

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041823\
 Data File : BM039517.D
 Acq On : 20 Apr 2023 04:53
 Operator : CG/JU
 Sample : 02296-15DL3 900X
 Misc :
 ALS Vial : 69 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DCBM0DL3

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: BM039517.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.966	155	158	168	rVB	15519	26652	0.40%	0.053%
2	6.484	580	586	598	rBV	547970	860518	12.76%	1.719%
3	6.754	625	632	634	rBV4	14924	22552	0.33%	0.045%
4	6.790	634	638	645	rVB	47100	75833	1.12%	0.151%
5	6.907	654	658	665	rBV2	11144	20320	0.30%	0.041%
6	7.043	676	681	687	rVB	19083	29913	0.44%	0.060%
7	7.248	710	716	724	rVB	140028	217253	3.22%	0.434%
8	7.472	748	754	760	rBV2	37490	68008	1.01%	0.136%
9	7.554	763	768	776	rVB8	7090	20134	0.30%	0.040%
10	7.825	807	814	829	rBV	659028	1208235	17.91%	2.413%
11	7.954	829	836	844	rVV	344250	557440	8.26%	1.113%
12	8.037	844	850	858	rVB	193353	297354	4.41%	0.594%
13	8.195	870	877	890	rVB	77024	132757	1.97%	0.265%
14	8.360	899	905	919	rBV	142145	282114	4.18%	0.563%
15	8.466	919	923	933	rVB2	13073	28142	0.42%	0.056%
16	8.942	998	1004	1011	rBV4	63155	120017	1.78%	0.240%
17	9.013	1011	1016	1022	rVV2	13803	28395	0.42%	0.057%
18	9.195	1040	1047	1054	rVB8	9116	20275	0.30%	0.040%
19	9.319	1061	1068	1075	rBV4	14196	36817	0.55%	0.074%
20	9.454	1087	1091	1097	rVB	11075	19420	0.29%	0.039%
21	9.519	1097	1102	1108	rVB	12561	21472	0.32%	0.043%
22	9.884	1158	1164	1172	rVB	24802	46897	0.70%	0.094%
23	9.995	1175	1183	1187	rBV	170273	300141	4.45%	0.599%
24	10.131	1201	1206	1210	rBV2	17230	34091	0.51%	0.068%
25	10.166	1210	1212	1221	rVV6	13612	21880	0.32%	0.044%
26	10.254	1221	1227	1236	rVB7	7309	17352	0.26%	0.035%
27	10.631	1281	1291	1295	rBV	1098209	1926597	28.56%	3.848%
28	10.683	1295	1300	1316	rVV	3875788	6746463	100.00%	13.473%
29	10.807	1316	1321	1335	rVB	70522	169633	2.51%	0.339%
30	11.083	1362	1368	1377	rVB	6280	16656	0.25%	0.033%
31	11.295	1399	1404	1410	rVB6	7694	15028	0.22%	0.030%
32	11.554	1441	1448	1456	rVB4	8467	18686	0.28%	0.037%
33	11.748	1473	1481	1484	rVB8	6045	12574	0.19%	0.025%
34	12.195	1550	1557	1563	rBV	32891	54850	0.81%	0.110%
35	12.295	1563	1574	1590	rBV	1985463	3410344	50.55%	6.811%
36	12.419	1590	1595	1602	rVV2	35414	93473	1.39%	0.187%

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If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
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37	12.519	1603	1612	1623	rVB	899496	1527246	22.64%	3.050%
38	13.313	1741	1747	1760	rBV	404646	704441	10.44%	1.407%
39	13.513	1774	1781	1791	rVB	148702	283190	4.20%	0.566%
40	13.660	1796	1806	1823	rBV4	161669	474175	7.03%	0.947%

41	13.801	1823	1830	1835	rBV	236938	437910	6.49%	0.875%
42	13.854	1835	1839	1854	rVB	155844	279135	4.14%	0.557%
43	14.042	1864	1871	1874	rBV	94163	156571	2.32%	0.313%
44	14.207	1894	1899	1907	rVB	58190	95984	1.42%	0.192%
45	14.377	1923	1928	1932	rBV2	8951	13220	0.20%	0.026%

46	14.483	1935	1946	1951	rVV	1841145	2922756	43.32%	5.837%
47	14.542	1951	1956	1967	rVV	2069661	3220602	47.74%	6.432%
48	14.630	1967	1971	1976	rVV	34812	73477	1.09%	0.147%
49	14.689	1976	1981	1991	rVB3	32384	80768	1.20%	0.161%
50	14.789	1991	1998	2003	rBV	80515	125474	1.86%	0.251%

51	14.883	2003	2014	2023	rBV	1143960	1772920	26.28%	3.541%
52	14.960	2024	2027	2045	rVB3	48706	108323	1.61%	0.216%
53	15.101	2045	2051	2056	rBV4	28570	50909	0.75%	0.102%
54	15.154	2056	2060	2071	rVB3	28340	49110	0.73%	0.098%
55	15.271	2073	2080	2083	rBV	22858	46583	0.69%	0.093%

56	15.413	2100	2104	2107	rBV2	12721	19749	0.29%	0.039%
57	15.454	2107	2111	2118	rVV4	19121	39028	0.58%	0.078%
58	15.530	2118	2124	2136	rVV	1215844	1846955	27.38%	3.688%
59	15.624	2136	2140	2142	rVV	49640	73896	1.10%	0.148%
60	15.654	2142	2145	2148	rVV	73557	115374	1.71%	0.230%

61	15.695	2148	2152	2157	rVV	153923	246672	3.66%	0.493%
62	15.742	2157	2160	2163	rVV2	36782	55194	0.82%	0.110%
63	15.783	2163	2167	2170	rVV	71859	114017	1.69%	0.228%
64	15.830	2170	2175	2185	rVB	155200	255893	3.79%	0.511%
65	15.977	2191	2200	2208	rBV	134475	315335	4.67%	0.630%

66	16.065	2211	2215	2222	rVB3	29140	57205	0.85%	0.114%
67	16.189	2230	2236	2237	rBV4	12358	16940	0.25%	0.034%
68	16.213	2237	2240	2246	rVB7	13275	24383	0.36%	0.049%
69	16.330	2252	2260	2268	rBV	65809	118681	1.76%	0.237%
70	16.542	2284	2296	2300	rBV2	79200	190962	2.83%	0.381%

71	16.589	2300	2304	2309	rVV3	28512	45638	0.68%	0.091%
72	16.689	2317	2321	2326	rVB2	32143	51416	0.76%	0.103%
73	16.765	2326	2334	2337	rBV3	32605	68056	1.01%	0.136%
74	16.807	2337	2341	2345	rVB3	29541	36604	0.54%	0.073%
75	16.889	2351	2355	2362	rVB8	13147	24813	0.37%	0.050%

76	16.983	2362	2371	2376	rBV2	46990	114188	1.69%	0.228%
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77	17.054	2378	2383	2395	rVB	169264	270002	4.00%	0.539%
78	17.230	2407	2413	2417	rBV	2158352	3242996	48.07%	6.476%
79	17.277	2417	2421	2430	rVV	2633083	3914607	58.02%	7.818%
80	17.371	2431	2437	2447	rVB	219567	413140	6.12%	0.825%
81	17.654	2479	2485	2495	rBV	59875	160493	2.38%	0.321%
82	17.754	2499	2502	2509	rVB2	30313	47522	0.70%	0.095%
83	17.959	2533	2537	2543	rBV5	9181	15099	0.22%	0.030%
84	18.112	2558	2563	2567	rBV	105658	165114	2.45%	0.330%
85	18.159	2567	2571	2579	rVB	121781	193286	2.86%	0.386%
86	18.248	2582	2586	2590	rBV2	23633	32525	0.48%	0.065%
87	18.312	2591	2597	2601	rBV	198365	334871	4.96%	0.669%
88	18.424	2613	2616	2620	rBV4	10505	12510	0.19%	0.025%
89	18.618	2644	2649	2654	rBV	65170	99114	1.47%	0.198%
90	18.801	2678	2680	2685	rVB6	7902	11183	0.17%	0.022%
91	18.859	2685	2690	2696	rBV4	17603	34309	0.51%	0.069%
92	18.918	2696	2700	2703	rVV3	11317	14858	0.22%	0.030%
93	19.030	2715	2719	2724	rBV2	19264	36726	0.54%	0.073%
94	19.295	2758	2764	2773	rBV	887610	1235210	18.31%	2.467%
95	19.542	2803	2806	2814	rVB2	19852	34248	0.51%	0.068%
96	19.659	2816	2826	2835	rBV	516431	1000025	14.82%	1.997%
97	20.159	2908	2911	2919	rVB3	50461	69859	1.04%	0.140%
98	20.259	2924	2928	2933	rBV	57084	88413	1.31%	0.177%
99	21.418	3120	3125	3143	rBV	1871713	2988586	44.30%	5.968%
100	23.783	3521	3527	3543	rVB2	1143578	2426971	35.97%	4.847%

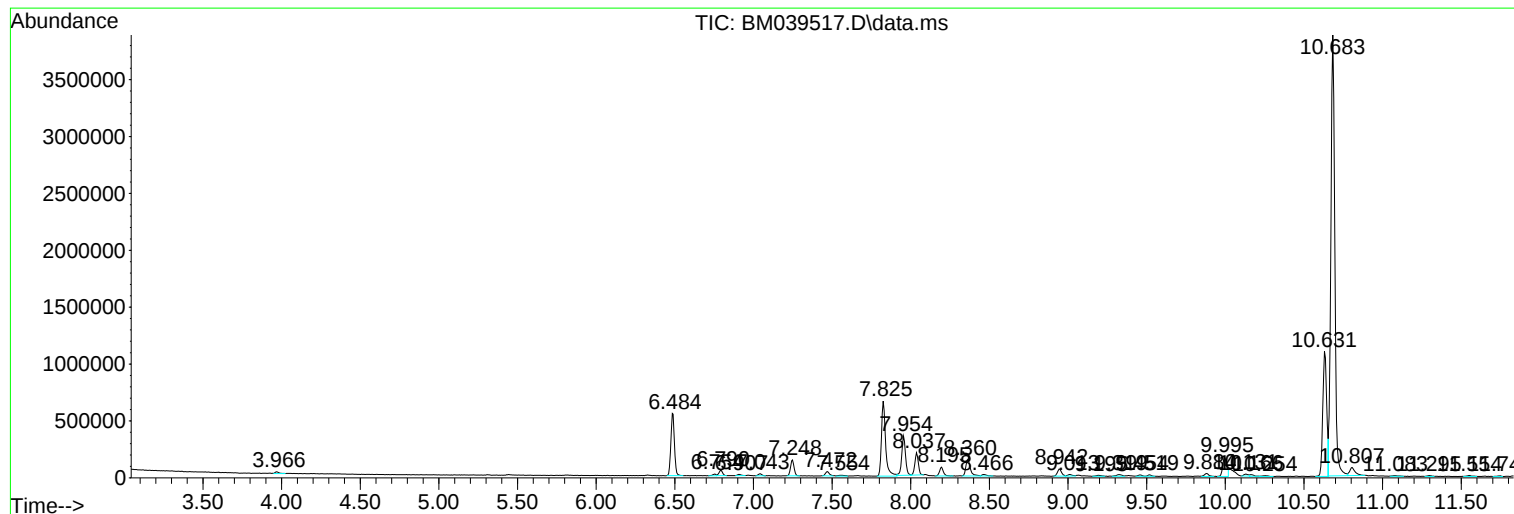
Sum of corrected areas: 50073746

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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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ALS Vial : 69 Sample Multiplier: 1

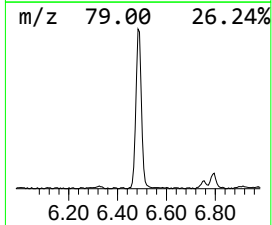
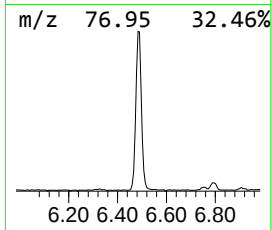
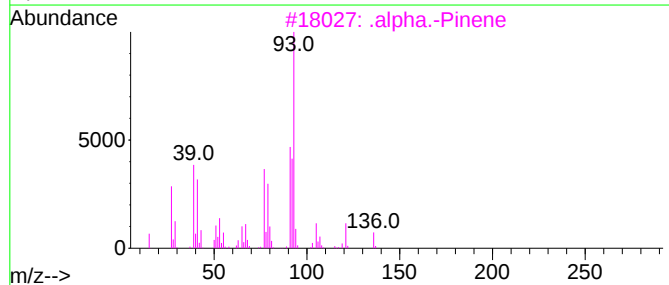
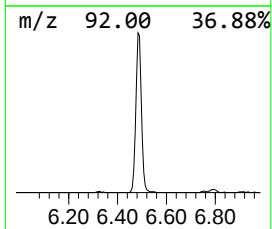
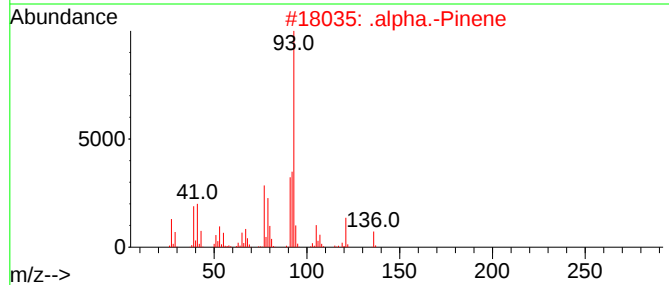
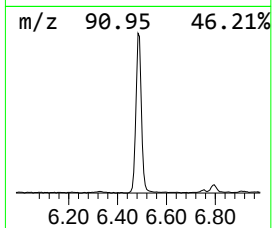
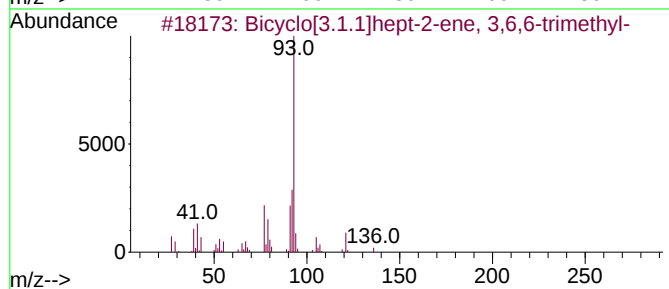
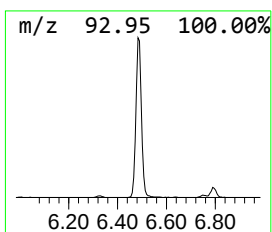
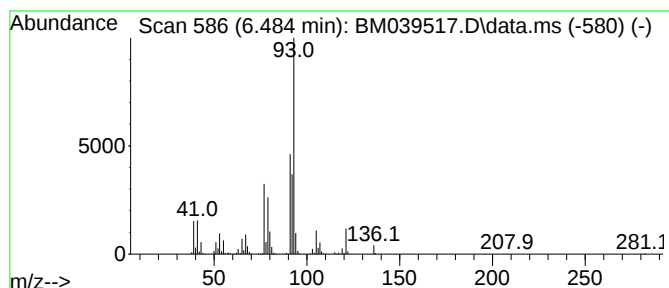
Instrument :
BNA_M
ClientSampleId :
DCBM0DL3

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 1 Bicyclo[3.1.1]hept-2-ene, 3... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
6.484	14.24 ng/ul	860518	1,4-Dichlorobenzene-d4	7.825	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[3.1.1]hept-2-ene, 3,6,6-...	136	C10H16	004889-83-2	94
2	.alpha.-Pinene	136	C10H16	000080-56-8	94
3	.alpha.-Pinene	136	C10H16	000080-56-8	94
4	(1R)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-70-8	94
5	3-Carene	136	C10H16	013466-78-9	91



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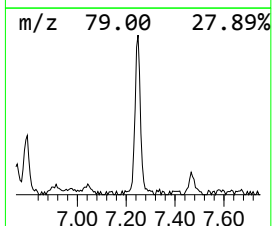
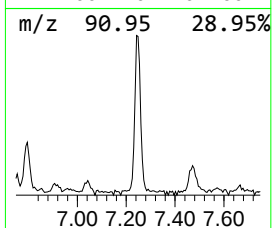
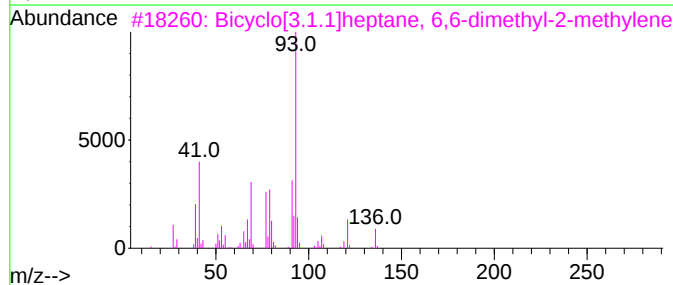
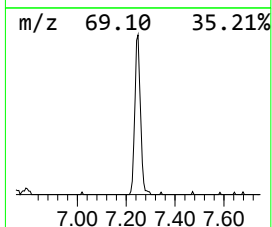
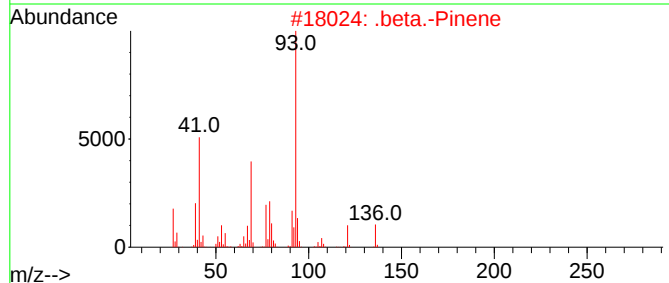
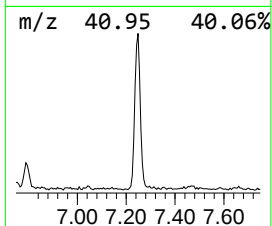
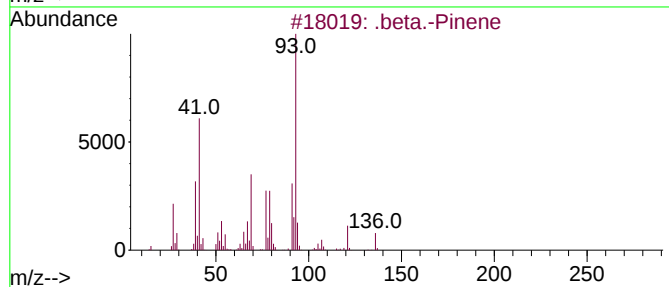
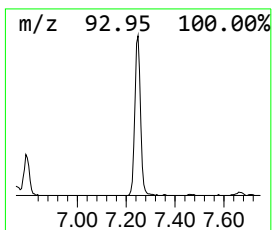
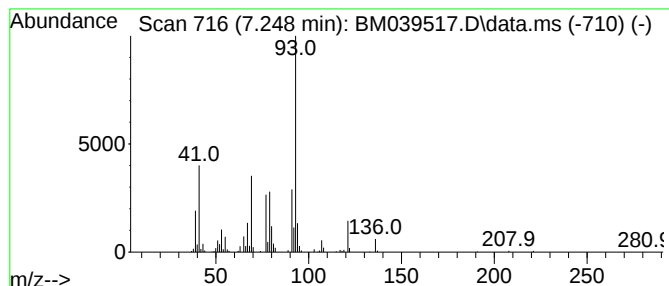
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 .beta.-Pinene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.248	3.60 ng/ul	217253	1,4-Dichlorobenzene-d4	7.825

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	beta.-Pinene	136	C10H16	000127-91-3	96	
2	.	beta.-Pinene	136	C10H16	000127-91-3	91	
3	Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	91		
4	.	beta.-Pinene	136	C10H16	000127-91-3	91	
5	Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	91		



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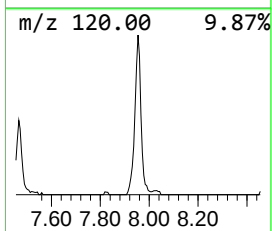
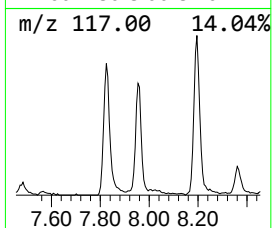
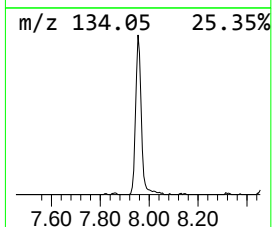
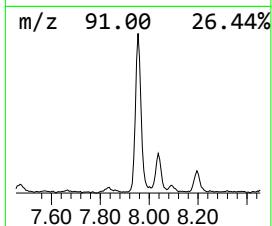
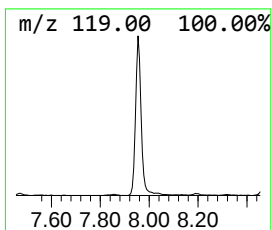
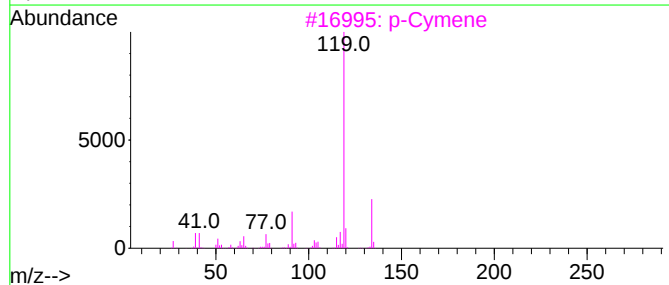
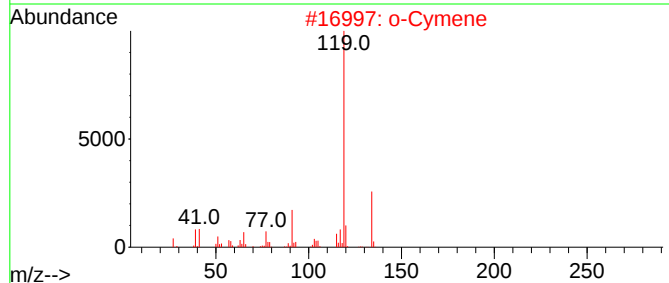
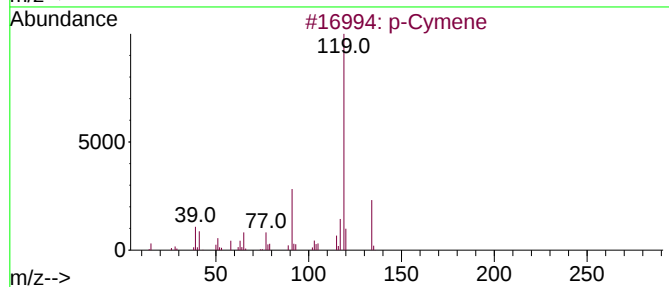
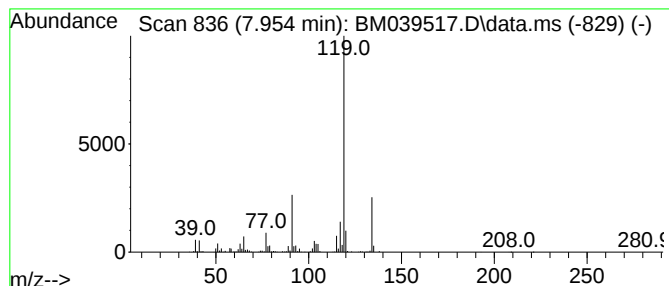
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 p-Cymene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.954	9.23 ng/ul	557440	1,4-Dichlorobenzene-d4	7.825

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene			134	C10H14	000099-87-6	97
2	o-Cymene			134	C10H14	000527-84-4	94
3	p-Cymene			134	C10H14	000099-87-6	94
4	o-Cymene			134	C10H14	000527-84-4	94
5	Benzene, 1-methyl-3-(1-methyleth...			134	C10H14	000535-77-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041823\
Data File : BM039517.D
Acq On : 20 Apr 2023 04:53
Operator : CG/JU
Sample : 02296-15DL3 900X
Misc :
ALS Vial : 69 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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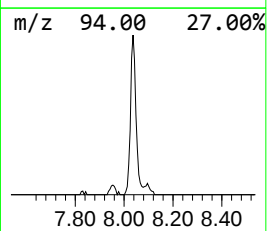
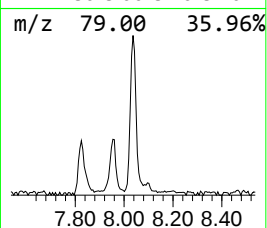
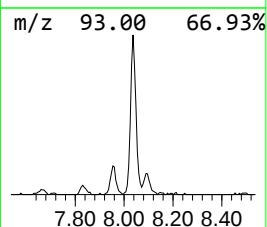
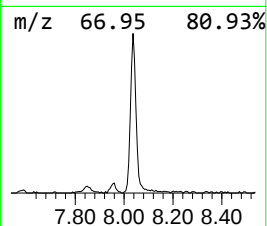
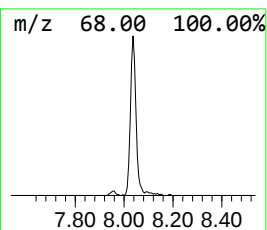
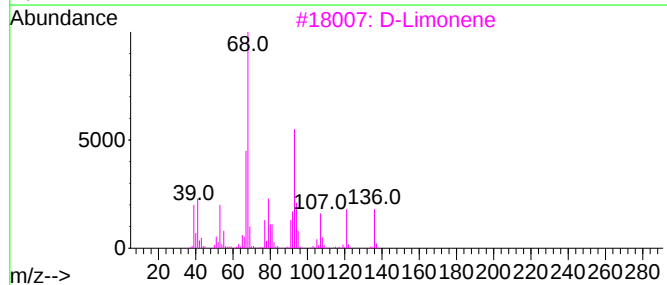
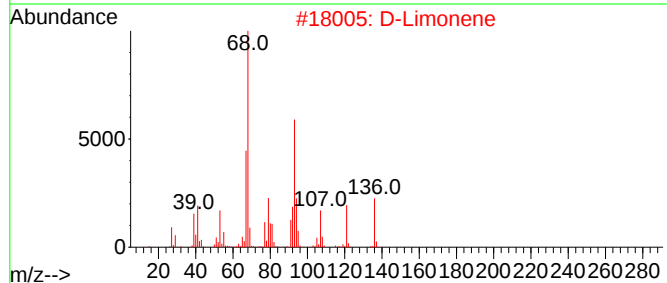
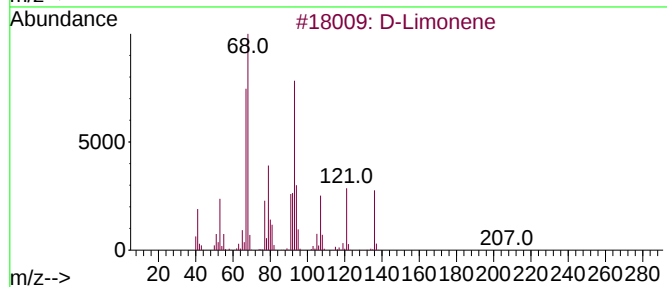
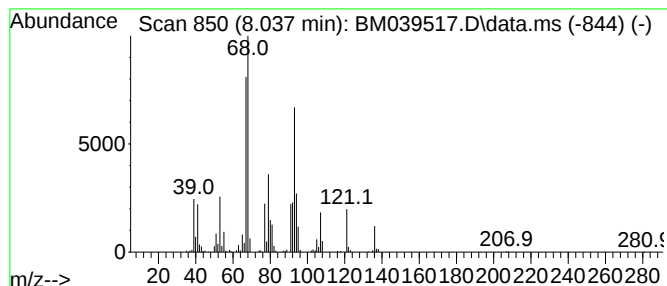
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 D-Limonene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.037	4.92 ng/ul	297354	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	D-Limonene			136	C10H16	005989-27-5	95
3	D-Limonene			136	C10H16	005989-27-5	94
4	D-Limonene			136	C10H16	005989-27-5	93
5	Limonene			136	C10H16	000138-86-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041823\
Data File : BM039517.D
Acq On : 20 Apr 2023 04:53
Operator : CG/JU
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ClientSampleId :
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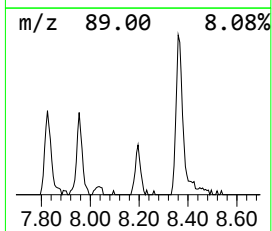
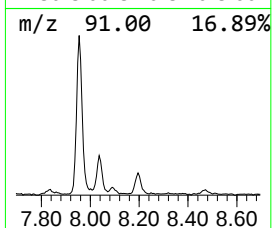
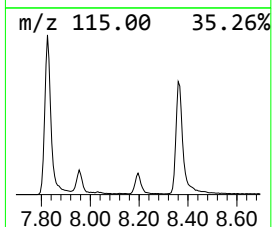
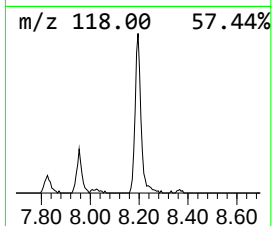
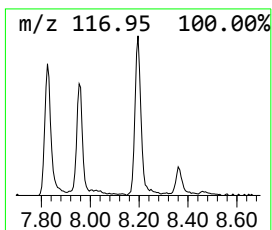
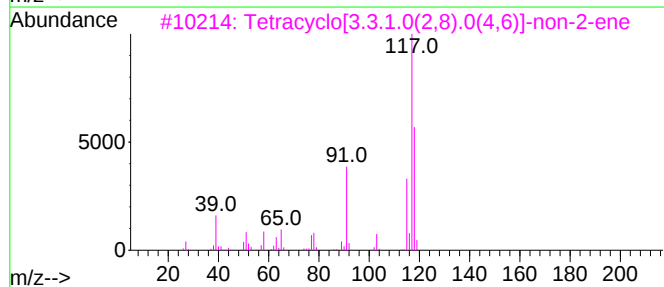
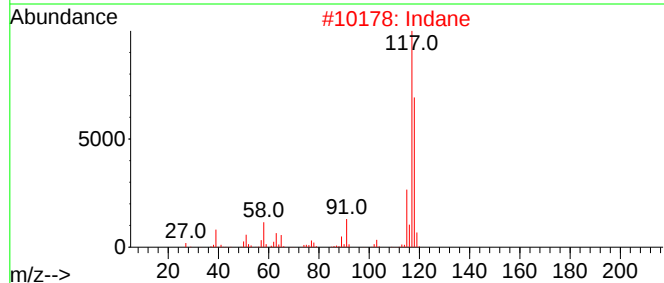
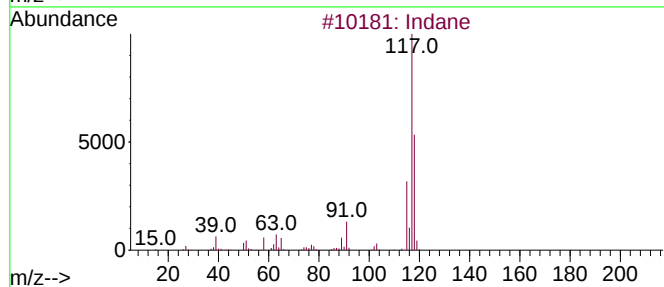
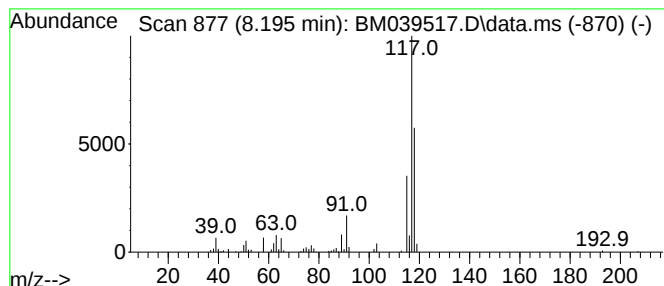
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Indane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.195	2.20 ng/ul	132757	1,4-Dichlorobenzene-d4	7.825

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	87
2			Indane	118	C9H10	000496-11-7	74
3			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	72
4			1,4:5,8-Dimethanobiphenylene, 1,...	184	C14H16	001624-13-1	72
5			Deltacyclene	118	C9H10	007785-10-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041823\
Data File : BM039517.D
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Operator : CG/JU
Sample : 02296-15DL3 900X
Misc :
ALS Vial : 69 Sample Multiplier: 1

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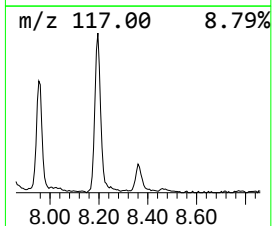
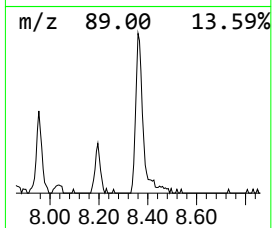
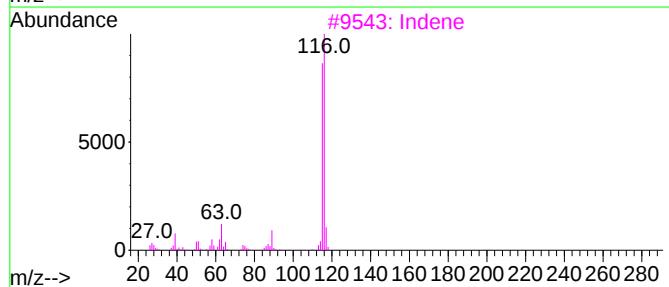
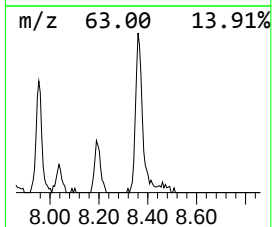
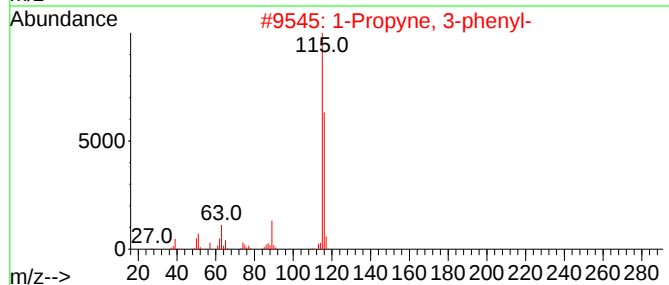
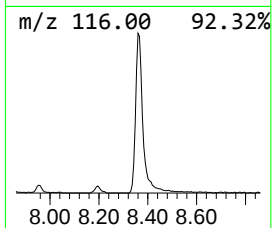
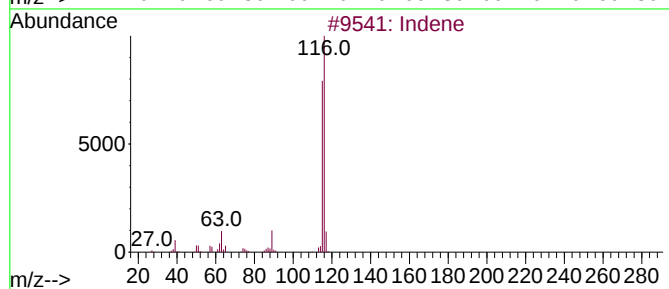
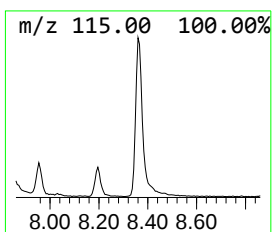
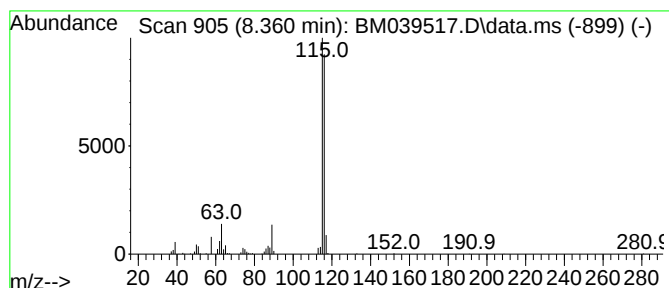
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Indene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.360	4.67 ng/ul	282114	1,4-Dichlorobenzene-d4	7.825

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene			116	C9H8	000095-13-6	97
2	1-Propyne, 3-phenyl-			116	C9H8	010147-11-2	95
3	Indene			116	C9H8	000095-13-6	95
4	Benzene, 1-propynyl-			116	C9H8	000673-32-5	94
5	Indene			116	C9H8	000095-13-6	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041823\
Data File : BM039517.D
Acq On : 20 Apr 2023 04:53
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Misc :
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Instrument :
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ClientSampleId :
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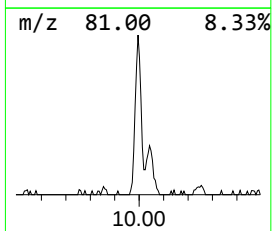
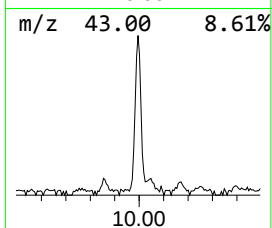
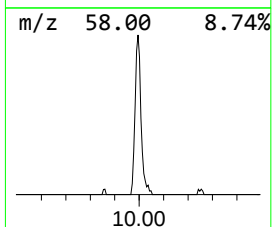
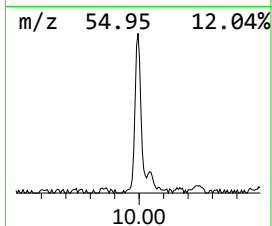
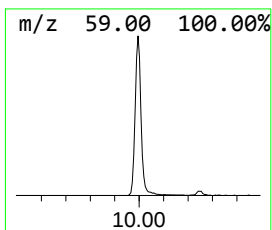
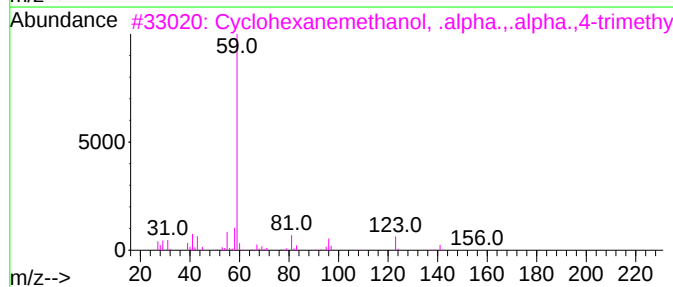
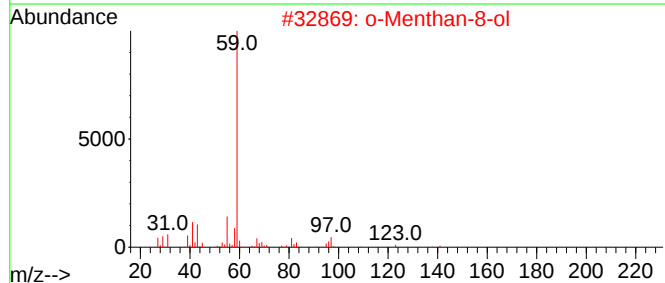
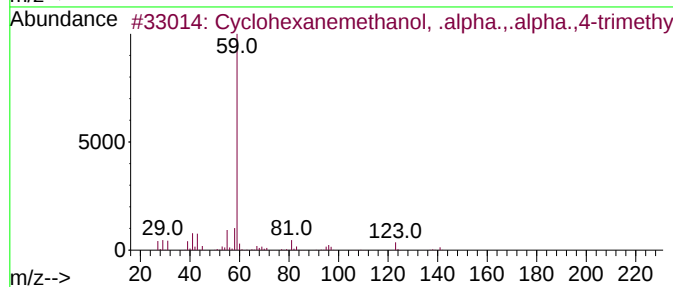
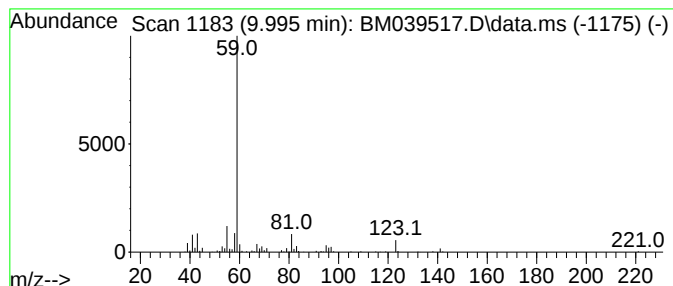
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Cyclohexanemethanol, .alpha... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.995	3.12 ng/ul	300141	Naphthalene-d8	10.631

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	83
2		o-Menthan-8-ol	156	C10H20O	343855-44-7	78
3		Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	005114-00-1	64
4		Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	007322-63-6	56
5		Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041823\
Data File : BM039517.D
Acq On : 20 Apr 2023 04:53
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Misc :
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ClientSampleId :
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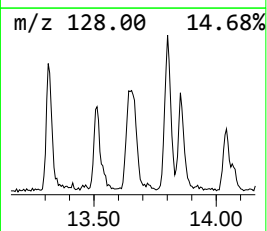
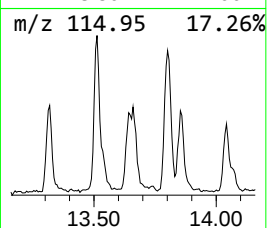
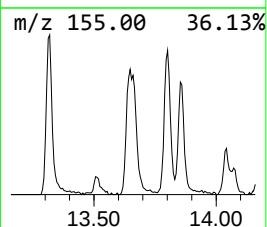
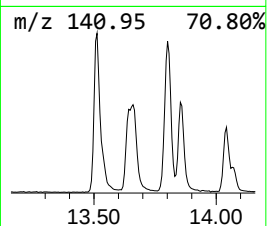
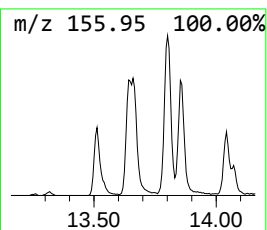
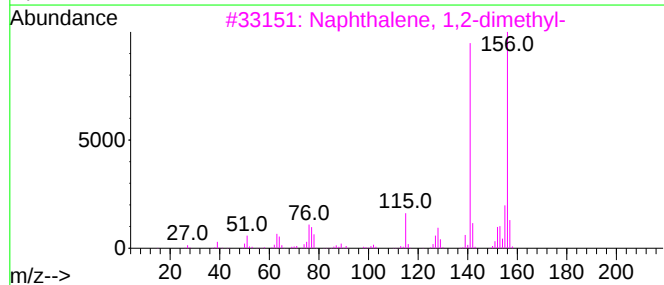
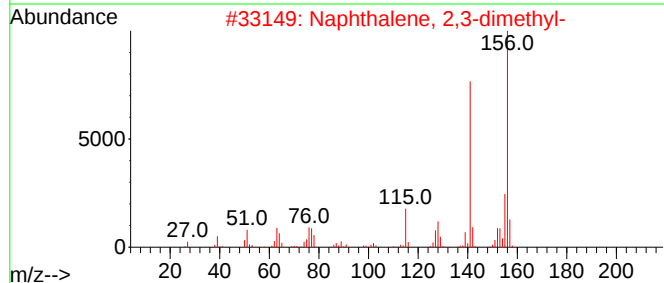
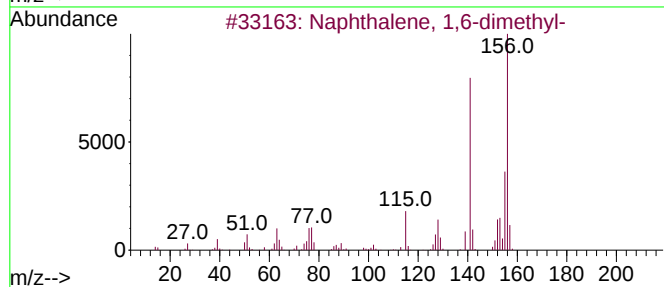
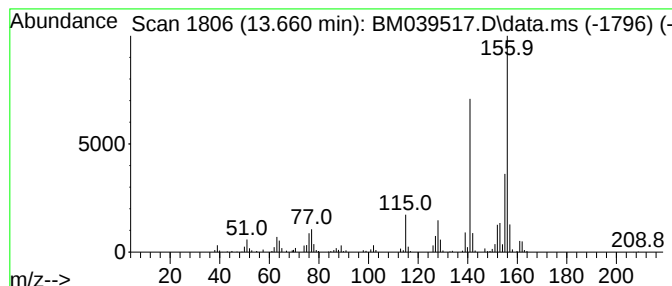
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Naphthalene, 1,6-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.660	3.24 ng/ul	474175	Acenaphthene-d10	14.483

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
3			Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	96
4			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041823\
Data File : BM039517.D
Acq On : 20 Apr 2023 04:53
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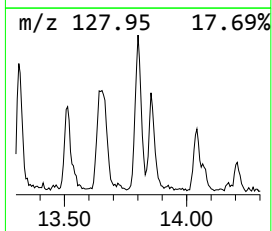
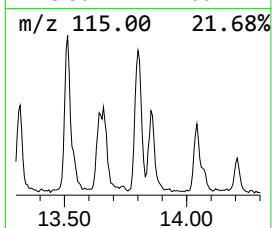
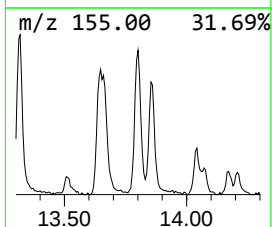
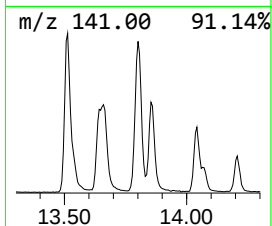
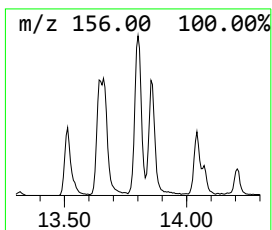
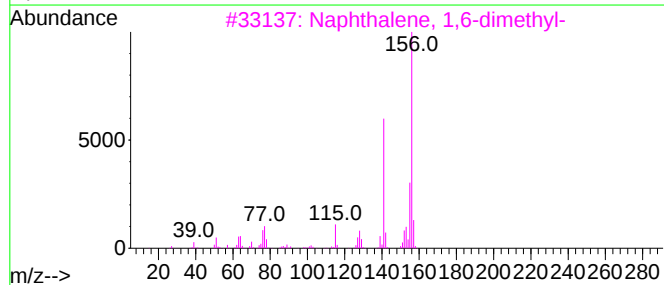
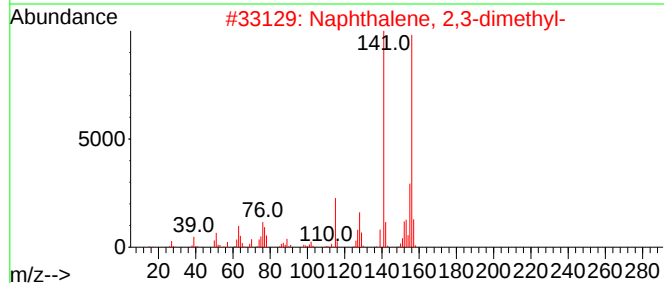
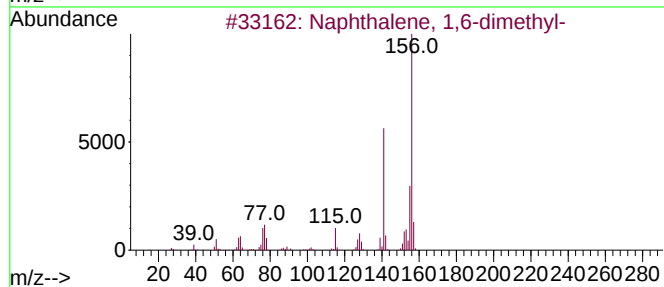
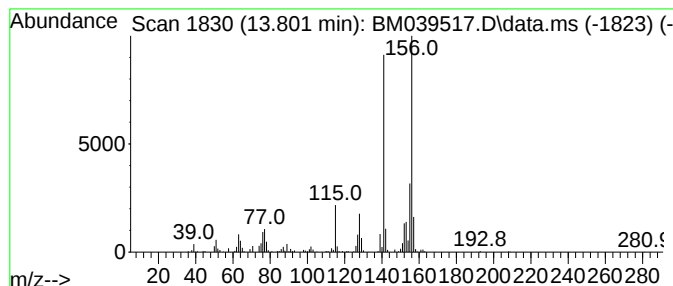
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Naphthalene, 2,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.801	3.00 ng/ul	437910	Acenaphthene-d10	14.483

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
5			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041823\
Data File : BM039517.D
Acq On : 20 Apr 2023 04:53
Operator : CG/JU
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ALS Vial : 69 Sample Multiplier: 1

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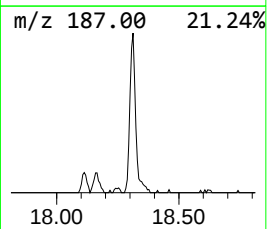
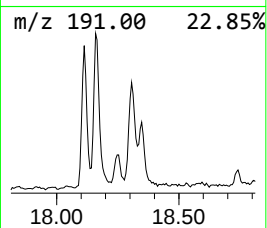
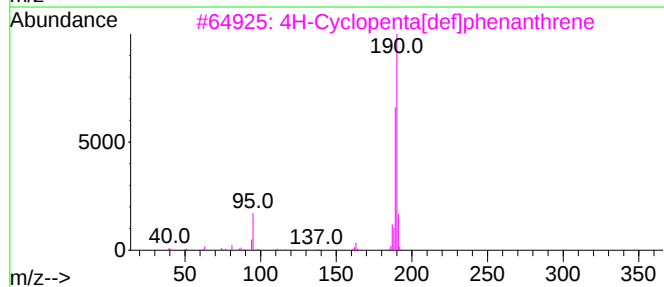
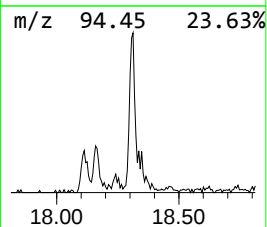
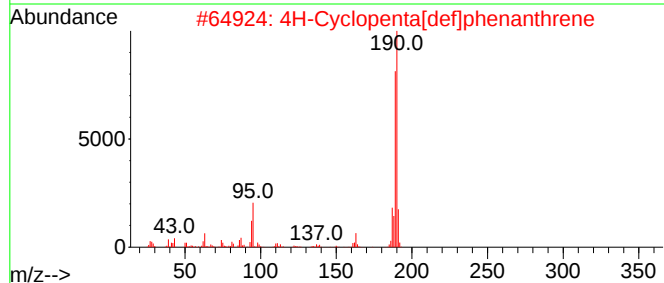
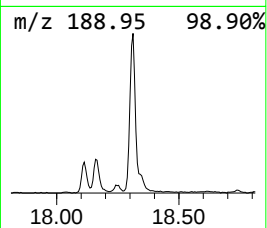
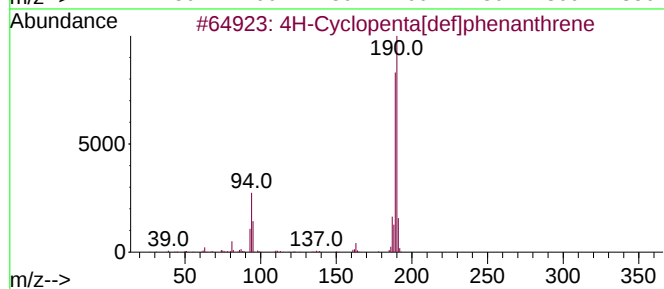
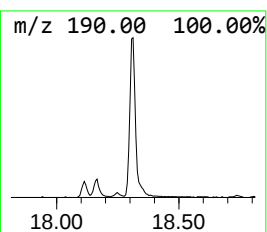
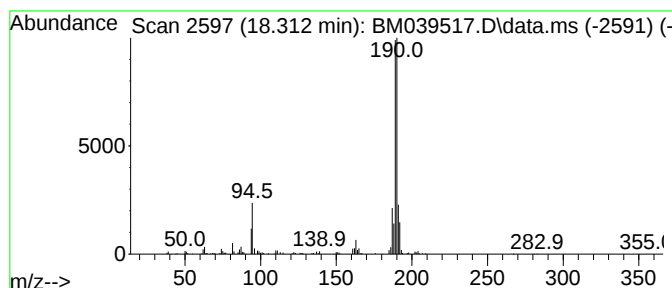
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 4H-Cyclopenta[def]phenanthrene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.312	2.07 ng/ul	334871	Phenanthrene-d10	17.230

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	95
2		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	90
3		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	83
4		6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	64
5		Methyl diselenide	190	C2H6Se2	007101-31-7	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041823\
Data File : BM039517.D
Acq On : 20 Apr 2023 04:53
Operator : CG/JU
Sample : 02296-15DL3 900X
Misc :
ALS Vial : 69 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCBM0DL3

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

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Bicyclo[3.1.1]h...	6.484	14.2	ng/ul	860518	1	7.825	1208240	20.0
.beta.-Pinene	7.248	3.6	ng/ul	217253	1	7.825	1208240	20.0
p-Cymene	7.954	9.2	ng/ul	557440	1	7.825	1208240	20.0
D-Limonene	8.037	4.9	ng/ul	297354	1	7.825	1208240	20.0
Indane	8.195	2.2	ng/ul	132757	1	7.825	1208240	20.0
Indene	8.360	4.7	ng/ul	282114	1	7.825	1208240	20.0
Cyclohexanemeth...	9.995	3.1	ng/ul	300141	2	10.631	1926600	20.0
Naphthalene, 1,...	13.660	3.2	ng/ul	474175	3	14.483	2922760	20.0
Naphthalene, 2,...	13.801	3.0	ng/ul	437910	3	14.483	2922760	20.0
4H-Cyclopenta[d...	18.312	2.1	ng/ul	334871	4	17.230	3243000	20.0