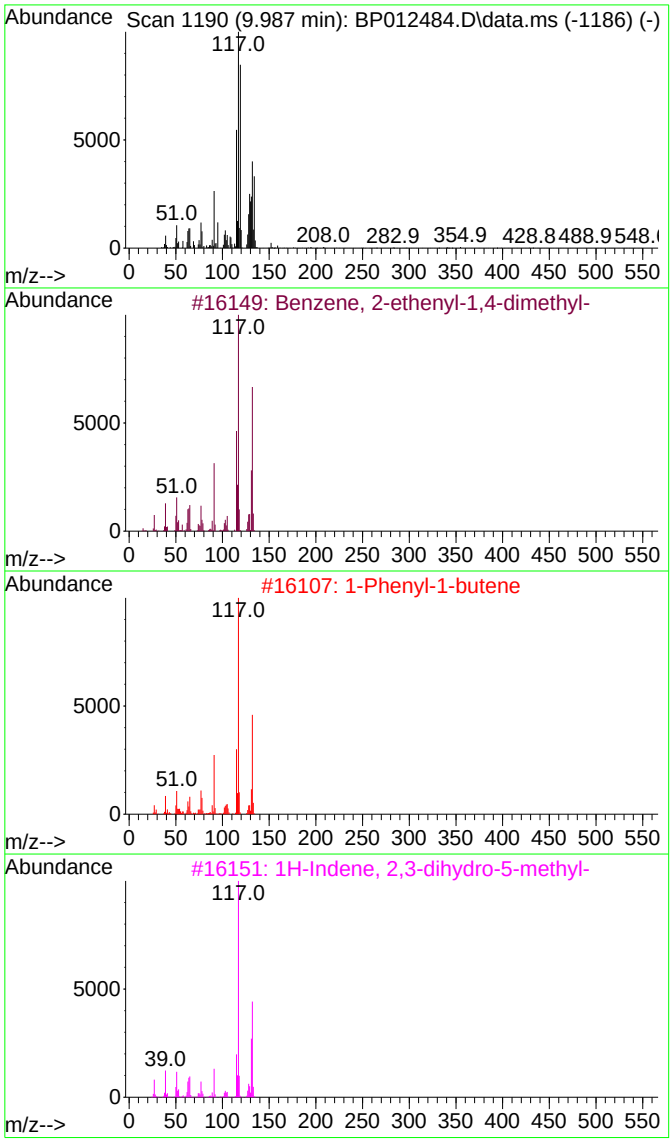
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

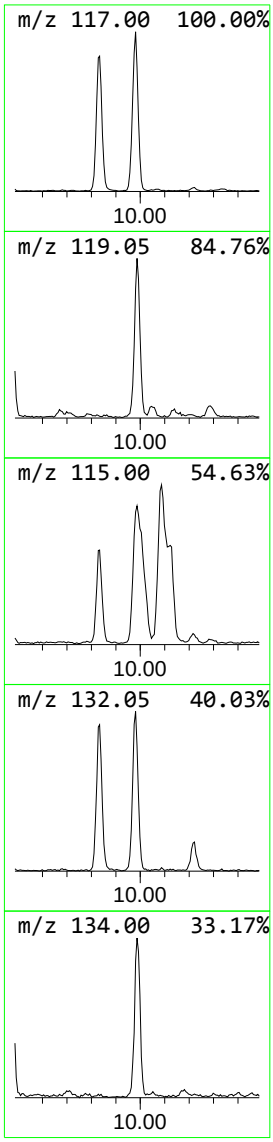
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.987	6.72 ng/ul	773261	Naphthalene-d8	10.593

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	70
2	1-Phenyl-1-butene	132	C10H12	000824-90-8	50
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	50
4	(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	49
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	46





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Data File : BP012484.D
Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

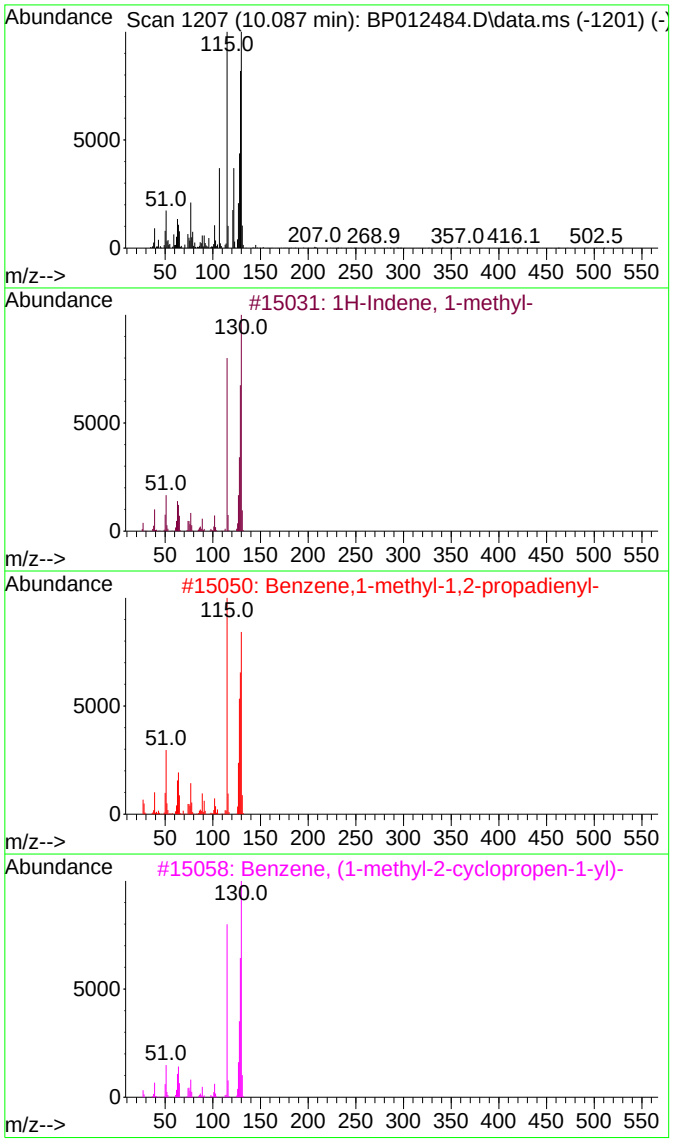
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

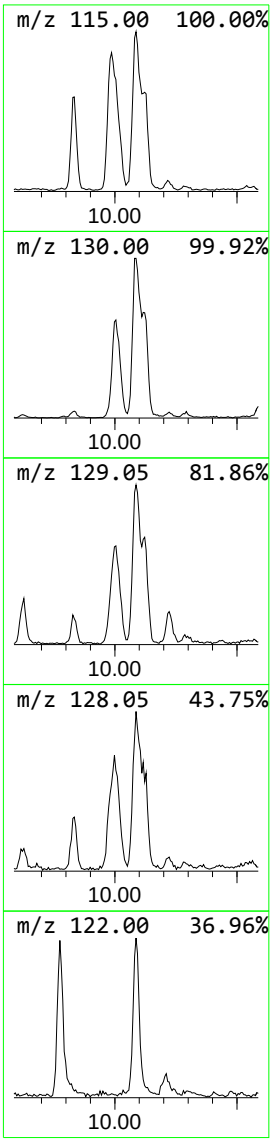
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 1H-Indene, 1-methyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.087	2.75 ng/ul	316004	Naphthalene-d8	10.593

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 1-methyl-	130	C10H10	000767-59-9	96
2	Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	95
3	Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	93
4	2-Methylindene	130	C10H10	002177-47-1	90
5	Benzene, 1-butynyl-	130	C10H10	000622-76-4	89





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Data File : BP012484.D
Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

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

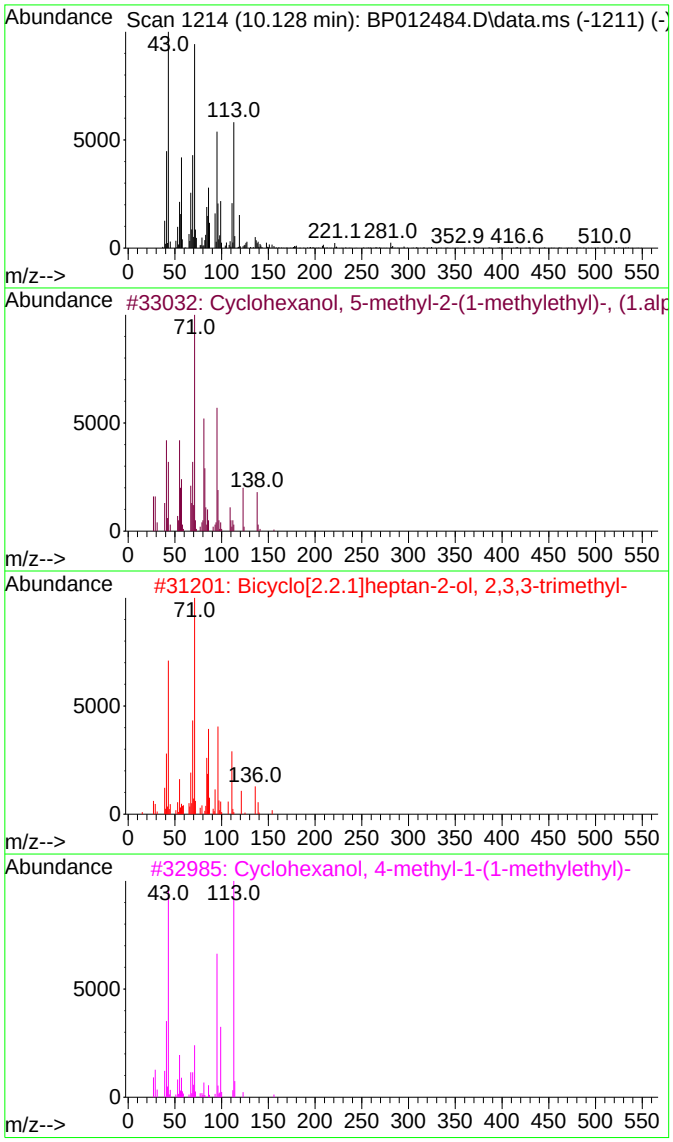
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

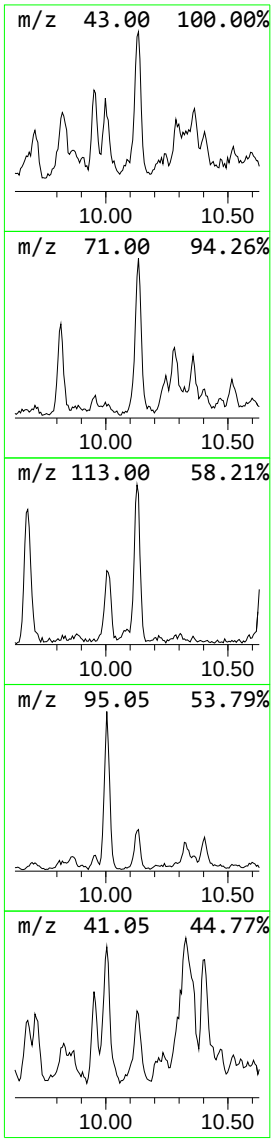
****

Peak Number 21 unknown-03 Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.128	2.05 ng/ul	235728	Naphthalene-d8	10.593

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanol, 5-methyl-2-(1-meth...	156	C10H20O	000491-02-1	46
2	Bicyclo[2.2.1]heptan-2-ol, 2,3,3...	154	C10H18O	000465-31-6	43
3	Cyclohexanol, 4-methyl-1-(1-meth...	156	C10H20O	000470-65-5	38
4	Heptane, 2,5,5-trimethyl-	142	C10H22	001189-99-7	35
5	1-Propene, 1-methoxy-2-methyl-	86	C5H10O	017574-84-4	30





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Data File : BP012484.D
Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

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

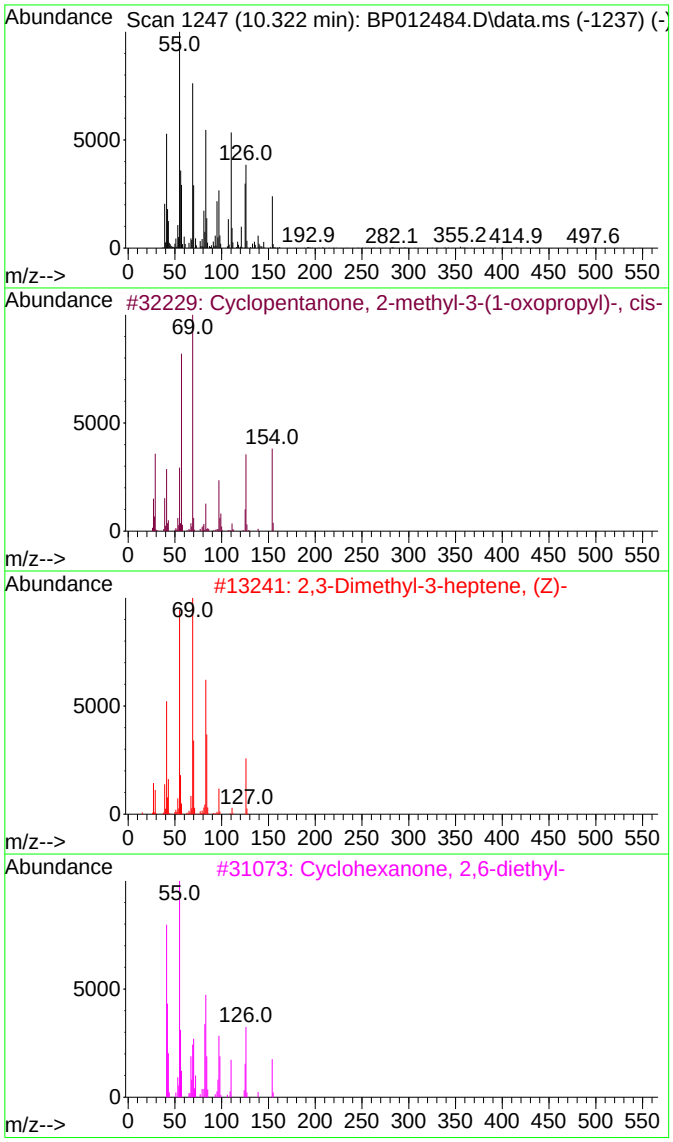
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

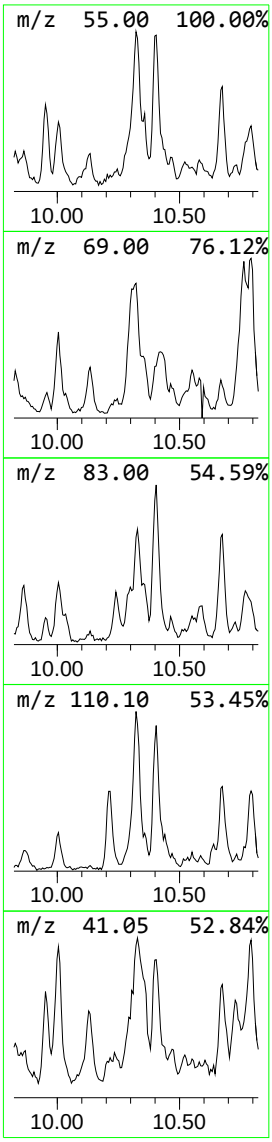
****

Peak Number 22 unknown-04 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.322	3.04 ng/ul	349282	Naphthalene-d8	10.593

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentanone, 2-methyl-3-(1-ox...	154	C9H14O2	079881-04-2	47
2	2,3-Dimethyl-3-heptene, (Z)-	126	C9H18	059643-73-1	43
3	Cyclohexanone, 2,6-diethyl-	154	C10H18O	016519-68-9	38
4	Cyclooctane, 1,4-dimethyl-, trans-	140	C10H20	013151-98-9	38
5	(1R,6S)-3-Methoxy-7-oxabicyclo[4...	154	C7H6O4	1000482-42-1	27





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Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

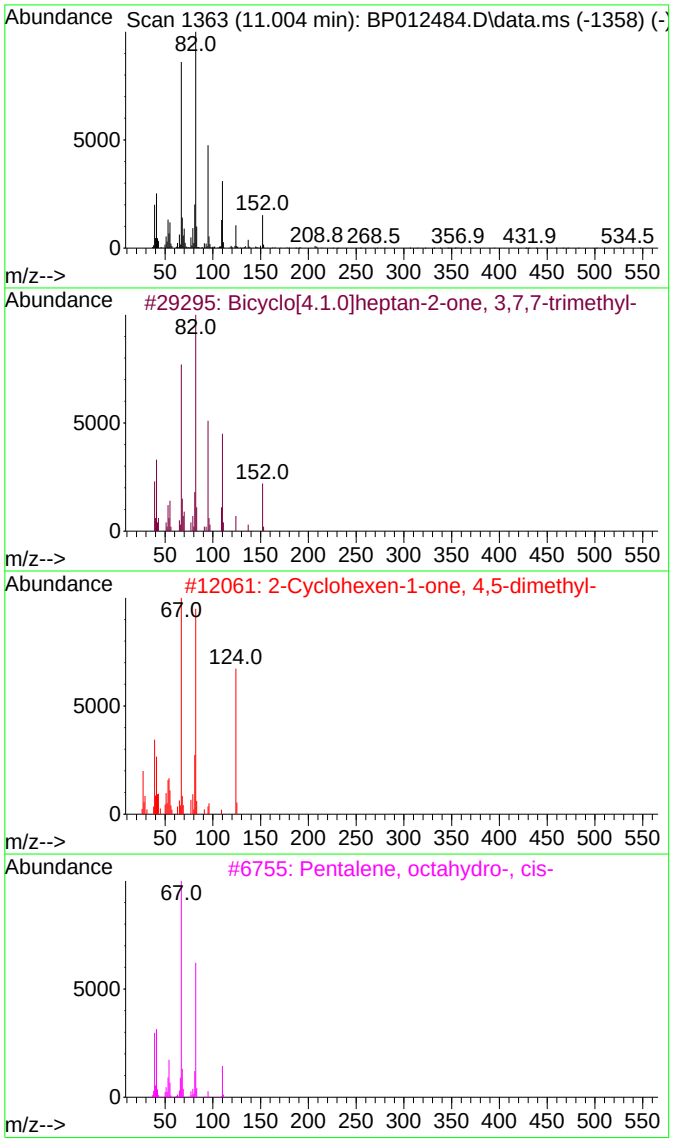
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

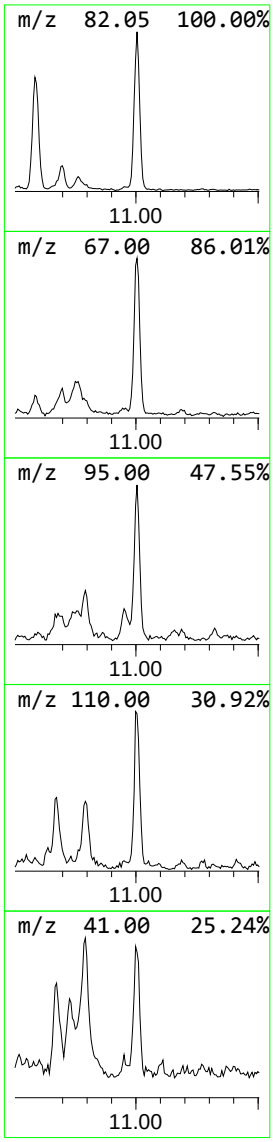
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Bicyclo[4.1.0]heptan-2-one,... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.004	2.28 ng/ul	262855	Naphthalene-d8	10.593

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[4.1.0]heptan-2-one, 3,7,...	152	C10H16O	000497-62-1	94
2	2-Cyclohexen-1-one, 4,5-dimethyl-	124	C8H12O	005715-25-3	62
3	Pentalene, octahydro-, cis-	110	C8H14	001755-05-1	62
4	Cyclopropene, 1-butyl-2-ethyl-	124	C9H16	050915-91-8	58
5	1-Methylcycloheptene	110	C8H14	055308-20-8	53





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Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

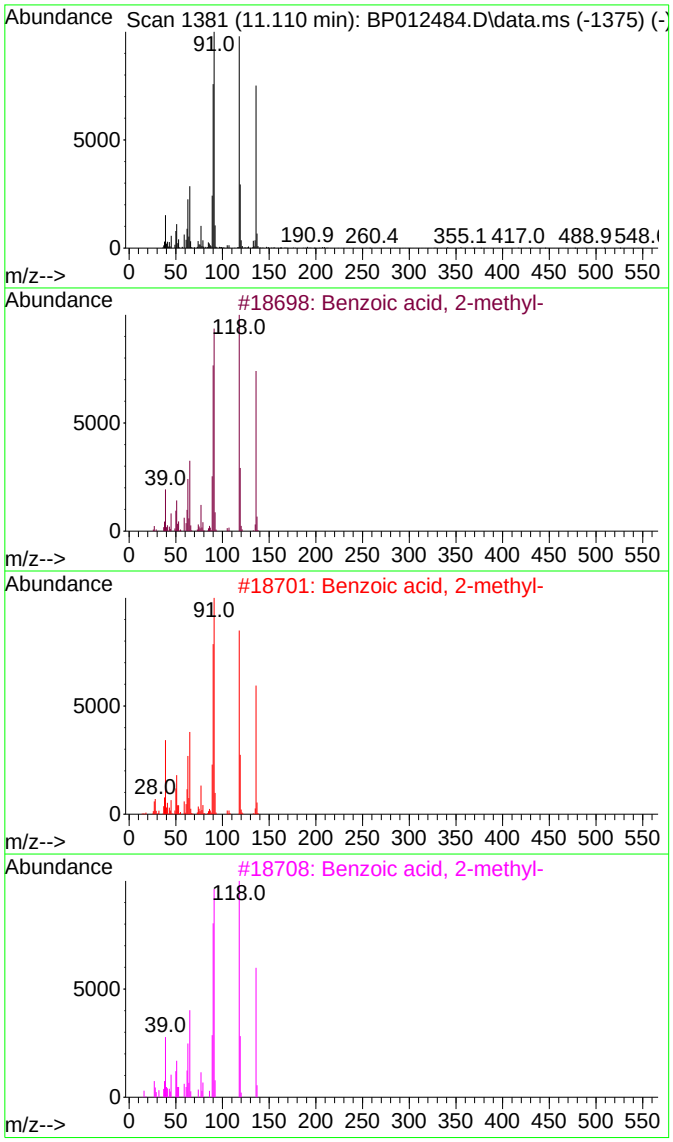
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

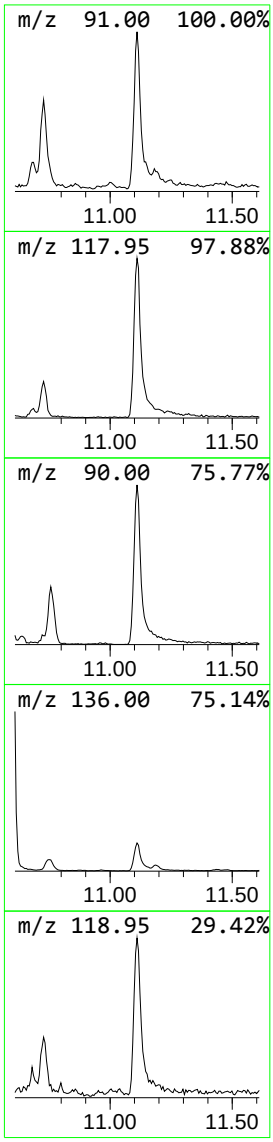
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Benzoic acid, 2-methyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.110	3.01 ng/ul	345845	Naphthalene-d8	10.593

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 2-methyl-	136	C8H8O2	000118-90-1	98
2	Benzoic acid, 2-methyl-	136	C8H8O2	000118-90-1	97
3	Benzoic acid, 2-methyl-	136	C8H8O2	000118-90-1	96
4	Benzoic acid, 2-methyl-	136	C8H8O2	000118-90-1	96
5	Bicyclo[4.1.0]hepta-1,3,5-triene	90	C7H6	004646-69-9	53





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Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

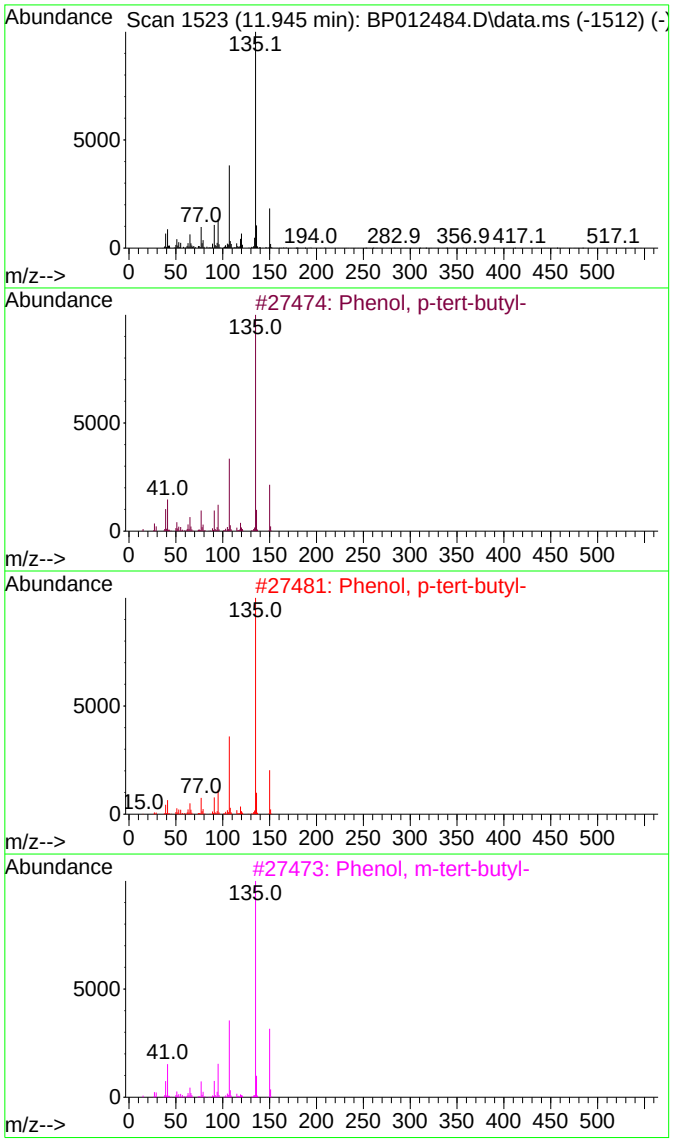
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

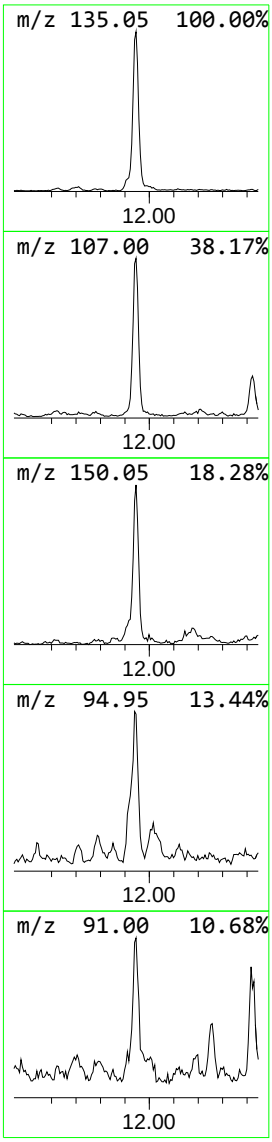
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 Phenol, p-tert-butyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.945	3.77 ng/ul	433977	Naphthalene-d8	10.593

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	94
3	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	93
4	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	91
5	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	90





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Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

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

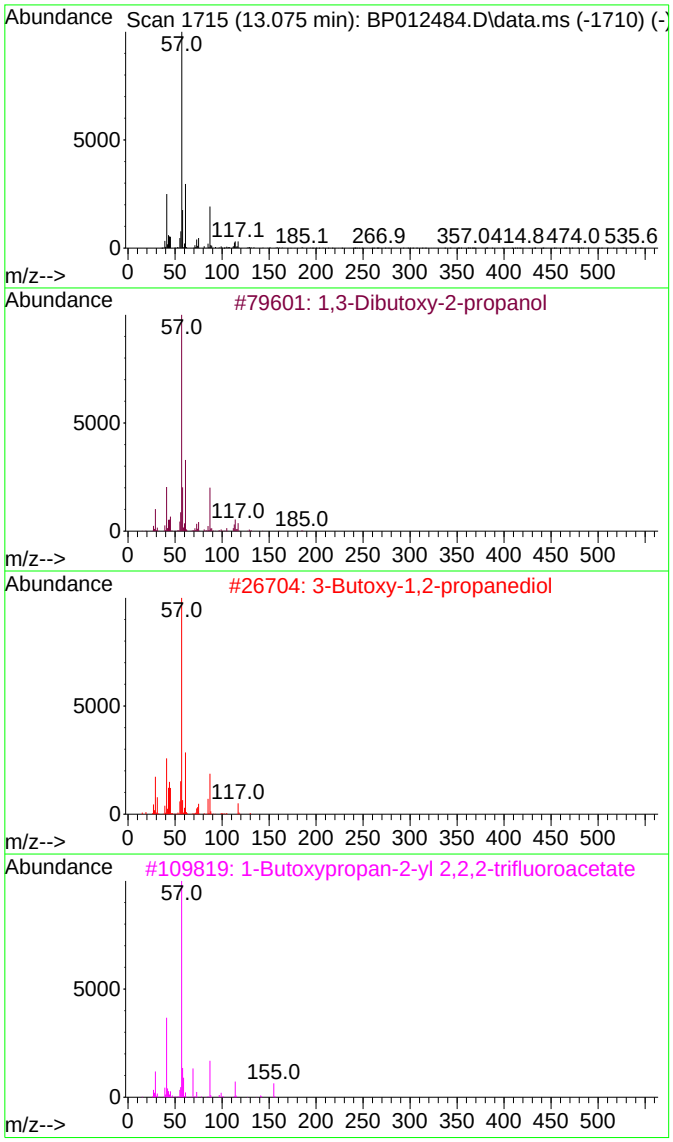
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

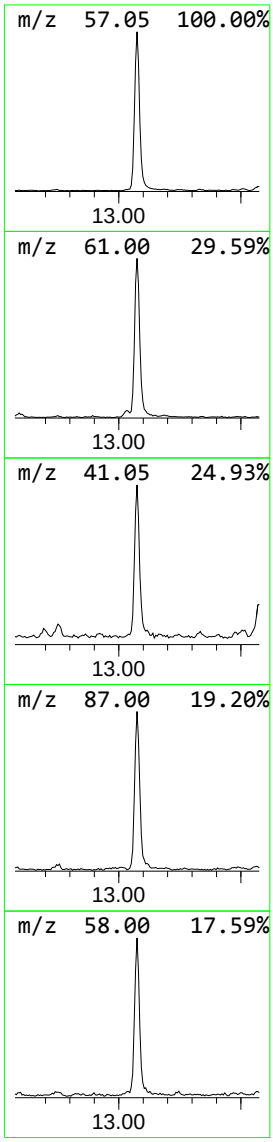
****

Peak Number 26 1,3-Dibutoxy-2-propanol Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.075	2.53 ng/ul	469267	Acenaphthene-d10	14.440

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Dibutoxy-2-propanol	204	C11H24O3	002216-77-5	59
2	3-Butoxy-1,2-propanediol	148	C7H16O3	000624-52-2	50
3	1-Butoxypropan-2-yl 2,2,2-triflu...	228	C9H15F3O3	1000367-08-1	40
4	n-Butyl ether	130	C8H18O	000142-96-1	28
5	S-(-)-2-Methylbutylamine	87	C5H13N	020626-52-2	10





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Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

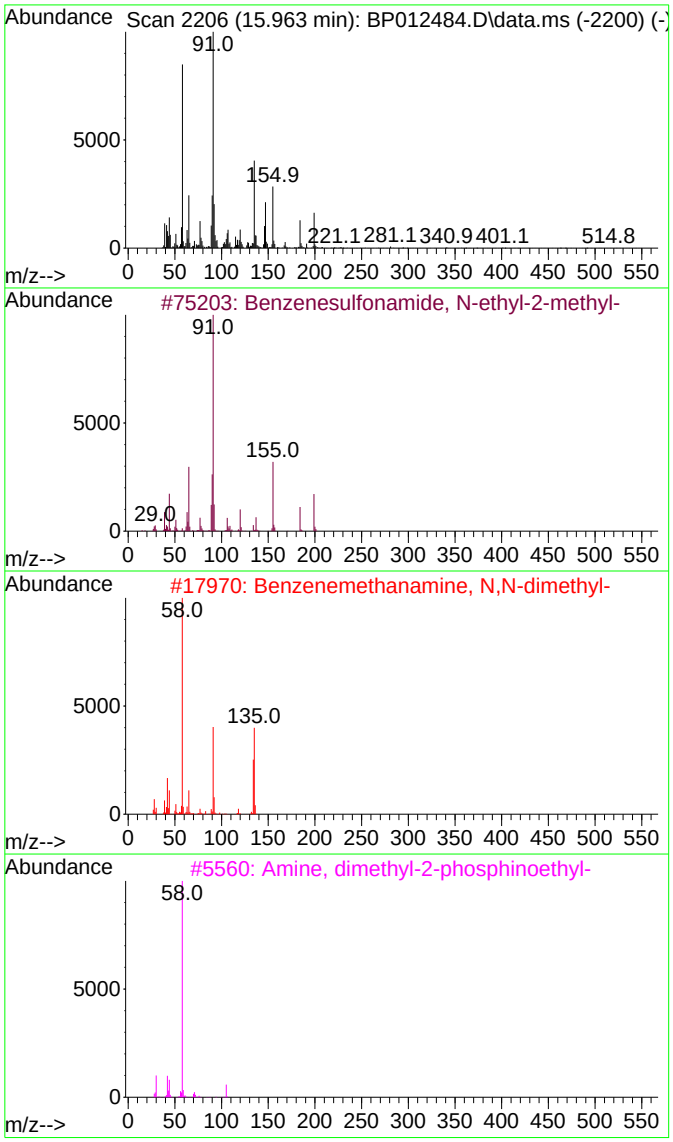
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

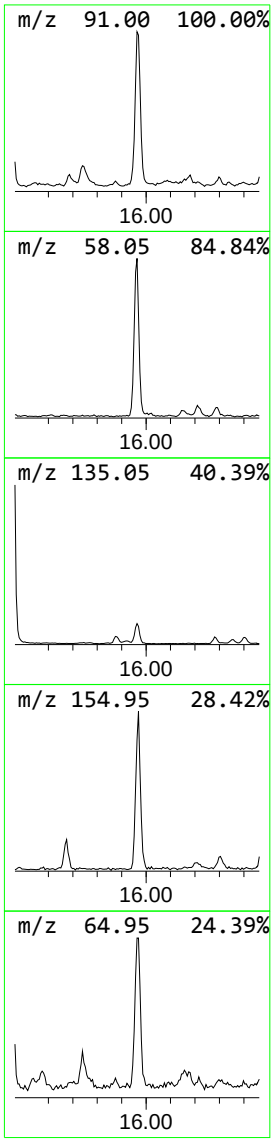
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 Benzenesulfonamide, N-ethyl... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.963	2.79 ng/ul	613416	Phenanthrene-d10	17.204

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	92
2	Benzenemethanamine, N,N-dimethyl-	135	C9H13N	000103-83-3	43
3	Amine, dimethyl-2-phosphinoethyl-	105	C4H12NP	161944-90-7	25
4	2,2'-Bis[(3-dimethylaminopropyl)...	474	C24H34N4O2S2	032276-24-7	25
5	5H-Dibenzo[a,d]cyclohepten-5-ol,...	295	C20H25NO	001159-03-1	25





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Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

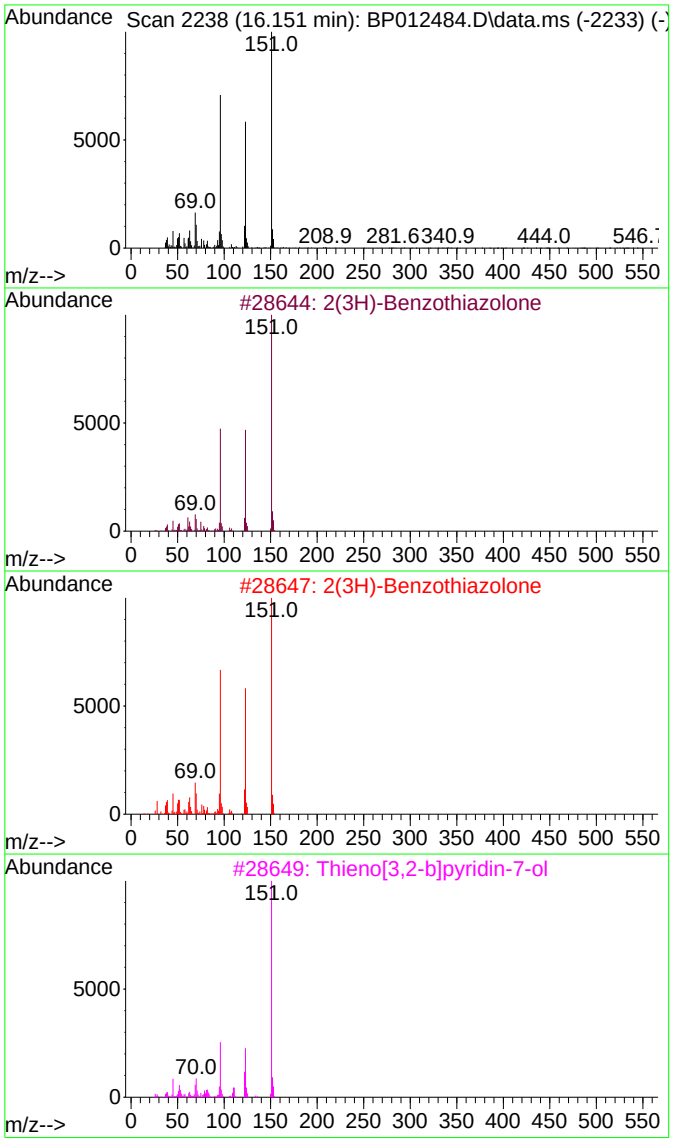
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

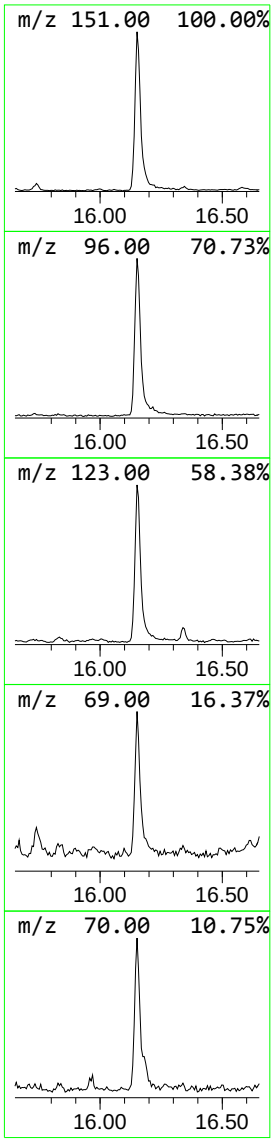
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 28 2(3H)-Benzothiazolone Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.151	2.68 ng/ul	589131	Phenanthrene-d10	17.204

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	94
2	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	94
3	Thieno[3,2-b]pyridin-7-ol	151	C7H5NOS	107818-20-2	58
4	1,2-Benzisothiazol-3(2H)-one	151	C7H5NOS	002634-33-5	53
5	6-Methyl-7-oxo-4,7-dihydro-triaz...	151	C5H5N5O	057250-39-2	50





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Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

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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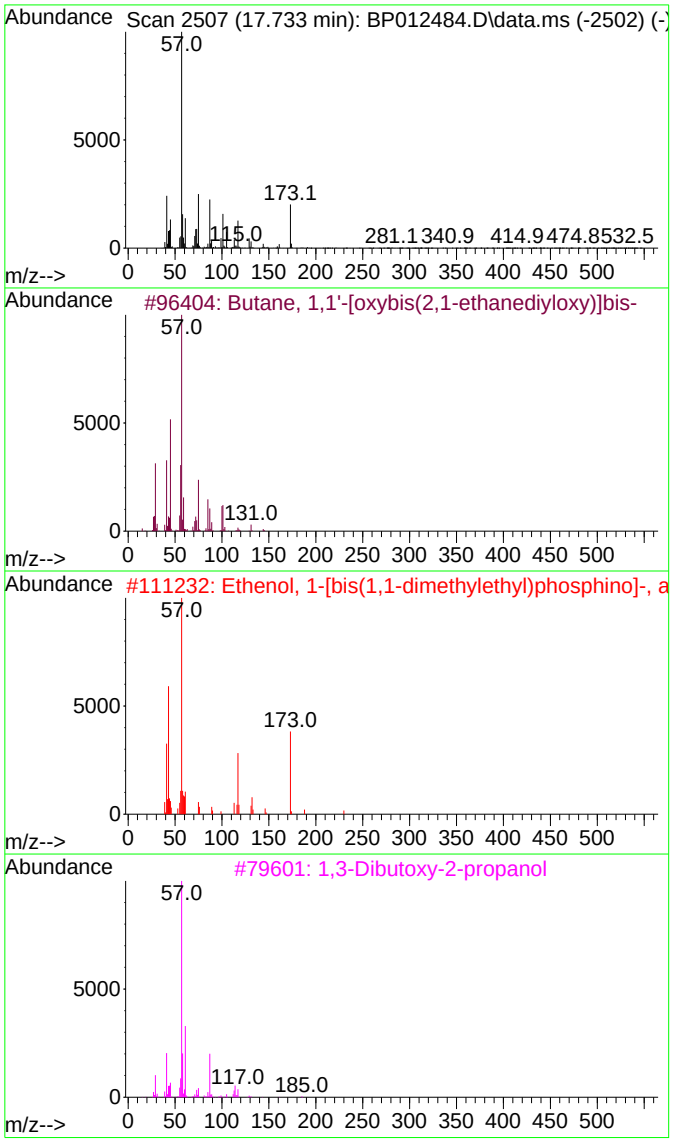
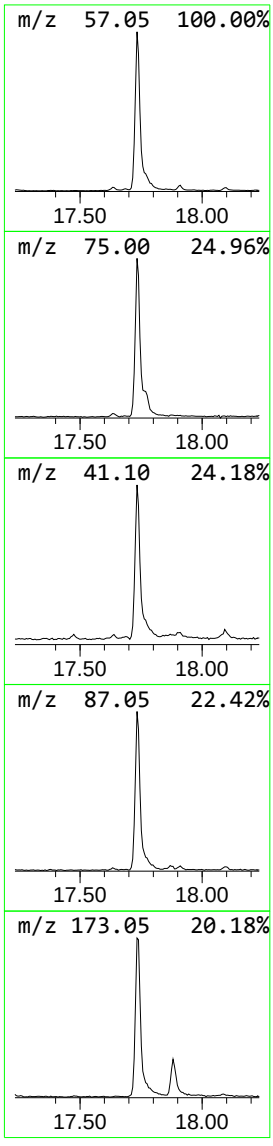
Peak Number 29 unknown-05 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.733	8.22 ng/ul	1808580	Phenanthrene-d10	17.204

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 1,1'-[oxybis(2,1-ethaned...	218	C12H26O3	000112-73-2	32
2	Ethenol, 1-[bis(1,1-dimethylethy...	230	C12H23O2P	050838-15-8	32
3	1,3-Dibutoxy-2-propanol	204	C11H24O3	002216-77-5	27
4	Butane, 1,1'-[oxybis(2,1-ethaned...	218	C12H26O3	000112-73-2	25
5	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	23





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Data File : BP012484.D
Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

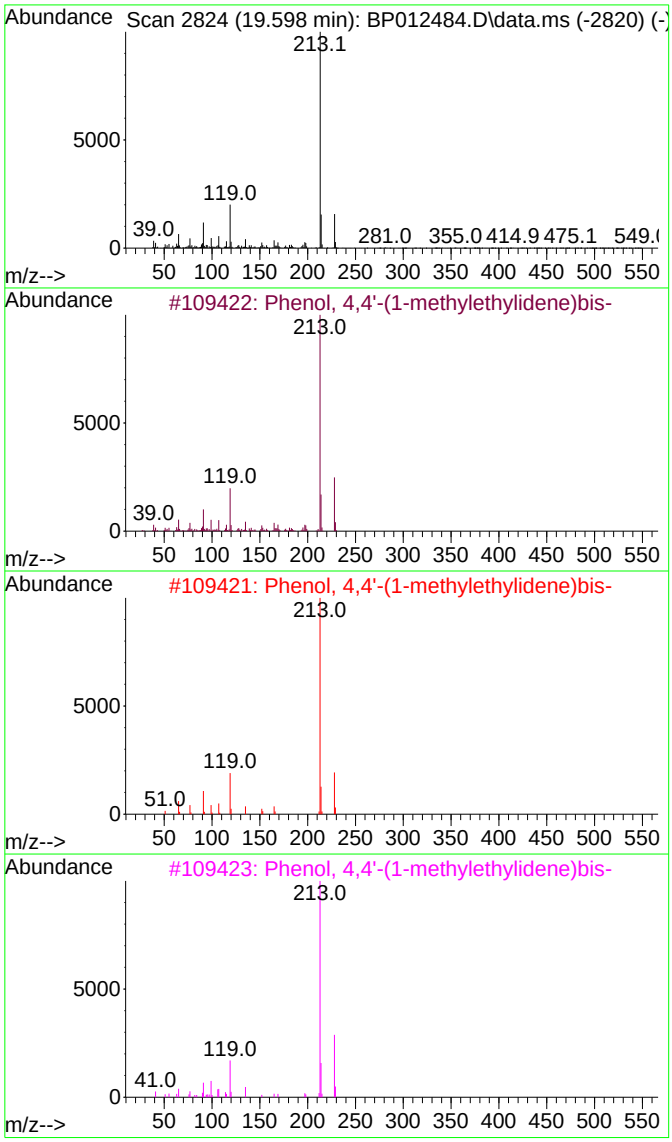
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

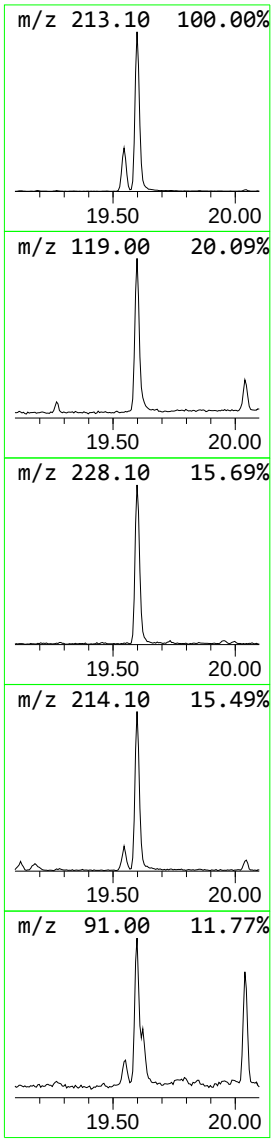
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 30 Phenol, 4,4'-(1-methylethyl... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.598	4.01 ng/ul	946555	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	96
2	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	95
3	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	90
4	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	87
5	4,4'-(1,3-Dimethylbutylidene)bis...	270	C18H22O2	006807-17-6	74





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Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
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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

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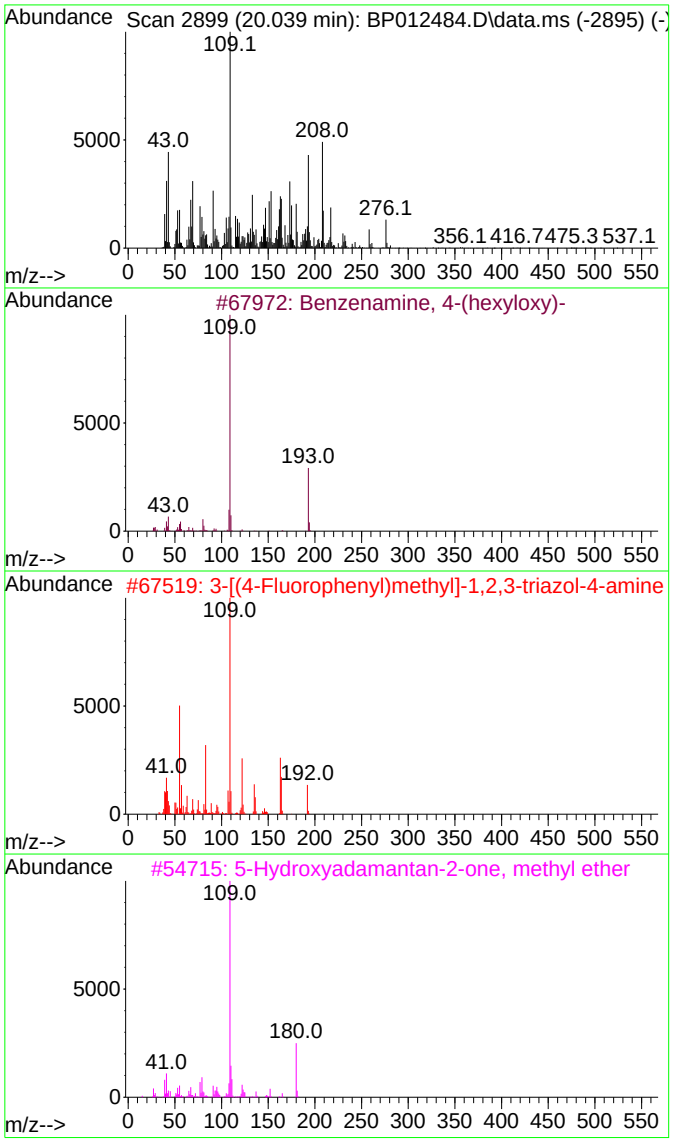
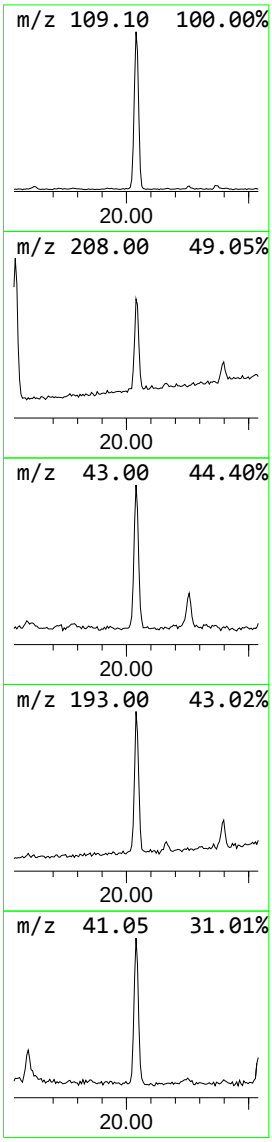
Peak Number 31 unknown-06 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.039	5.37 ng/ul	1269180	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 4-(hexyloxy)-	193	C12H19NO	039905-57-2	38
2	3-[(4-Fluorophenyl)methyl]-1,2,3...	192	C9H9FN4	1000435-39-8	30
3	5-Hydroxyadamantan-2-one, methyl...	180	C11H16O2	115009-22-8	25
4	1,2,4-Oxadiazol-5-amine, 3-(4-am...	276	C10H12N8O2	1000351-27-1	18
5	2-(1-Cyclopent-1-enyl-1-methylet...	192	C13H20O	1000190-57-4	18





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
Data File : BP012484.D
Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

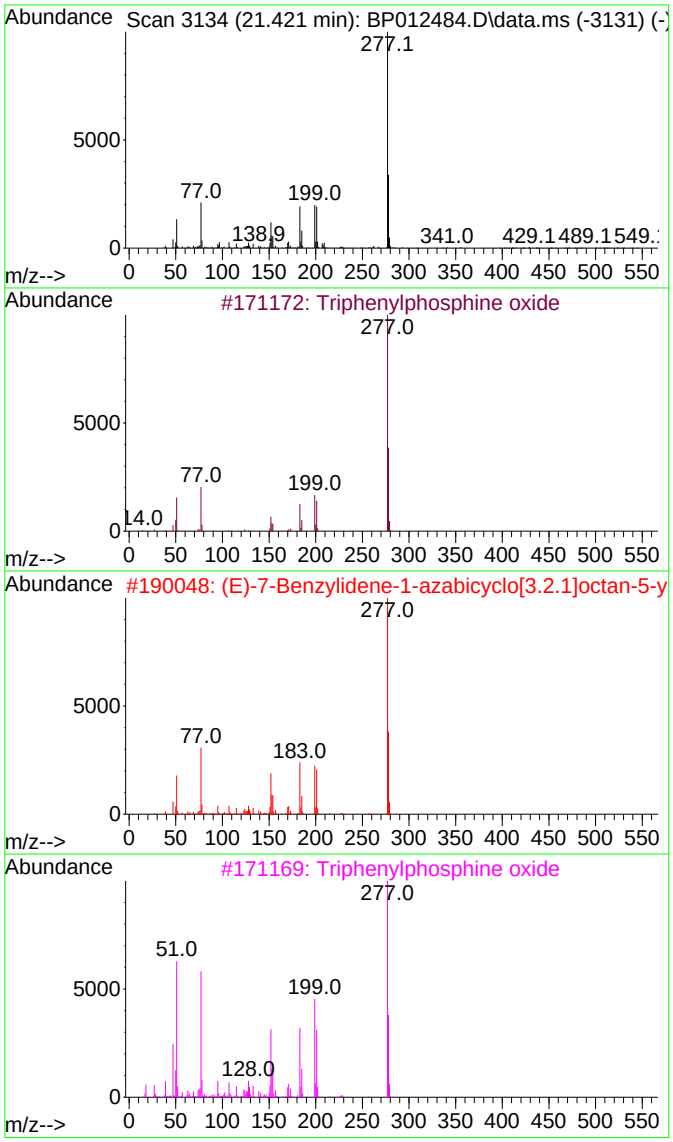
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

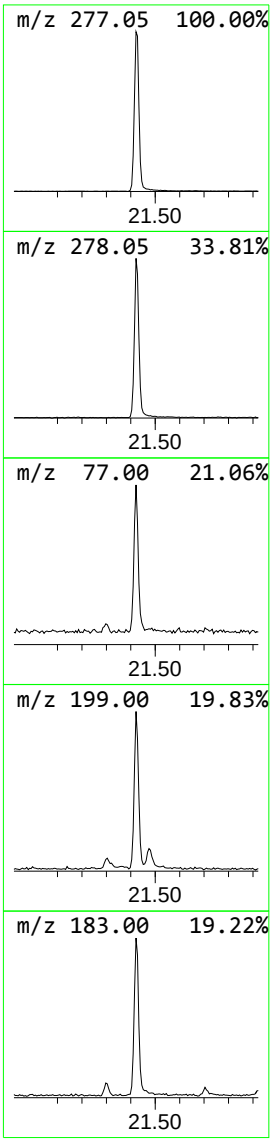
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 32 Triphenylphosphine oxide Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.421	2.23 ng/ul	527266	Chrysene-d12	21.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	96
2	(E)-7-Benzylidene-1-azabicyclo[3...	293	C15H19NO3S	1000483-83-4	90
3	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	87
4	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	58
5	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	53





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
Data File : BP012484.D
Acq On : 03 Nov 2022 11:25
Operator : CG/JU
Sample : N5300-02DL2 10X
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73DL2

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
Benzene, 1,3-di...	5.416	2.6	ng/ul	203118	1	7.787	1578530	20.0
1-[1,2,4]Triazo...	5.499	10.1	ng/ul	799168	1	7.787	1578530	20.0
3-Heptanone, 5-...	6.452	3.6	ng/ul	282725	1	7.787	1578530	20.0
Benzene, propyl-	6.758	2.6	ng/ul	203484	1	7.787	1578530	20.0
Benzene, 1,2,3-...	7.428	18.5	ng/ul	1458080	1	7.787	1578530	20.0
Indane	8.152	28.6	ng/ul	2257860	1	7.787	1578530	20.0
Cyclohexanone, ...	8.222	8.6	ng/ul	678375	1	7.787	1578530	20.0
Cyclohexanol, 3...	8.416	9.9	ng/ul	781709	1	7.787	1578530	20.0
unknown-01	8.693	2.1	ng/ul	166282	1	7.787	1578530	20.0
Benzene, 1-ethy...	8.887	3.4	ng/ul	270127	1	7.787	1578530	20.0
unknown-02	9.016	6.1	ng/ul	479086	1	7.787	1578530	20.0
1,2-Dihydrolina...	9.575	3.3	ng/ul	377981	2	10.593	2300740	20.0
1H-Indene, 2,3-...	9.828	3.4	ng/ul	394653	2	10.593	2300740	20.0
Cyclohexanemeth...	9.957	2.1	ng/ul	239107	2	10.593	2300740	20.0
Benzene, 2-ethe...	9.987	6.7	ng/ul	773261	2	10.593	2300740	20.0
1H-Indene, 1-me...	10.087	2.8	ng/ul	316004	2	10.593	2300740	20.0
unknown-03	10.128	2.0	ng/ul	235728	2	10.593	2300740	20.0
unknown-04	10.322	3.0	ng/ul	349282	2	10.593	2300740	20.0
Bicyclo[4.1.0]h...	11.004	2.3	ng/ul	262855	2	10.593	2300740	20.0
Benzoic acid, 2...	11.110	3.0	ng/ul	345845	2	10.593	2300740	20.0
Phenol, p-tert-...	11.945	3.8	ng/ul	433977	2	10.593	2300740	20.0
1,3-Dibutoxy-2-...	13.075	2.5	ng/ul	469267	3	14.440	3708950	20.0
Benzenesulfonam...	15.963	2.8	ng/ul	613416	4	17.204	4400760	20.0
2(3H)-Benzothia...	16.151	2.7	ng/ul	589131	4	17.204	4400760	20.0
unknown-05	17.733	8.2	ng/ul	1808580	4	17.204	4400760	20.0
Phenol, 4,4'-(1...	19.598	4.0	ng/ul	946555	5	21.298	4723340	20.0
unknown-06	20.039	5.4	ng/ul	1269180	5	21.298	4723340	20.0
Triphenylphosph...	21.421	2.2	ng/ul	527266	5	21.298	4723340	20.0