LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042023\
Data File : BM039528.D
Acq On : 20 Apr 2023 17:12
Operator : CG/JU
Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBN2

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_M\Methods\SFAM-EPA-BM041823.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BM039528.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.307	380	386	396	rBV	483523	703246	0.52%	0.094%
2	5.442	402	409	425	rBV	1027305	1997698	1.47%	0.268%
3	5.813	463	472	495	rVB2	558287	1052812	0.77%	0.141%
4	6.484	580	586	595	rVB	670947	1015484	0.74%	0.136%
5	6.789	634	638	648	rVB	358626	573455	0.42%	0.077%
6	6.901	648	657	663	rBV	1509198	2293739	1.68%	0.308%
7	6.960	663	667	672	rVV	624405	938525	0.69%	0.126%
8	7.037	672	680	690	rVB	1626531	2720068	2.00%	0.365%
9	7.160	695	701	705	rVV	764876	1264266	0.93%	0.170%
10	7.201	705	708	712	rVV	305622	494072	0.36%	0.066%
11	7.360	730	735	742	rVV	764660	1301517	0.95%	0.175%
12	7.466	746	753	762	rVV	4414142	7154226	5.25%	0.960%
13	7.560	762	769	778	rVB4	221865	479093	0.35%	0.064%
14	7.825	805	814	825	rBV2	648352	1198072	0.88%	0.161%
15	7.936	826	833	841	rVV2	1442463	3575043	2.62%	0.480%
16	8.036	841	850	862	rVB2	1474736	2632887	1.93%	0.353%
17	8.201	870	878	893	rBV	14358255	24070455	17.66%	3.229%
18	8.372	899	907	917	rVV	18339666	32113043	23.56%	4.308%
19	8.460	917	922	933	rVB	413096	682976	0.50%	0.092%
20	8.560	933	939	949	rBV	563063	1063732	0.78%	0.143%
21	8.931	996	1002	1008	rBV	477570	824191	0.60%	0.111%
22	9.007	1008	1015	1022	rVV2	1678851	3491998	2.56%	0.469%
23	9.072	1022	1026	1038	rVB	775932	1328280	0.97%	0.178%
24	9.183	1038	1045	1054	rBV	2013445	3476720	2.55%	0.466%
25	9.313	1054	1067	1073	rVV	4429679	9369746	6.87%	1.257%
26	9.372	1073	1077	1086	rVV	1359442	2334811	1.71%	0.313%
27	9.454	1086	1091	1096	rVV	364815	577338	0.42%	0.077%
28	9.513	1096	1101	1110	rVB	370168	623244	0.46%	0.084%
29	9.760	1132	1143	1151	rBV3	1045696	2684829	1.97%	0.360%
30	9.878	1156	1163	1170	rVB	1209835	2029104	1.49%	0.272%
31	9.978	1170	1180	1184	rBV2	930663	2100034	1.54%	0.282%
32	10.060	1184	1194	1202	rVV4	3180417	10452158	7.67%	1.402%
33	10.142	1202	1208	1210	rVV	1255253	2585861	1.90%	0.347%
34	10.177	1210	1214	1224	rVB	2149538	4278562	3.14%	0.574%
35	10.319	1224	1238	1251	rVB3	634275	2310441	1.69%	0.310%
36	10.719	1287	1306	1314	rBV2	27546744	136329109	100.00%	18.291%

LSC Area Percent Report

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Operator : CG/JU
Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBN2

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_M\Methods\SFAM-EPA-BM041823.MA.M
Title : SVOA CALIBRATION

37	10.883	1329	1334	1342	rVB	21446874	32460304	23.81%	4.355%
38	11.066	1360	1365	1371	rVV	369450	692942	0.51%	0.093%
39	11.419	1413	1425	1432	rBV2	623193	1248337	0.92%	0.167%
40	11.754	1474	1482	1487	rBV2	253375	511907	0.38%	0.069%
41	12.195	1548	1557	1568	rVB	1692816	3177951	2.33%	0.426%
42	12.336	1568	1581	1591	rBV2	30129742	85420856	62.66%	11.461%
43	12.430	1591	1597	1604	rVV2	2811894	5397590	3.96%	0.724%
44	12.548	1604	1617	1628	rVB	26200262	54302012	39.83%	7.285%
45	12.954	1679	1686	1693	rBV	639051	1021692	0.75%	0.137%
46	13.318	1741	1748	1761	rBV	10694660	16343529	11.99%	2.193%
47	13.513	1775	1781	1793	rVB	2118178	4236248	3.11%	0.568%
48	13.642	1793	1803	1804	rBV2	1952513	2701750	1.98%	0.362%
49	13.801	1823	1830	1835	rVV	3248309	5842667	4.29%	0.784%
50	13.854	1835	1839	1843	rVV	2200280	3429667	2.52%	0.460%
51	13.901	1843	1847	1858	rVB	2313010	3463065	2.54%	0.465%
52	14.036	1858	1870	1874	rBV	1225597	2004205	1.47%	0.269%
53	14.071	1874	1876	1882	rVB	518707	596981	0.44%	0.080%
54	14.177	1887	1894	1897	rBV	2810552	4340695	3.18%	0.582%
55	14.483	1937	1946	1951	rBV2	1360751	3431465	2.52%	0.460%
56	14.565	1951	1960	1966	rVV	27848488	53299131	39.10%	7.151%
57	14.630	1966	1971	1978	rVB	8416113	12744521	9.35%	1.710%
58	14.760	1989	1993	1996	rBV	378281	534933	0.39%	0.072%
59	14.795	1996	1999	2004	rVB	495758	661573	0.49%	0.089%
60	14.895	2008	2016	2025	rBV2	19651080	37560986	27.55%	5.039%
61	15.477	2110	2115	2120	rBV	3175635	4505667	3.30%	0.605%
62	15.536	2120	2125	2133	rVB	11511216	17687099	12.97%	2.373%
63	15.612	2133	2138	2143	rBV2	1033370	1710141	1.25%	0.229%
64	15.654	2143	2145	2149	rVV	458980	628796	0.46%	0.084%
65	15.695	2149	2152	2156	rVV2	610500	849264	0.62%	0.114%
66	15.783	2162	2167	2170	rVV2	483160	822755	0.60%	0.110%
67	15.824	2170	2174	2185	rVB3	825333	1452296	1.07%	0.195%
68	15.971	2195	2199	2209	rVV	1020300	1983492	1.45%	0.266%
69	16.212	2234	2240	2254	rBV	3475877	5658462	4.15%	0.759%
70	16.395	2266	2271	2274	rBV	384023	692837	0.51%	0.093%
71	16.436	2275	2278	2284	rVB	469990	767583	0.56%	0.103%
72	16.754	2328	2332	2335	rVV	355530	584190	0.43%	0.078%
73	16.806	2335	2341	2347	rVV	701429	1525124	1.12%	0.205%
74	16.895	2347	2356	2369	rVB3	2148194	4950563	3.63%	0.664%
75	17.054	2374	2383	2391	rVB	1037175	1694814	1.24%	0.227%
76	17.201	2403	2408	2410	rVV	321372	458205	0.34%	0.061%

LSC Area Percent Report

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Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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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
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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Title : SVOA CALIBRATION

77	17.236	2410	2414	2417	rVV	1626491	2498473	1.83%	0.335%
78	17.277	2417	2421	2427	rVV	10832946	16689023	12.24%	2.239%
79	17.336	2427	2431	2435	rVV	3385481	5099382	3.74%	0.684%
80	17.448	2436	2450	2462	rVV3	1539558	6448328	4.73%	0.865%
81	17.571	2462	2471	2476	rVV	1191517	2248508	1.65%	0.302%
82	17.659	2476	2486	2496	rVV	17702385	31038332	22.77%	4.164%
83	17.759	2500	2503	2507	rVV	518018	776018	0.57%	0.104%
84	17.801	2507	2510	2516	rVV3	306256	560019	0.41%	0.075%
85	17.895	2520	2526	2531	rVB	363076	799374	0.59%	0.107%
86	17.948	2531	2535	2540	rBV	376861	512191	0.38%	0.069%
87	18.165	2564	2572	2577	rVV3	1615610	3489578	2.56%	0.468%
88	18.306	2591	2596	2603	rBV2	417732	705953	0.52%	0.095%
89	18.400	2608	2612	2617	rBV2	492476	1051648	0.77%	0.141%
90	18.465	2620	2623	2627	rVV	461894	610939	0.45%	0.082%
91	18.559	2634	2639	2646	rBV2	669488	1168779	0.86%	0.157%
92	19.412	2780	2784	2795	rBV	601928	878202	0.64%	0.118%
93	19.630	2815	2821	2827	rBV	3901282	5350474	3.92%	0.718%
94	19.689	2827	2831	2842	rVB	387061	809924	0.59%	0.109%
95	20.047	2887	2892	2899	rBV	325409	474033	0.35%	0.064%
96	20.118	2900	2904	2915	rVV	336679	597244	0.44%	0.080%
97	20.771	3011	3015	3031	rVB2	460402	814182	0.60%	0.109%
98	21.424	3121	3126	3134	rBV	1729022	2247323	1.65%	0.302%
99	23.641	3496	3503	3520	rVB2	2315994	4805807	3.53%	0.645%
100	23.794	3522	3529	3539	rVB2	1029246	2158134	1.58%	0.290%

Sum of corrected areas: 745348998

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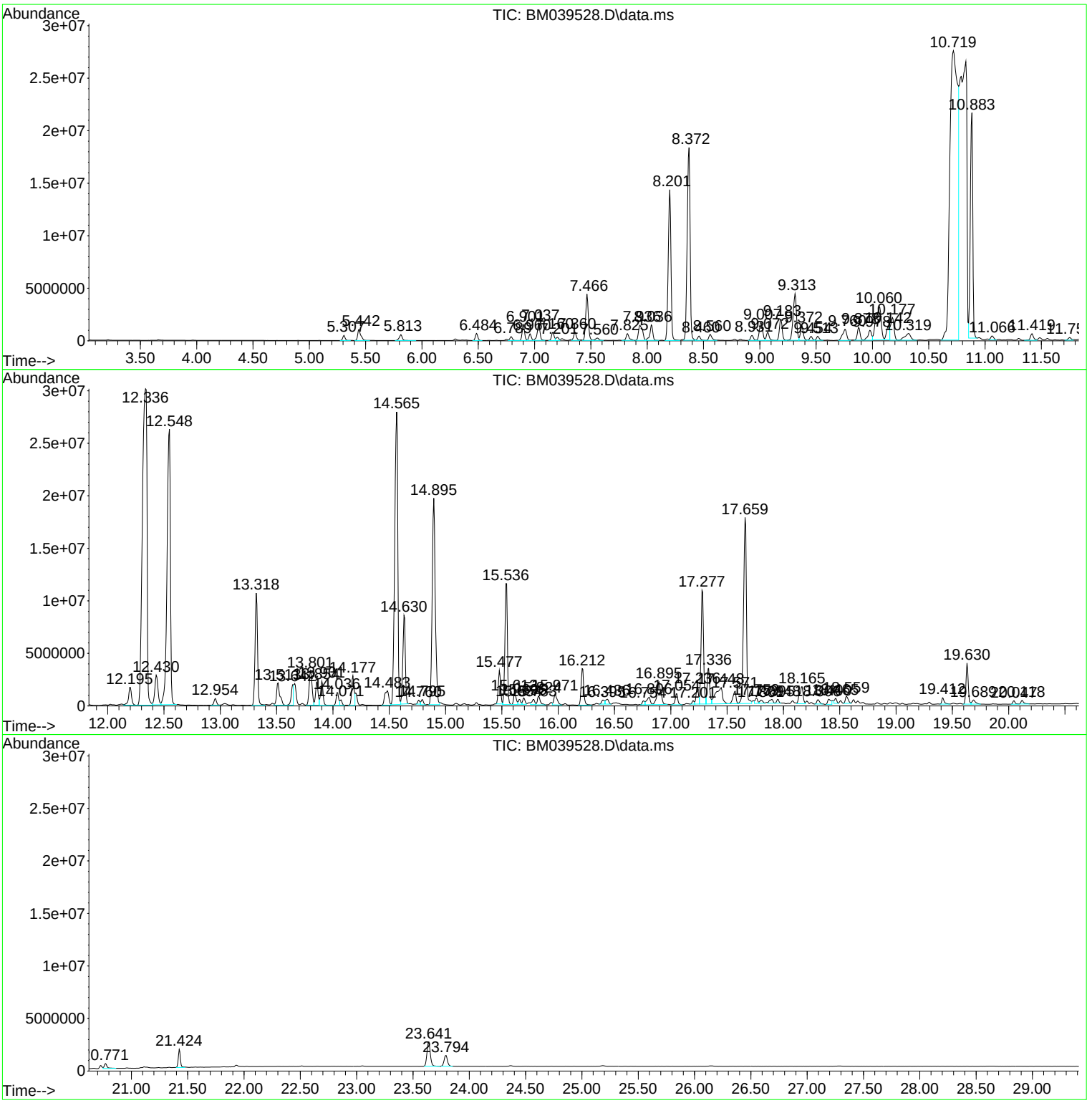
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBN2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042023\
Data File : BM039528.D
Acq On : 20 Apr 2023 17:12
Operator : CG/JU
Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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

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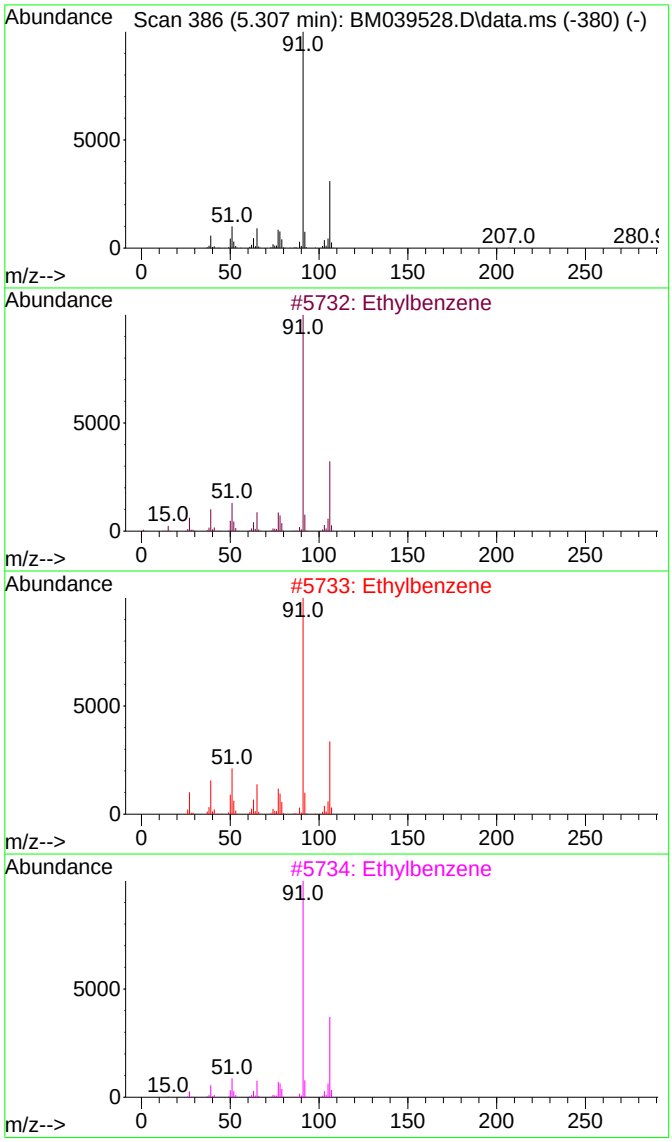
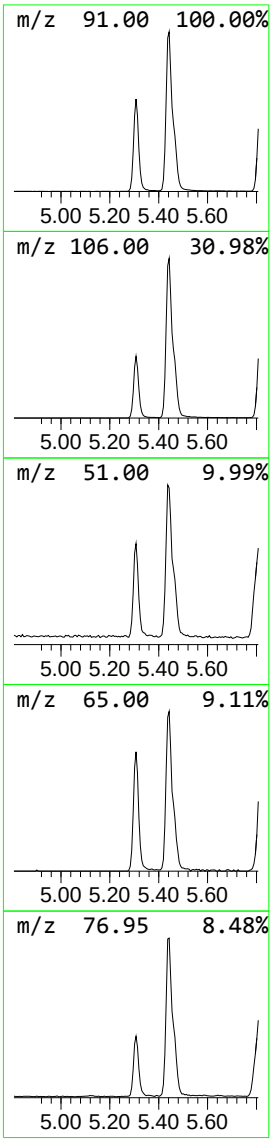
Peak Number 1 Ethylbenzene Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.307	11.74 ng/ul	703246	1,4-Dichlorobenzene-d4	7.825

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethylbenzene	106	C8H10	000100-41-4	94
2	Ethylbenzene	106	C8H10	000100-41-4	91
3	Ethylbenzene	106	C8H10	000100-41-4	90
4	Ethylbenzene	106	C8H10	000100-41-4	87
5	o-Xylene	106	C8H10	000095-47-6	80





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042023\
Data File : BM039528.D
Acq On : 20 Apr 2023 17:12
Operator : CG/JU
Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

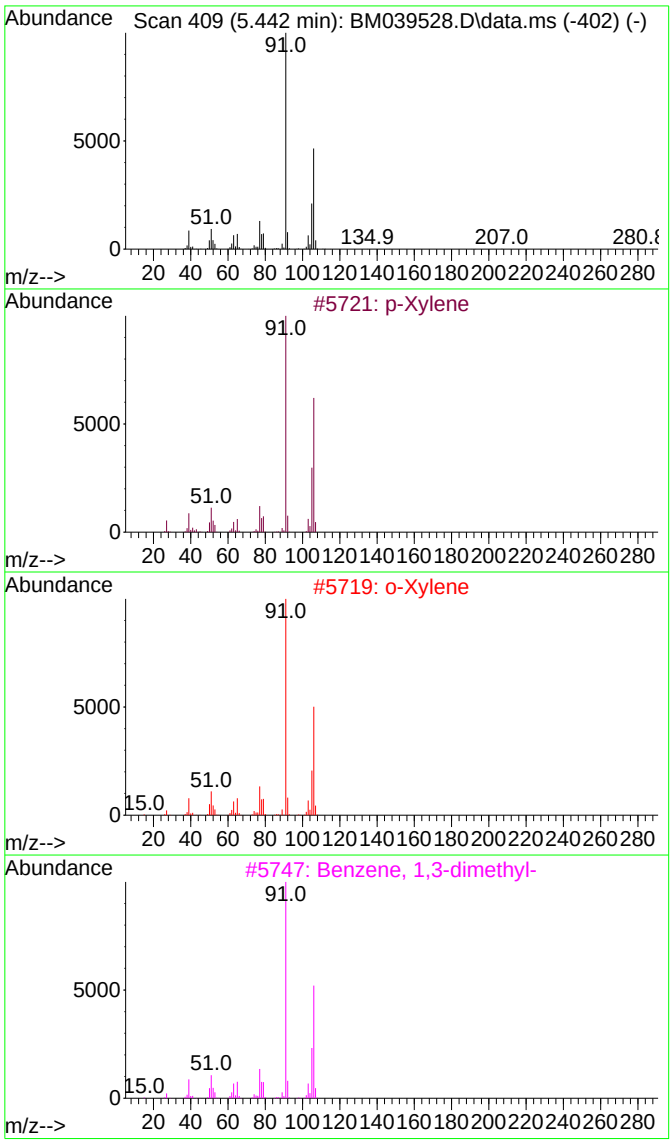
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

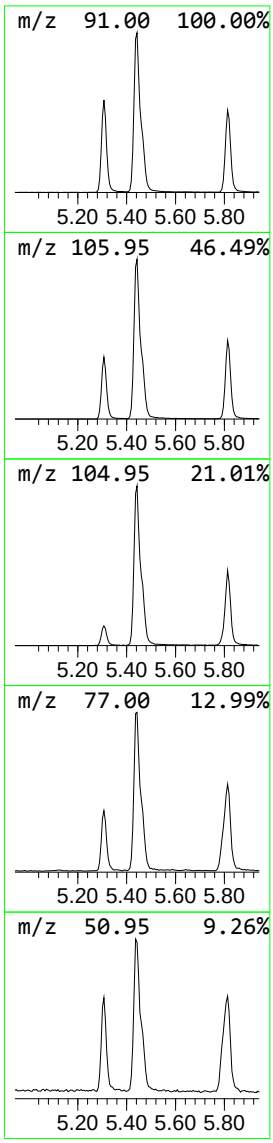
****

Peak Number 2 p-Xylene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.442	33.35 ng/ul	1997700	1,4-Dichlorobenzene-d4	7.825

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





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Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBN2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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

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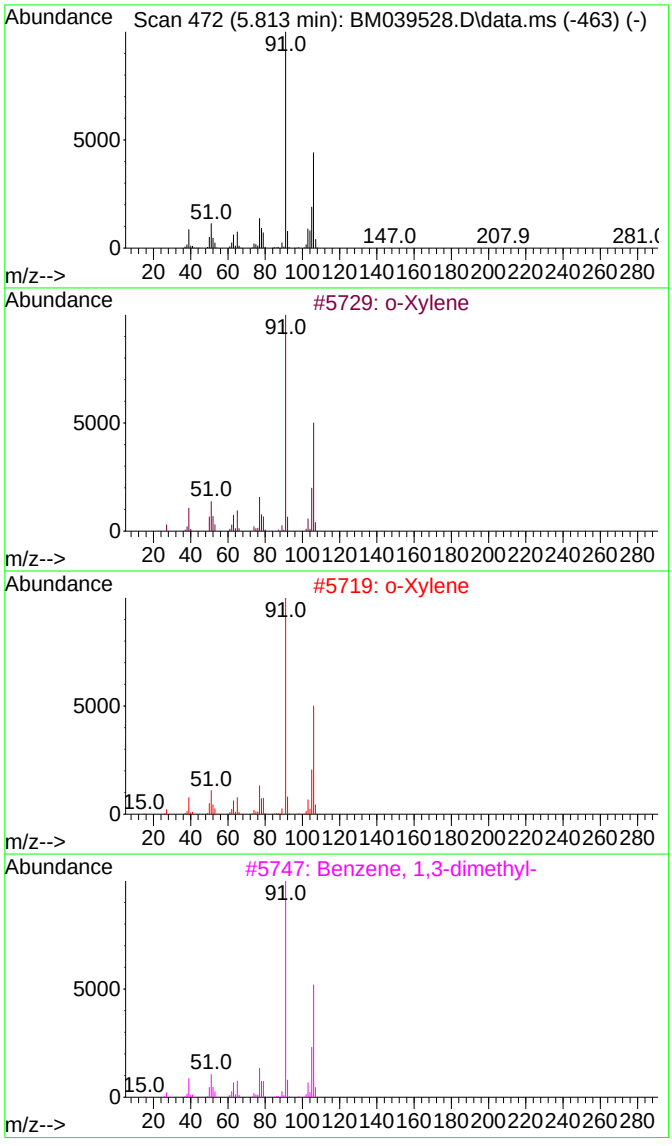
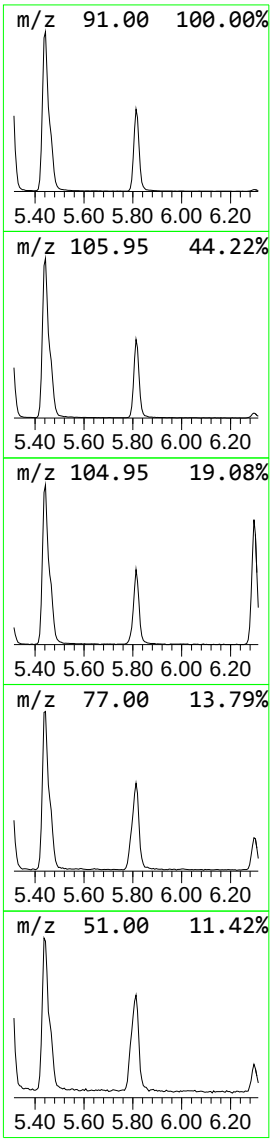
Peak Number 3 o-Xylene Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.813	17.58 ng/ul	1052810	1,4-Dichlorobenzene-d4	7.825

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





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Data File : BM039528.D
Acq On : 20 Apr 2023 17:12
Operator : CG/JU
Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBN2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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

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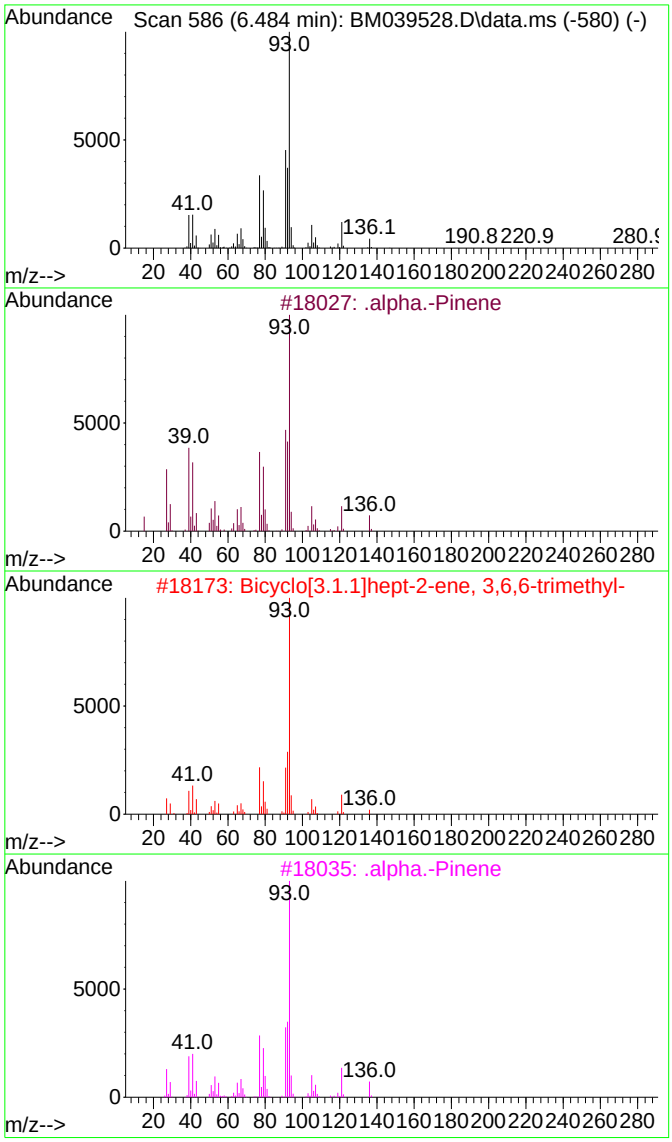
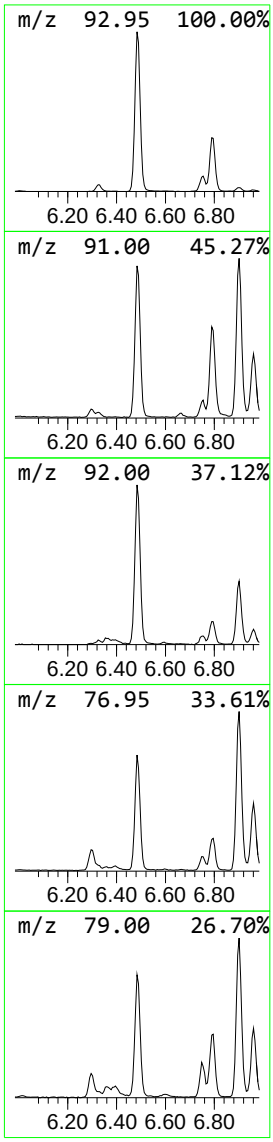
Peak Number 4 .alpha.-Pinene Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.484	16.95 ng/ul	1015480	1,4-Dichlorobenzene-d4	7.825

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-Pinene	136	C10H16	000080-56-8	95
2	Bicyclo[3.1.1]hept-2-ene, 3,6,6-...	136	C10H16	004889-83-2	94
3	.alpha.-Pinene	136	C10H16	000080-56-8	94
4	.alpha.-Pinene	136	C10H16	000080-56-8	91
5	3-Carene	136	C10H16	013466-78-9	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042023\
Data File : BM039528.D
Acq On : 20 Apr 2023 17:12
Operator : CG/JU
Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBN2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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

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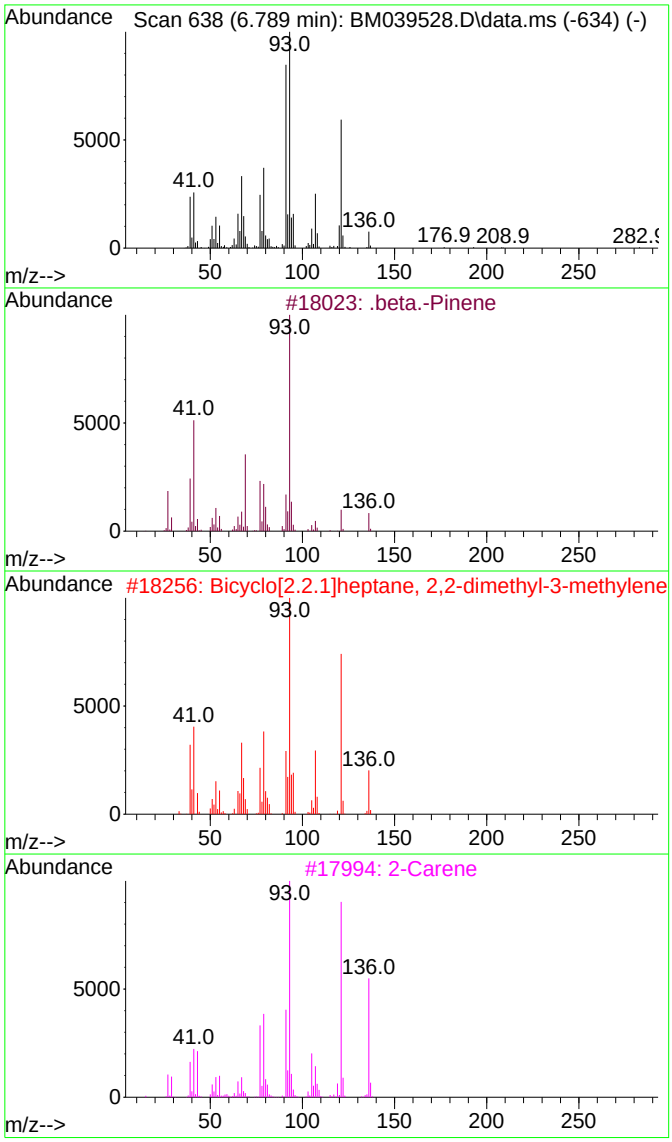
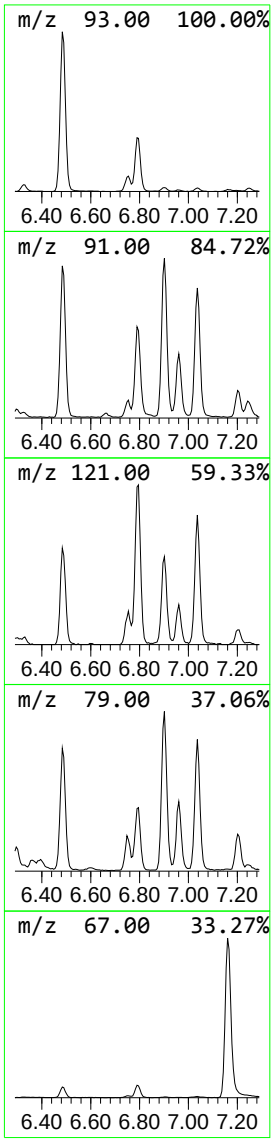
Peak Number 5 .beta.-Pinene Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.789	9.57 ng/ul	573455	1,4-Dichlorobenzene-d4	7.825

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.beta.-Pinene	136	C10H16	000127-91-3	72
2	Bicyclo[2.2.1]heptane, 2,2-dimet...	136	C10H16	005794-04-7	72
3	2-Carene	136	C10H16	000554-61-0	68
4	Cyclohexene, 1-methyl-4-(1-methy...	136	C10H16	000586-62-9	64
5	Camphene	136	C10H16	000079-92-5	60





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Acq On : 20 Apr 2023 17:12
Operator : CG/JU
Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBN2

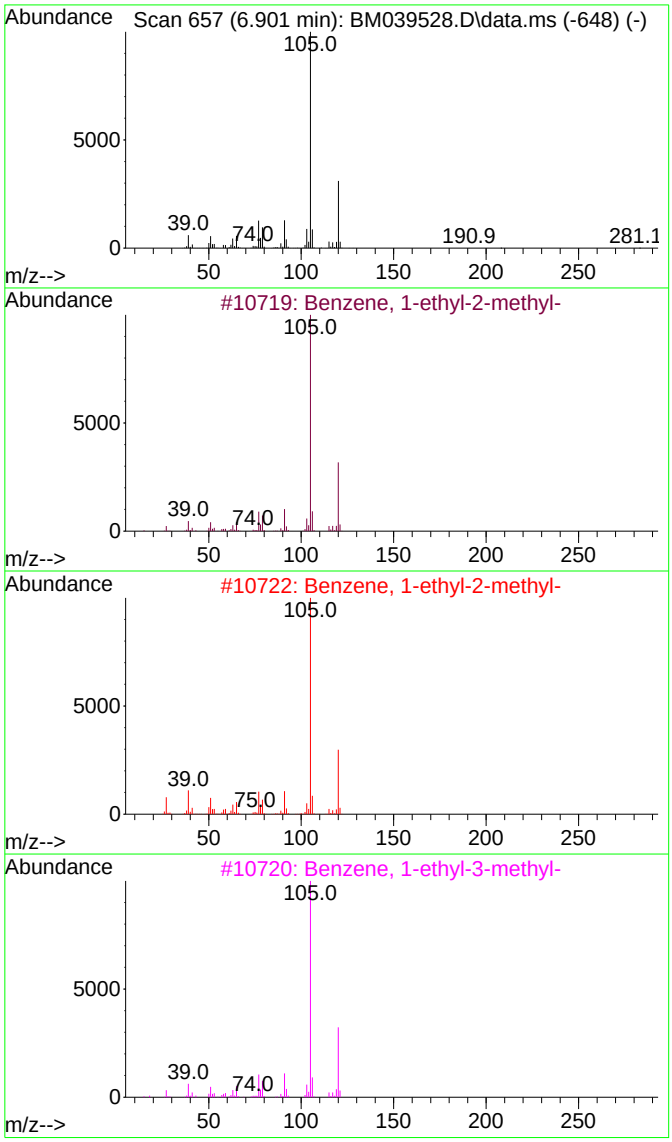
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

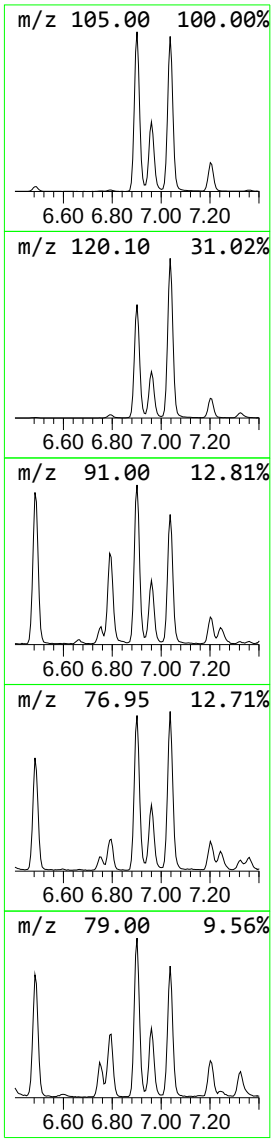
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1-ethyl-2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.901	38.29 ng/ul	2293740	1,4-Dichlorobenzene-d4	7.825

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





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Operator : CG/JU
Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBN2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

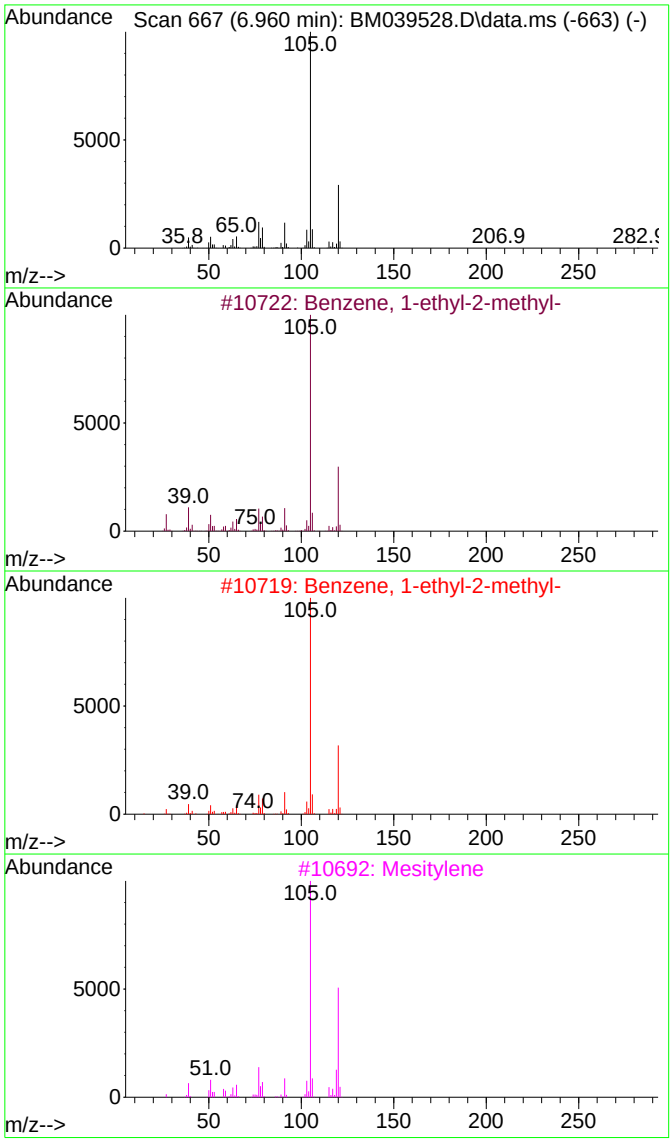
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

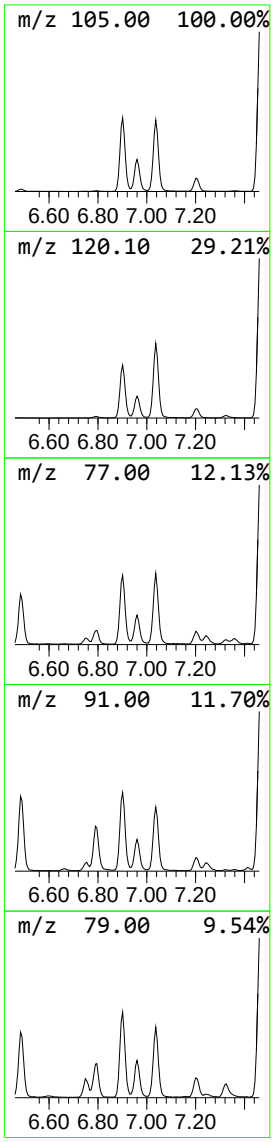
****

Peak Number 7 Mesitylene Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.960	15.67 ng/ul	938525	1,4-Dichlorobenzene-d4	7.825

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	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91





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Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBN2

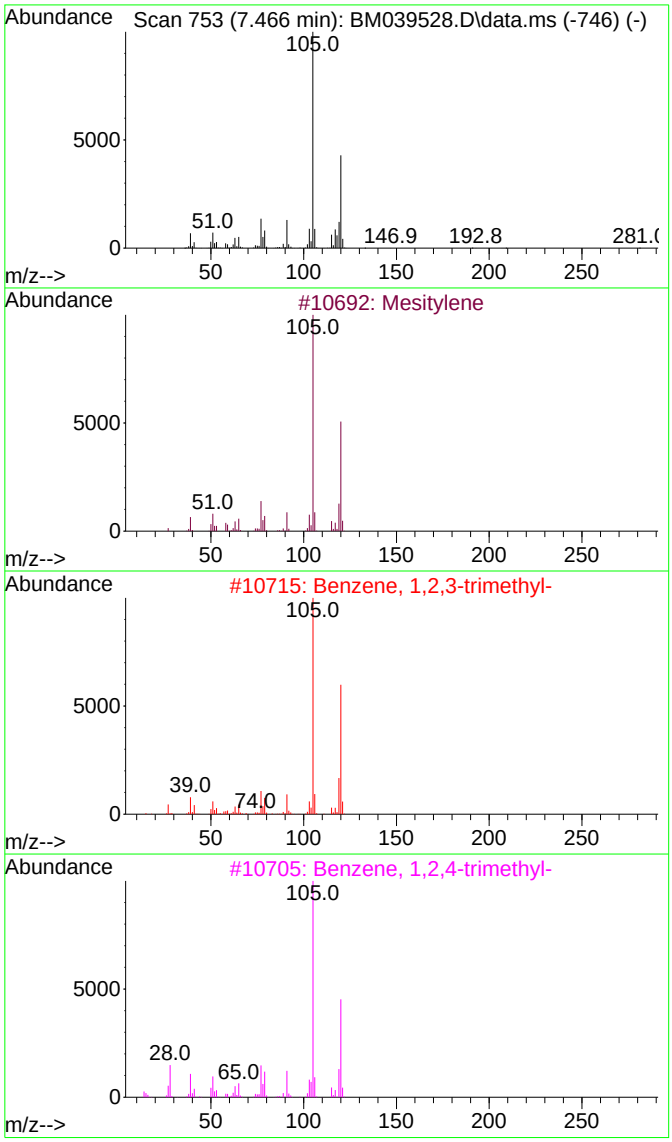
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

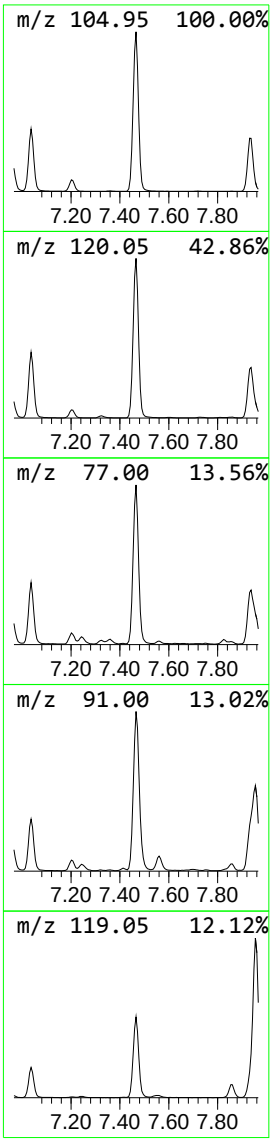
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.466	119.43 ng/ul	7154230	1,4-Dichlorobenzene-d4	7.825

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	94
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	93
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	93
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90





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Operator : CG/JU
Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBN2

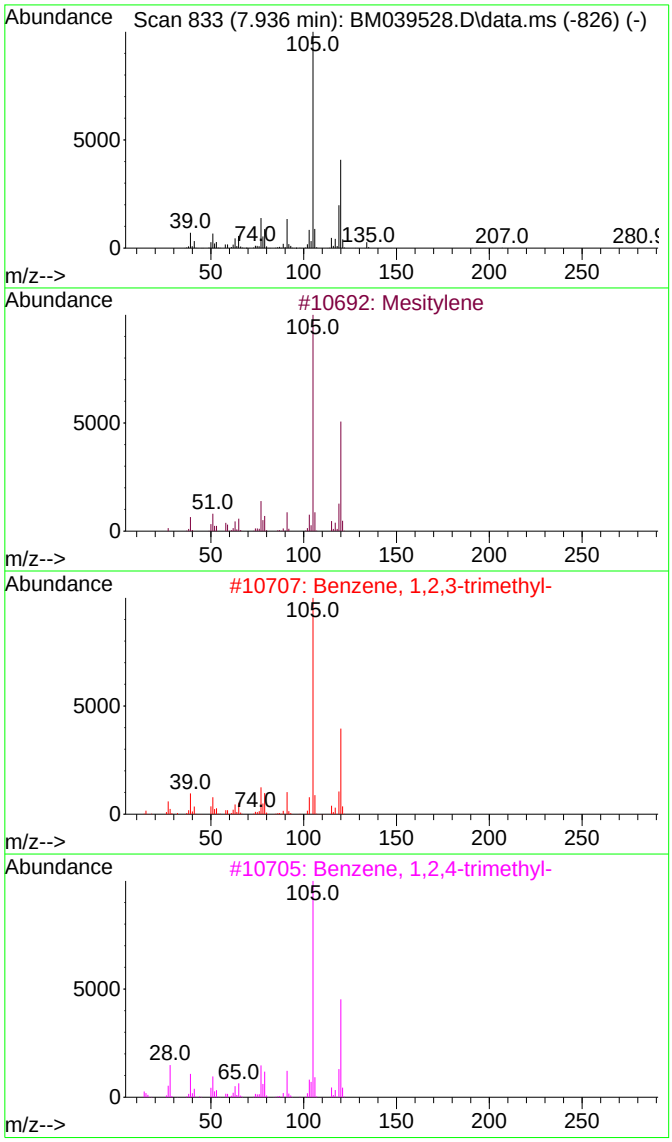
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

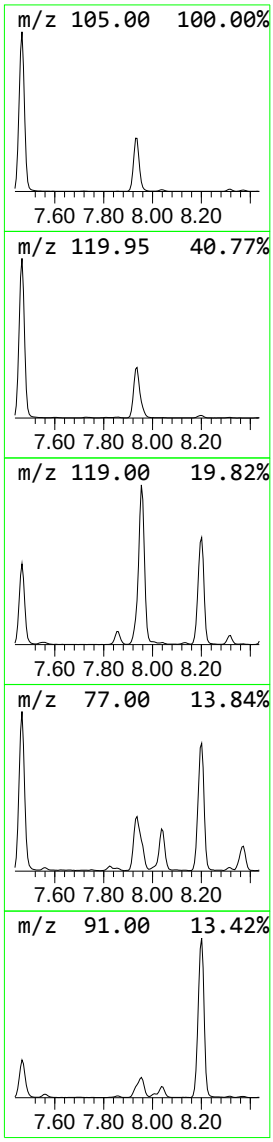
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.936	59.68 ng/ul	3575040	1,4-Dichlorobenzene-d4	7.825

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	93
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	93
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	93
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90





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Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBN2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

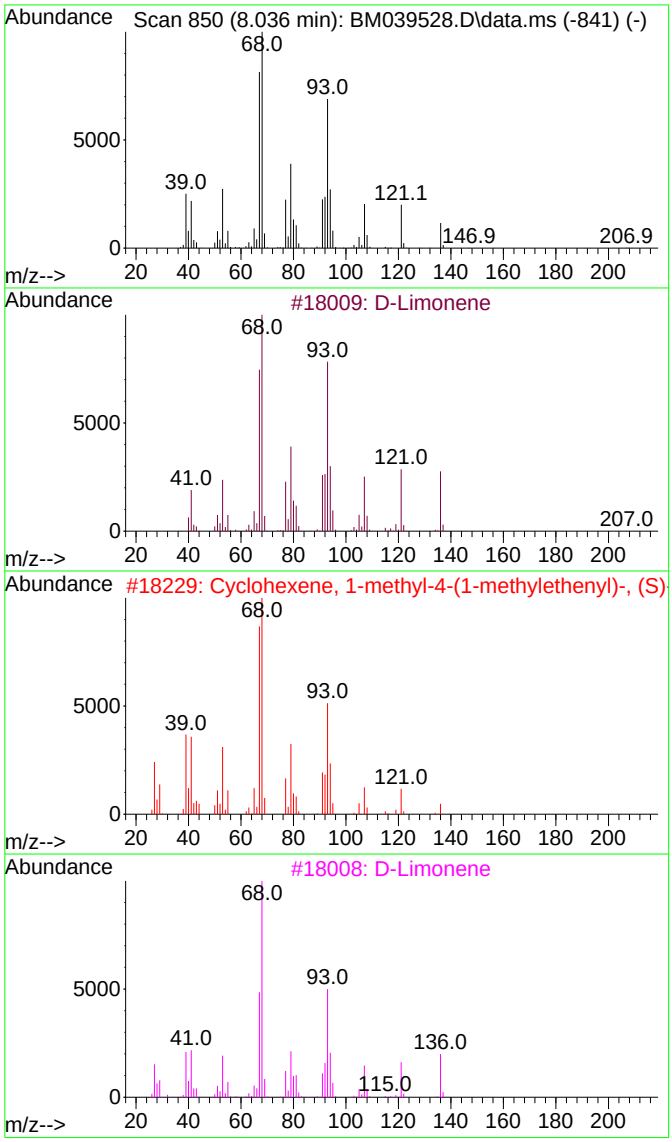
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

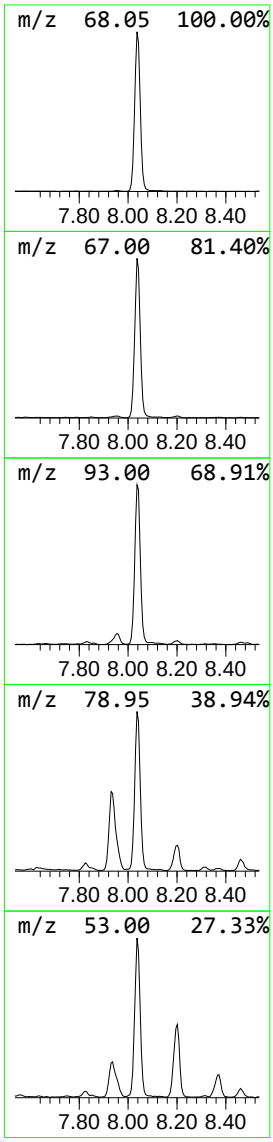
****

Peak Number 11 D-Limonene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.036	43.95 ng/ul	2632890	1,4-Dichlorobenzene-d4	7.825

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	D-Limonene	136	C10H16	005989-27-5	98
2	Cyclohexene, 1-methyl-4-(1-methy...	136	C10H16	005989-54-8	95
3	D-Limonene	136	C10H16	005989-27-5	93
4	Limonene	136	C10H16	000138-86-3	91
5	Cyclohexene, 1-methyl-5-(1-methy...	136	C10H16	013898-73-2	90





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Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBN2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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

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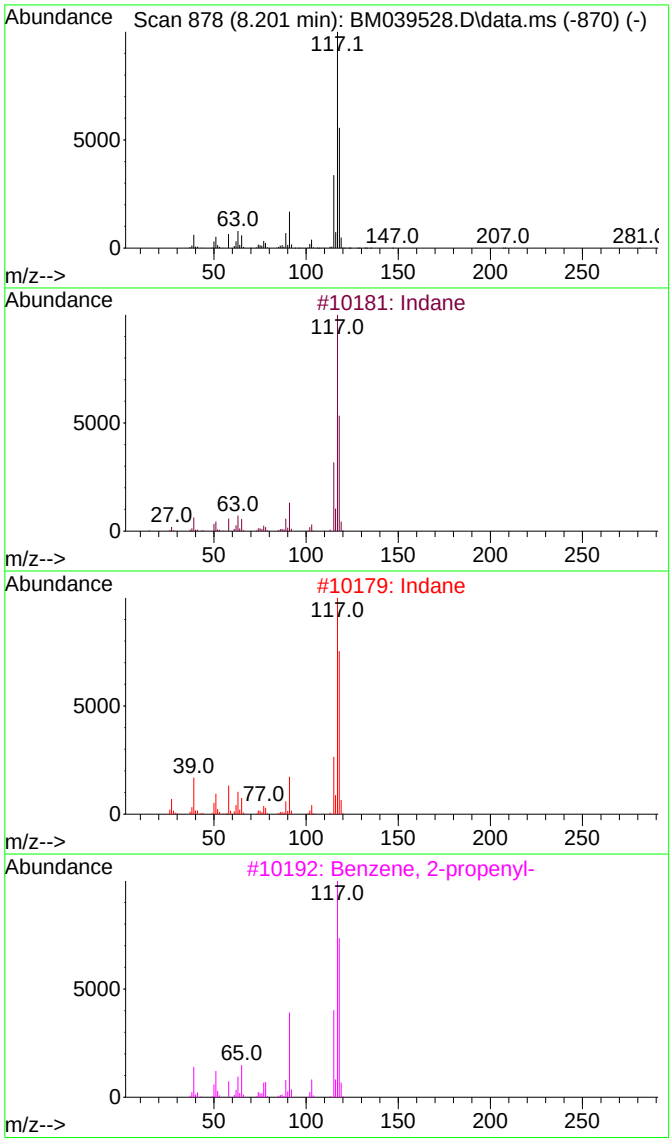
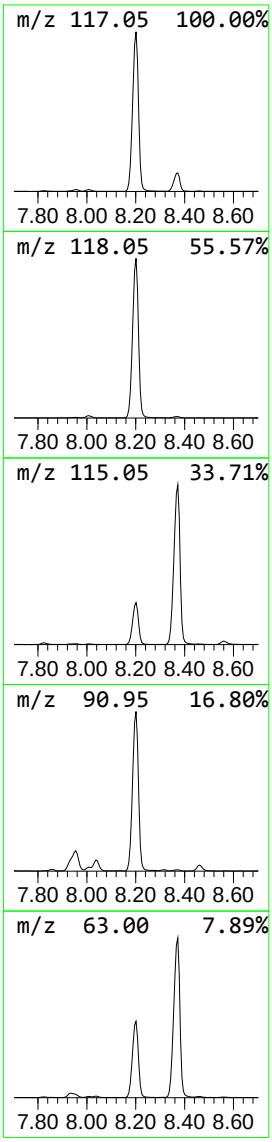
Peak Number 12 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.201	401.82 ng/ul	24070500	1,4-Dichlorobenzene-d4	7.825

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, 2-propenyl-	118	C9H10	000300-57-2	83
4	Deltacyclene	118	C9H10	007785-10-6	72
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	72





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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBN2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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

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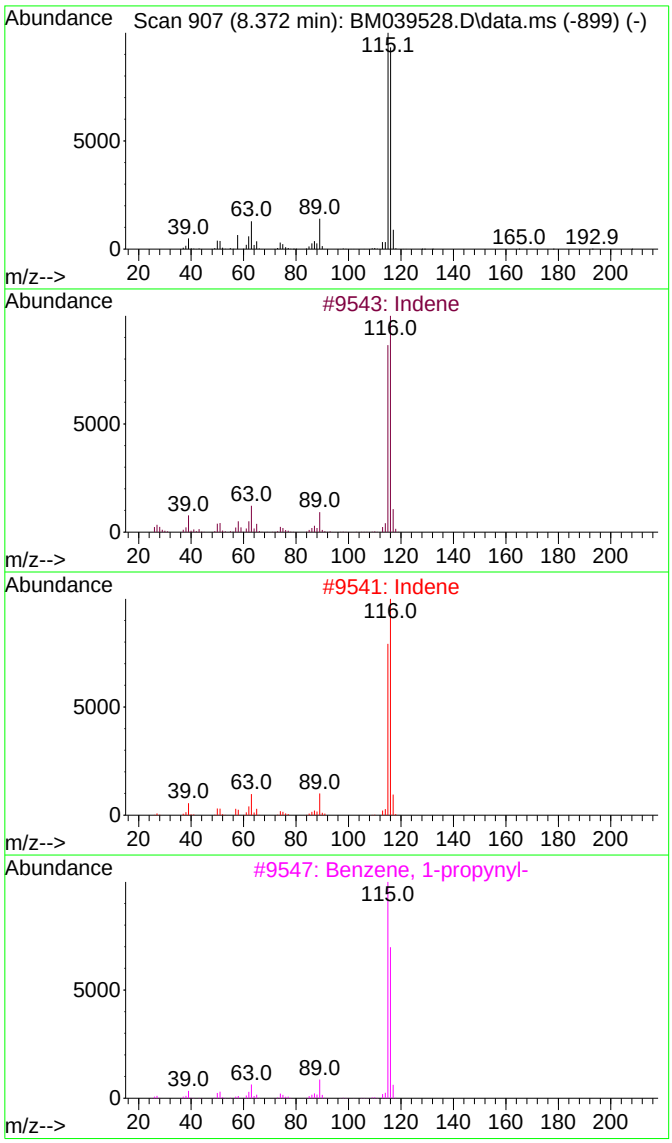
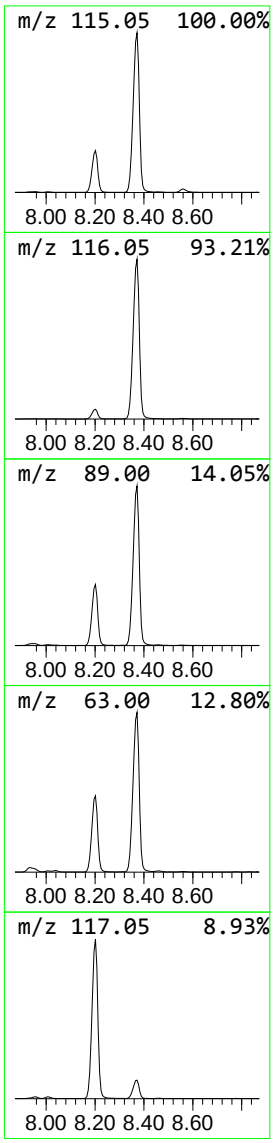
Peak Number 13 Indene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.372	536.08 ng/ul	32113000	1,4-Dichlorobenzene-d4	7.825

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene	116	C9H8	000095-13-6	95
2	Indene	116	C9H8	000095-13-6	95
3	Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
4	Indene	116	C9H8	000095-13-6	94
5	1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	94





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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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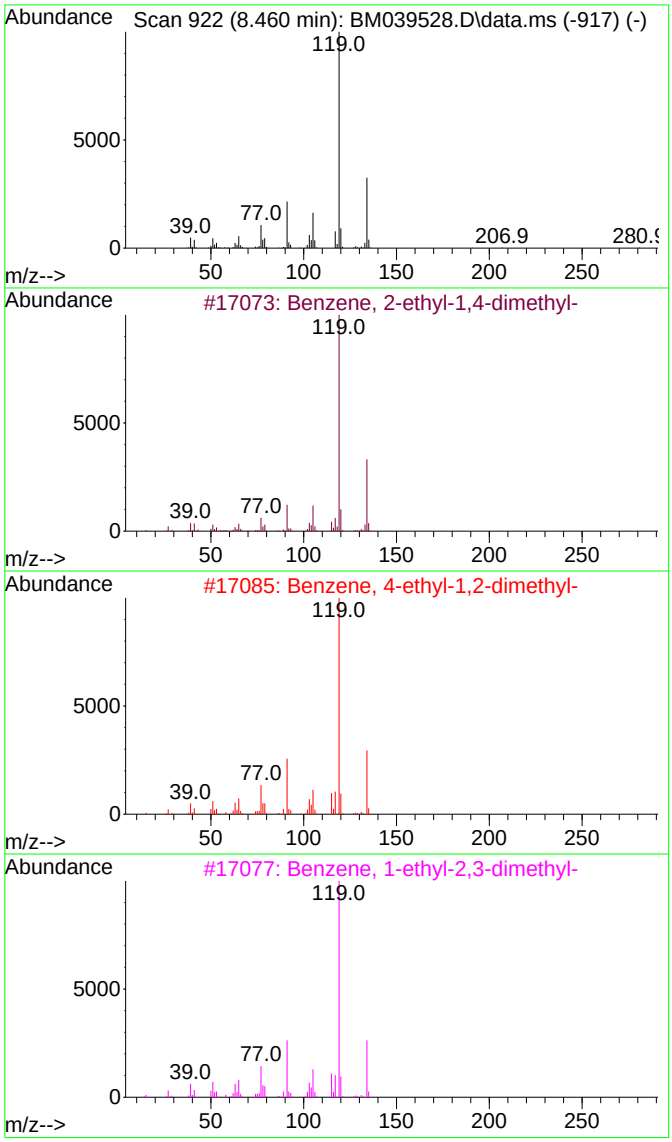
Quant Title : SVOA CALIBRATION

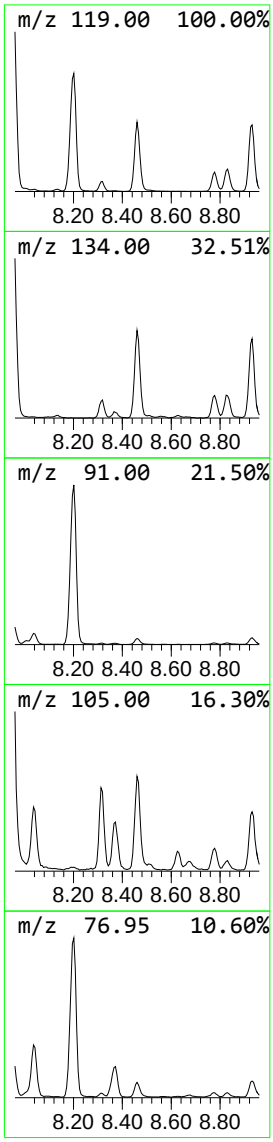
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.460	11.40 ng/ul	682976	1,4-Dichlorobenzene-d4	7.825

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	91
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91





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Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DCBN2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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

****
Peak Number 15 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.931	13.76 ng/ul	824191	1,4-Dichlorobenzene-d4	7.825

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
3	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93
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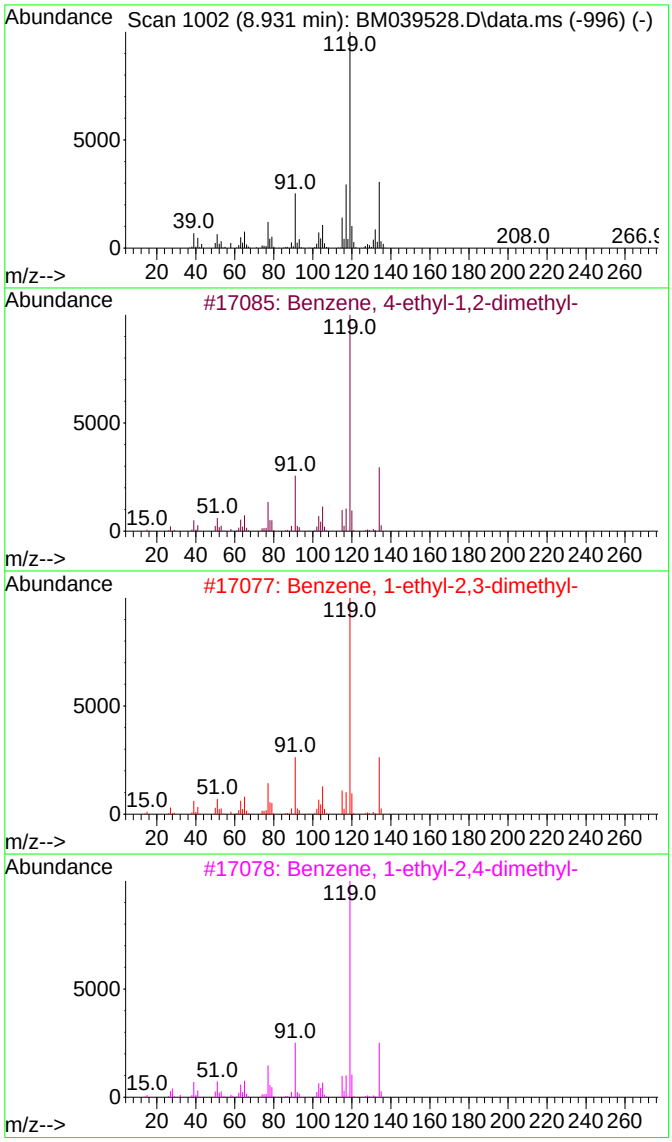

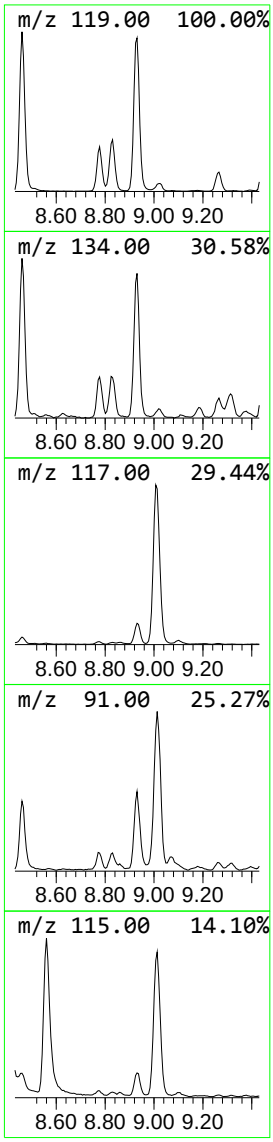

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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

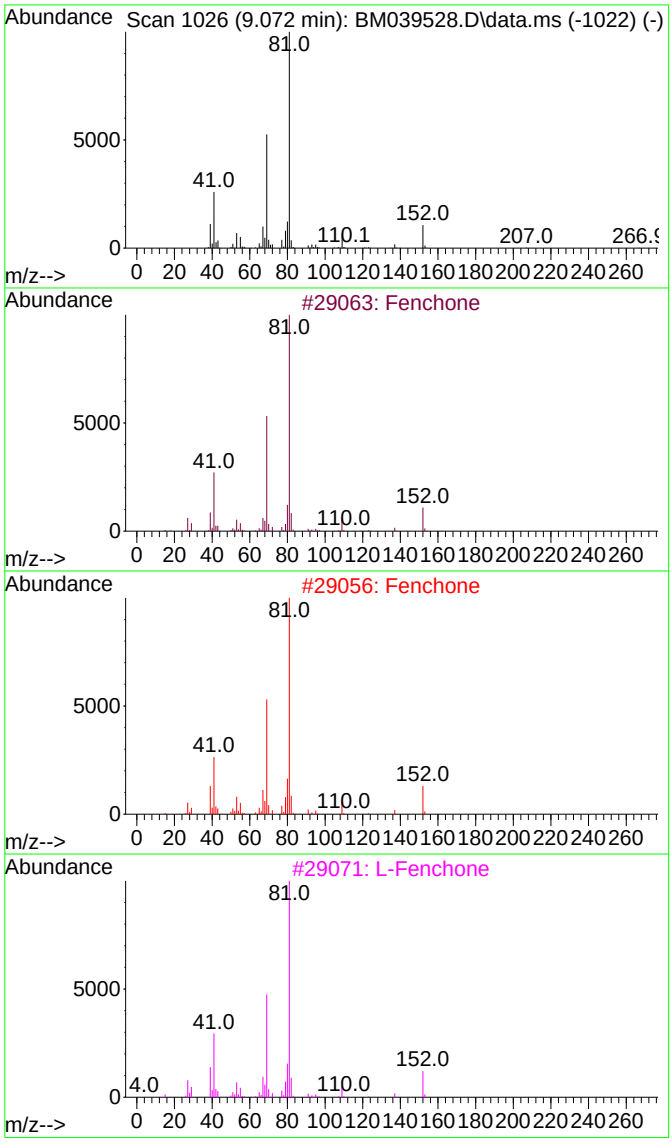
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

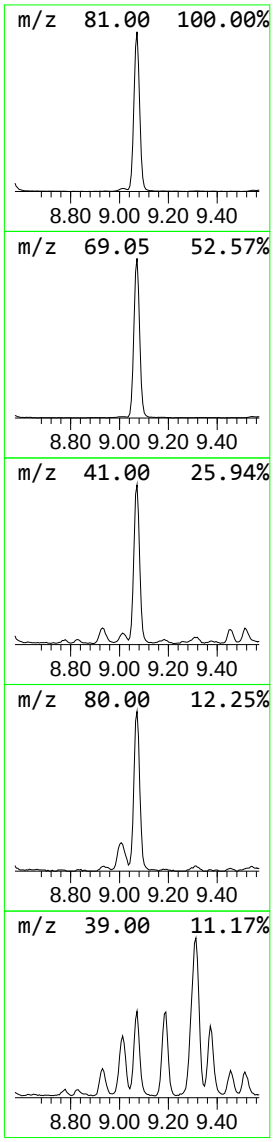
****

Peak Number 16 Fenchone Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.072	22.17 ng/ul	1328280	1,4-Dichlorobenzene-d4	7.825

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fenchone	152	C10H16O	001195-79-5	94
2	Fenchone	152	C10H16O	001195-79-5	94
3	L-Fenchone	152	C10H16O	007787-20-4	94
4	D-Fenchone	152	C10H16O	004695-62-9	91
5	L-Fenchone	152	C10H16O	007787-20-4	91





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ALS Vial : 9 Sample Multiplier: 1

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ClientSampleId :
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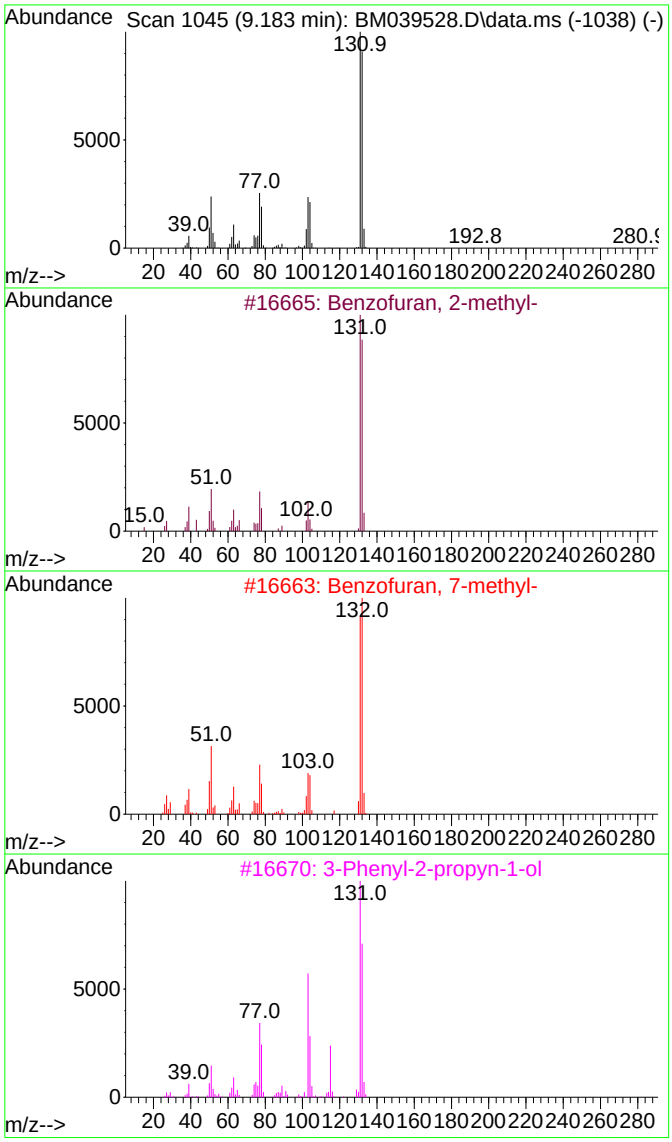
Quant Title : SVOA CALIBRATION

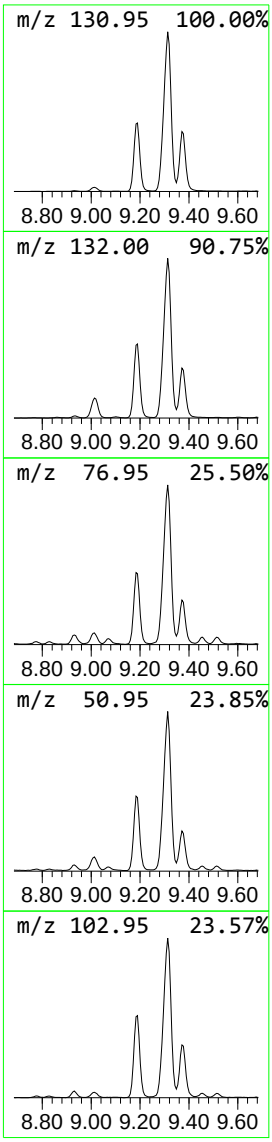
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Benzofuran, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.183	58.04 ng/ul	3476720	1,4-Dichlorobenzene-d4	7.825

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
2	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	94
3	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	91
4	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
5	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91





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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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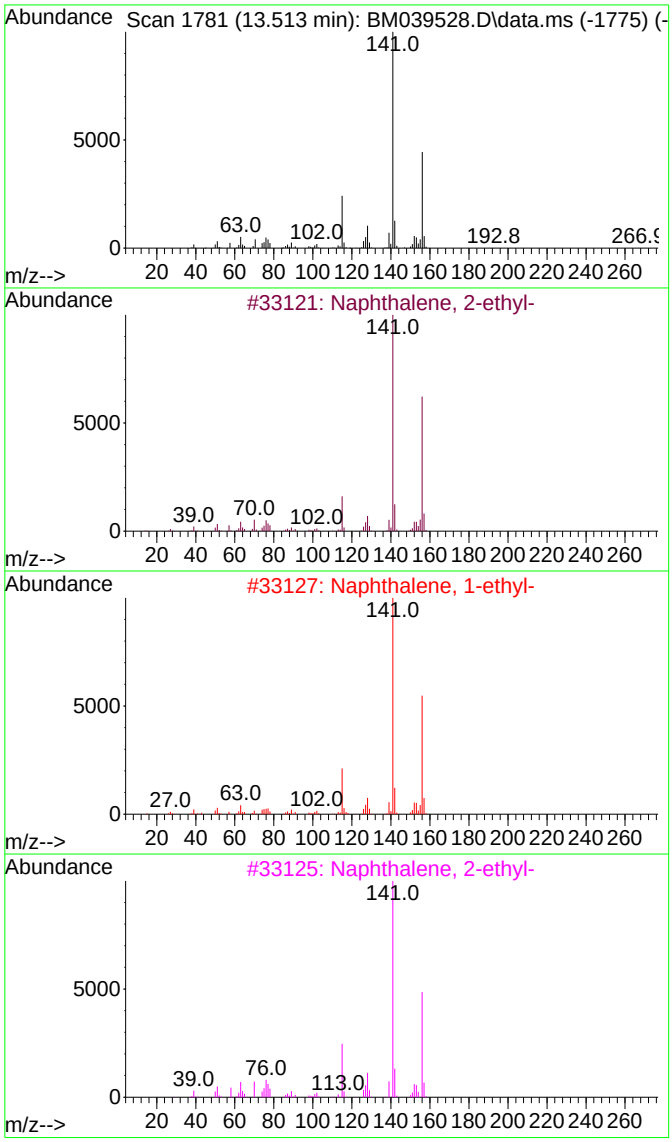
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

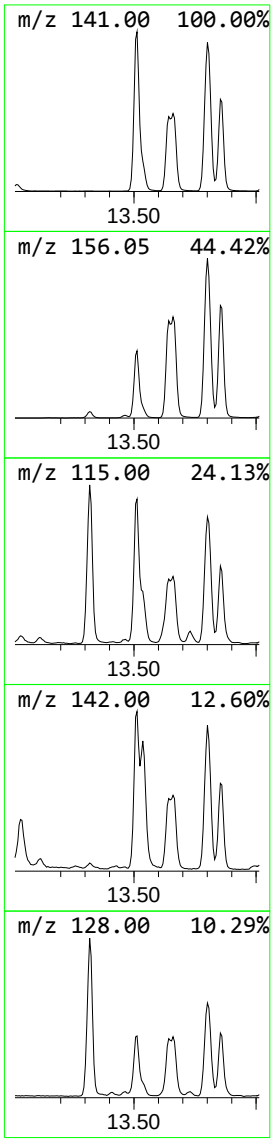
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Naphthalene, 2-ethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.513	24.69 ng/ul	4236250	Acenaphthene-d10	14.483

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	95
2	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	95
3	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	94
4	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	94
5	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	94





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Acq On : 20 Apr 2023 17:12
Operator : CG/JU
Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBN2

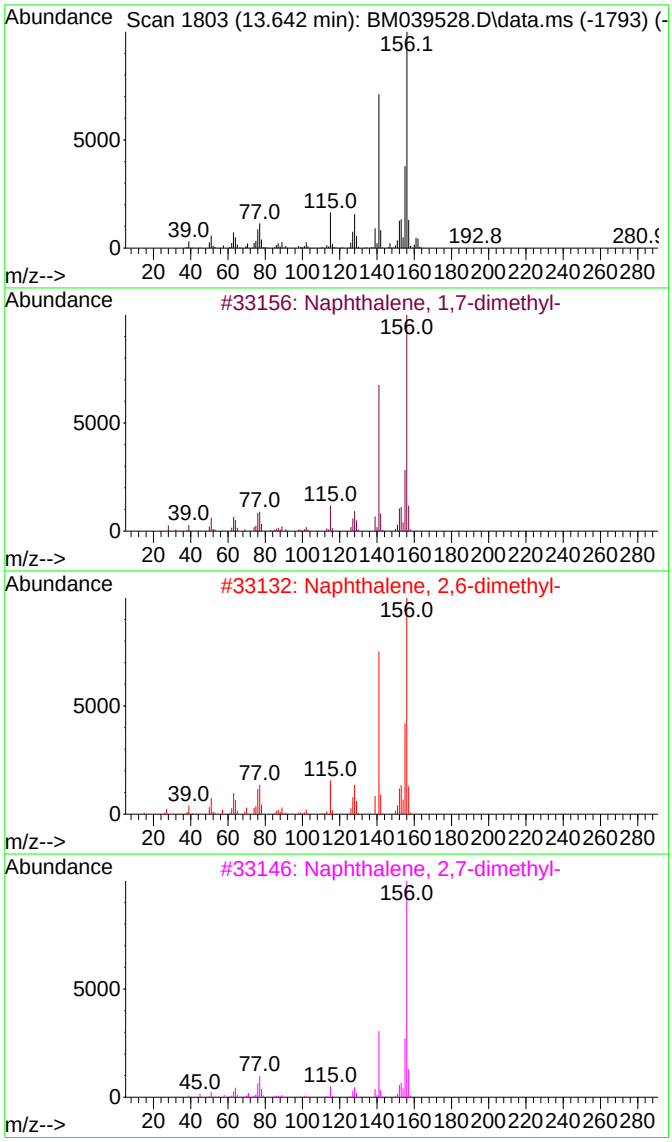
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

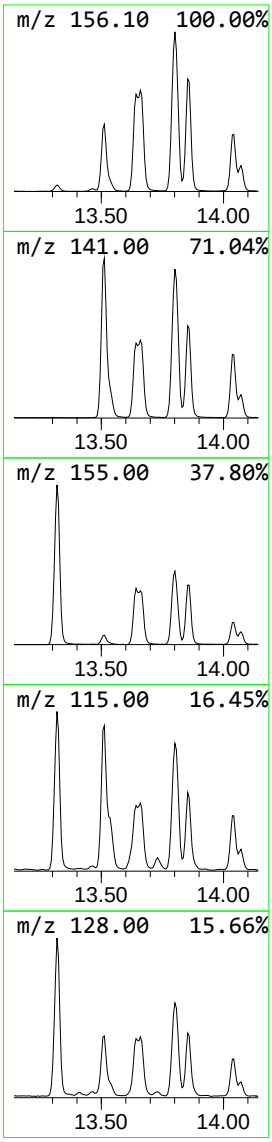
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Naphthalene, 1,7-dimethyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.642	15.75 ng/ul	2701750	Acenaphthene-d10	14.483

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
2	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96





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Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBN2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

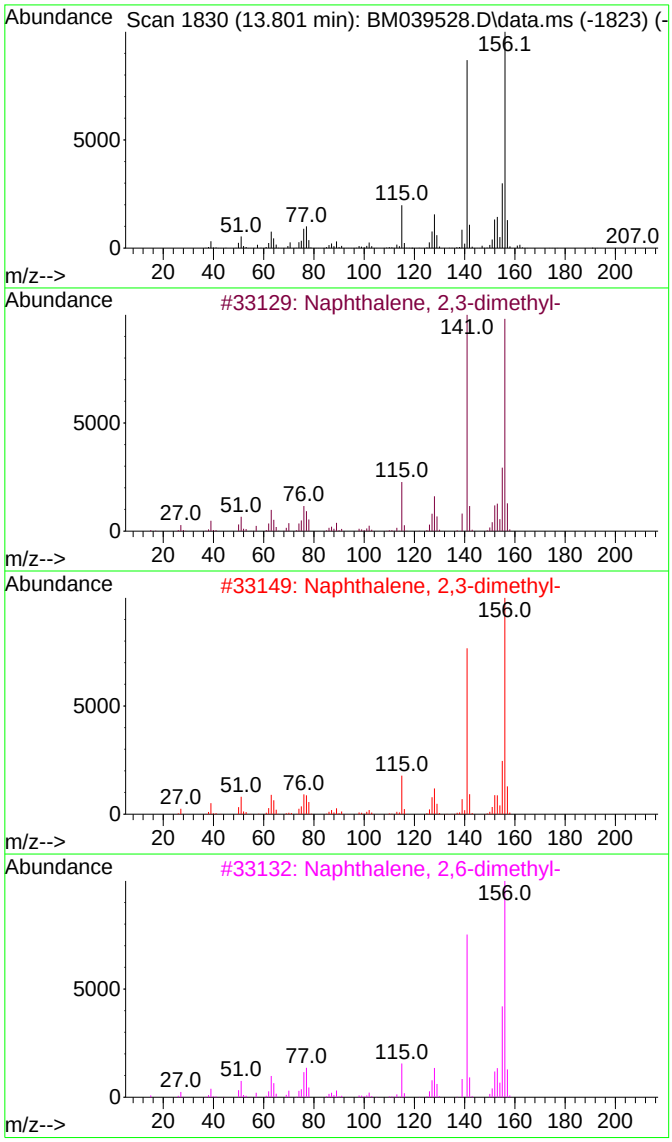
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

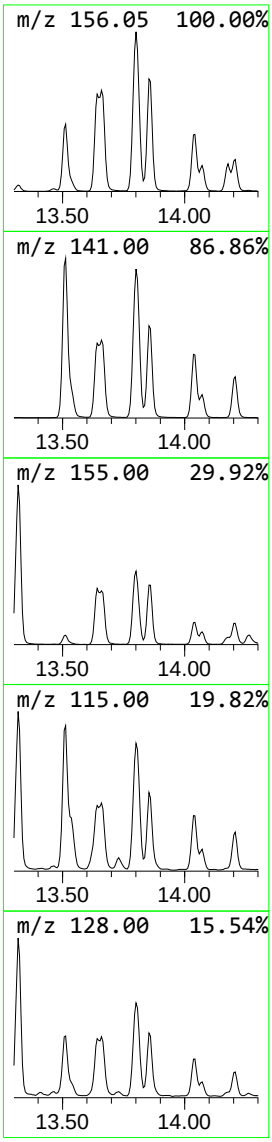
****

Peak Number 21 Naphthalene, 2,3-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.801	34.05 ng/ul	5842670	Acenaphthene-d10	14.483

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
4	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
5	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97





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Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBN2

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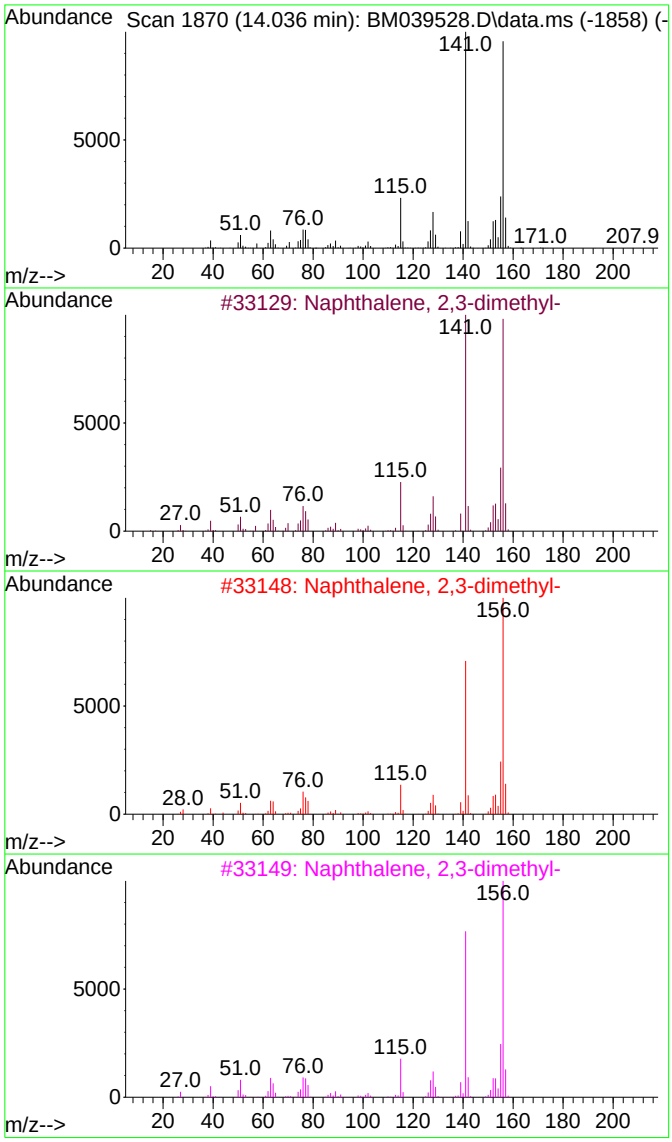
Quant Title : SVOA CALIBRATION

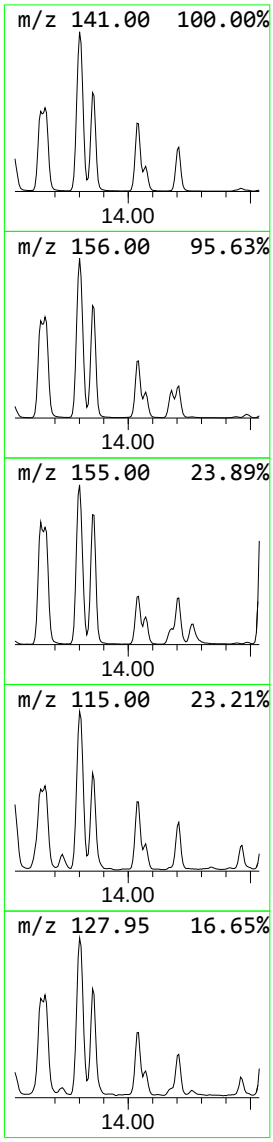
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 Naphthalene, 1,6-dimethyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.036	11.68 ng/ul	2004210	Acenaphthene-d10	14.483

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97





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Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBN2

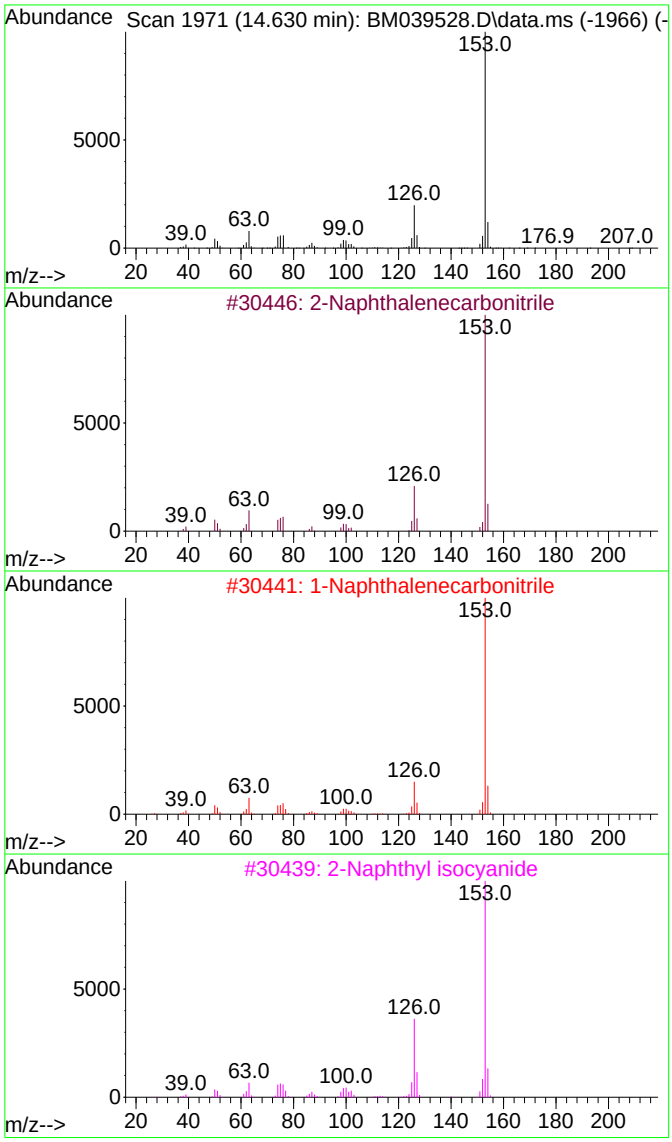
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

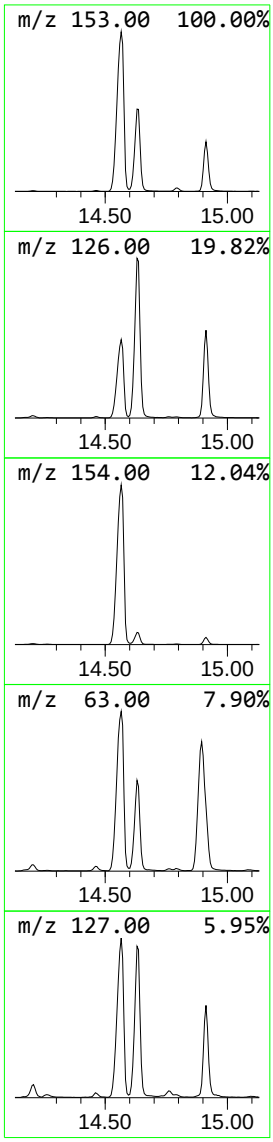
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 2-Naphthalenecarbonitrile Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.630	74.28 ng/ul	12744500	Acenaphthene-d10	14.483

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	97
2	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	97
3	2-Naphthyl isocyanide	153	C11H7N	010124-78-4	94
4	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	91
5	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	91





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Instrument :
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ClientSampleId :
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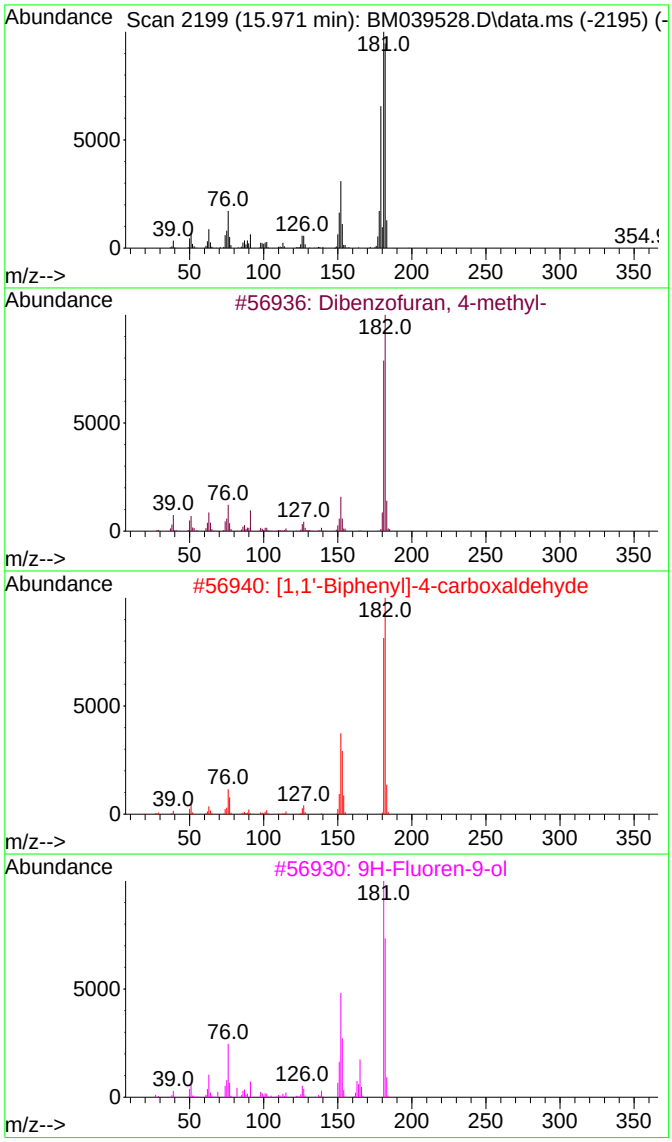
Quant Title : SVOA CALIBRATION

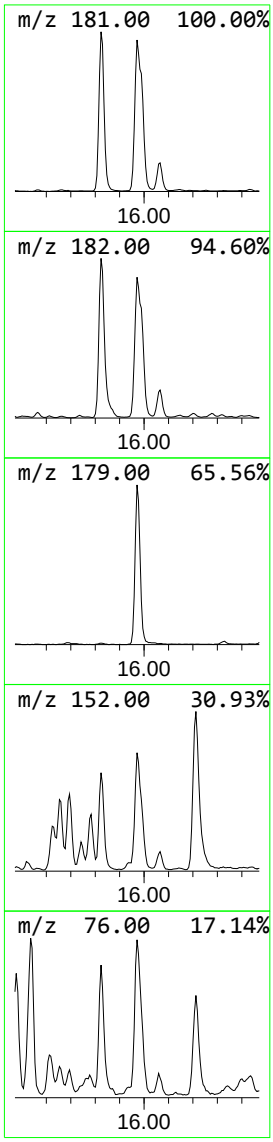
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 29 Dibenzofuran, 4-methyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.971	15.88 ng/ul	1983490	Phenanthrene-d10	17.236

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	81
2	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	76
3	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	68
4	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	68
5	9H-Fluoren-3-ol	182	C13H10O	006344-67-8	64





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Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DCBN2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

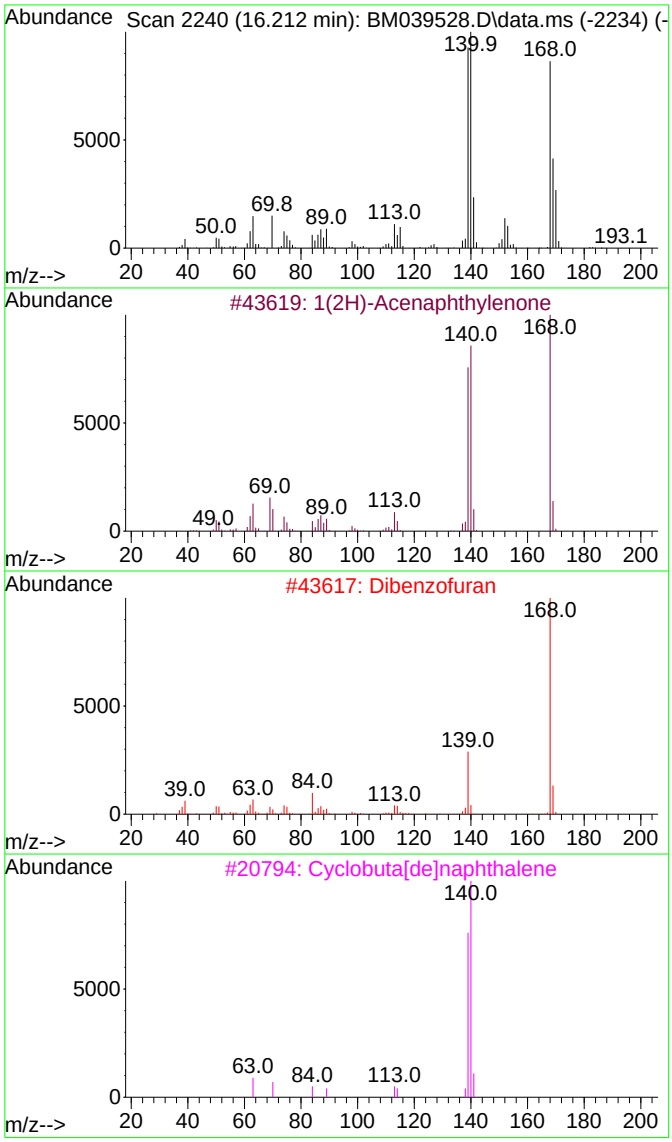
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

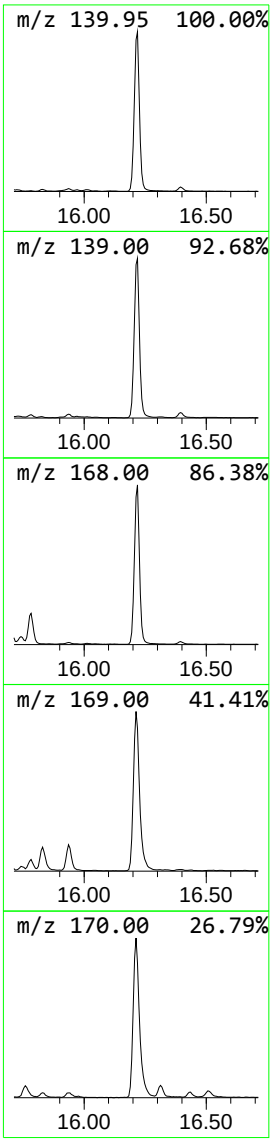
****

Peak Number 30 1(2H)-Acenaphthylenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.212	45.30 ng/ul	5658460	Phenanthrene-d10	17.236

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1(2H)-Acenaphthylenone	168	C12H8O	002235-15-6	94
2	Dibenzofuran	168	C12H8O	000132-64-9	43
3	Cyclobuta[de]naphthalene	140	C11H8	024973-91-9	43
4	1-Isoquinolinecarbonitrile, 3-me...	168	C11H8N2	022381-52-8	38
5	Benzoic acid, 2-chloro-, methyl ...	170	C8H7ClO2	000610-96-8	25





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Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DCBN2

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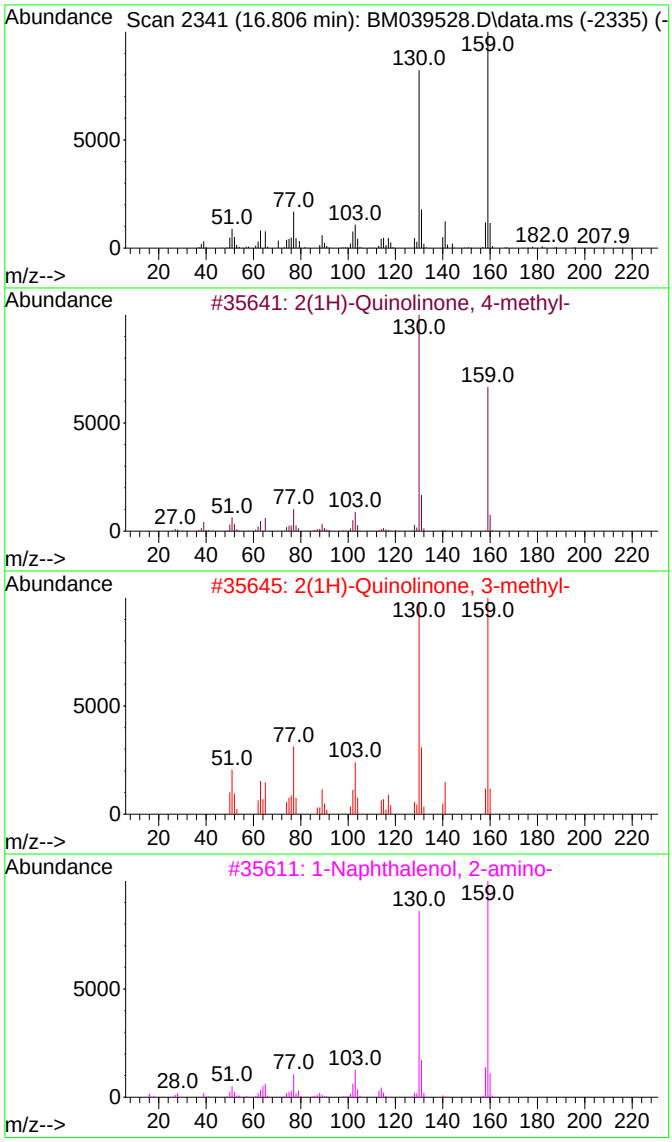
Quant Title : SVOA CALIBRATION

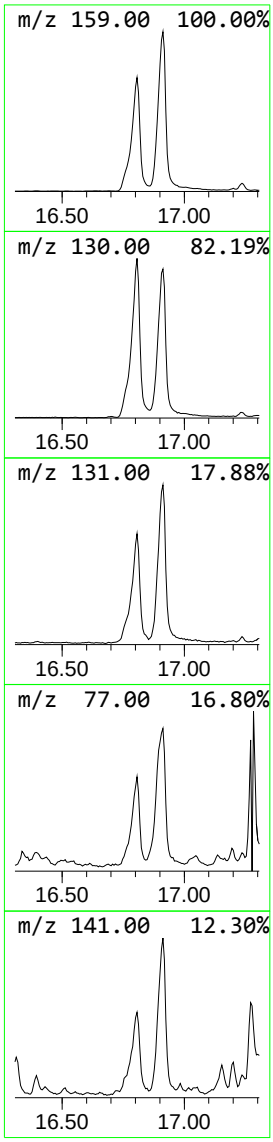
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 34 2(1H)-Quinolinone, 4-methyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.807	12.21 ng/ul	1525120	Phenanthrene-d10	17.236

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(1H)-Quinolinone, 4-methyl-	159	C10H9NO	000607-66-9	94
2	2(1H)-Quinolinone, 3-methyl-	159	C10H9NO	002721-59-7	93
3	1-Naphthalenol, 2-amino-	159	C10H9NO	000606-41-7	93
4	5-Amino-1-naphthol	159	C10H9NO	000083-55-6	91
5	2-(Aminomethylidene)-2,3-dihydro...	159	C10H9NO	053394-92-6	90





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Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBN2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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

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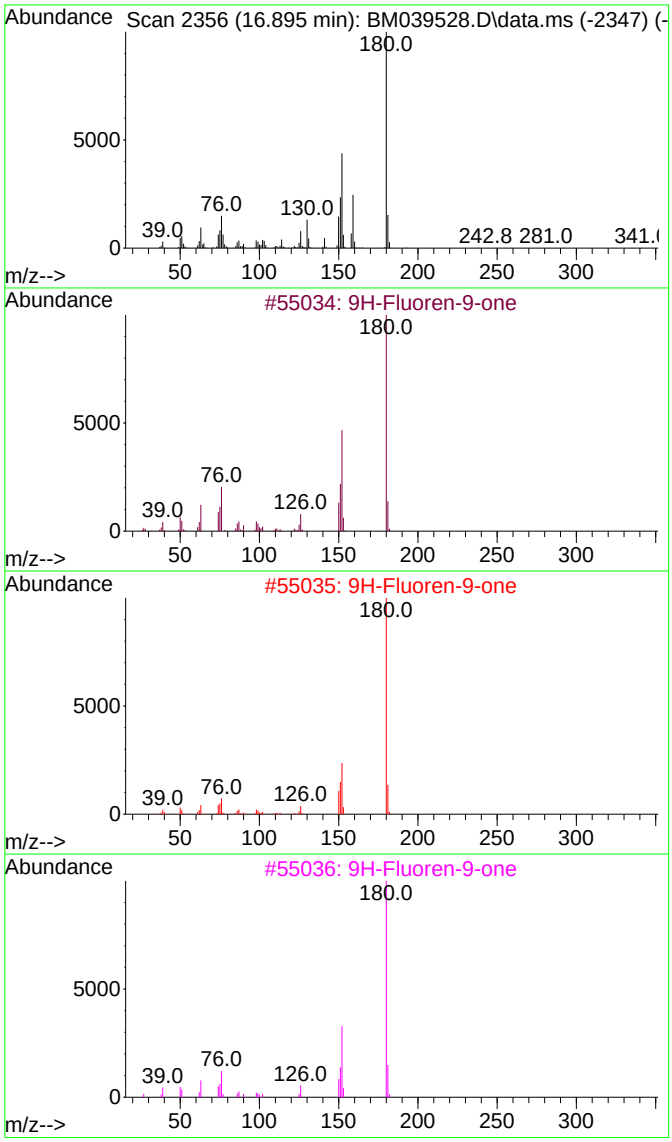
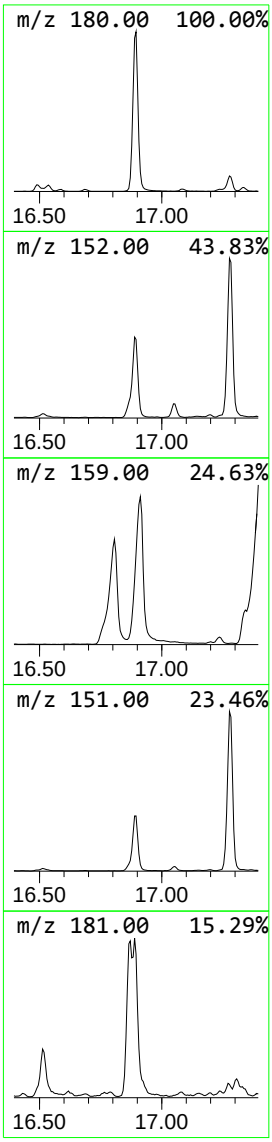
Peak Number 35 9H-Fluoren-9-one Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.895	39.63 ng/ul	4950560	Phenanthrene-d10	17.236

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluoren-9-one	180	C13H8O	000486-25-9	95
2	9H-Fluoren-9-one	180	C13H8O	000486-25-9	95
3	9H-Fluoren-9-one	180	C13H8O	000486-25-9	94
4	9H-Fluoren-9-one	180	C13H8O	000486-25-9	93
5	9H-Fluoren-9-one	180	C13H8O	000486-25-9	93





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Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

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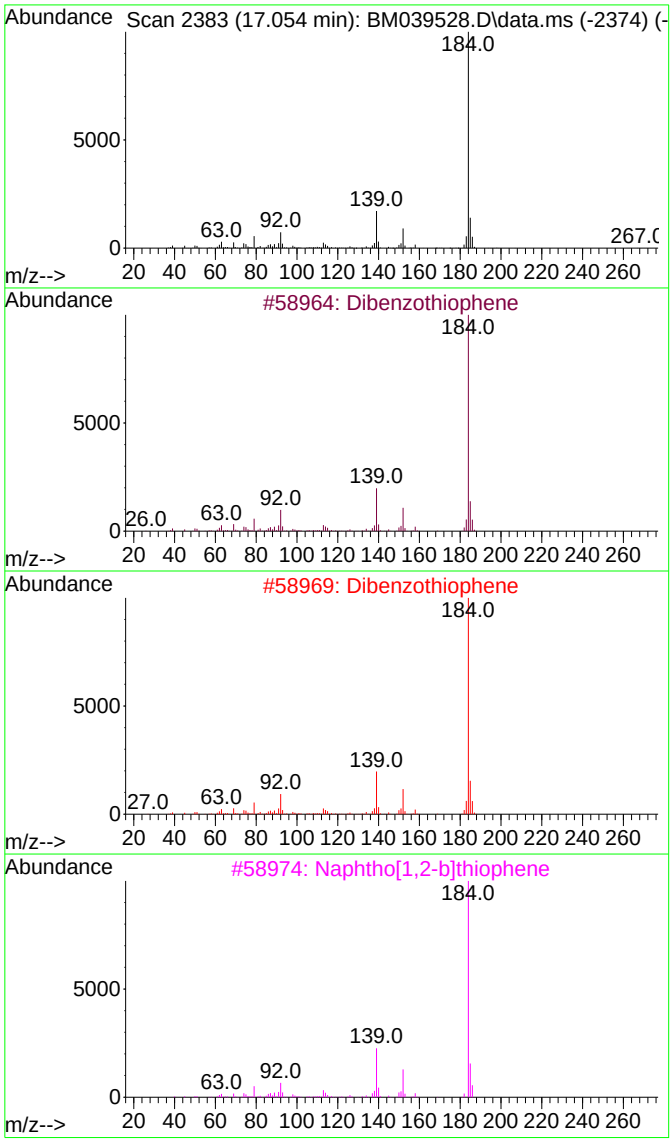
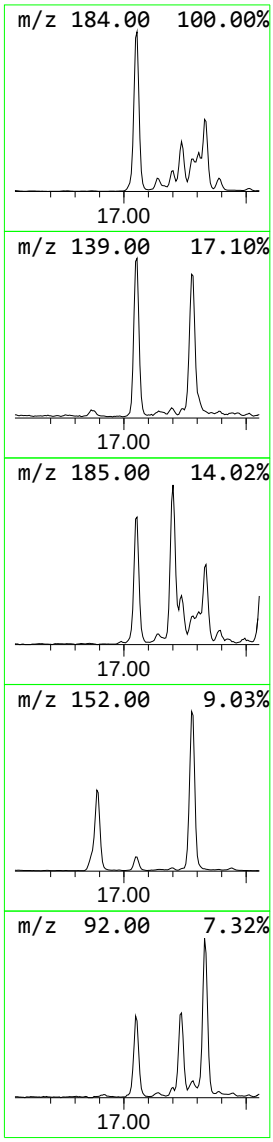
Peak Number 36 Dibenzothiophene Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.054	13.57 ng/ul	1694810	Phenanthrene-d10	17.236

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzothiophene	184	C12H8S	000132-65-0	97
2	Dibenzothiophene	184	C12H8S	000132-65-0	97
3	Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	97
4	Dibenzothiophene	184	C12H8S	000132-65-0	97
5	Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	96





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Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DCBN2

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Quant Title : SVOA CALIBRATION

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

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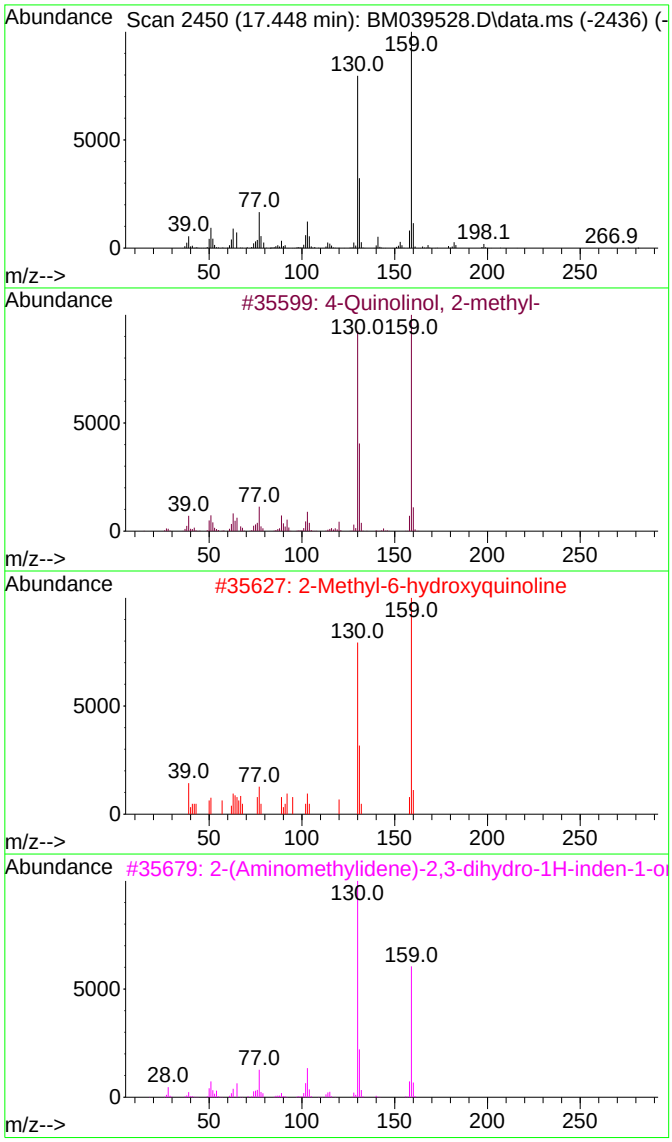
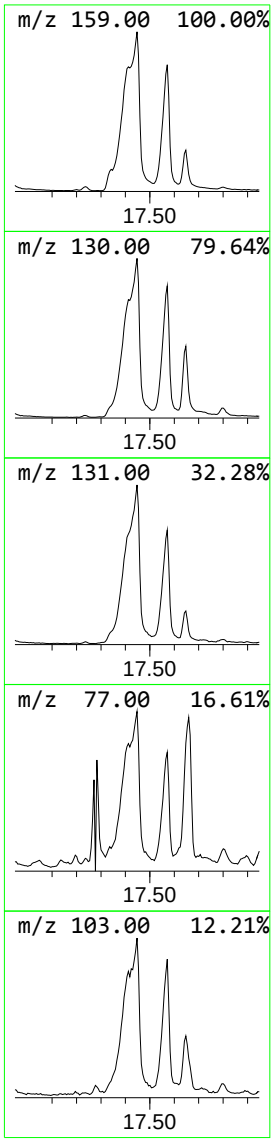
Peak Number 37 4-Quinolinol, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.448	51.62 ng/ul	6448330	Phenanthrene-d10	17.236

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Quinolinol, 2-methyl-	159	C10H9NO	000607-67-0	94
2	2-Methyl-6-hydroxyquinoline	159	C10H9NO	000613-21-8	91
3	2-(Aminomethylidene)-2,3-dihydro...	159	C10H9NO	053394-92-6	90
4	2-Methyl-6-hydroxyquinoline	159	C10H9NO	000613-21-8	90
5	6-Methyl-4(1H)-quinolinone	159	C10H9NO	023432-40-8	87





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ALS Vial : 9 Sample Multiplier: 1

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ClientSampleId :
DCBN2

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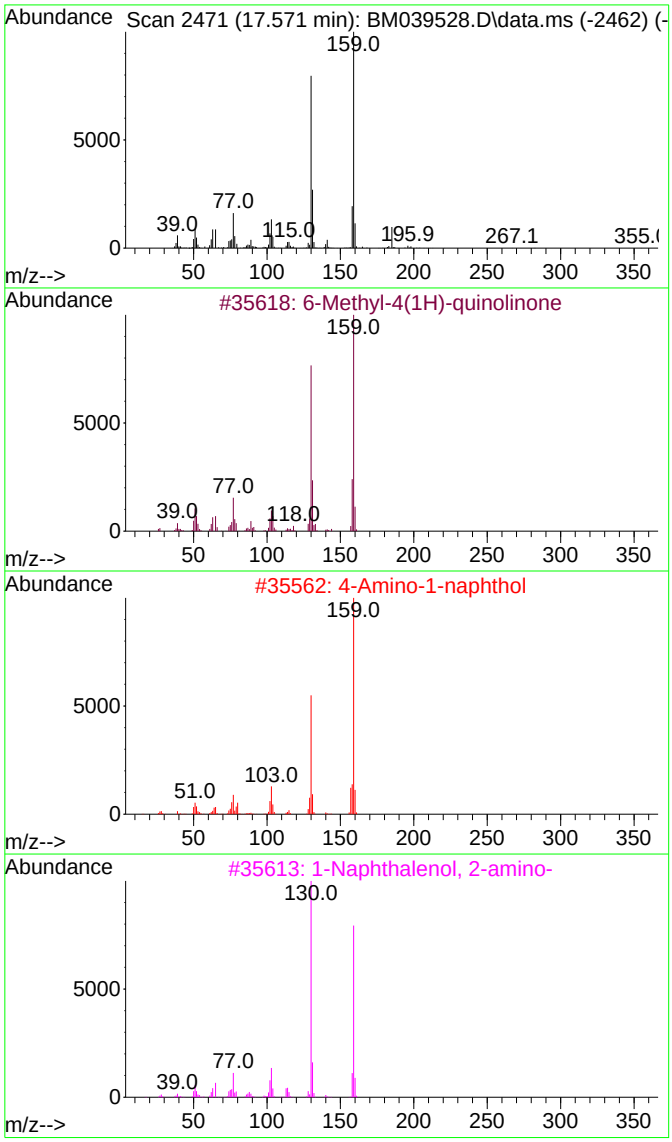
Quant Title : SVOA CALIBRATION

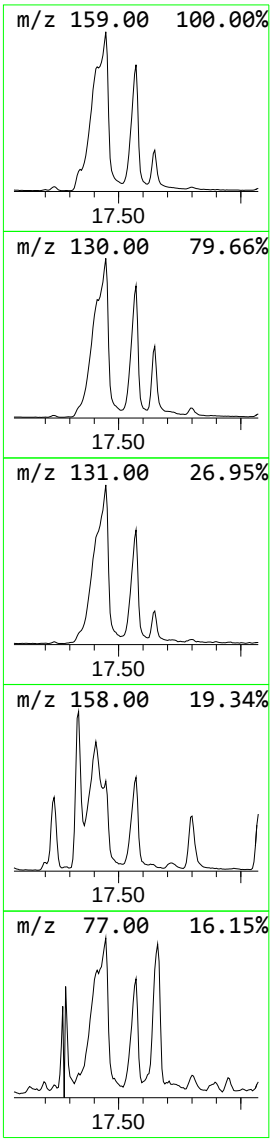
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 38 6-Methyl-4(1H)-quinolinone Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.571	18.00 ng/ul	2248510	Phenanthrene-d10	17.236

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Methyl-4(1H)-quinolinone	159	C10H9NO	023432-40-8	91
2	4-Amino-1-naphthol	159	C10H9NO	002834-90-4	87
3	1-Naphthalenol, 2-amino-	159	C10H9NO	000606-41-7	87
4	2-(Aminomethylidene)-2,3-dihydro...	159	C10H9NO	053394-92-6	87
5	1-Naphthalenol, 6-amino-	159	C10H9NO	023894-12-4	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042023\
Data File : BM039528.D
Acq On : 20 Apr 2023 17:12
Operator : CG/JU
Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBN2

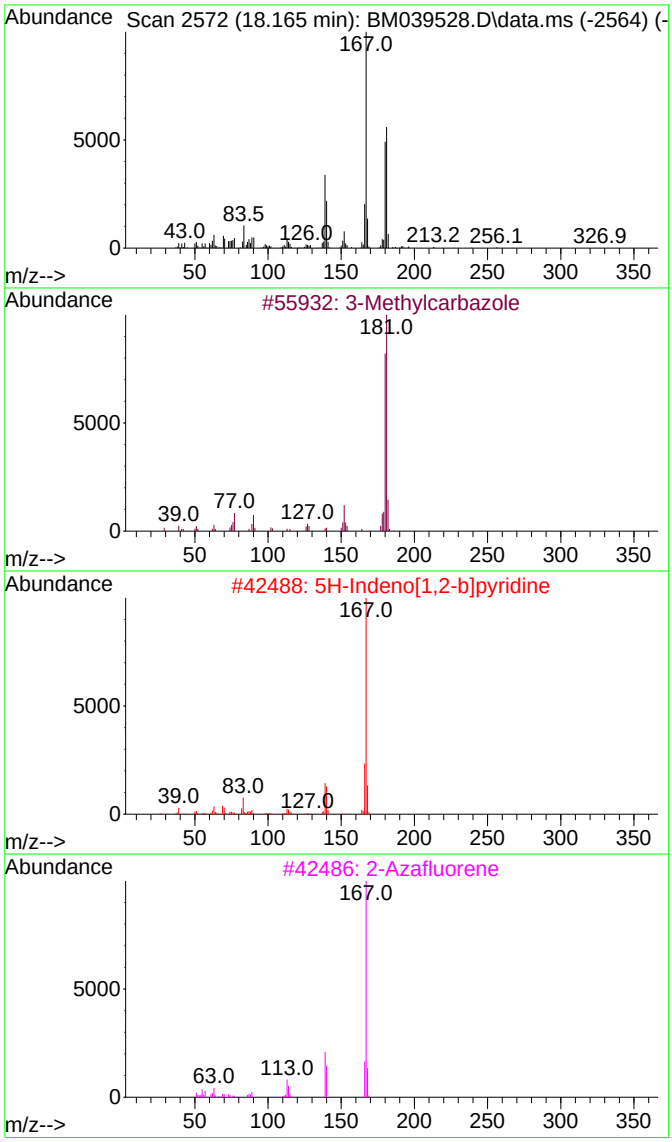
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

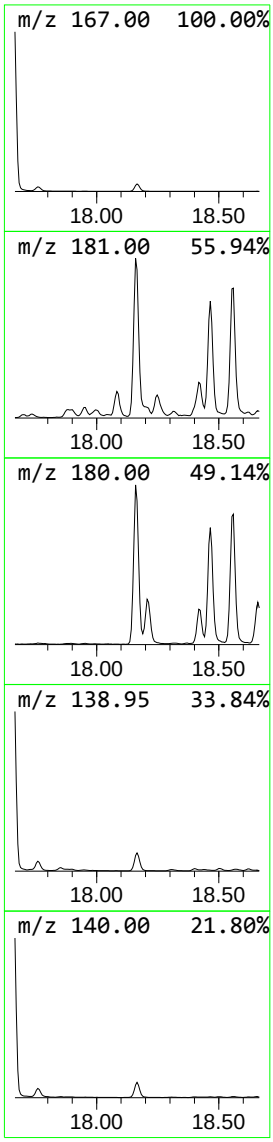
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 43 3-Methylcarbazole Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.165	27.93 ng/ul	3489580	Phenanthrene-d10	17.236

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Methylcarbazole	181	C13H11N	004630-20-0	70
2	5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	60
3	2-Azafluorene	167	C12H9N	000244-40-6	53
4	Carbazole	167	C12H9N	000086-74-8	53
5	Carbazole	167	C12H9N	000086-74-8	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042023\
Data File : BM039528.D
Acq On : 20 Apr 2023 17:12
Operator : CG/JU
Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBN2

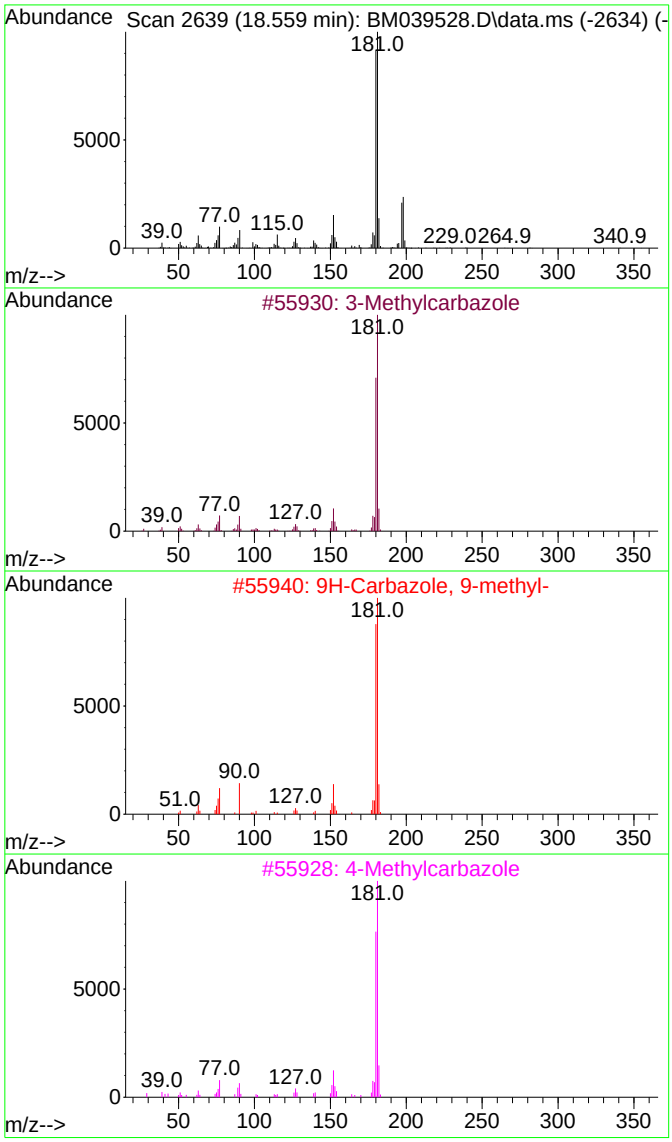
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

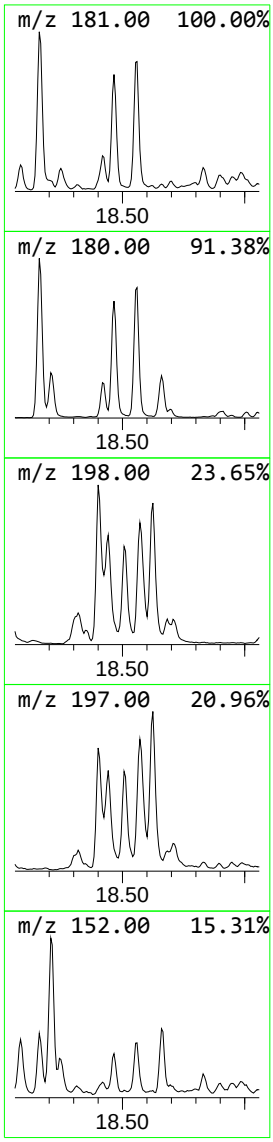
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 47 9H-Carbazole, 9-methyl- Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.559	9.36 ng/ul	1168780	Phenanthrene-d10	17.236

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Methylcarbazole	181	C13H11N	004630-20-0	95
2	9H-Carbazole, 9-methyl-	181	C13H11N	001484-12-4	93
3	4-Methylcarbazole	181	C13H11N	003770-48-7	92
4	9H-Carbazole, 2-methyl-	181	C13H11N	003652-91-3	92
5	1-Methylcarbazole	181	C13H11N	006510-65-2	92





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042023\
Data File : BM039528.D
Acq On : 20 Apr 2023 17:12
Operator : CG/JU
Sample : 02296-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBN2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.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
Ethylbenzene	5.307	11.7	ng/ul	703246	1	7.825	1198070	20.0
p-Xylene	5.442	33.4	ng/ul	1997700	1	7.825	1198070	20.0
o-Xylene	5.813	17.6	ng/ul	1052810	1	7.825	1198070	20.0
.alpha.-Pinene	6.484	16.9	ng/ul	1015480	1	7.825	1198070	20.0
.beta.-Pinene	6.789	9.6	ng/ul	573455	1	7.825	1198070	20.0
Benzene, 1-ethy...	6.901	38.3	ng/ul	2293740	1	7.825	1198070	20.0
Mesitylene	6.960	15.7	ng/ul	938525	1	7.825	1198070	20.0
Benzene, 1,2,3-...	7.466	119.4	ng/ul	7154230	1	7.825	1198070	20.0
Benzene, 1,2,4-...	7.936	59.7	ng/ul	3575040	1	7.825	1198070	20.0
D-Limonene	8.036	44.0	ng/ul	2632890	1	7.825	1198070	20.0
Indane	8.201	401.8	ng/ul	24070500	1	7.825	1198070	20.0
Indene	8.372	536.1	ng/ul	32113000	1	7.825	1198070	20.0
Benzene, 2-ethy...	8.460	11.4	ng/ul	682976	1	7.825	1198070	20.0
Benzene, 4-ethy...	8.931	13.8	ng/ul	824191	1	7.825	1198070	20.0
Fenchone	9.072	22.2	ng/ul	1328280	1	7.825	1198070	20.0
Benzofuran, 2-m...	9.183	58.0	ng/ul	3476720	1	7.825	1198070	20.0
Naphthalene, 2-...	13.513	24.7	ng/ul	4236250	3	14.483	3431470	20.0
Naphthalene, 1,...	13.642	15.8	ng/ul	2701750	3	14.483	3431470	20.0
Naphthalene, 2,...	13.801	34.0	ng/ul	5842670	3	14.483	3431470	20.0
Naphthalene, 1,...	14.036	11.7	ng/ul	2004210	3	14.483	3431470	20.0
2-Naphthaleneca...	14.630	74.3	ng/ul	12744500	3	14.483	3431470	20.0
Dibenzofuran, 4...	15.971	15.9	ng/ul	1983490	4	17.236	2498470	20.0
1(2H)-Acenaphth...	16.212	45.3	ng/ul	5658460	4	17.236	2498470	20.0
2(1H)-Quinolino...	16.807	12.2	ng/ul	1525120	4	17.236	2498470	20.0
9H-Fluoren-9-one	16.895	39.6	ng/ul	4950560	4	17.236	2498470	20.0
Dibenzothiophene	17.054	13.6	ng/ul	1694810	4	17.236	2498470	20.0
4-Quinolinol, 2...	17.448	51.6	ng/ul	6448330	4	17.236	2498470	20.0
6-Methyl-4(1H)-...	17.571	18.0	ng/ul	2248510	4	17.236	2498470	20.0
3-Methylcarbazole	18.165	27.9	ng/ul	3489580	4	17.236	2498470	20.0
9H-Carbazole, 9...	18.559	9.4	ng/ul	1168780	4	17.236	2498470	20.0