LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011924.D
Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

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-BP093022.M
Title : SVOA CALIBRATION

Signal : TIC: BP011924.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.034	3	8	20	rVV	4847897	6046314	86.35%	6.540%
2	3.128	21	24	30	rVB	49080	60448	0.86%	0.065%
3	3.299	49	53	60	rVB	50320	59361	0.85%	0.064%
4	3.728	123	126	134	rBV	240495	347392	4.96%	0.376%
5	3.799	134	138	149	rVB2	75123	135315	1.93%	0.146%
6	3.899	150	155	160	rVV	66728	91793	1.31%	0.099%
7	3.952	160	164	173	rVB	36229	54884	0.78%	0.059%
8	5.534	424	433	437	rBV2	41537	80666	1.15%	0.087%
9	5.881	484	492	497	rBV2	31147	59139	0.84%	0.064%
10	6.964	670	676	681	rBV	49088	73975	1.06%	0.080%
11	7.046	681	690	696	rVV2	207217	490453	7.00%	0.531%
12	7.099	696	699	705	rVB	53643	77899	1.11%	0.084%
13	7.211	712	718	723	rBV	753227	1237905	17.68%	1.339%
14	7.269	723	728	736	rVB	1325874	2096034	29.93%	2.267%
15	7.411	745	752	764	rBV	743084	1203959	17.19%	1.302%
16	7.528	764	772	787	rVB	1778572	2844470	40.62%	3.077%
17	7.881	825	832	845	rVV	766211	1316021	18.79%	1.424%
18	7.993	845	851	866	rVV	995986	1563483	22.33%	1.691%
19	8.187	875	884	890	rVV3	19905	54319	0.78%	0.059%
20	8.258	890	896	906	rVB	95935	159797	2.28%	0.173%
21	8.375	909	916	921	rBV	304703	472711	6.75%	0.511%
22	8.522	935	941	946	rBV	139367	230642	3.29%	0.249%
23	8.593	946	953	963	rVV	581415	1095343	15.64%	1.185%
24	8.687	963	969	977	rVB	240241	380637	5.44%	0.412%
25	8.834	988	994	999	rBV	449925	711325	10.16%	0.769%
26	8.887	999	1003	1012	rVB	539416	816483	11.66%	0.883%
27	8.987	1012	1020	1025	rBV	840567	1313342	18.76%	1.421%
28	9.052	1025	1031	1046	rVB2	883909	2071142	29.58%	2.240%
29	9.240	1054	1063	1070	rVB4	45067	120974	1.73%	0.131%
30	9.322	1070	1077	1084	rBV	226733	382459	5.46%	0.414%
31	9.510	1103	1109	1114	rVV	711877	1122255	16.03%	1.214%
32	9.575	1114	1120	1127	rVB	807121	1298253	18.54%	1.404%
33	9.769	1145	1153	1159	rBV	659131	1144188	16.34%	1.238%
34	9.822	1159	1162	1166	rVB	66518	95289	1.36%	0.103%
35	9.881	1166	1172	1176	rBV3	70816	126133	1.80%	0.136%
36	9.934	1176	1181	1194	rVB	340184	634012	9.05%	0.686%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011924.D
Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

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-BP093022.M
Title : SVOA CALIBRATION

37	10.087	1194	1207	1215	rBV2	1032865	2009878	28.70%	2.174%
38	10.157	1215	1219	1223	rBV	103294	166295	2.37%	0.180%
39	10.246	1228	1234	1235	rVV2	66493	128092	1.83%	0.139%
40	10.269	1235	1238	1240	rVV3	86911	137934	1.97%	0.149%
41	10.310	1240	1245	1253	rVV	1039985	1802730	25.75%	1.950%
42	10.387	1253	1258	1267	rVV	105420	206031	2.94%	0.223%
43	10.569	1281	1289	1292	rVV2	63482	125862	1.80%	0.136%
44	10.610	1292	1296	1298	rVV2	70581	121191	1.73%	0.131%
45	10.693	1298	1310	1314	rVV2	1106808	2465352	35.21%	2.667%
46	10.740	1314	1318	1325	rVV4	397555	894713	12.78%	0.968%
47	10.834	1325	1334	1342	rVV	883235	1856375	26.51%	2.008%
48	10.904	1342	1346	1354	rVV	121138	194674	2.78%	0.211%
49	11.057	1367	1372	1375	rBV	36646	54726	0.78%	0.059%
50	11.104	1375	1380	1383	rVV3	36651	73424	1.05%	0.079%
51	11.146	1383	1387	1395	rVV3	77104	166557	2.38%	0.180%
52	11.310	1410	1415	1419	rVV2	183784	346448	4.95%	0.375%
53	11.352	1419	1422	1427	rVV	170556	272277	3.89%	0.295%
54	11.410	1427	1432	1437	rVV3	73142	143084	2.04%	0.155%
55	11.610	1459	1466	1471	rBV	196291	323686	4.62%	0.350%
56	11.669	1471	1476	1483	rBV	307553	610387	8.72%	0.660%
57	11.799	1491	1498	1504	rBV	172992	291582	4.16%	0.315%
58	11.881	1507	1512	1515	rVV	97120	172118	2.46%	0.186%
59	11.916	1515	1518	1527	rVV6	116561	238896	3.41%	0.258%
60	12.010	1528	1534	1539	rVV3	121239	232883	3.33%	0.252%
61	12.081	1539	1546	1553	rVB2	285831	493235	7.04%	0.534%
62	12.246	1569	1574	1578	rVV2	68161	105892	1.51%	0.115%
63	12.328	1582	1588	1598	rVB3	100805	209810	3.00%	0.227%
64	12.416	1598	1603	1615	rVB2	57568	109018	1.56%	0.118%
65	12.510	1615	1619	1623	rBV	107717	173212	2.47%	0.187%
66	12.569	1623	1629	1636	rVV	512092	881071	12.58%	0.953%
67	12.651	1637	1643	1652	rVB6	81078	240561	3.44%	0.260%
68	12.740	1652	1658	1663	rBV4	42185	86262	1.23%	0.093%
69	12.869	1675	1680	1686	rVB3	43667	87357	1.25%	0.094%
70	12.993	1696	1701	1707	rVB	62185	91941	1.31%	0.099%
71	13.263	1739	1747	1753	rBV4	40262	91627	1.31%	0.099%
72	13.404	1766	1771	1775	rBV2	37123	56544	0.81%	0.061%
73	13.498	1782	1787	1794	rVB	193425	306221	4.37%	0.331%
74	13.575	1794	1800	1809	rVB7	20739	55749	0.80%	0.060%
75	13.663	1809	1815	1817	rBV2	84886	147224	2.10%	0.159%
76	13.698	1817	1821	1830	rVV4	161202	350744	5.01%	0.379%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011924.D
Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

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-BP093022.M
Title : SVOA CALIBRATION

77	13.846	1841	1846	1851	rBV	111060	201606	2.88%	0.218%
78	13.898	1851	1855	1857	rVV2	79531	115884	1.65%	0.125%
79	13.940	1857	1862	1870	rVB	1950060	2734243	39.05%	2.958%
80	14.222	1904	1910	1921	rBV	2409650	3473523	49.61%	3.757%
81	14.345	1928	1931	1939	rVB2	51413	73552	1.05%	0.080%
82	14.528	1953	1962	1977	rBV	1634295	2540461	36.28%	2.748%
83	14.740	1993	1998	2010	rVB	101941	229984	3.28%	0.249%
84	14.851	2012	2017	2025	rBV2	58171	89837	1.28%	0.097%
85	15.157	2060	2069	2078	rVB3	40772	86844	1.24%	0.094%
86	15.522	2121	2131	2145	rVB	3179729	4662093	66.58%	5.043%
87	15.640	2145	2151	2160	rBV	856594	1182247	16.88%	1.279%
88	16.034	2213	2218	2231	rBV7	20846	58416	0.83%	0.063%
89	17.287	2424	2431	2441	rBV	2097710	2942787	42.03%	3.183%
90	17.387	2441	2448	2458	rBV	3771275	5323641	76.03%	5.759%
91	17.951	2540	2544	2549	rBV	86175	103000	1.47%	0.111%
92	18.169	2576	2581	2589	rBV2	230630	365919	5.23%	0.396%
93	19.198	2752	2756	2762	rBV4	54833	95062	1.36%	0.103%
94	19.263	2764	2767	2774	rVV	52401	78417	1.12%	0.085%
95	19.398	2787	2790	2796	rBV	95926	138932	1.98%	0.150%
96	19.622	2822	2828	2841	rBV	5046767	6656863	95.07%	7.201%
97	21.080	3072	3076	3084	rBV2	288065	512889	7.32%	0.555%
98	21.369	3121	3125	3133	rBV	2697164	3519920	50.27%	3.808%
99	23.645	3505	3512	3524	rVB2	3367266	7002097	100.00%	7.574%
100	23.804	3533	3539	3555	rVB2	1790313	3667246	52.37%	3.967%

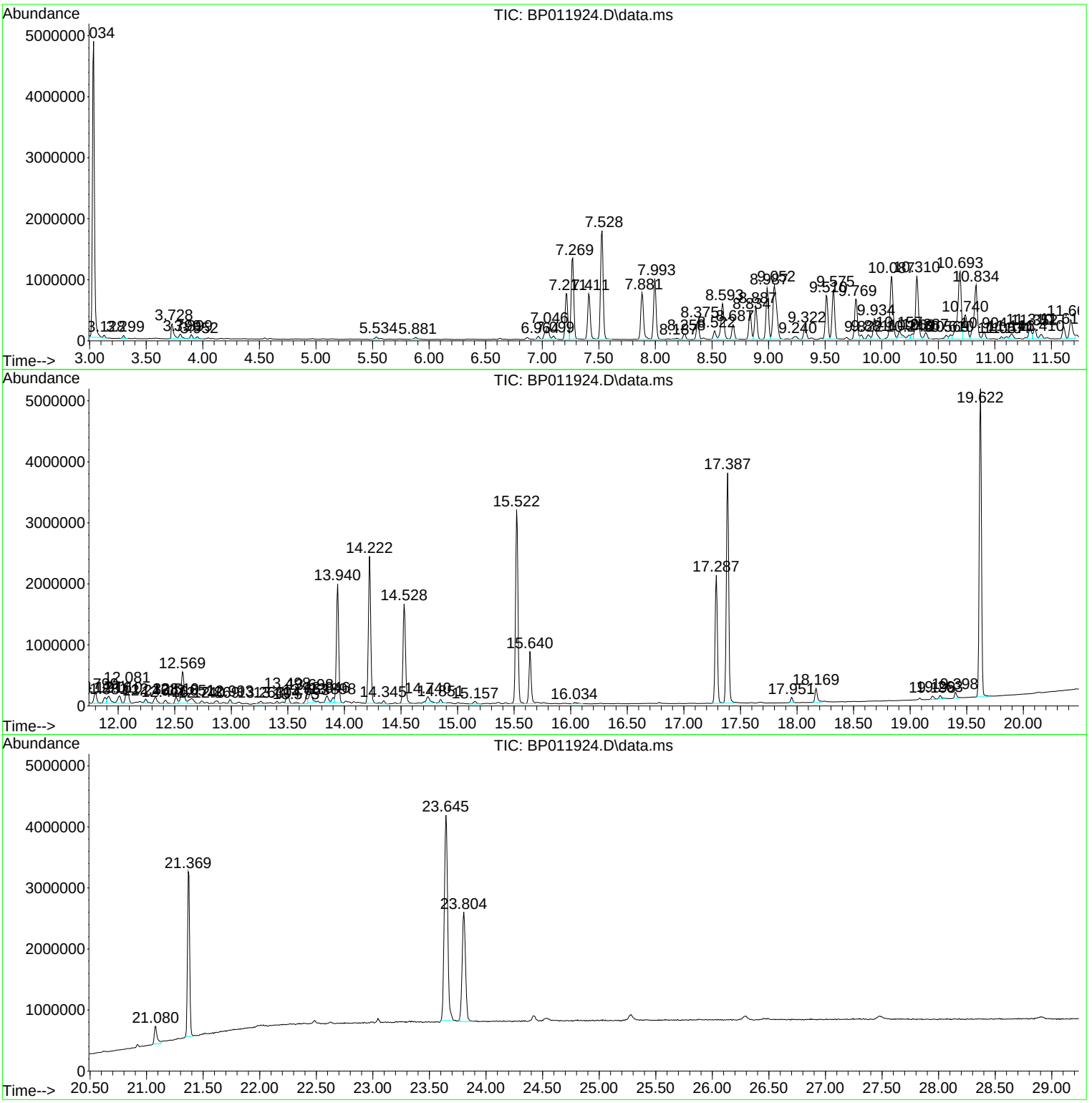
Sum of corrected areas: 92445941

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011924.D
Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011924.D
Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

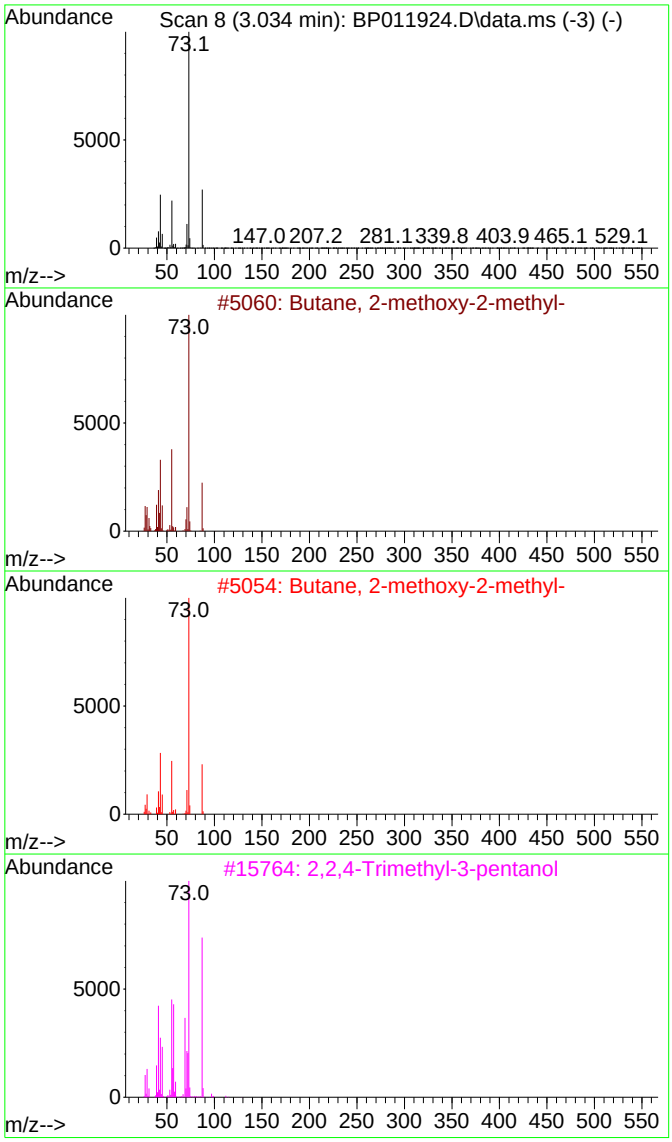
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

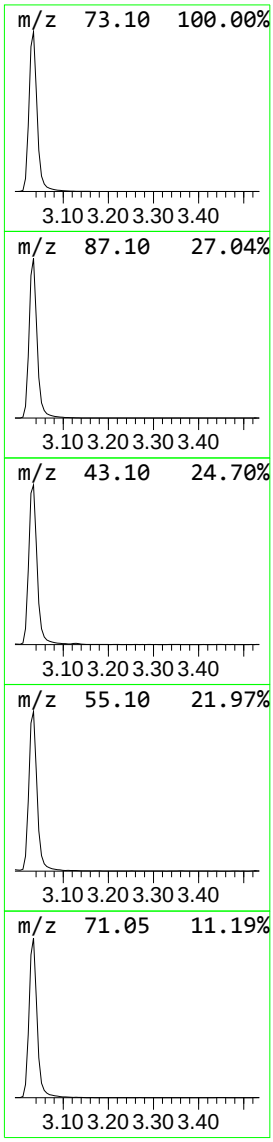
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.034	91.89 ng/ul	6046310	1,4-Dichlorobenzene-d4	7.881

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5	N-Ethylformamide	73	C3H7NO	000627-45-2	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011924.D
Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

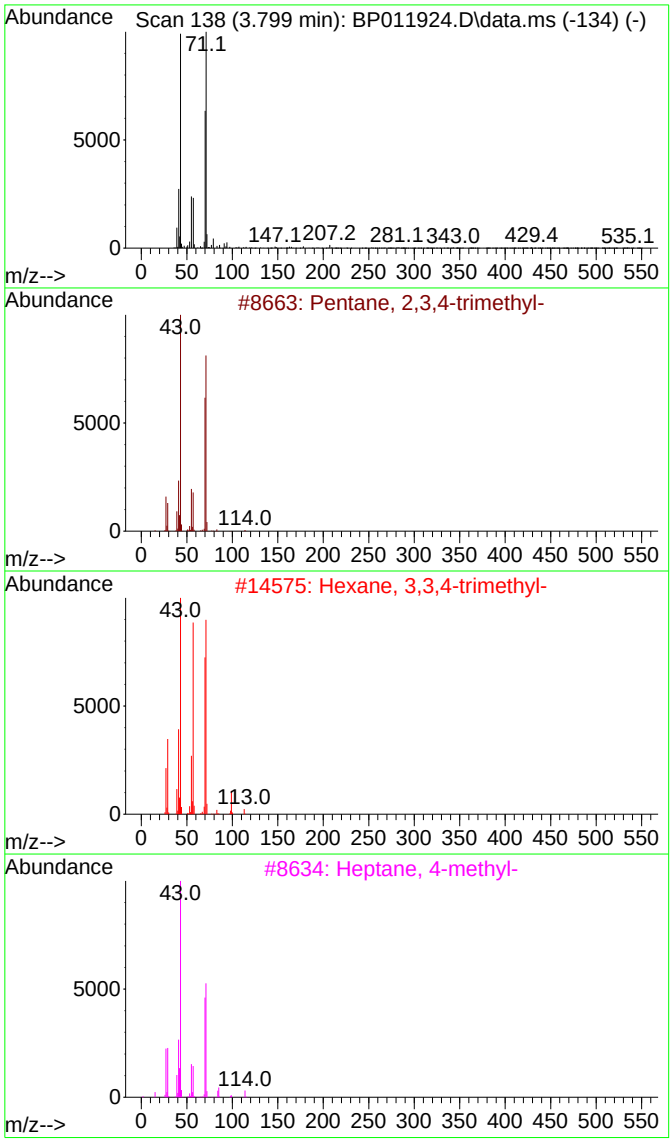
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

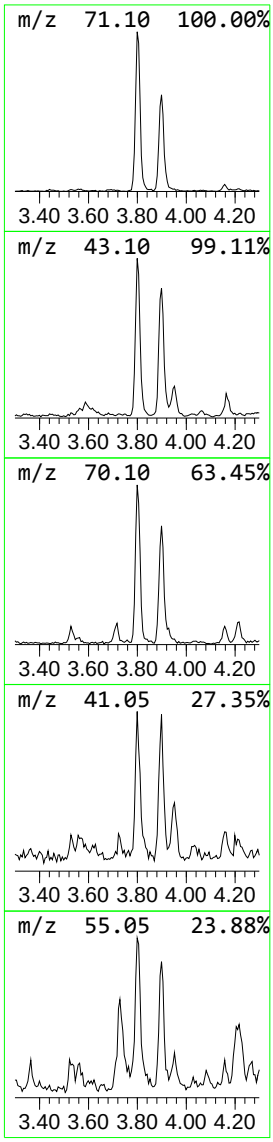
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.799	2.06 ng/ul	135315	1,4-Dichlorobenzene-d4	7.881

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	78
2	Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	78
3	Heptane, 4-methyl-	114	C8H18	000589-53-7	78
4	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	78
5	Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011924.D
Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

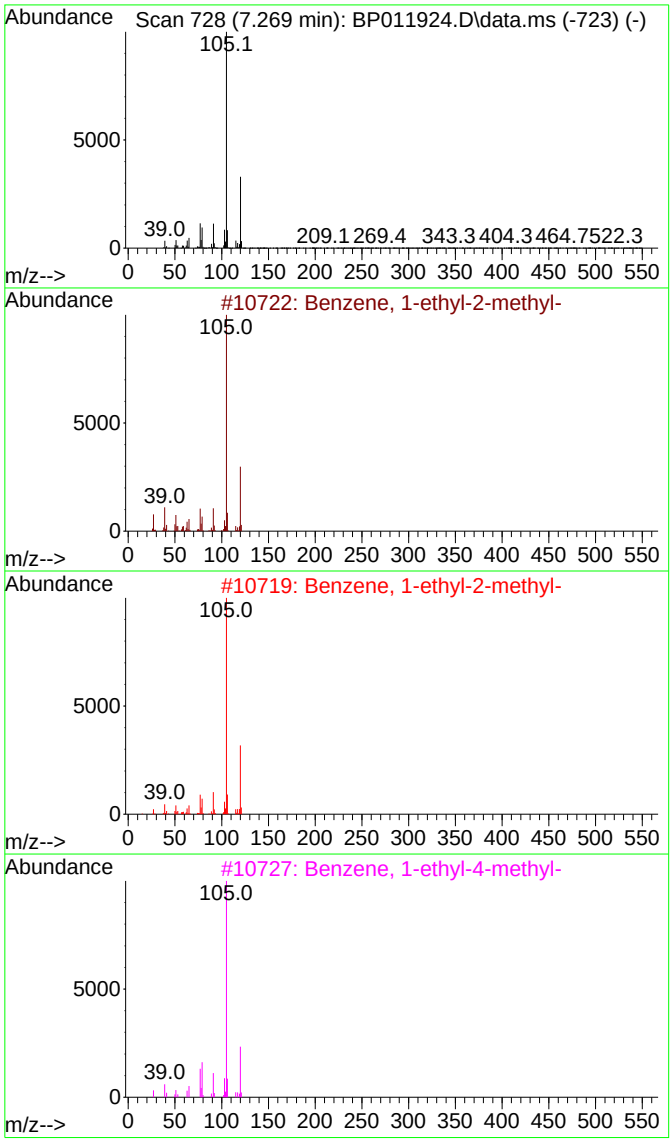
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

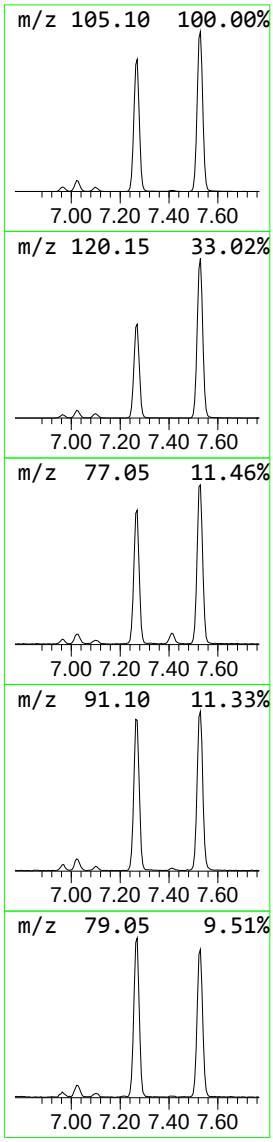
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.269	31.85 ng/ul	2096030	1,4-Dichlorobenzene-d4	7.881

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-4-methyl-	120	C9H12	000622-96-8	94
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011924.D
Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

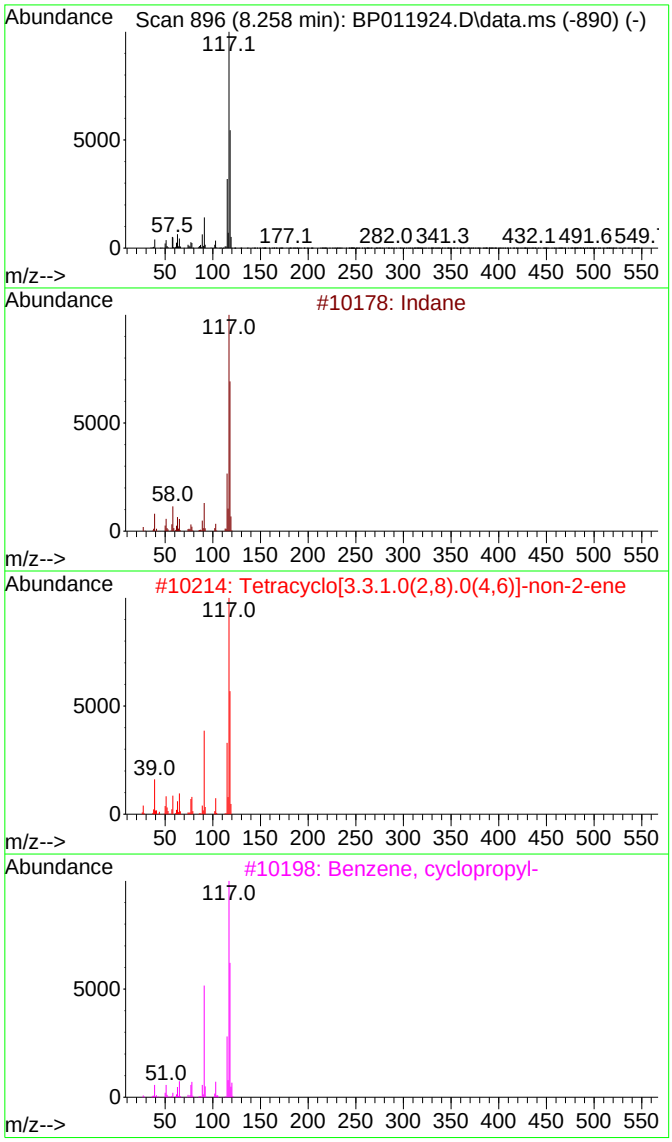
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

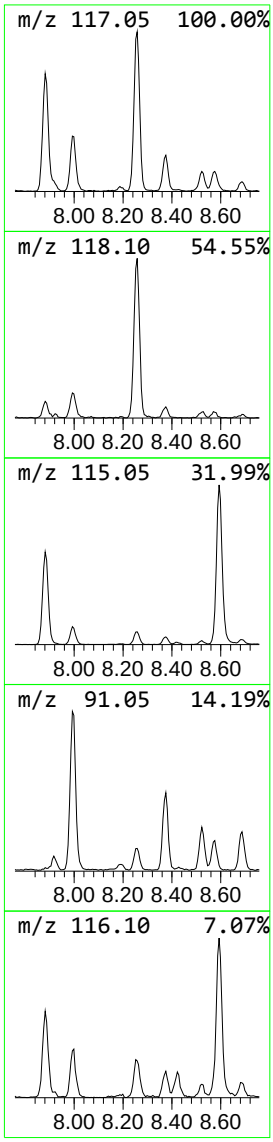
****

Peak Number 6 Indane Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.258	2.43 ng/ul	159797	1,4-Dichlorobenzene-d4	7.881

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	91
2	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	87
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	87
4	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	72
5	Indane	118	C9H10	000496-11-7	60





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011924.D
Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

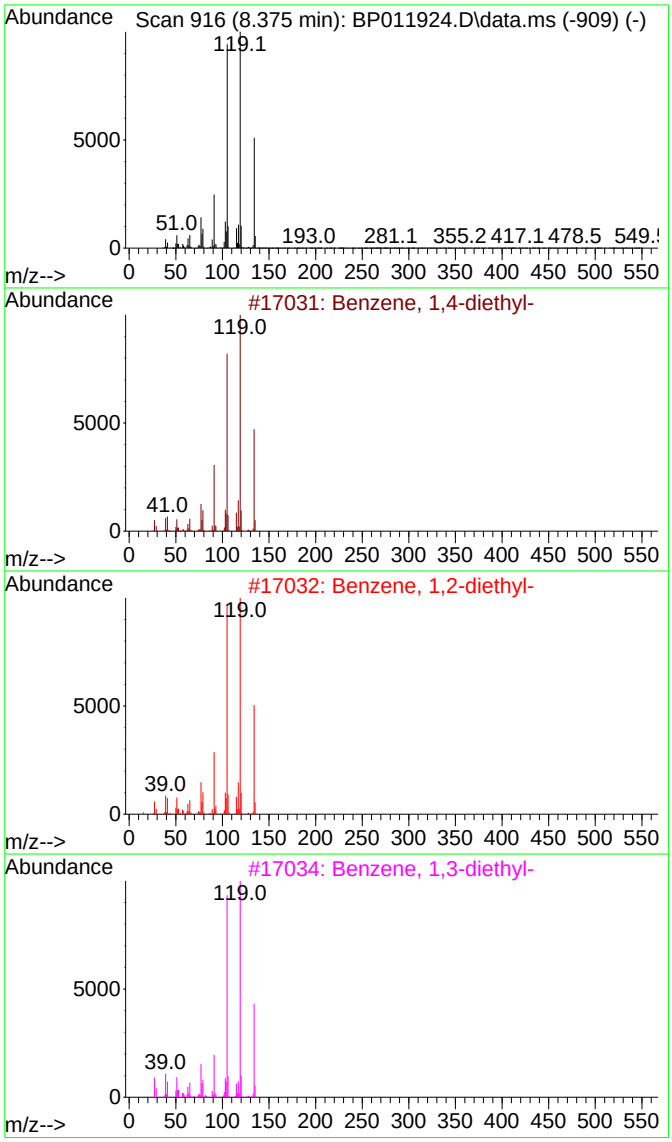
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

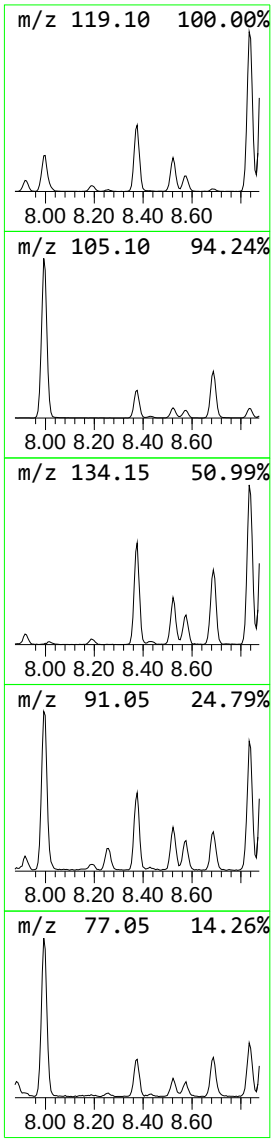
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1,4-diethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.375	7.18 ng/ul	472711	1,4-Dichlorobenzene-d4	7.881

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	97
2	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
4	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
5	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94





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Data File : BP011924.D
Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

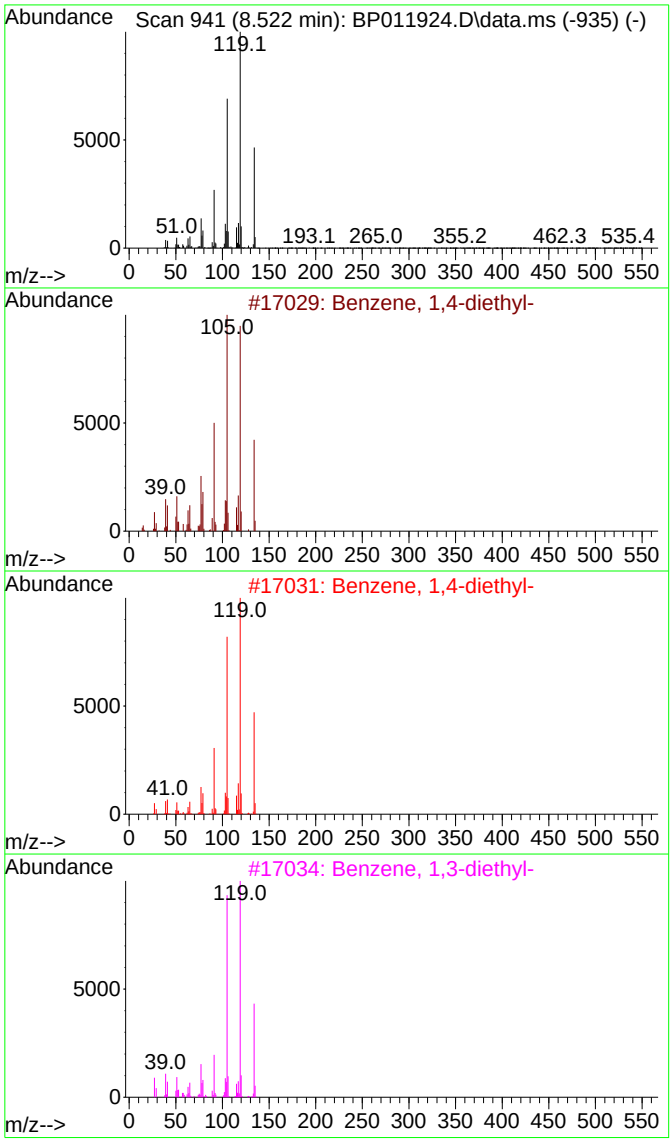
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

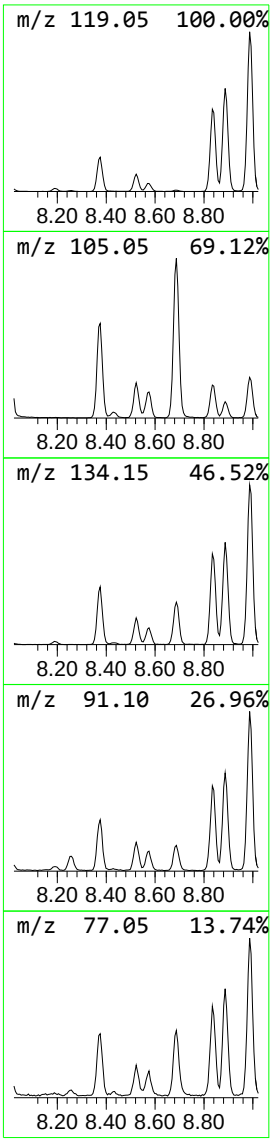
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1,3-diethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.522	3.51 ng/ul	230642	1,4-Dichlorobenzene-d4	7.881

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	97
2	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	96
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
4	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95
5	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	95





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Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

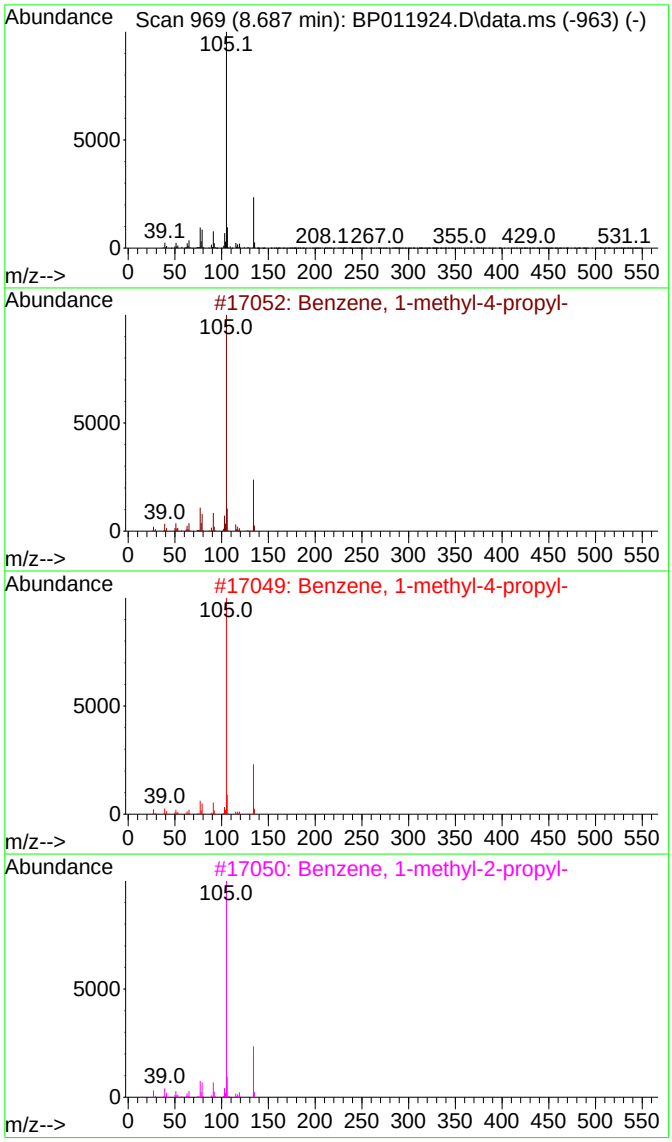
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

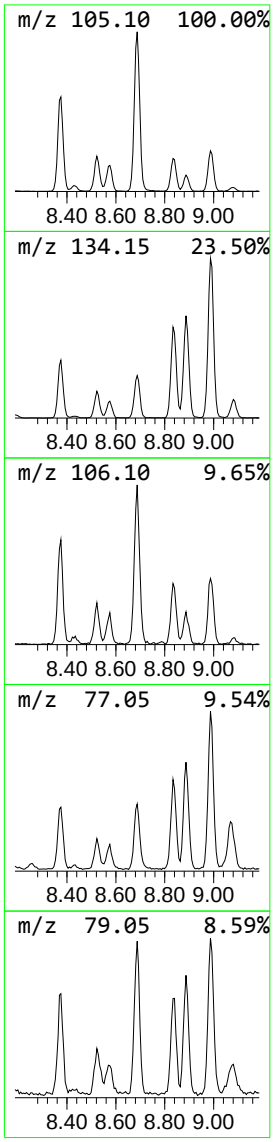
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1-methyl-4-propyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.687	5.78 ng/ul	380637	1,4-Dichlorobenzene-d4	7.881

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	95
2	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	94
3	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	94
4	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	91
5	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91





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Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

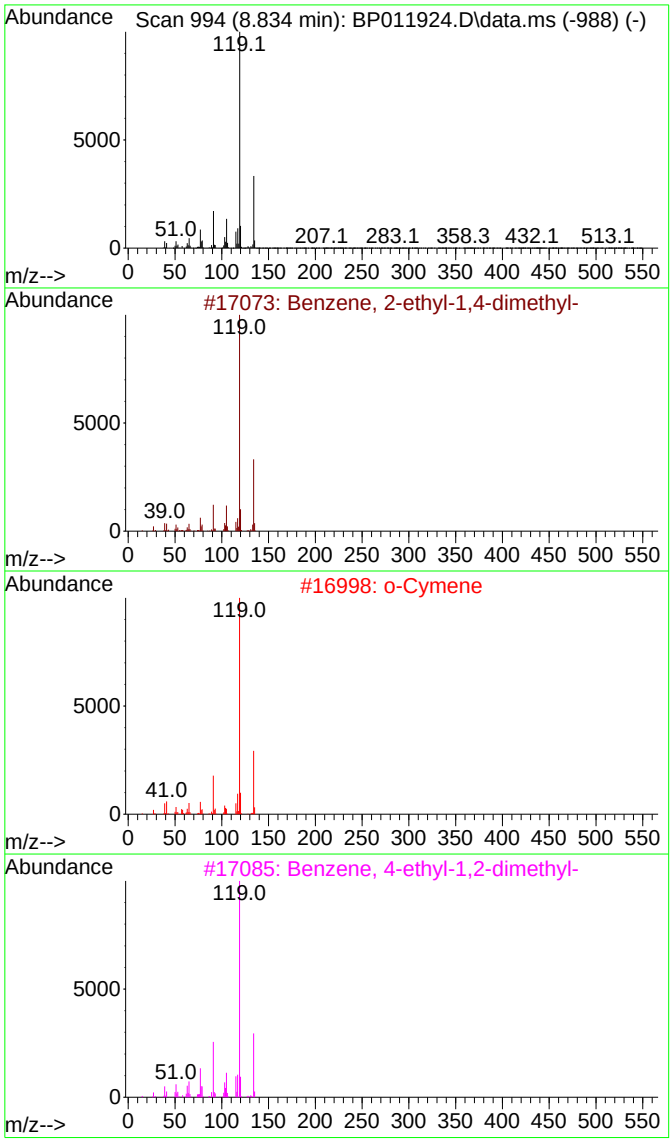
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

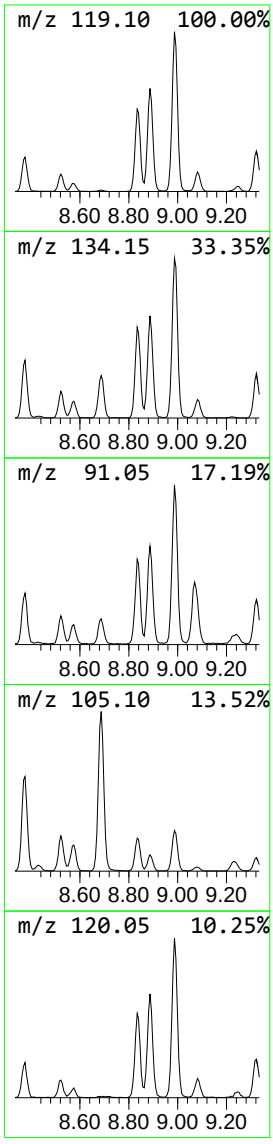
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.834	10.81 ng/ul	711325	1,4-Dichlorobenzene-d4	7.881

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2	o-Cymene	134	C10H14	000527-84-4	95
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
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Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

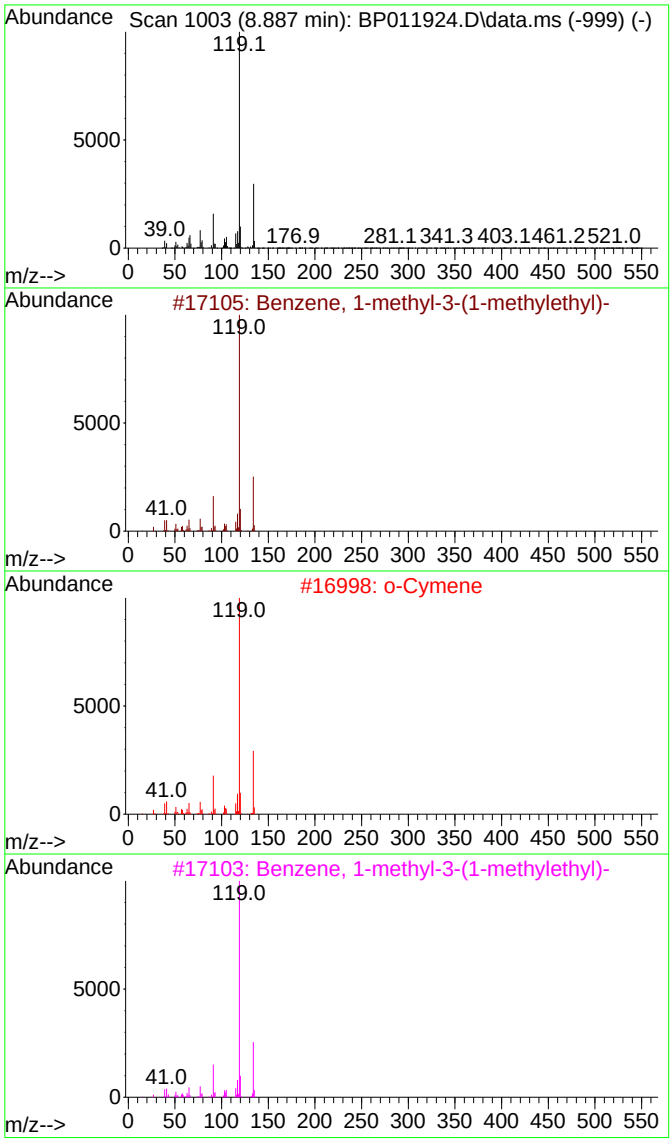
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

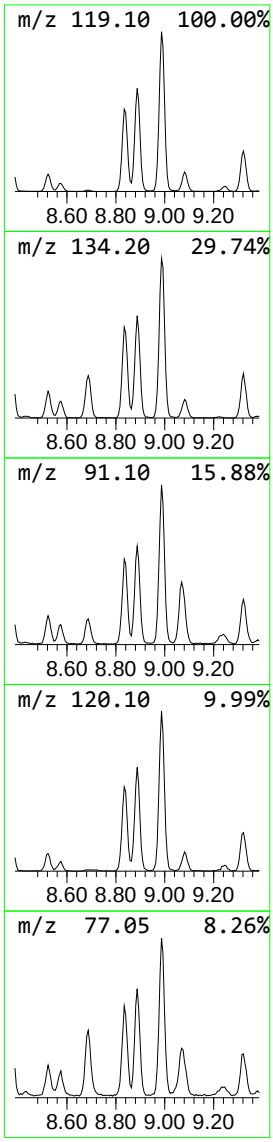
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzene, 1-methyl-3-(1-meth... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.887	12.41 ng/ul	816483	1,4-Dichlorobenzene-d4	7.881

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	97
2	o-Cymene	134	C10H14	000527-84-4	97
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	o-Cymene	134	C10H14	000527-84-4	95





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Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

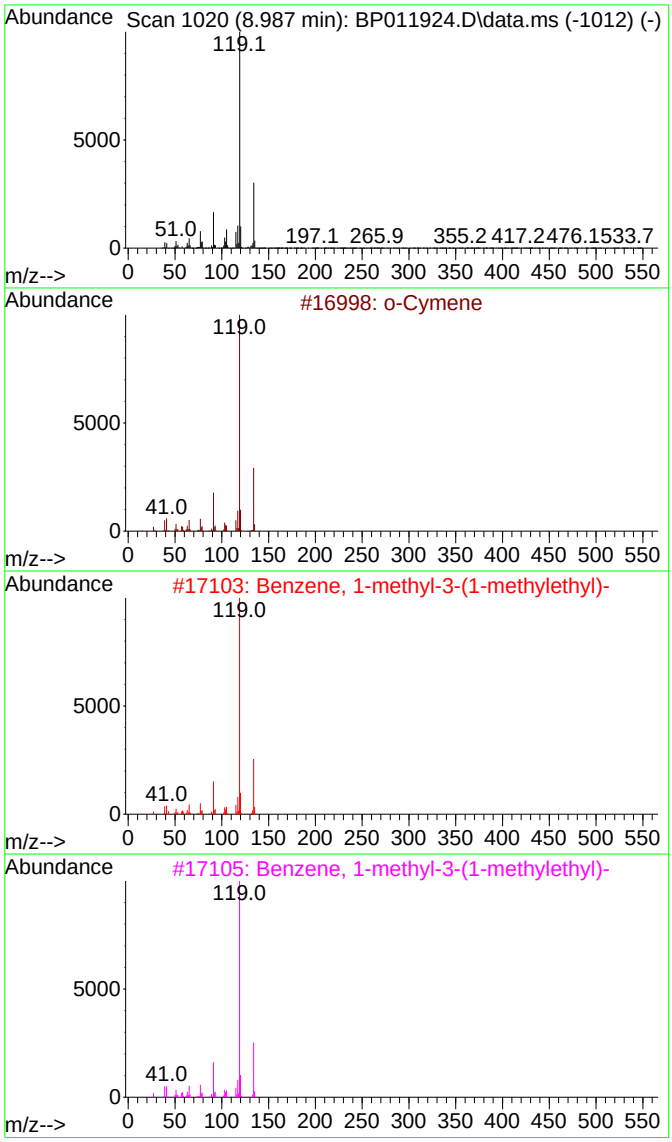
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

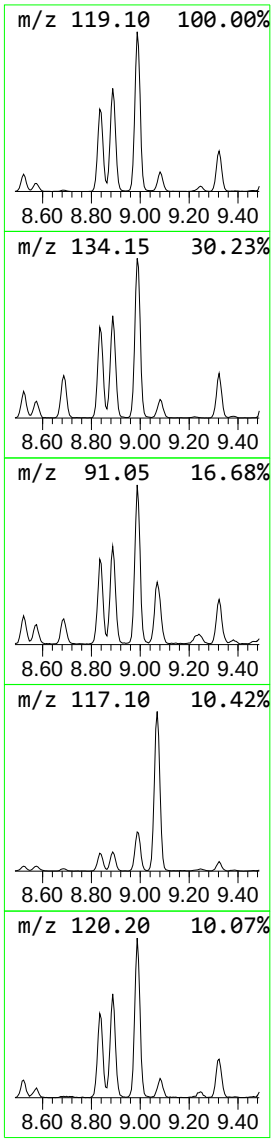
****

Peak Number 12 o-Cymene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.987	19.96 ng/ul	1313340	1,4-Dichlorobenzene-d4	7.881

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	97
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94





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Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

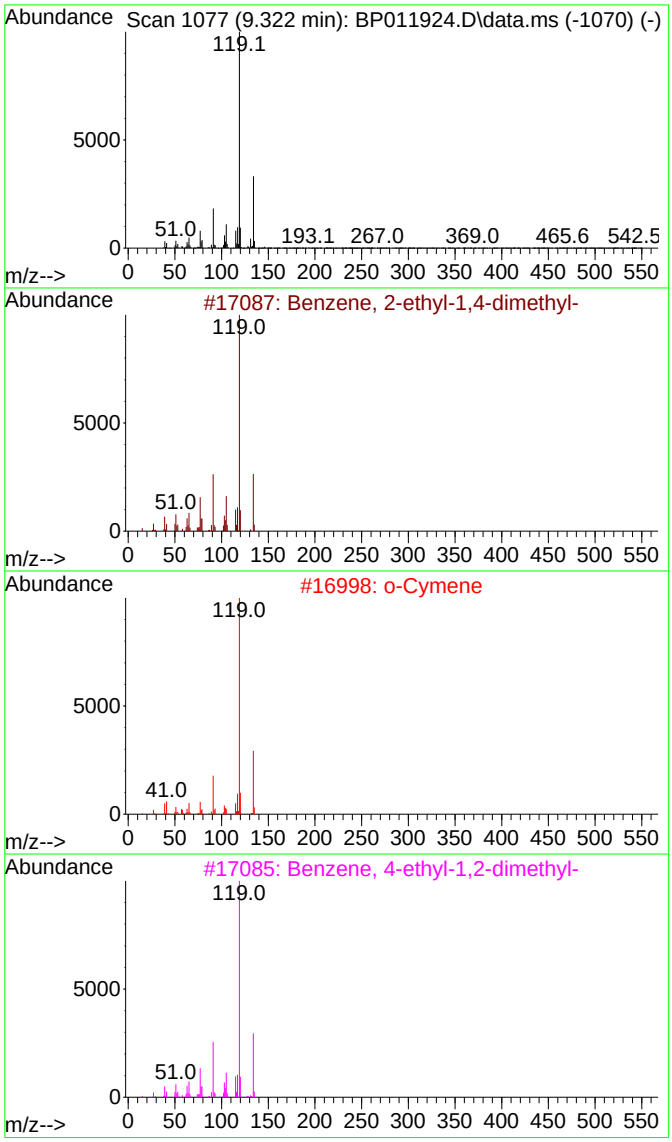
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

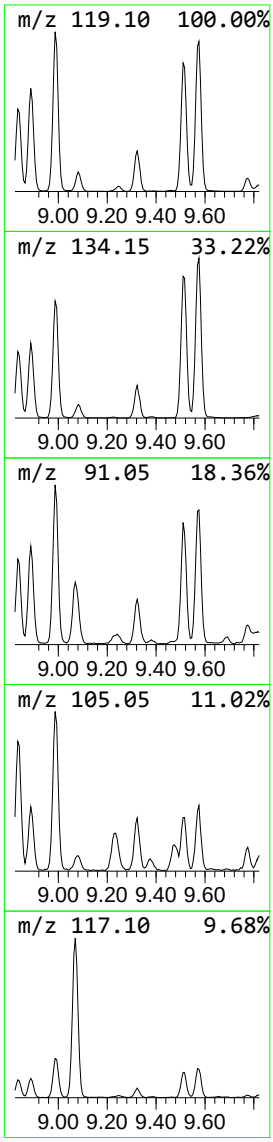
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.322	3.10 ng/ul	382459	Naphthalene-d8	10.693

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2	o-Cymene	134	C10H14	000527-84-4	95
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94





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Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

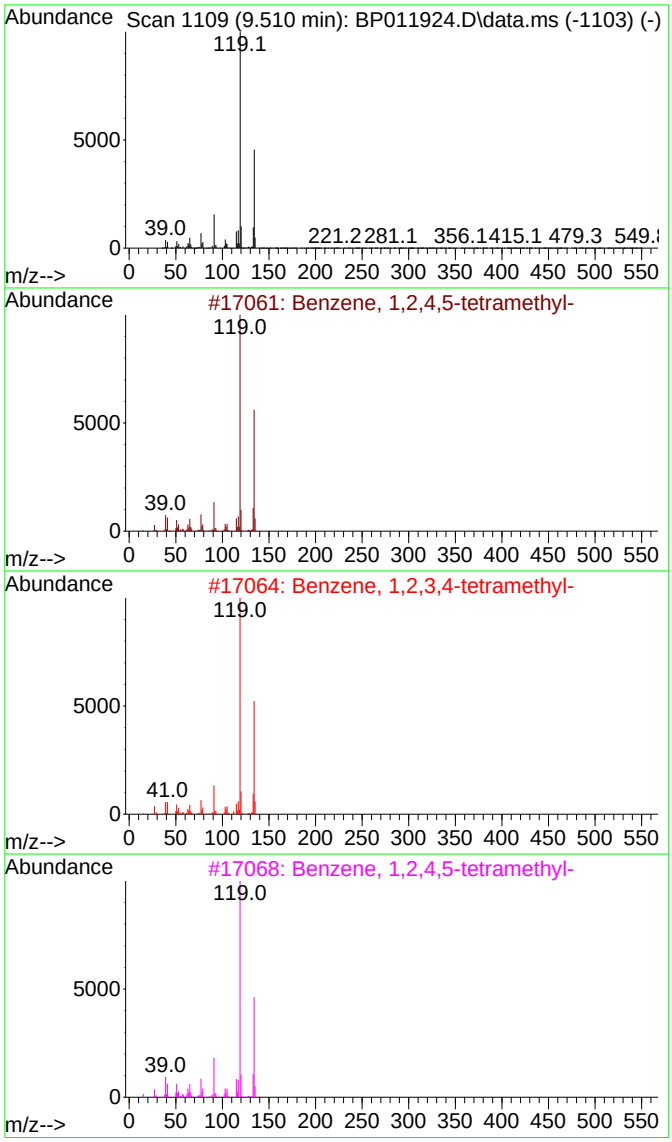
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

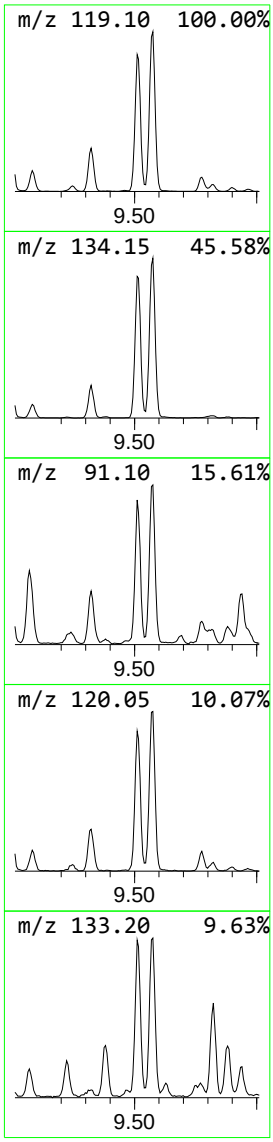
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.510	9.10 ng/ul	1122260	Naphthalene-d8	10.693

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
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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Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

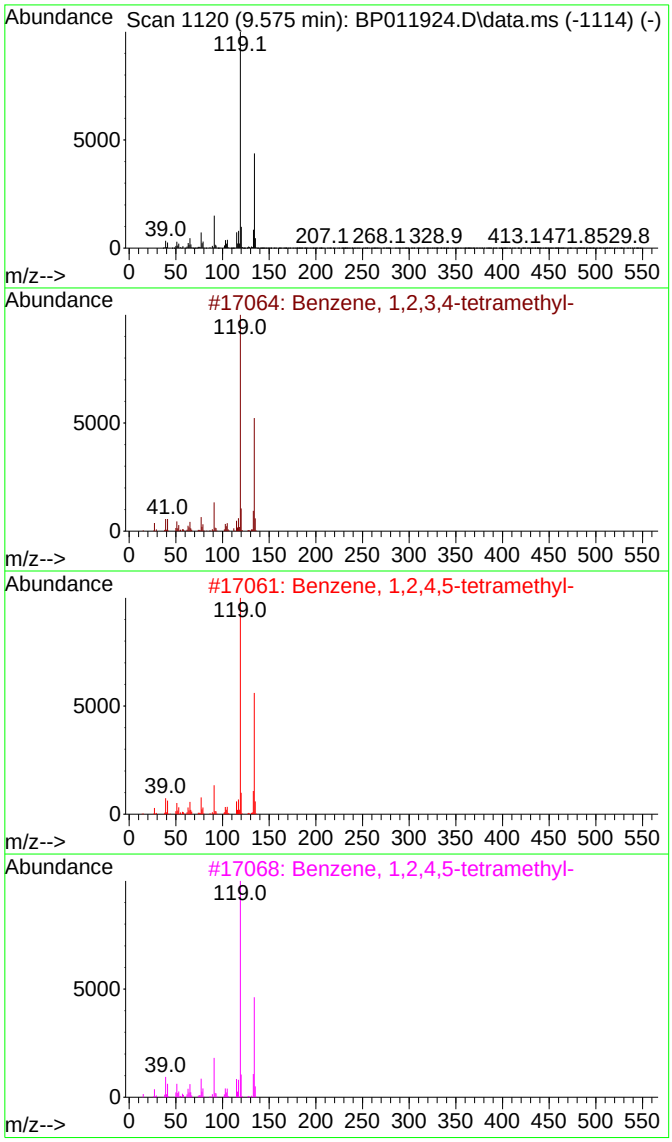
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

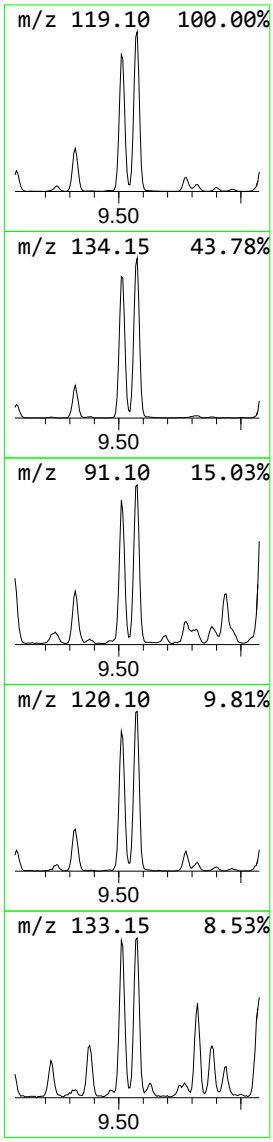
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.575	10.53 ng/ul	1298250	Naphthalene-d8	10.693

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96





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Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

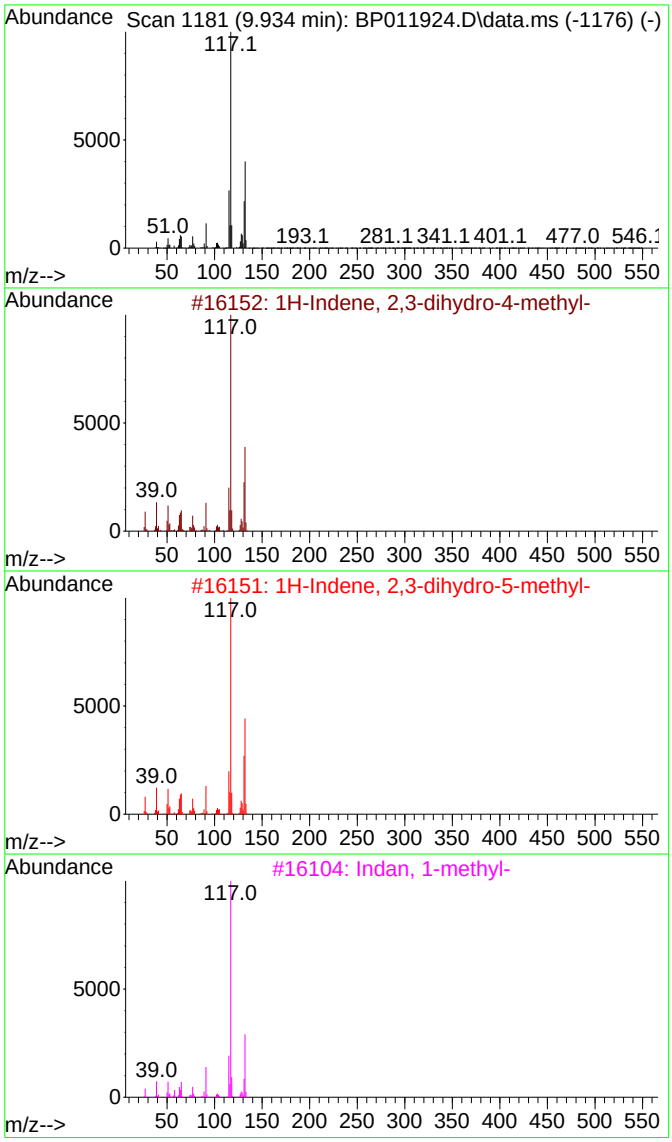
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

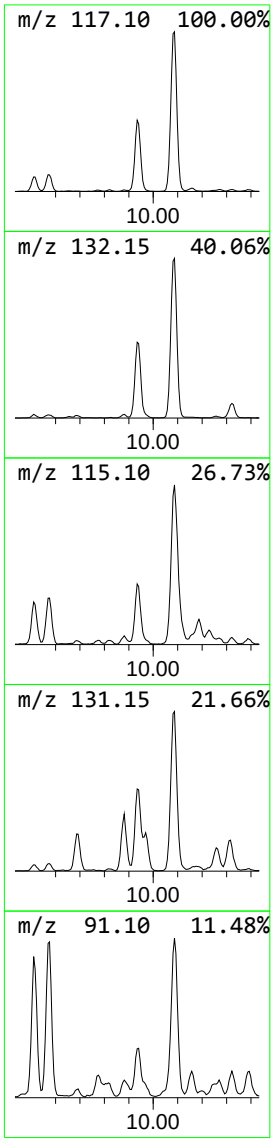
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.934	5.14 ng/ul	634012	Naphthalene-d8	10.693

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
3	Indan, 1-methyl-	132	C10H12	000767-58-8	87
4	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	86
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	81





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Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

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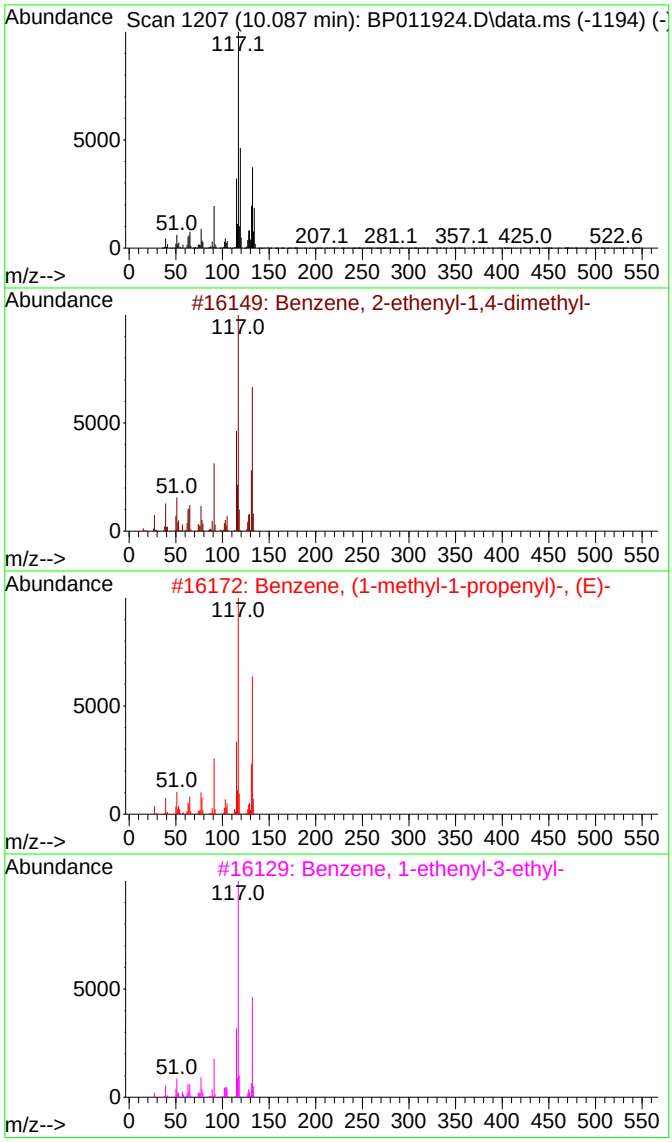
Quant Title : SVOA CALIBRATION

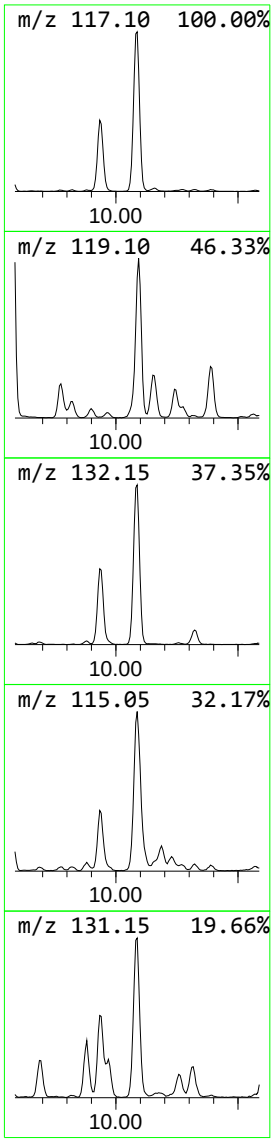
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.087	16.31 ng/ul	2009880	Naphthalene-d8	10.693

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	68
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	64
4	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	64
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	62





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Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

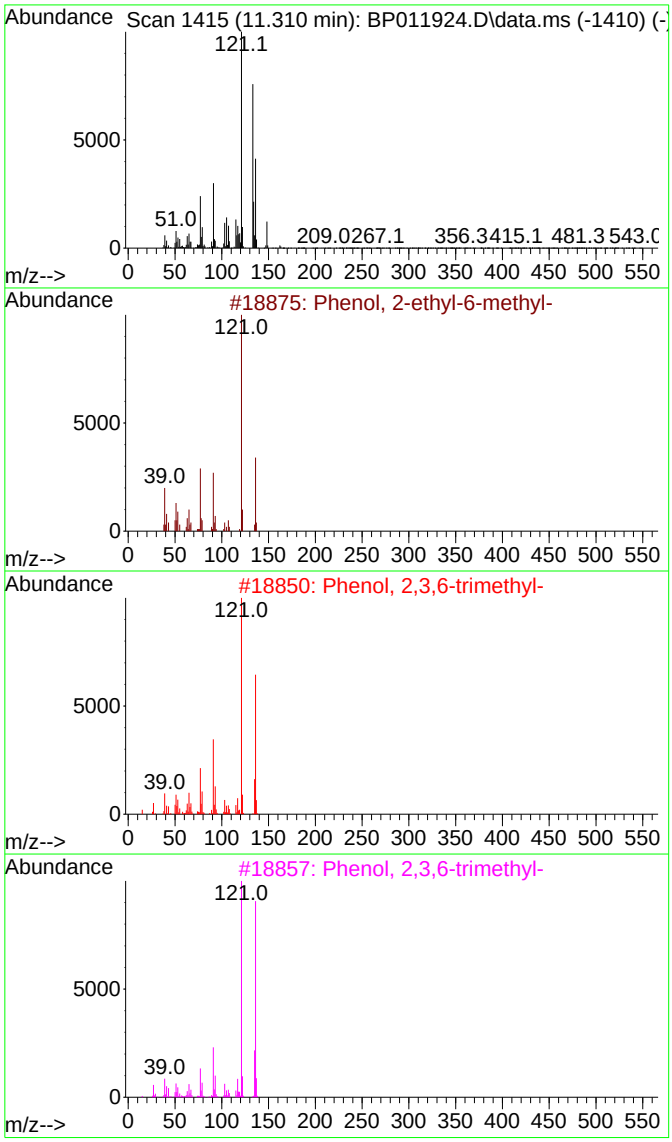
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

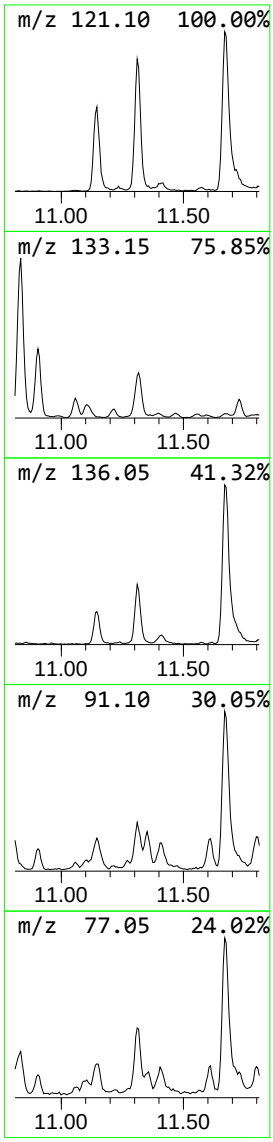
****

Peak Number 18 unknown-01 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.310	2.81 ng/ul	346448	Naphthalene-d8	10.693

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-ethyl-6-methyl-	136	C9H12O	001687-64-5	49
2	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	49
3	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	49
4	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	46
5	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	46





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Data File : BP011924.D
Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

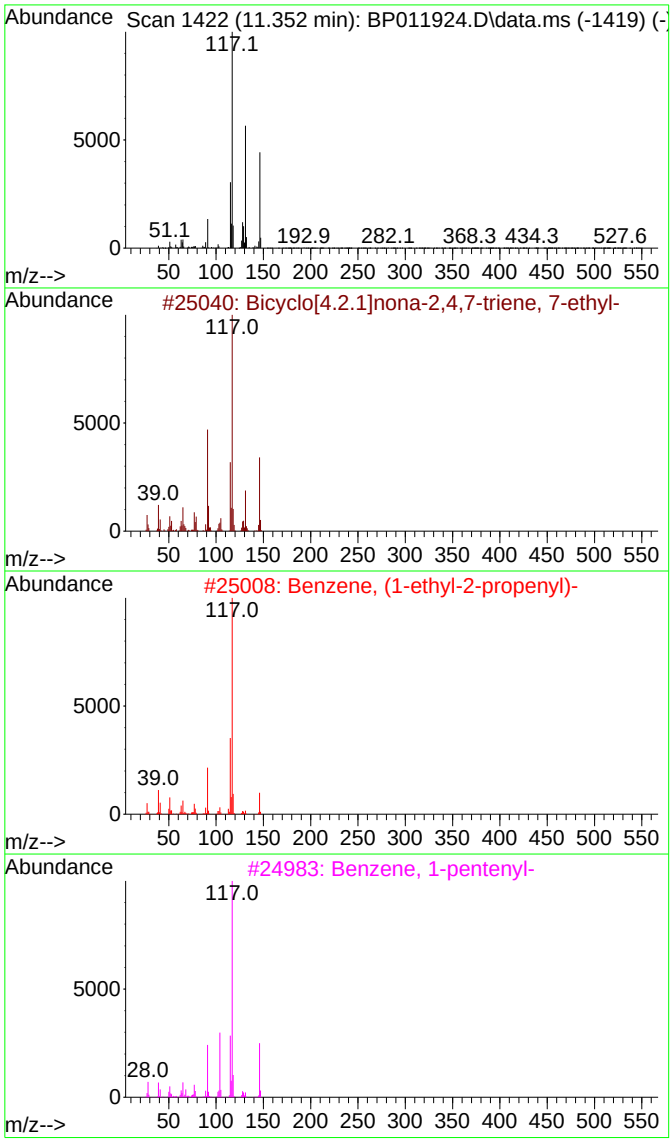
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

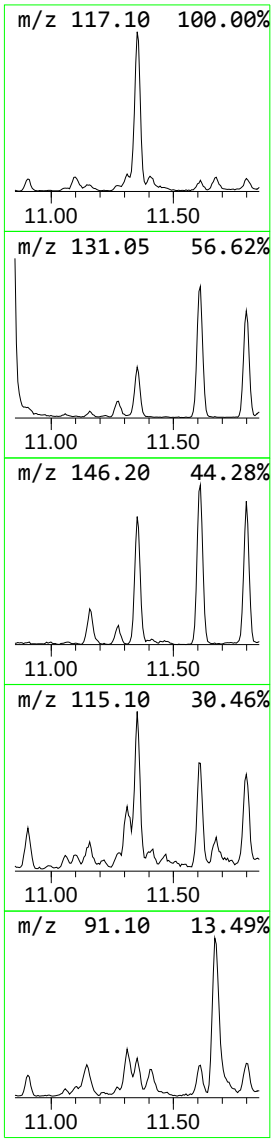
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Bicyclo[4.2.1]nona-2,4,7-tr... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.352	2.21 ng/ul	272277	Naphthalene-d8	10.693

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[4.2.1]nona-2,4,7-triene,...	146	C11H14	1000164-42-6	72
2	Benzene, (1-ethyl-2-propenyl)-	146	C11H14	019947-22-9	59
3	Benzene, 1-pentenyl-	146	C11H14	000826-18-6	59
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	58
5	1H-Indene, 1-ethyl-2,3-dihydro-	146	C11H14	004830-99-3	58





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Data File : BP011924.D
Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

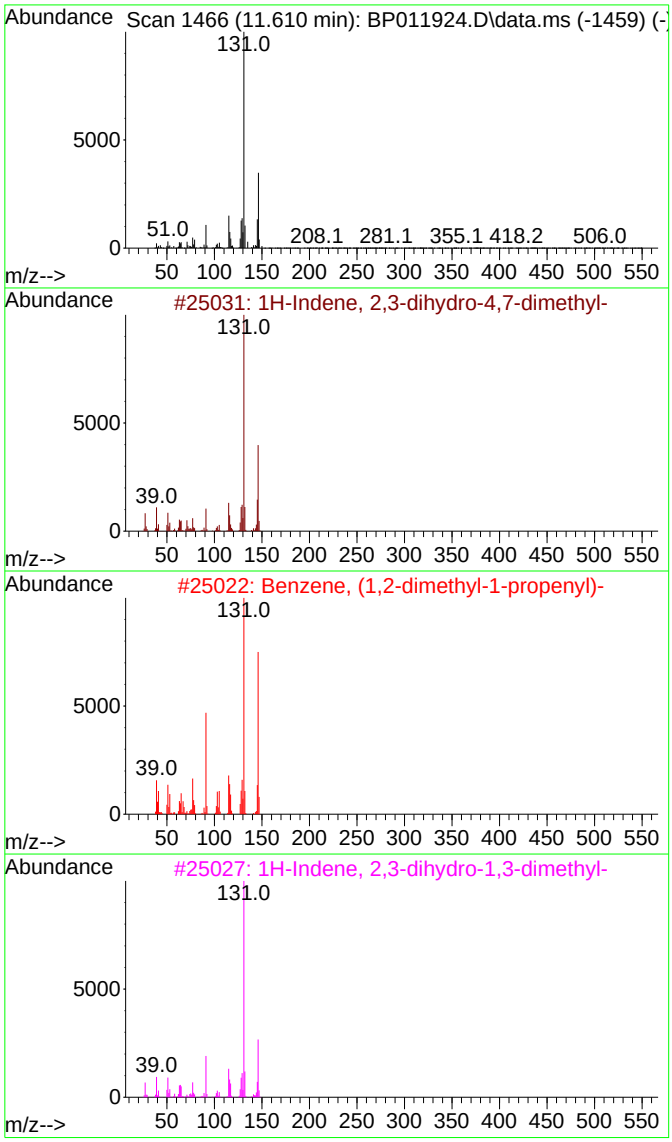
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

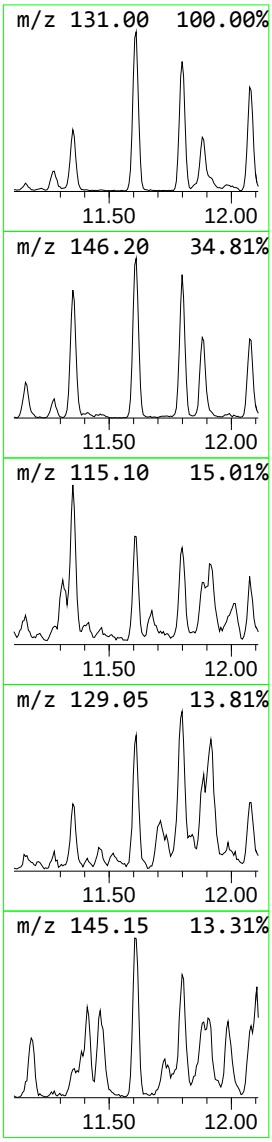
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.610	2.63 ng/ul	323686	Naphthalene-d8	10.693

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
2	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	94
3	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	94
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93
5	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	93





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Data File : BP011924.D
Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

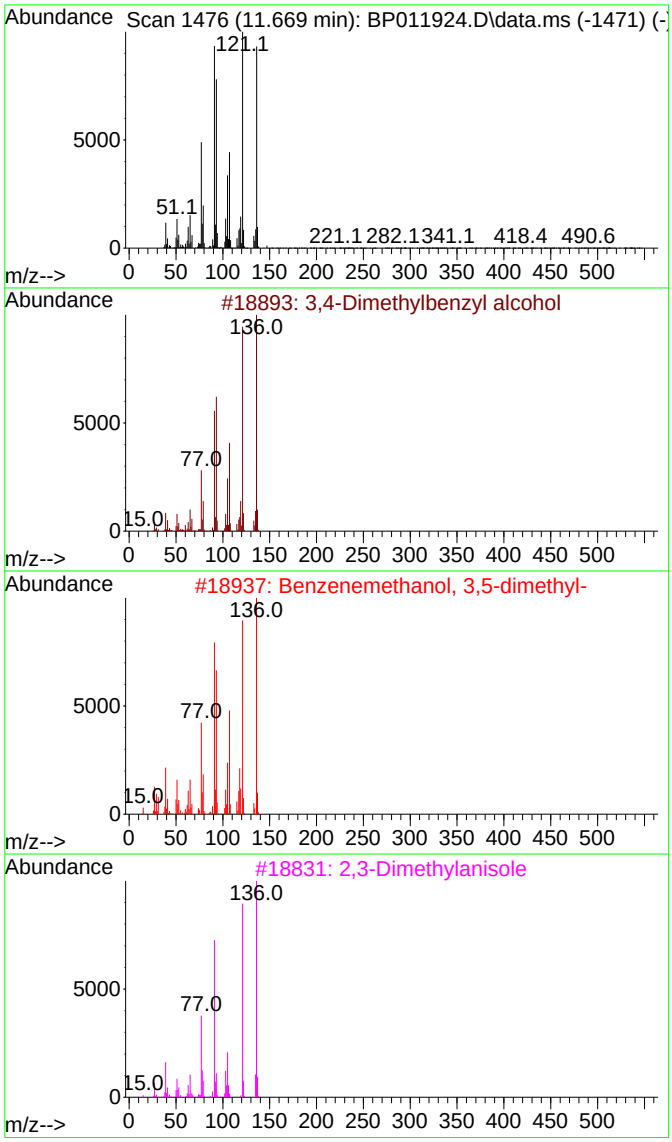
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

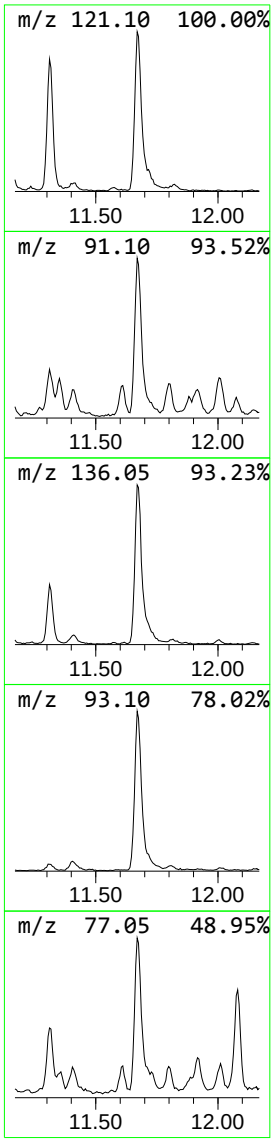
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 3,4-Dimethylbenzyl alcohol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.669	4.95 ng/ul	610387	Naphthalene-d8	10.693

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,4-Dimethylbenzyl alcohol	136	C9H12O	006966-10-5	97
2	Benzenemethanol, 3,5-dimethyl-	136	C9H12O	027129-87-9	95
3	2,3-Dimethylanisole	136	C9H12O	002944-49-2	87
4	Cyclohexene, 1-methyl-4-(1-methy...	136	C10H16	000586-62-9	83
5	1,3-Cyclohexadiene, 1-methyl-4-(...	136	C10H16	000099-86-5	81





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Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

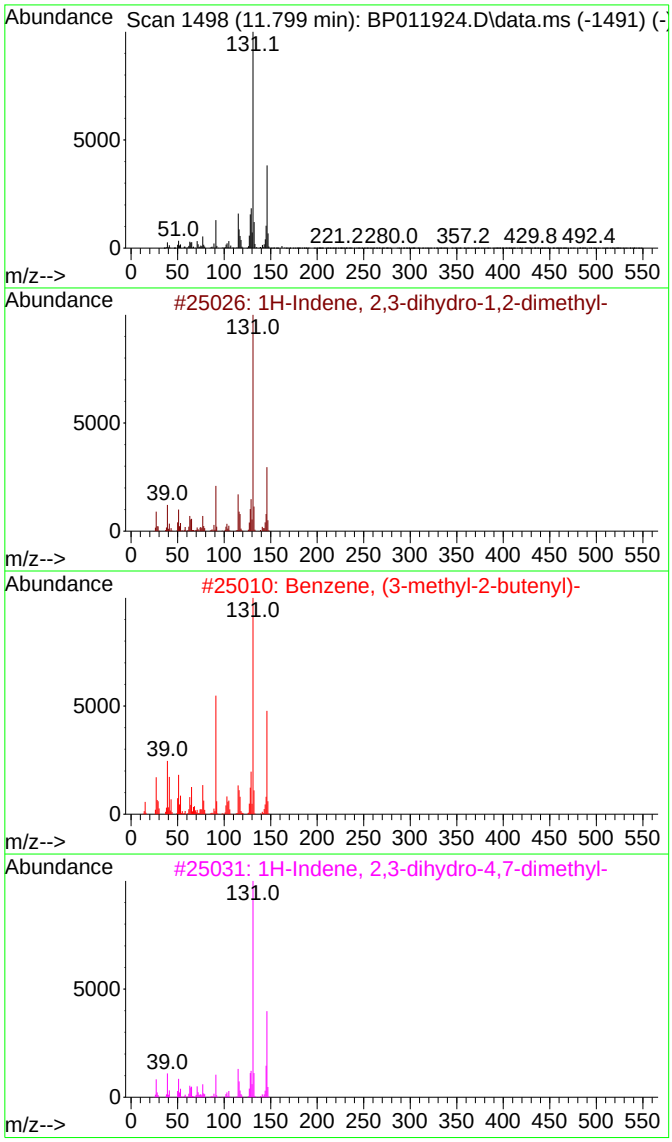
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

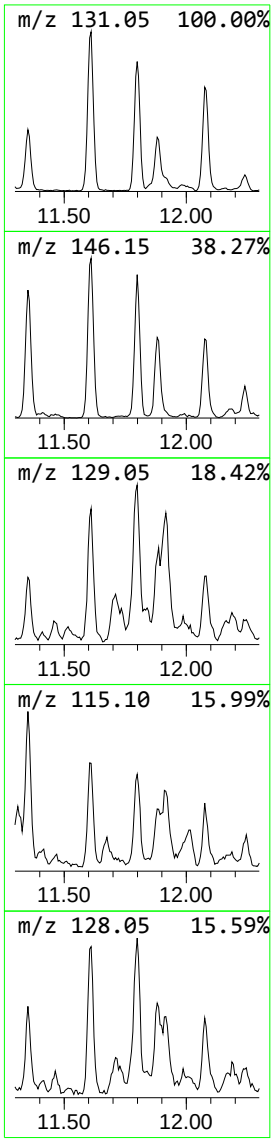
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.799	2.37 ng/ul	291582	Naphthalene-d8	10.693

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	95
2	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	94
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
4	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	94
5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93





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Data File : BP011924.D
Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

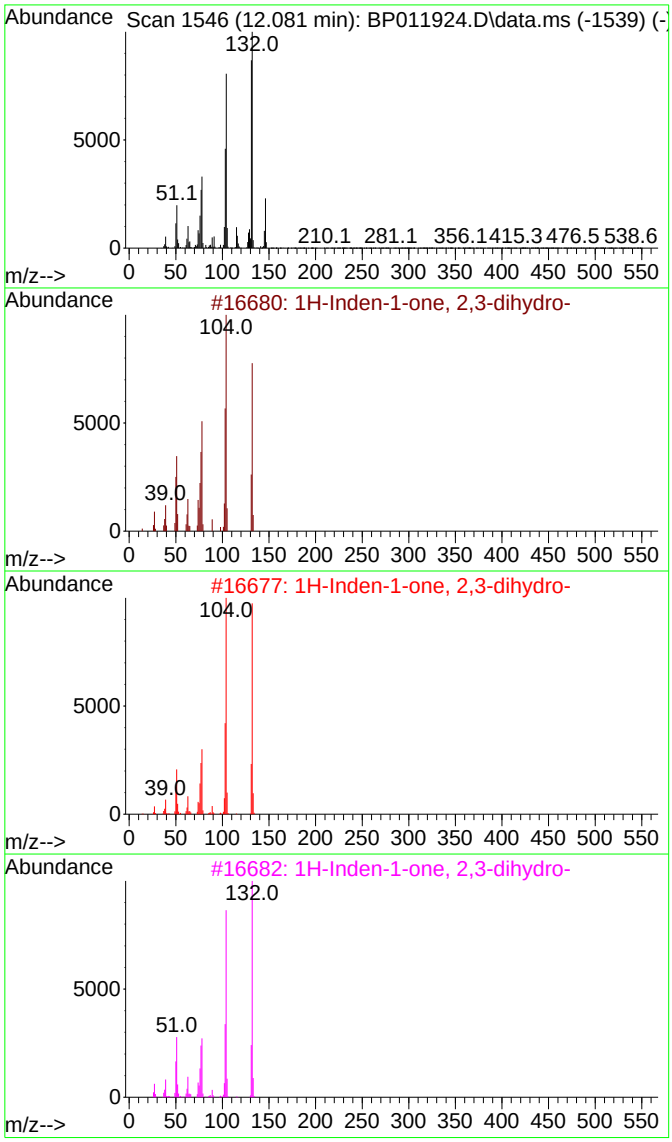
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

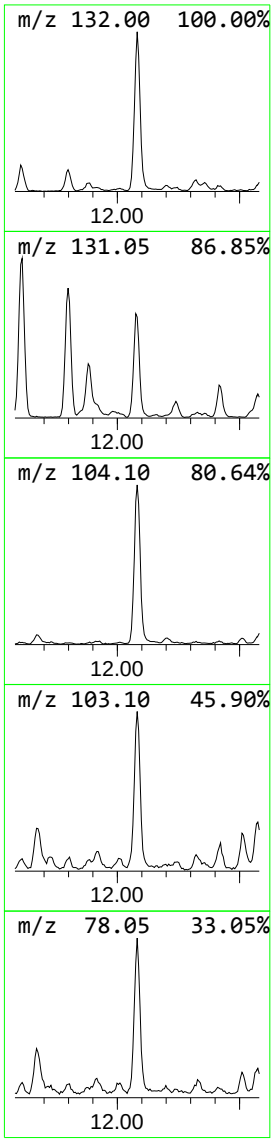
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 1H-Inden-1-one, 2,3-dihydro- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.081	4.00 ng/ul	493235	Naphthalene-d8	10.693

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	89
2	1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	70
3	1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	70
4	Benzene, (2-methyl-1-methylenepr...	146	C11H14	017498-71-4	50
5	1H-Indol-3-amine	132	C8H8N2	007250-19-3	47





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Data File : BP011924.D
Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

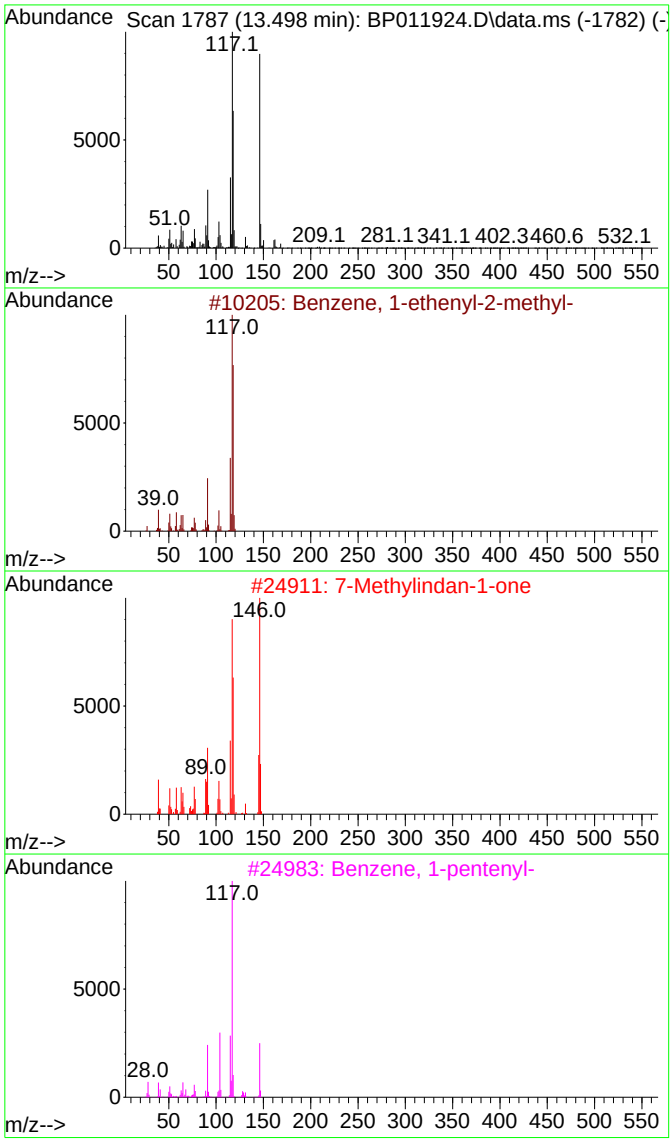
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

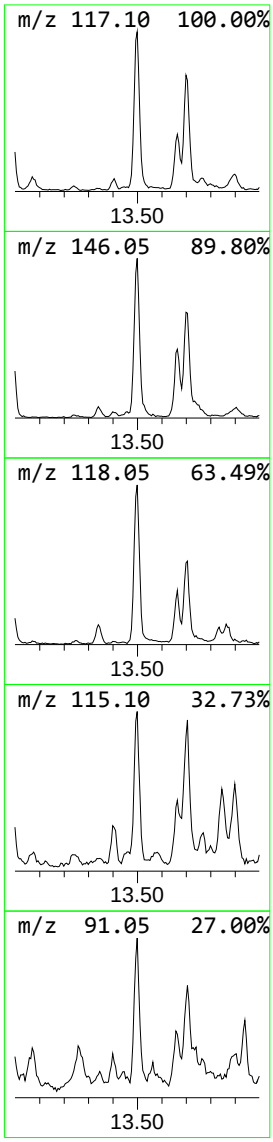
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Benzene, 1-ethenyl-2-methyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.499	2.41 ng/ul	306221	Acenaphthene-d10	14.528

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	86
2	7-Methylindan-1-one	146	C10H10O	039627-61-7	68
3	Benzene, 1-pentenyl-	146	C11H14	000826-18-6	68
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	64





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Data File : BP011924.D
Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

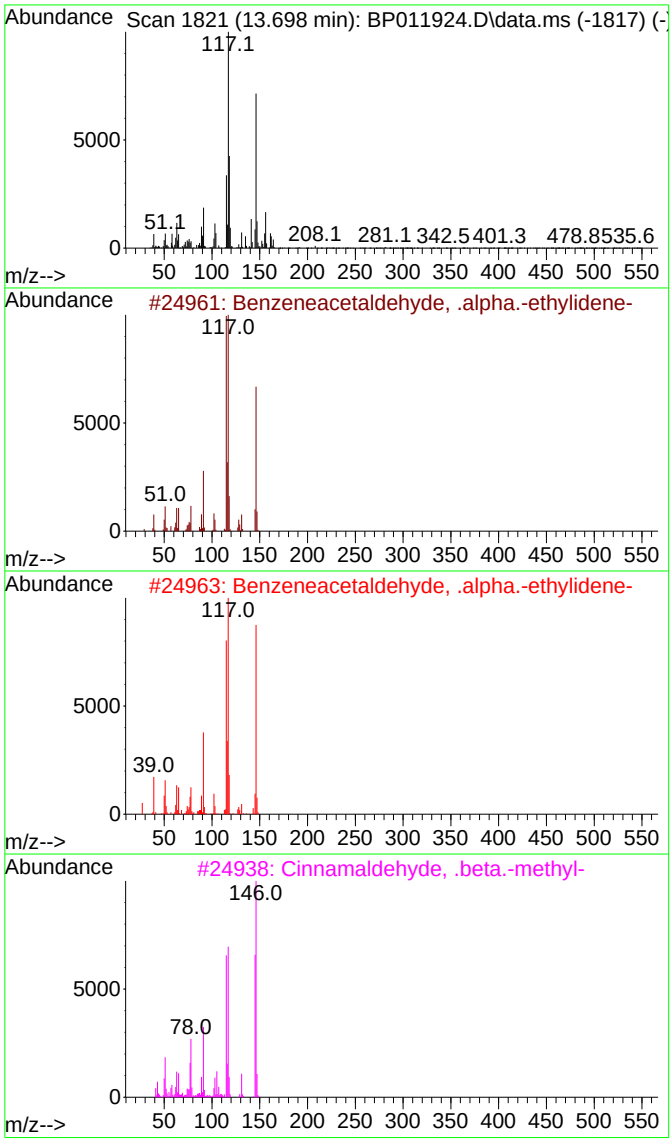
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

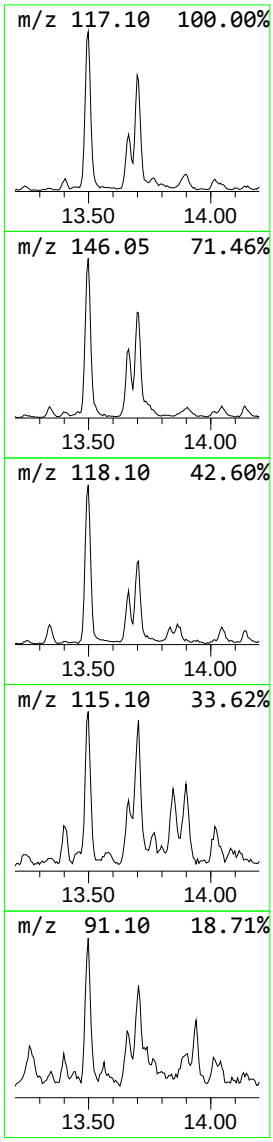
****

Peak Number 25 unknown-02 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.698	2.76 ng/ul	350744	Acenaphthene-d10	14.528

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzeneacetaldehyde, .alpha.-eth...	146	C10H10O	004411-89-6	49
2	Benzeneacetaldehyde, .alpha.-eth...	146	C10H10O	004411-89-6	49
3	Cinnamaldehyde, .beta.-methyl-	146	C10H10O	001196-67-4	47
4	Indane	118	C9H10	000496-11-7	46
5	Benzene, 1-(chloromethyl)-3-ethe...	152	C9H9Cl	039833-65-3	46





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Data File : BP011924.D
Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

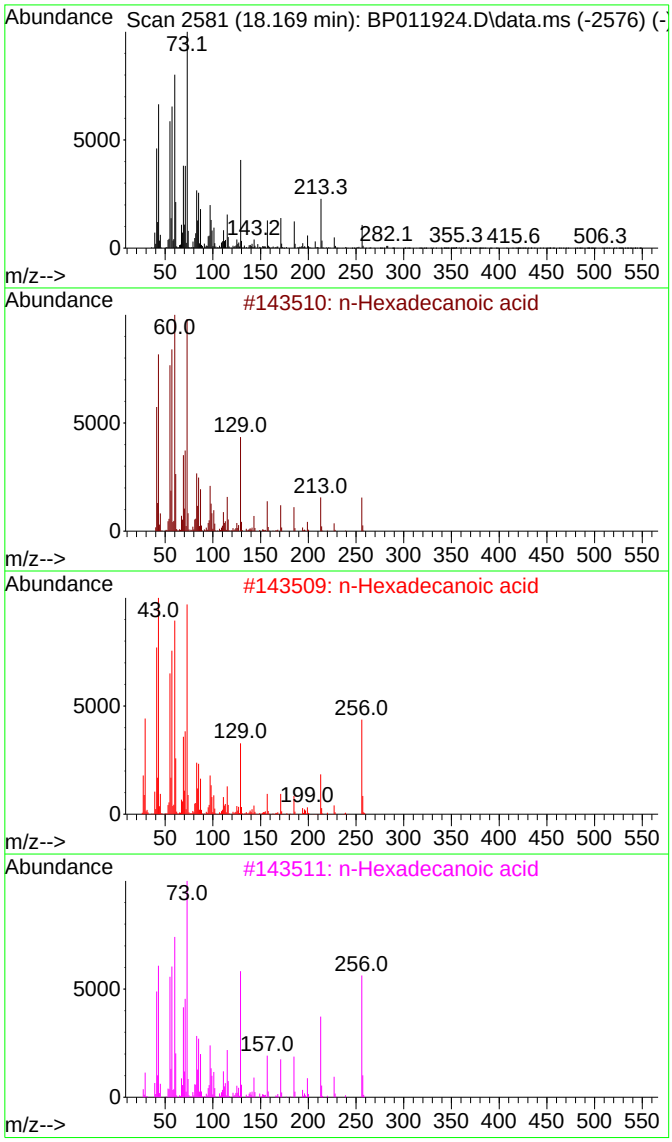
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

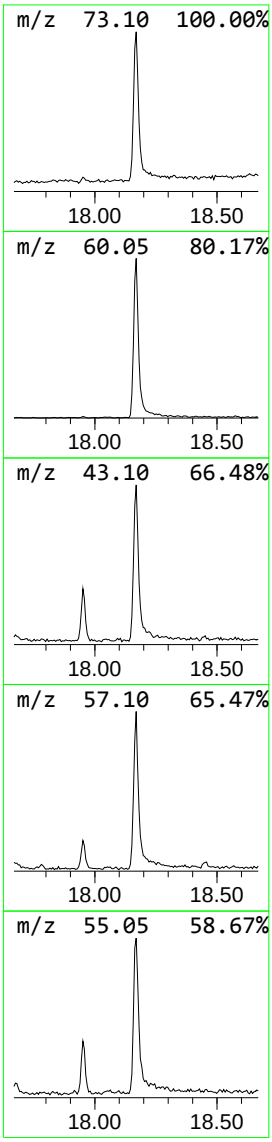
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 n-Hexadecanoic acid Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.169	2.49 ng/ul	365919	Phenanthrene-d10	17.287

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	93
5	Octadecanoic acid	284	C18H36O2	000057-11-4	91





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Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

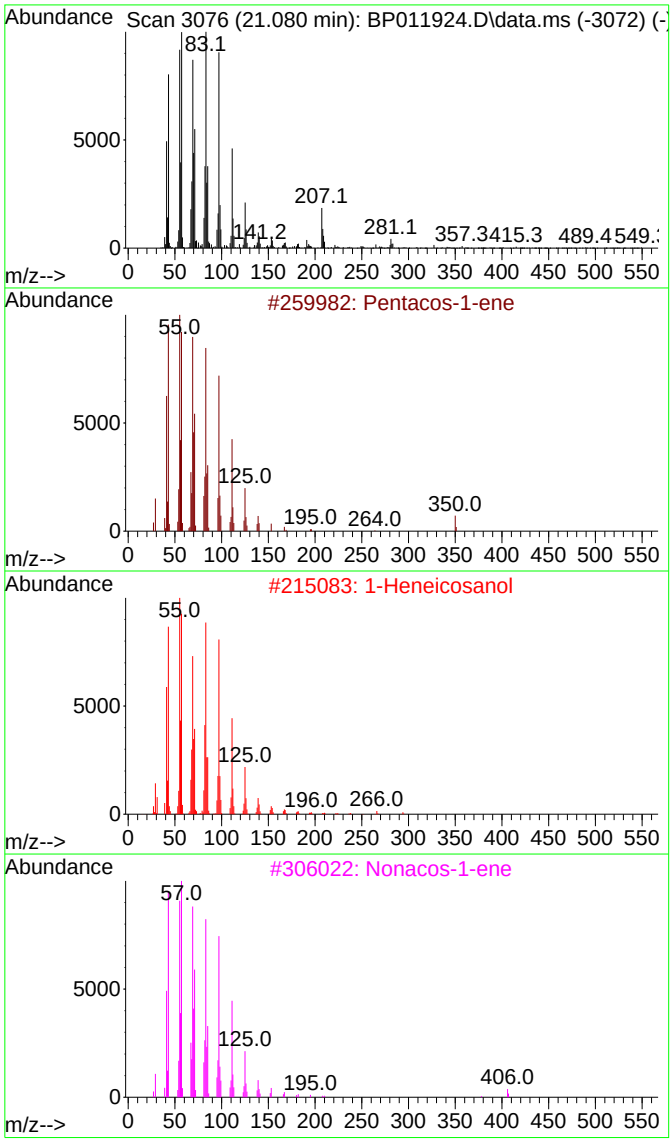
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION

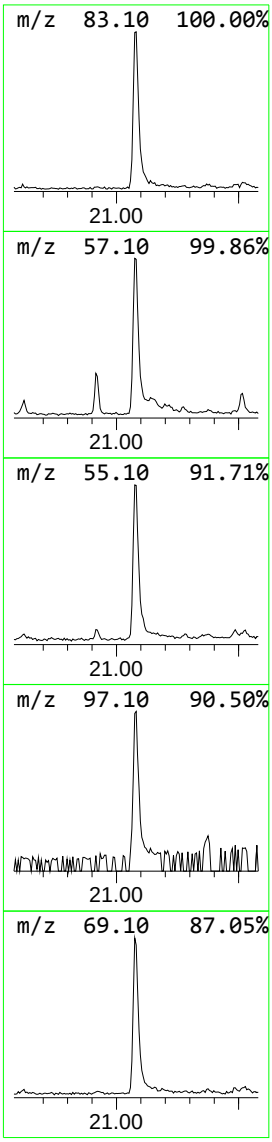
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.080	2.91 ng/ul	512889	Chrysene-d12	21.369

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentacos-1-ene	350	C25H50	016980-85-1	94
2	1-Heneicosanol	312	C21H44O	015594-90-8	94
3	Nonacos-1-ene	406	C29H58	018835-35-3	94
4	1-Hexacosene	364	C26H52	018835-33-1	94
5	n-Nonadecanol-1	284	C19H40O	001454-84-8	94





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011924.D
Acq On : 05 Oct 2022 11:39
Operator : CG/JU
Sample : N4890-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EY0D8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.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
Butane, 2-metho...	3.034	91.9	ng/ul	6046310	1	7.881	1316020	20.0
(DEL) Alkane: S...	3.799	2.1	ng/ul	135315	1	7.881	1316020	20.0
Benzene, 1-ethy...	7.269	31.9	ng/ul	2096030	1	7.881	1316020	20.0
Indane	8.258	2.4	ng/ul	159797	1	7.881	1316020	20.0
Benzene, 1,4-di...	8.375	7.2	ng/ul	472711	1	7.881	1316020	20.0
Benzene, 1,3-di...	8.522	3.5	ng/ul	230642	1	7.881	1316020	20.0
Benzene, 1-meth...	8.687	5.8	ng/ul	380637	1	7.881	1316020	20.0
Benzene, 2-ethy...	8.834	10.8	ng/ul	711325	1	7.881	1316020	20.0
Benzene, 1-meth...	8.887	12.4	ng/ul	816483	1	7.881	1316020	20.0
o-Cymene	8.987	20.0	ng/ul	1313340	1	7.881	1316020	20.0
Benzene, 4-ethy...	9.322	3.1	ng/ul	382459	2	10.693	2465350	20.0
Benzene, 1,2,4,...	9.510	9.1	ng/ul	1122260	2	10.693	2465350	20.0
Benzene, 1,2,3,...	9.575	10.5	ng/ul	1298250	2	10.693	2465350	20.0
1H-Indene, 2,3-...	9.934	5.1	ng/ul	634012	2	10.693	2465350	20.0
Benzene, 2-ethe...	10.087	16.3	ng/ul	2009880	2	10.693	2465350	20.0
unknown-01	11.310	2.8	ng/ul	346448	2	10.693	2465350	20.0
Bicyclo[4.2.1]n...	11.352	2.2	ng/ul	272277	2	10.693	2465350	20.0
1H-Indene, 2,3-...	11.610	2.6	ng/ul	323686	2	10.693	2465350	20.0
3,4-Dimethylben...	11.669	5.0	ng/ul	610387	2	10.693	2465350	20.0
1H-Indene, 2,3-...	11.799	2.4	ng/ul	291582	2	10.693	2465350	20.0
1H-Inden-1-one,...	12.081	4.0	ng/ul	493235	2	10.693	2465350	20.0
Benzene, 1-ethe...	13.499	2.4	ng/ul	306221	3	14.528	2540460	20.0
unknown-02	13.698	2.8	ng/ul	350744	3	14.528	2540460	20.0
n-Hexadecanoic ...	18.169	2.5	ng/ul	365919	4	17.287	2942790	20.0
(DEL) Alkane: S...	21.080	2.9	ng/ul	512889	5	21.369	3519920	20.0