

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030223\
 Data File : BM038841.D
 Acq On : 02 Mar 2023 18:06
 Operator : CG/JU
 Sample : 01710-08
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBAA7

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-BM022823.M
 Title : SVOA CALIBRATION

Signal : TIC: BM038841.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.116	3	5	25	rVB	5276732	6448383	100.00%	10.295%
2	3.381	46	50	61	rVB	31065	50827	0.79%	0.081%
3	4.052	159	164	169	rBV	32728	53105	0.82%	0.085%
4	4.252	194	198	204	rVB3	20609	31815	0.49%	0.051%
5	5.081	333	339	347	rBV	70538	116750	1.81%	0.186%
6	6.387	556	561	567	rVB	20746	31875	0.49%	0.051%
7	6.687	604	612	620	rVB	72942	113199	1.76%	0.181%
8	6.998	659	665	673	rVB	90624	146305	2.27%	0.234%
9	7.157	685	692	704	rVB	504077	816105	12.66%	1.303%
10	7.334	715	722	732	rBV	426798	653699	10.14%	1.044%
11	7.451	736	742	747	rVV	67554	109047	1.69%	0.174%
12	7.534	749	756	767	rVV	488016	765148	11.87%	1.222%
13	8.010	830	837	845	rBV	881367	1351590	20.96%	2.158%
14	8.239	867	876	882	rBV6	19272	38225	0.59%	0.061%
15	8.322	882	890	898	rVB	773156	1231409	19.10%	1.966%
16	8.710	947	956	967	rBV	607626	946104	14.67%	1.510%
17	8.834	971	977	982	rVV	91967	149353	2.32%	0.238%
18	8.887	982	986	992	rVV3	23922	37435	0.58%	0.060%
19	9.175	1023	1035	1045	rBV	429423	732514	11.36%	1.169%
20	9.898	1149	1158	1165	rBV	279646	463132	7.18%	0.739%
21	10.239	1210	1216	1225	rBV4	65401	120660	1.87%	0.193%
22	10.369	1230	1238	1243	rBV4	32454	60896	0.94%	0.097%
23	10.439	1243	1250	1255	rVV	583177	980542	15.21%	1.565%
24	10.481	1255	1257	1270	rVB	64032	114700	1.78%	0.183%
25	10.598	1270	1277	1287	rBV	217143	375244	5.82%	0.599%
26	10.828	1309	1316	1323	rVV	1220791	2024106	31.39%	3.232%
27	10.957	1330	1338	1350	rBV	429704	754909	11.71%	1.205%
28	11.128	1362	1367	1372	rBV3	33658	54583	0.85%	0.087%
29	11.210	1379	1381	1394	rVB10	12347	29380	0.46%	0.047%
30	11.369	1401	1408	1414	rVB2	45638	78290	1.21%	0.125%
31	11.575	1437	1443	1448	rBV	47548	83256	1.29%	0.133%
32	12.945	1667	1676	1682	rBV8	20722	40211	0.62%	0.064%
33	13.351	1739	1745	1748	rBV5	16479	26651	0.41%	0.043%
34	13.398	1748	1753	1763	rVV8	62988	134889	2.09%	0.215%
35	13.480	1763	1767	1771	rVV3	20352	33168	0.51%	0.053%
36	13.533	1771	1776	1785	rVV6	49414	89728	1.39%	0.143%

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Integrator: RTE

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Start Thrs: 0.2

Stop Thrs : 0

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Max Peaks: 100

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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-BM022823.M
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37	13.904	1835	1839	1842	rVV5	16077	27110	0.42%	0.043%
38	13.980	1842	1852	1858	rVB2	148140	228604	3.55%	0.365%
39	14.057	1858	1865	1876	rBV	1074503	1518147	23.54%	2.424%
40	14.180	1882	1886	1890	rVB4	21378	26689	0.41%	0.043%

41	14.339	1907	1913	1921	rVB2	1228897	1752018	27.17%	2.797%
42	14.416	1921	1926	1931	rVB2	66859	95079	1.47%	0.152%
43	14.492	1934	1939	1945	rBV2	54660	78998	1.23%	0.126%
44	14.645	1952	1965	1972	rBV	1941244	2932619	45.48%	4.682%
45	14.716	1972	1977	1983	rBV3	77045	124198	1.93%	0.198%

46	14.821	1989	1995	2002	rBV	401164	565683	8.77%	0.903%
47	14.898	2002	2008	2015	rVV2	144099	229512	3.56%	0.366%
48	14.968	2015	2020	2026	rVB	55312	92678	1.44%	0.148%
49	15.057	2031	2035	2041	rVB4	28013	50070	0.78%	0.080%
50	15.192	2052	2058	2062	rBV4	25842	46748	0.72%	0.075%

51	15.633	2126	2133	2144	rBV	1688919	2383047	36.96%	3.805%
52	15.739	2145	2151	2159	rBV2	317341	546877	8.48%	0.873%
53	15.798	2159	2161	2164	rVV3	43312	52967	0.82%	0.085%
54	15.839	2165	2168	2172	rVV6	25626	37967	0.59%	0.061%
55	15.880	2172	2175	2181	rVB7	16692	29543	0.46%	0.047%

56	15.998	2188	2195	2205	rBV	650115	973580	15.10%	1.554%
57	16.086	2205	2210	2213	rVV	143205	224075	3.47%	0.358%
58	16.121	2213	2216	2221	rVB3	71281	99585	1.54%	0.159%
59	16.210	2226	2231	2235	rBV3	69358	109795	1.70%	0.175%
60	16.268	2236	2241	2247	rVV	183393	256112	3.97%	0.409%

61	16.363	2251	2257	2261	rBV	117841	158687	2.46%	0.253%
62	16.962	2352	2359	2363	rBV9	21485	56674	0.88%	0.090%
63	17.021	2365	2369	2375	rVB2	52080	78403	1.22%	0.125%
64	17.098	2377	2382	2388	rBV	493170	685610	10.63%	1.095%
65	17.286	2410	2414	2417	rBV5	32005	48833	0.76%	0.078%

66	17.386	2417	2431	2439	rVV	2726976	4016838	62.29%	6.413%
67	17.486	2442	2448	2457	rVB	1839891	2564276	39.77%	4.094%
68	17.586	2460	2465	2472	rVV4	187254	321090	4.98%	0.513%
69	17.651	2472	2476	2482	rVV6	50431	111663	1.73%	0.178%
70	17.709	2482	2486	2494	rVB3	92341	160887	2.49%	0.257%

71	18.145	2555	2560	2567	rVB2	288983	409002	6.34%	0.653%
72	18.280	2577	2583	2592	rBV2	500511	821717	12.74%	1.312%
73	18.357	2593	2596	2602	rVV6	75752	142937	2.22%	0.228%
74	18.421	2602	2607	2611	rVV	166710	229287	3.56%	0.366%
75	18.509	2619	2622	2627	rVB3	69949	93709	1.45%	0.150%

76	18.715	2655	2657	2670	rVB7	46886	94735	1.47%	0.151%
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77	18.833	2673	2677	2679	rVV2	73720	103107	1.60%	0.165%
78	18.862	2679	2682	2691	rVB	167427	242395	3.76%	0.387%
79	18.974	2698	2701	2705	rVB6	24661	28252	0.44%	0.045%
80	19.086	2712	2720	2725	rBV2	228845	348637	5.41%	0.557%
81	19.215	2738	2742	2746	rVB3	67409	86353	1.34%	0.138%
82	19.345	2759	2764	2770	rBV4	87580	150712	2.34%	0.241%
83	19.409	2772	2775	2778	rBV3	53876	64397	1.00%	0.103%
84	19.556	2797	2800	2810	rVB3	73805	109164	1.69%	0.174%
85	19.774	2831	2837	2844	rBV	2432228	3237218	50.20%	5.168%
86	19.868	2849	2853	2857	rBV2	147161	172275	2.67%	0.275%
87	19.921	2857	2862	2866	rBV	563287	701981	10.89%	1.121%
88	20.145	2895	2900	2904	rBV	236566	308125	4.78%	0.492%
89	20.221	2910	2913	2915	rBV2	29939	39087	0.61%	0.062%
90	20.327	2927	2931	2936	rVB3	143264	171500	2.66%	0.274%
91	20.462	2951	2954	2959	rVB3	97484	117970	1.83%	0.188%
92	20.527	2961	2965	2972	rVB	277661	334111	5.18%	0.533%
93	20.645	2980	2985	2993	rBV	496880	624155	9.68%	0.996%
94	21.115	3061	3065	3070	rBV	731243	795975	12.34%	1.271%
95	21.280	3089	3093	3104	rVB3	318169	531440	8.24%	0.848%
96	21.562	3136	3141	3154	rBV	3086134	4120774	63.90%	6.579%
97	22.215	3249	3252	3258	rVB3	96572	157091	2.44%	0.251%
98	23.303	3433	3437	3441	rBV	106971	139215	2.16%	0.222%
99	23.856	3524	3531	3541	rBV2	1799266	3541542	54.92%	5.654%
100	24.021	3552	3559	3576	rVB	2332852	4746853	73.61%	7.579%

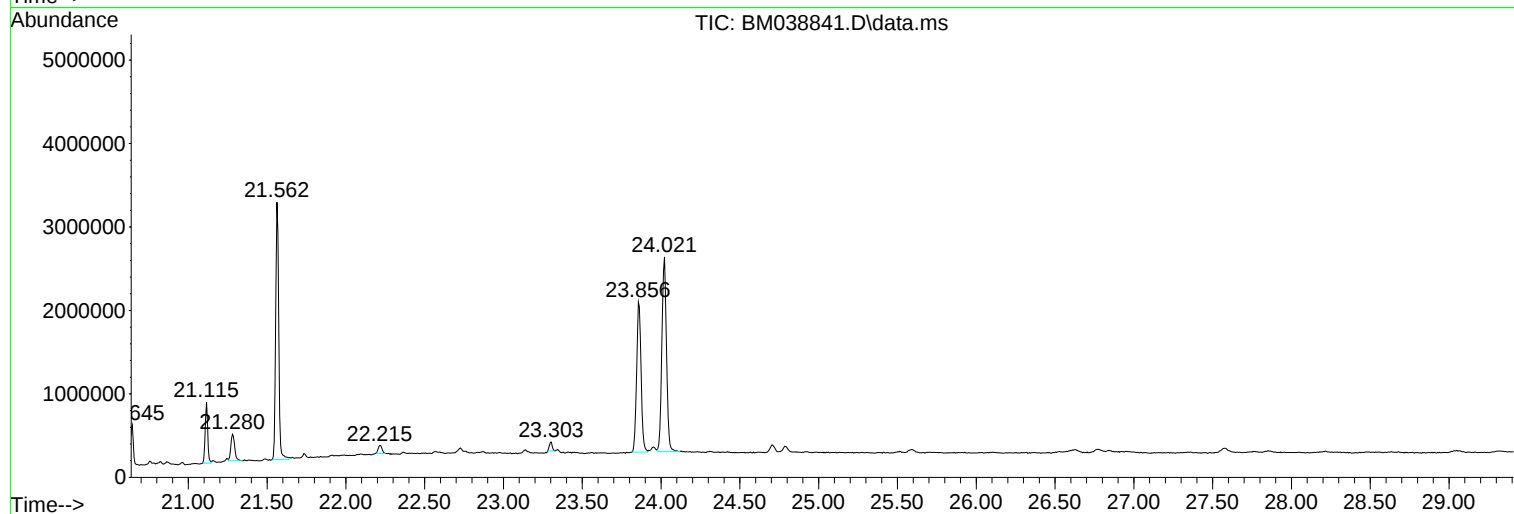
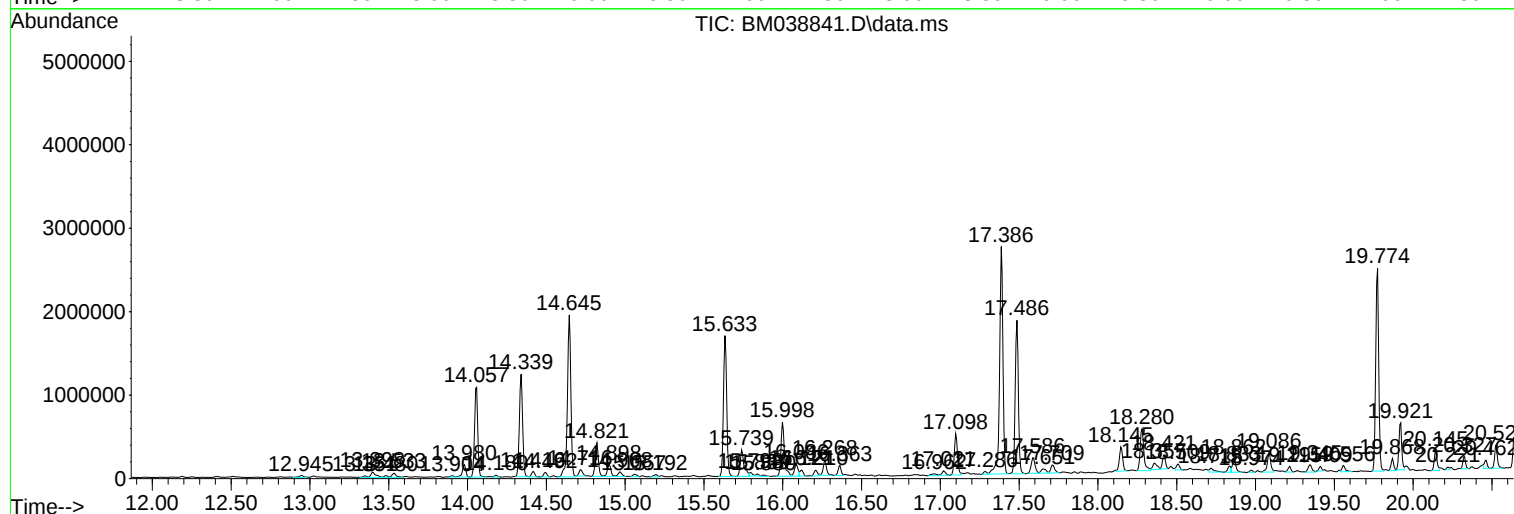
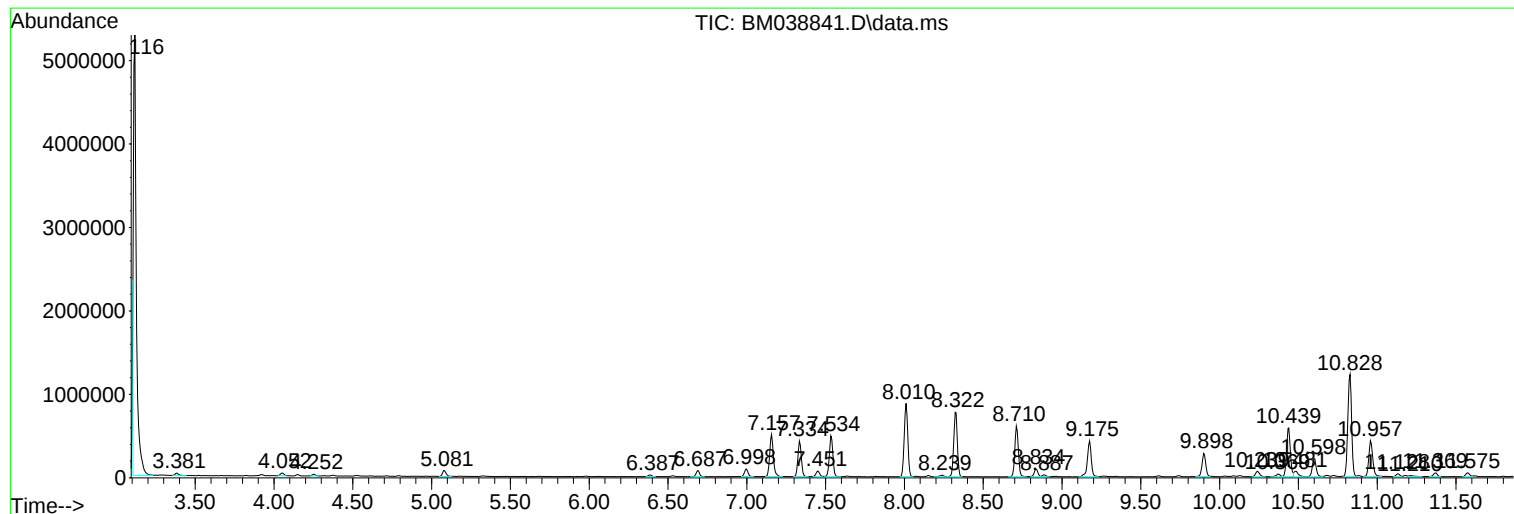
Sum of corrected areas: 62635611

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022823.M
Quant Title : SVOA CALIBRATION

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



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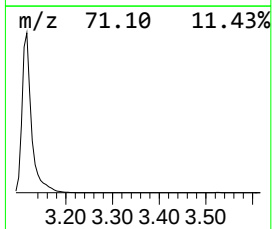
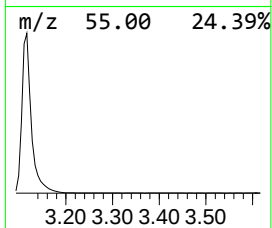
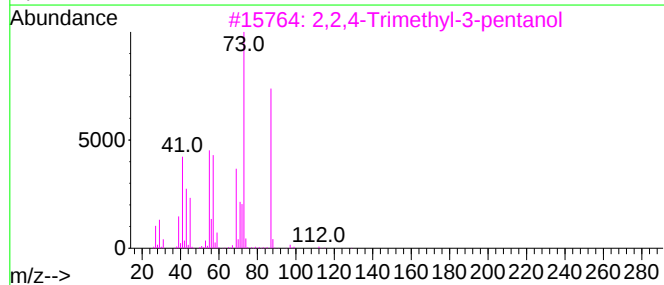
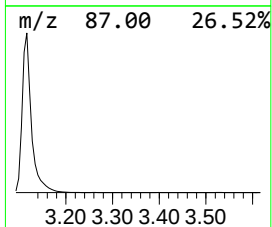
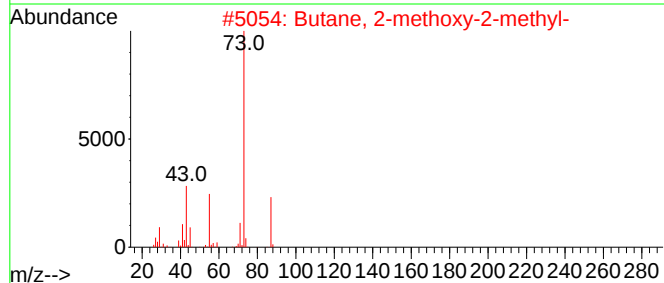
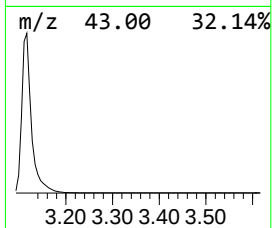
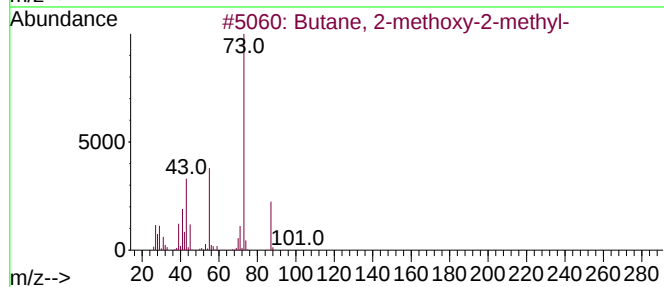
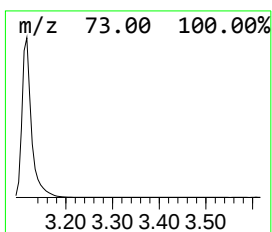
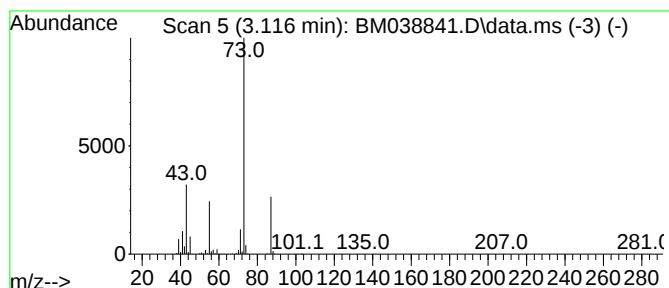
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022823.M
Quant Title : SVOA CALIBRATION

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.
3.116	95.42 ng/ul	6448380	1,4-Dichlorobenzene-d4	8.010

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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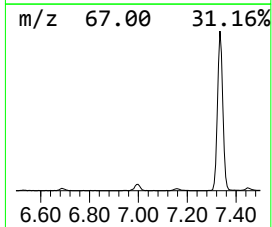
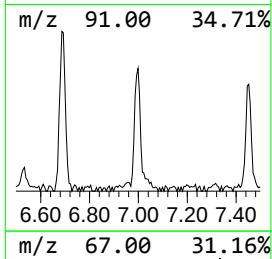
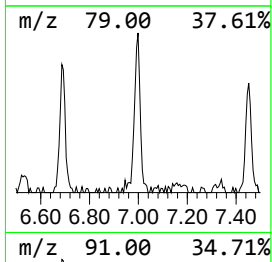
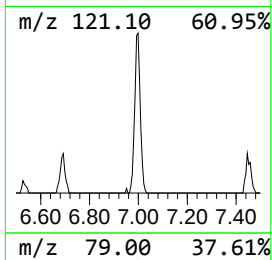
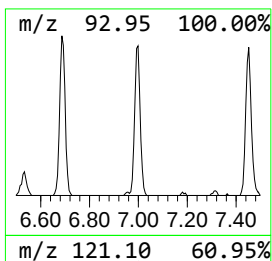
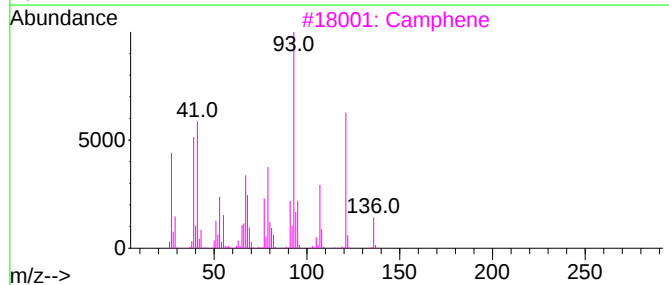
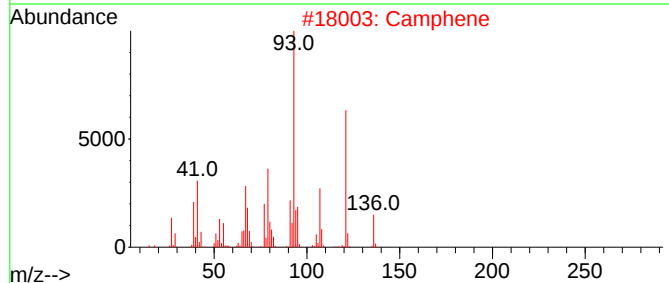
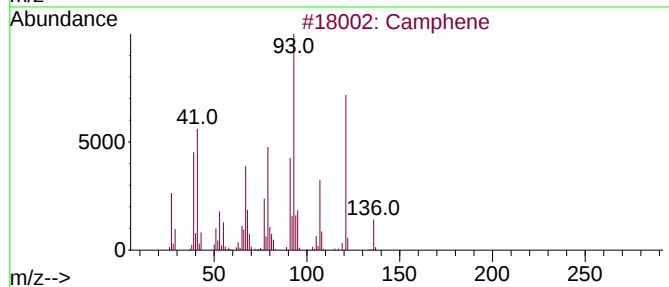
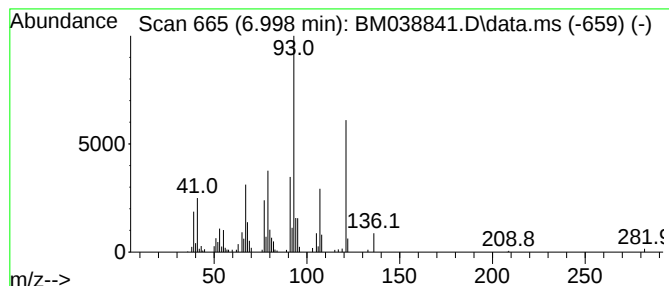
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022823.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Camphene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.998	2.16 ng/ul	146305	1,4-Dichlorobenzene-d4	8.010

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Camphene	136	C10H16	000079-92-5	95
2			Camphene	136	C10H16	000079-92-5	93
3			Camphene	136	C10H16	000079-92-5	90
4			Camphene	136	C10H16	000079-92-5	87
5			Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	80



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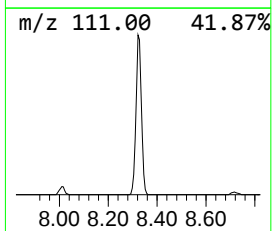
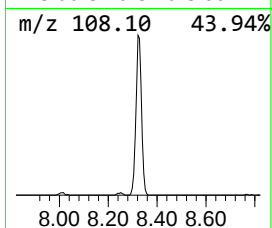
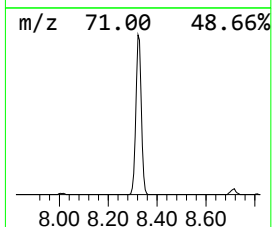
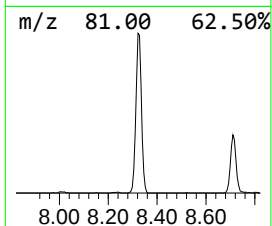
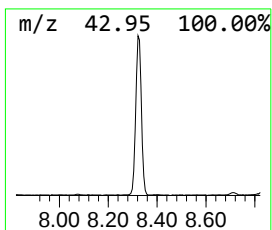
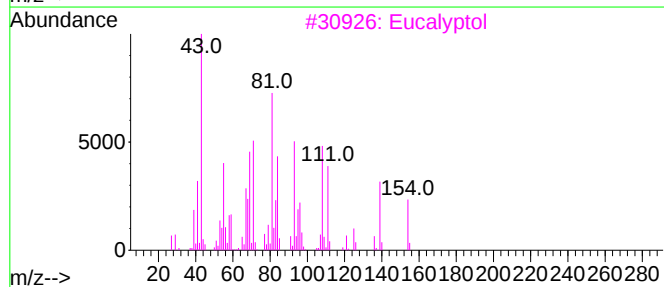
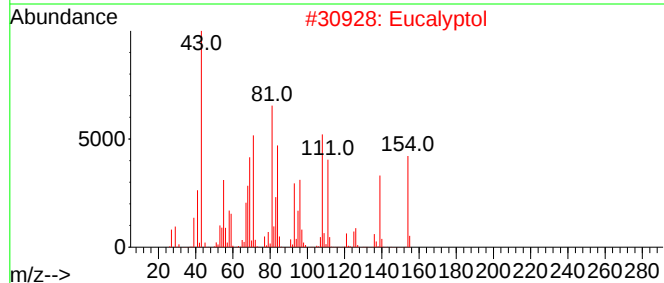
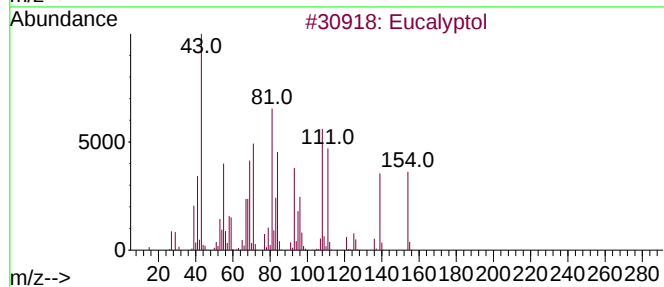
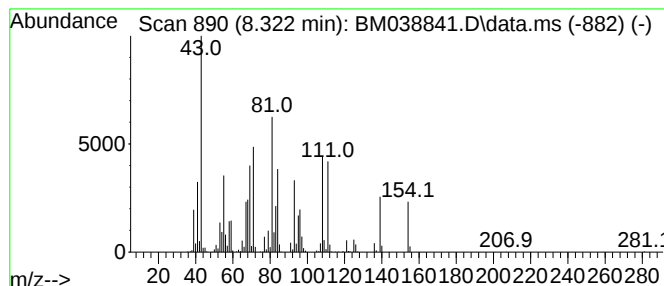
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Eucalyptol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.322	18.22 ng/ul	1231410	1,4-Dichlorobenzene-d4	8.010

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eucalyptol			154	C10H18O	000470-82-6	98
2	Eucalyptol			154	C10H18O	000470-82-6	93
3	Eucalyptol			154	C10H18O	000470-82-6	93
4	Eucalyptol			154	C10H18O	000470-82-6	91
5	Eucalyptol			154	C10H18O	000470-82-6	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030223\
Data File : BM038841.D
Acq On : 02 Mar 2023 18:06
Operator : CG/JU
Sample : 01710-08
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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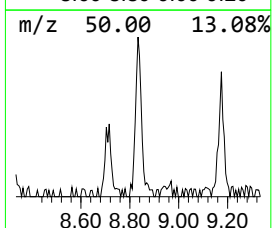
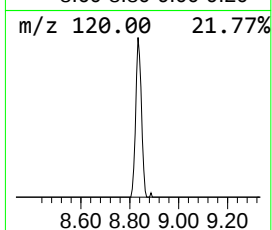
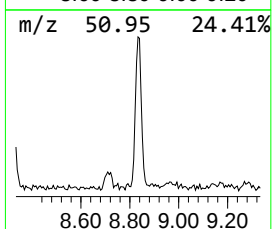
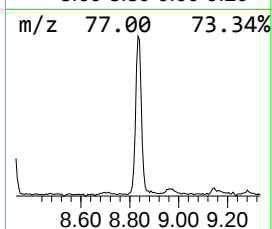
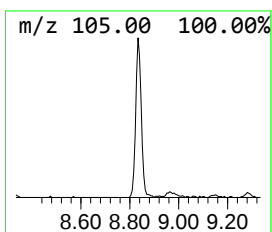
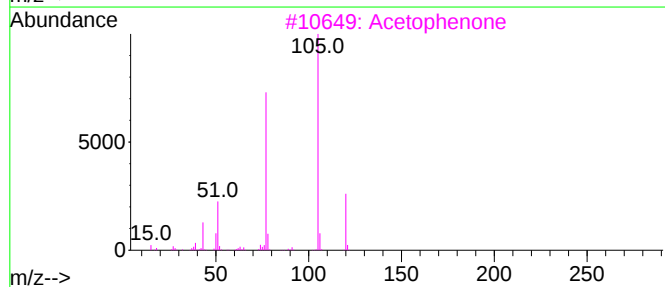
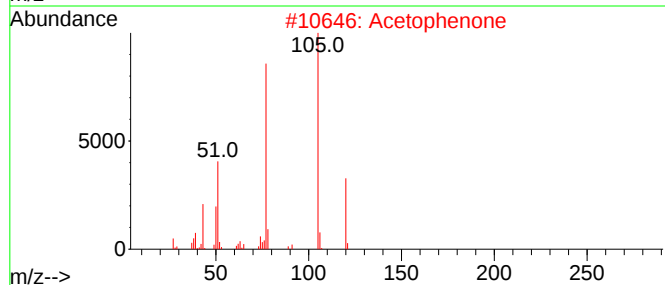
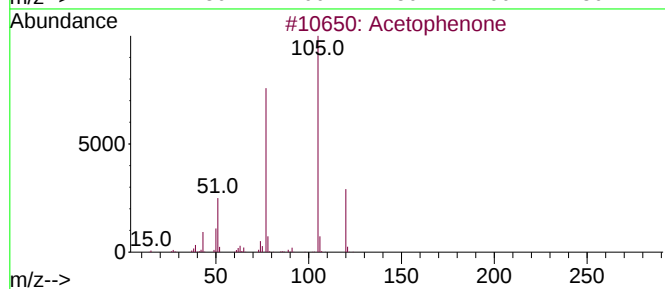
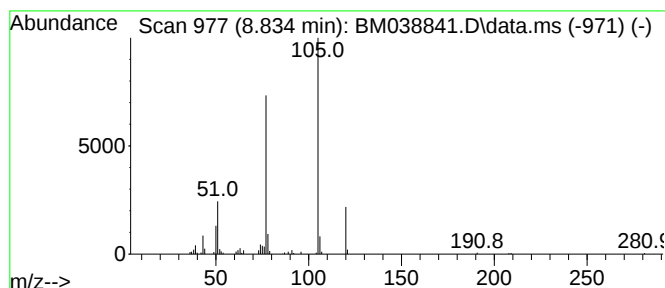
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Acetophe none Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.834	2.21 ng/ul	149353	1,4-Dichlorobenzene-d4	8.010

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetophenone	120	C8H8O	000098-86-2	95
2			Acetophenone	120	C8H8O	000098-86-2	95
3			Acetophenone	120	C8H8O	000098-86-2	94
4			Acetophenone	120	C8H8O	000098-86-2	91
5			Acetophenone	120	C8H8O	000098-86-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030223\
Data File : BM038841.D
Acq On : 02 Mar 2023 18:06
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Sample : 01710-08
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ClientSampleId :
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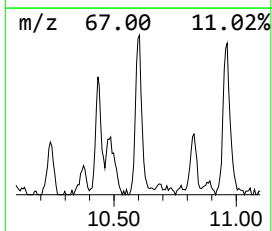
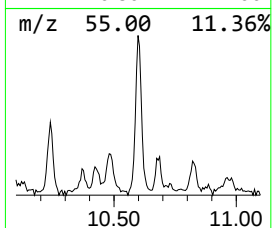
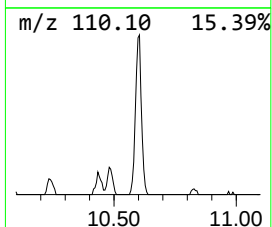
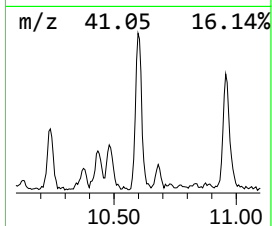
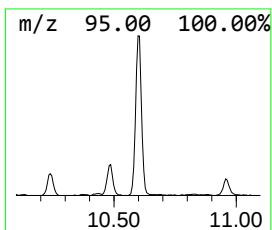
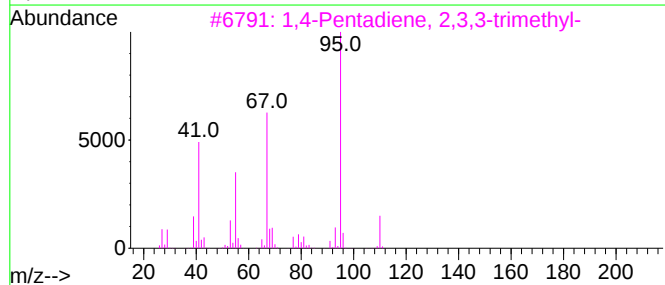
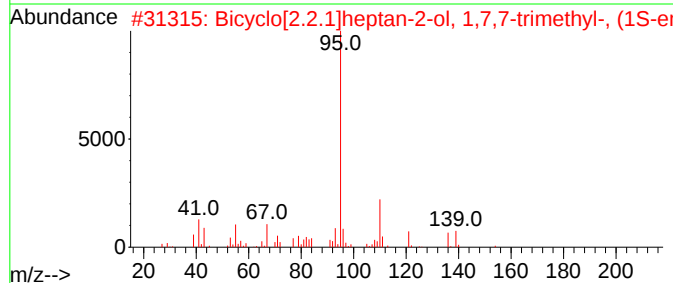
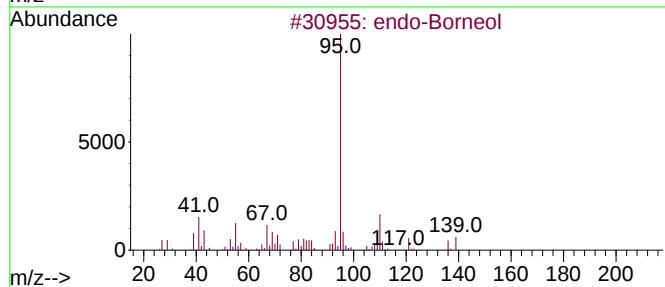
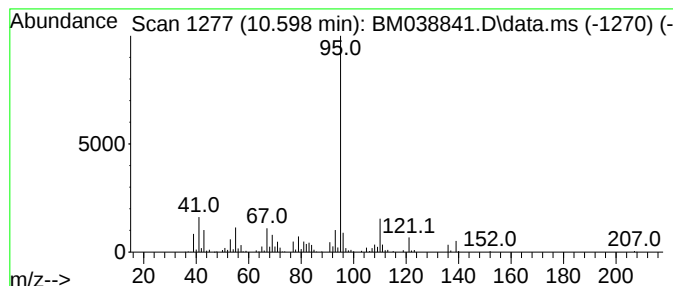
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 endo-Borneol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.598	3.71 ng/ul	375244	Naphthalene-d8	10.828

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			endo-Borneol	154	C10H18O	000507-70-0	86
2			Bicyclo[2.2.1]heptan-2-ol, 1,7,7...	154	C10H18O	000464-45-9	78
3			1,4-Pentadiene, 2,3,3-trimethyl-	110	C8H14	000756-02-5	64
4			endo-Borneol	154	C10H18O	000507-70-0	64
5			5,6-Dibromo-5-methyl-hex-1-ene	254	C7H12Br2	1000192-24-6	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030223\
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Sample : 01710-08
Misc :
ALS Vial : 13 Sample Multiplier: 1

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ClientSampleId :
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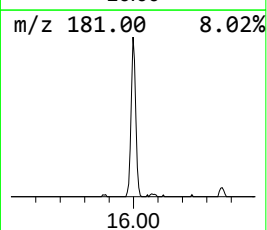
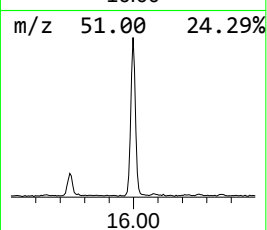
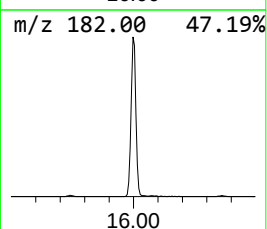
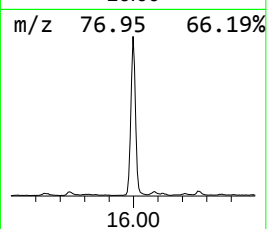
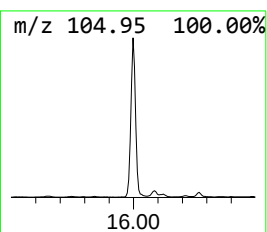
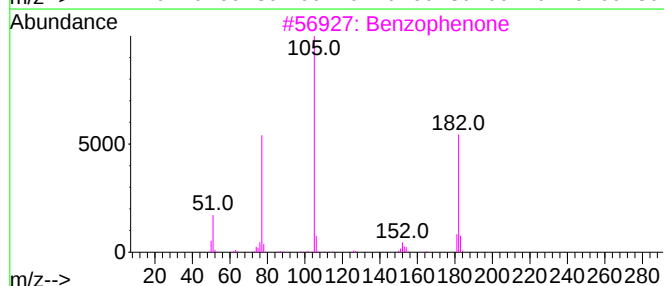
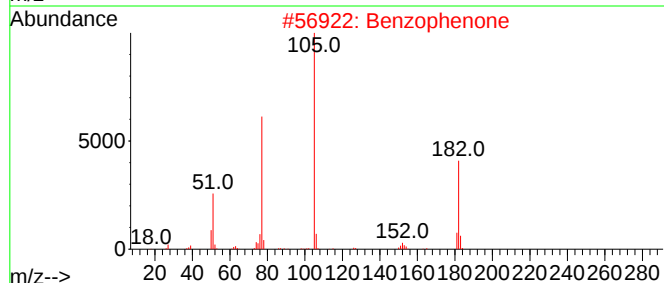
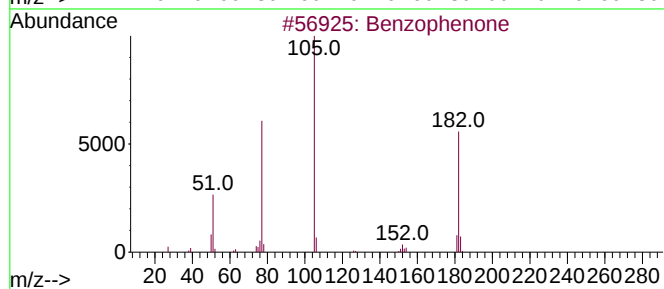
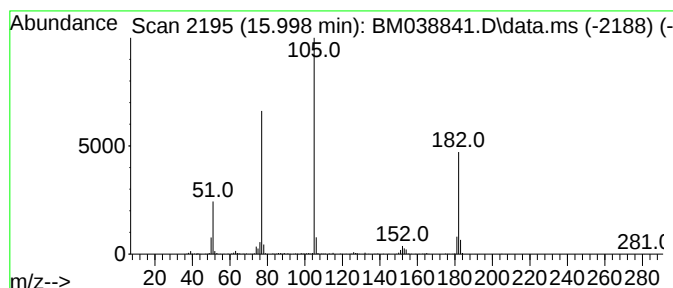
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.998	6.64 ng/ul	973580	Acenaphthene-d10	14.645

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030223\
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Acq On : 02 Mar 2023 18:06
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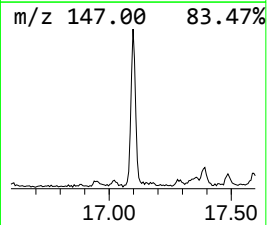
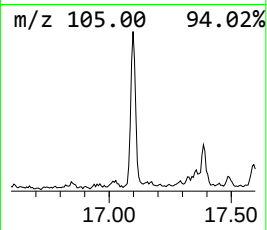
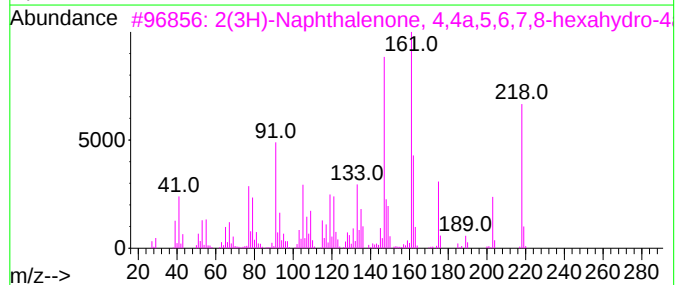
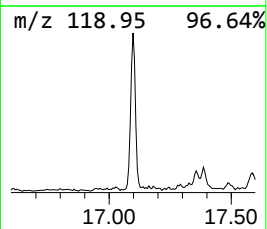
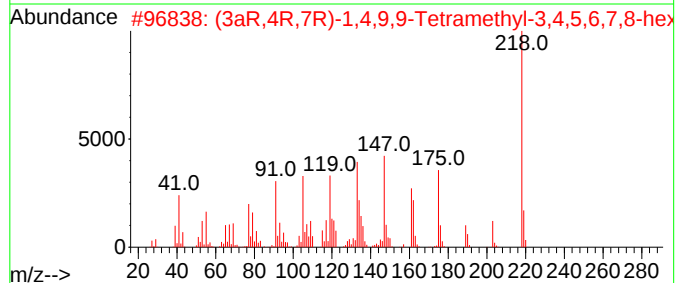
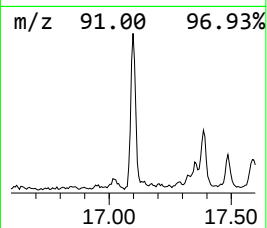
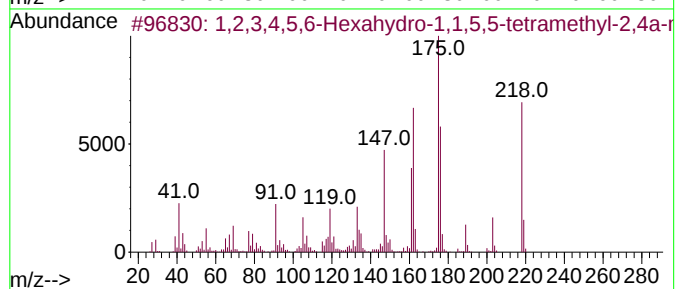
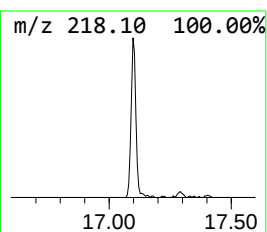
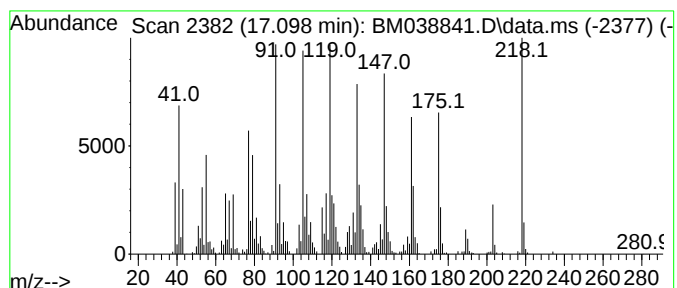
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1,2,3,4,5,6-Hexahydro-1,1,5... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.098	3.41 ng/ul	685610	Phenanthrene-d10	17.386	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,3,4,5,6-Hexahydro-1,1,5,5-te...	218	C15H22O	023747-14-0	91
2	(3aR,4R,7R)-1,4,9,9-Tetramethyl-...	218	C15H22O	003466-15-7	83
3	2(3H)-Naphthalenone, 4,4a,5,6,7,...	218	C15H22O	019598-45-9	83
4	5(1H)-Azulenone, 2,4,6,7,8,8a-he...	218	C15H22O	006754-66-1	83
5	Isolongifolen-5-one	218	C15H22O	1000159-37-1	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030223\
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Misc :
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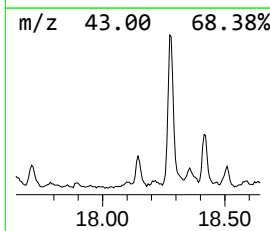
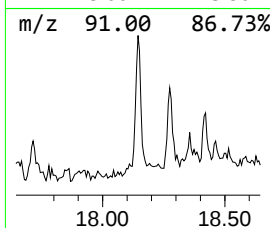
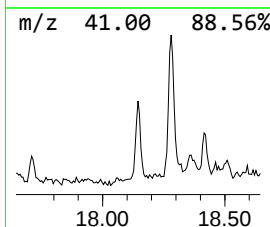
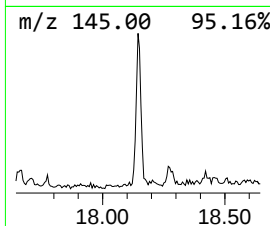
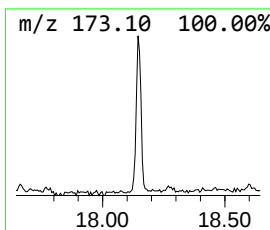
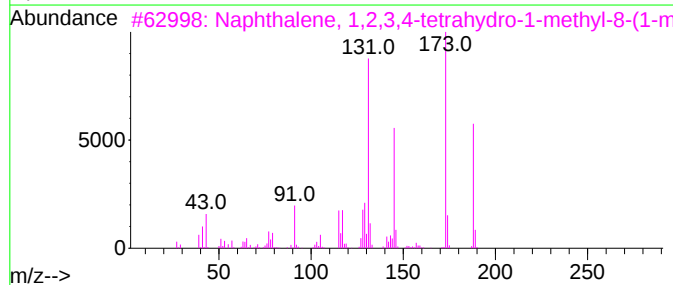
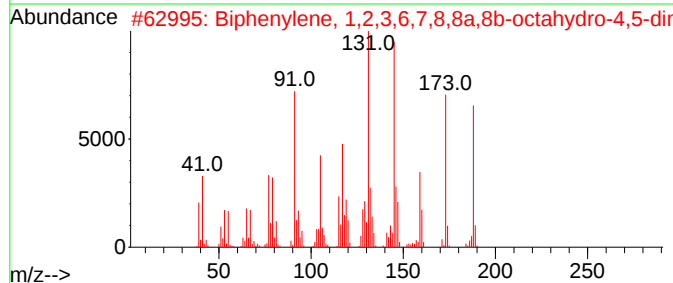
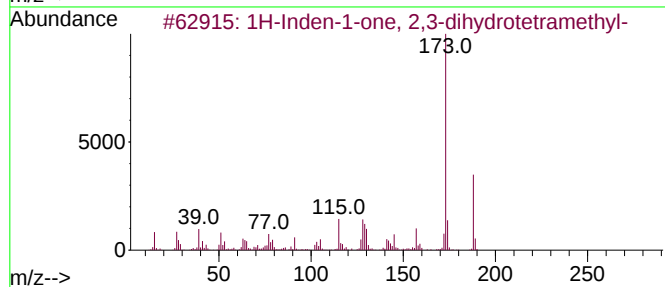
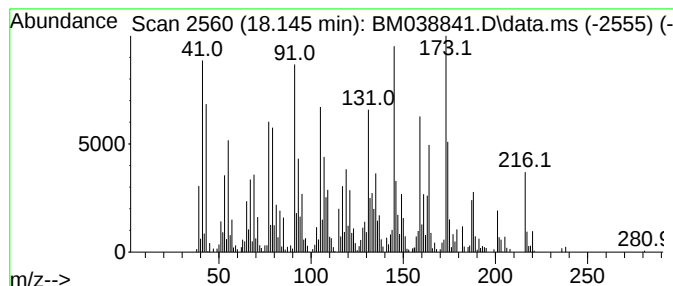
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 1H-Inden-1-one, 2,3-dihydro... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.145	2.04 ng/ul	409002	Phenanthrene-d10	17.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Inden-1-one, 2,3-dihydro-1-methyl-8-(1-methyl-1H-inden-5-yl)-1,2,3,4-tetrahydro-5H-benzo[5,6-b]indole	188	C13H16O	089907-33-5	56
2			Biphenylene, 1,2,3,6,7,8,8a,8b-octahydro-4,5-dimethyl-	188	C14H20	106988-87-8	46
3			Naphthalene, 1,2,3,4-tetrahydro-1-methyl-8-(1-methyl-1H-inden-5-yl)-1,2,3,4-tetrahydro-5H-benzo[5,6-b]indole	188	C14H20	081603-43-2	42
4			8-Amino-6-methoxyquinoline	174	C10H10N2O	000090-52-8	38
5			Benzene, 1-tert-butyl-4-cyclopropyl-	188	C14H20	058249-45-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030223\
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Sample : 01710-08
Misc :
ALS Vial : 13 Sample Multiplier: 1

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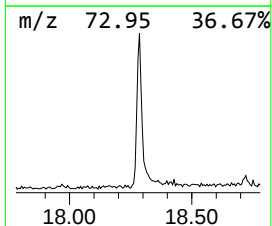
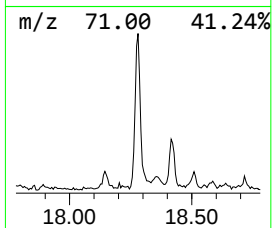
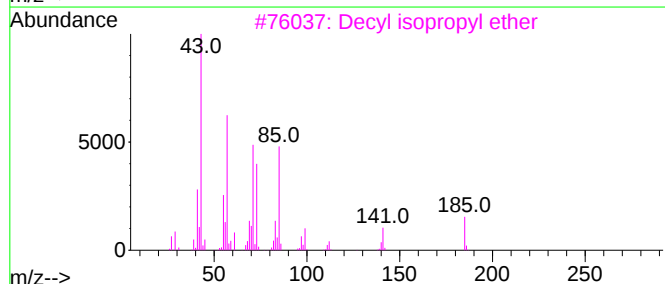
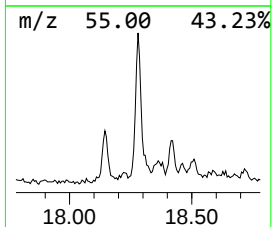
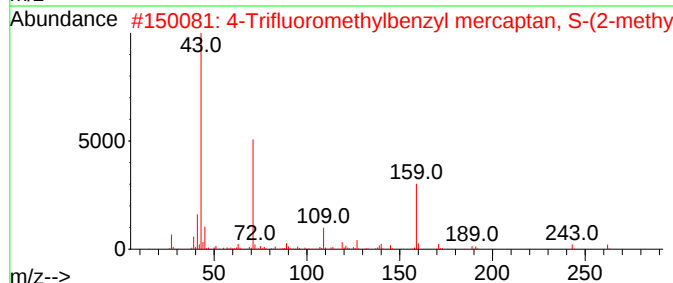
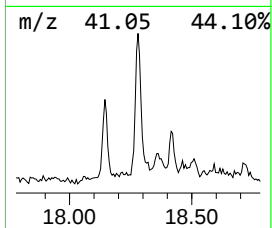
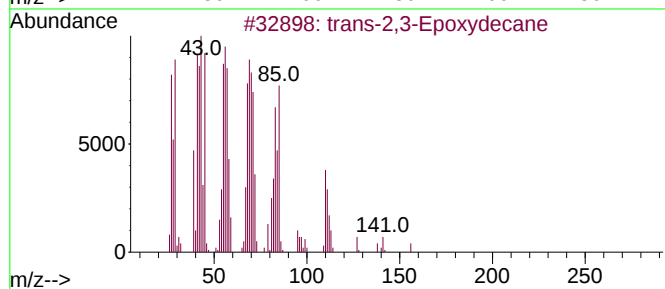
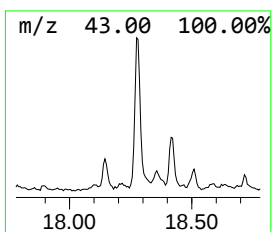
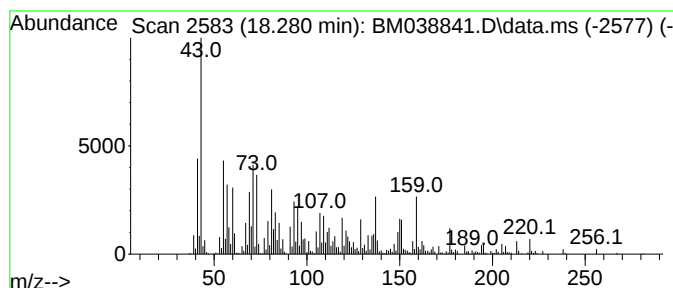
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 trans-2,3-Epoxydecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.280	4.09 ng/ul	821717	Phenanthrene-d10	17.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-2,3-Epoxydecane	156	C10H20O	054125-39-2	59
2			4-Trifluoromethylbenzyl mercapta...	262	C12H13F3OS	1000470-57-2	22
3			Decyl isopropyl ether	200	C13H28O	1000406-33-8	22
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	18
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030223\
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ALS Vial : 13 Sample Multiplier: 1

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ClientSampleId :
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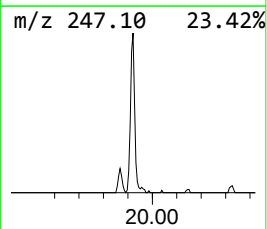
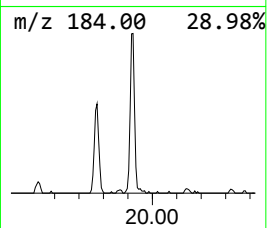
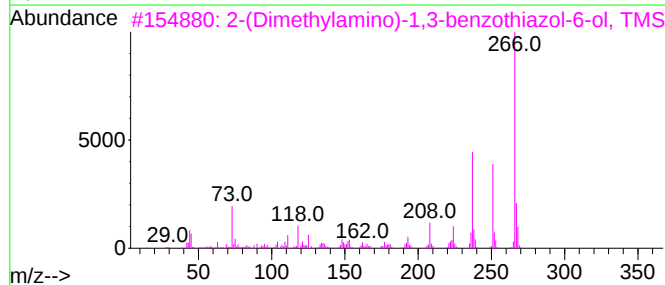
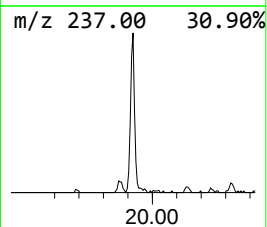
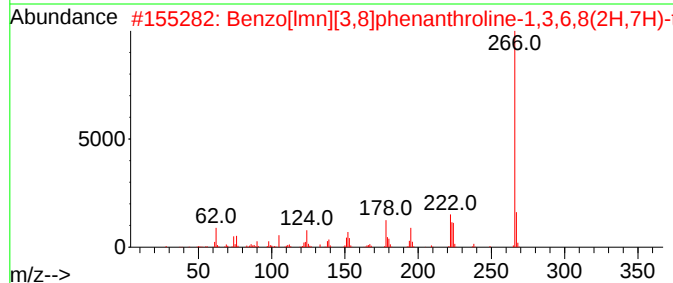
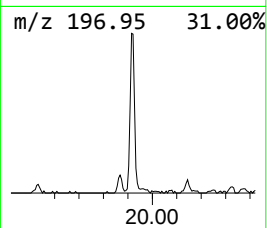
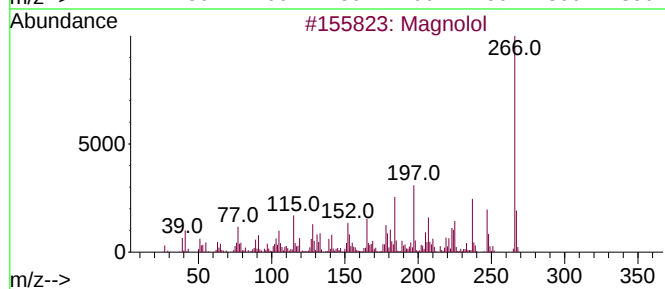
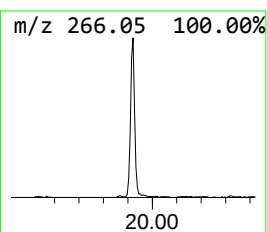
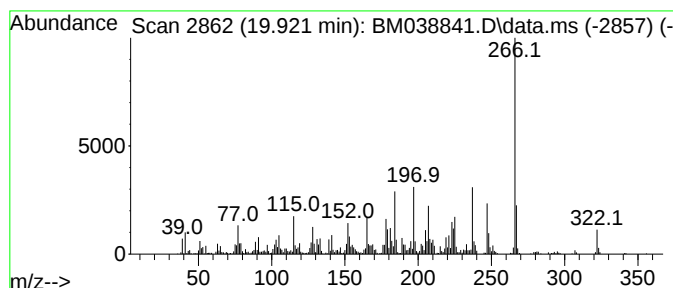
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Magnolol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.921	3.41 ng/ul	701981	Chrysene-d12	21.562

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Magnolol	266	C18H18O2	000528-43-8	99
2			Benzo[1mn][3,8]phenanthroline-1,...	266	C14H6N2O4	005690-24-4	56
3			2-(Dimethylamino)-1,3-benzothiaz...	266	C12H18N2OSSi	1000474-64-0	49
4			Equilenin	266	C18H18O2	000517-09-9	46
5			10-Nitro-15-oxa-14-azatetracyclo...	266	C14H6N2O4	085192-91-2	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030223\
Data File : BM038841.D
Acq On : 02 Mar 2023 18:06
Operator : CG/JU
Sample : 01710-08
Misc :
ALS Vial : 13 Sample Multiplier: 1

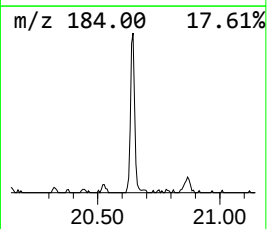
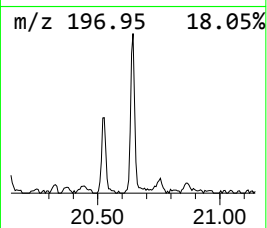
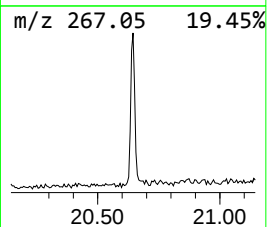
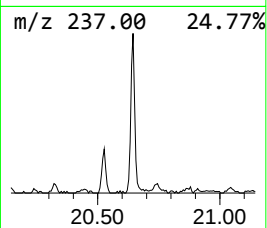
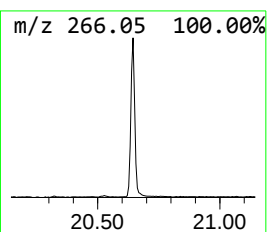
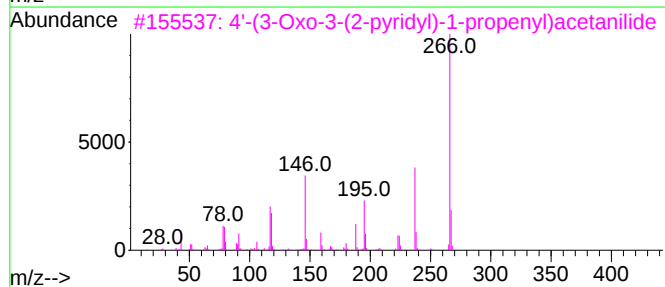
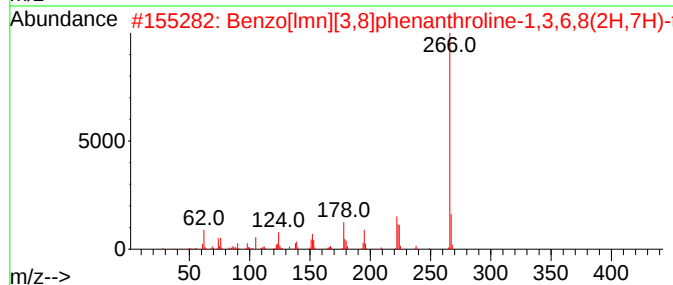
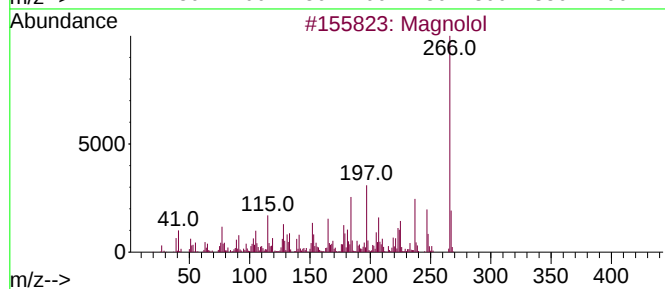
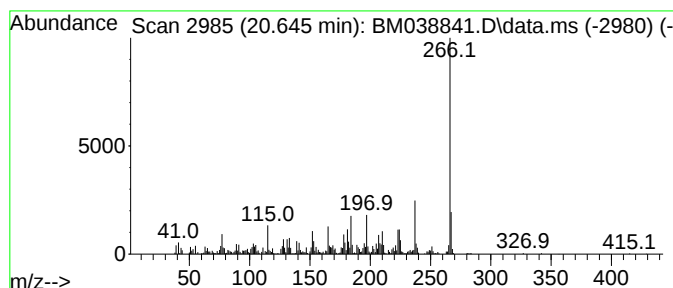
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022823.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 Benzo[lmn][3,8]phenanthroli... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.645	3.03 ng/ul	624155	Chrysene-d12		21.562	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Magnolol		266	C18H18O2	000528-43-8	91
2	Benzo[lmn][3,8]phenanthroline-1,...		266	C14H6N2O4	005690-24-4	83
3	4'-(3-Oxo-3-(2-pyridyl)-1-propen...		266	C16H14N2O2	013309-10-9	76
4	Equilenin		266	C18H18O2	000517-09-9	76
5	6-Methoxy-4-methyl-3-phenylcoumarin		266	C17H14O3	115059-27-3	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030223\
Data File : BM038841.D
Acq On : 02 Mar 2023 18:06
Operator : CG/JU
Sample : 01710-08
Misc :
ALS Vial : 13 Sample Multiplier: 1

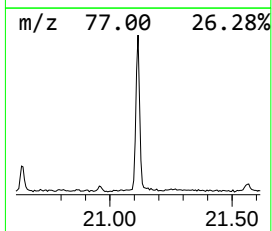
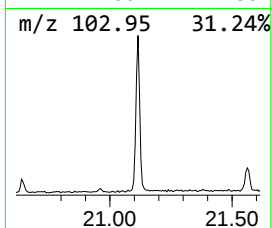
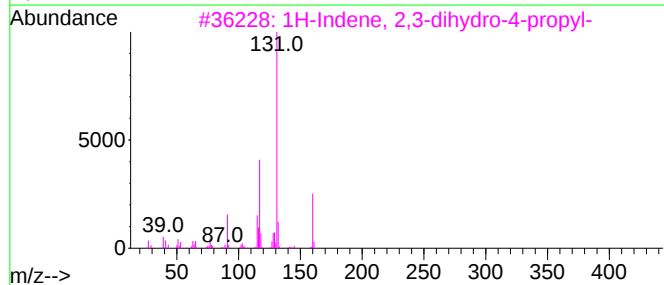
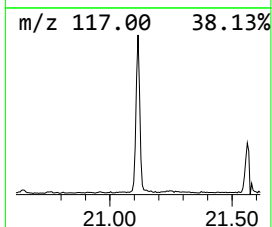
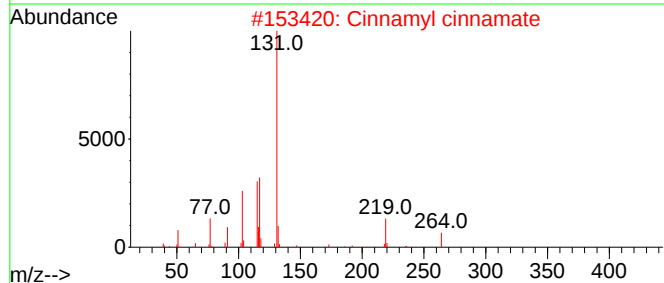
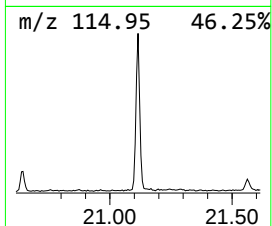
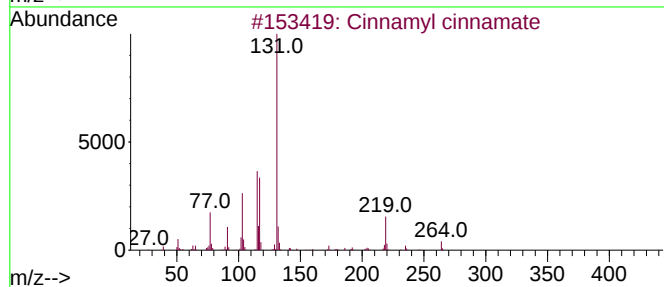
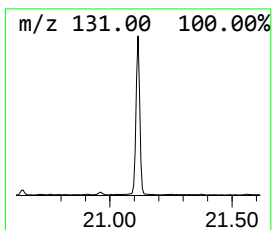
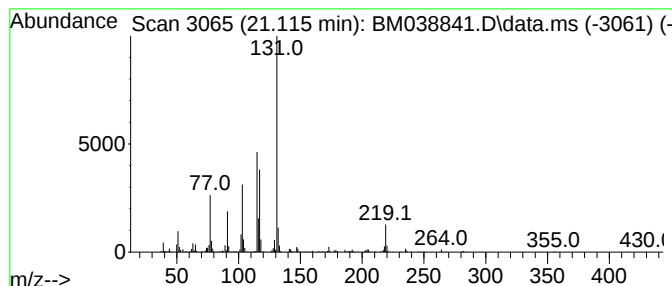
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022823.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 Cinnamyl cinnamate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.115	3.86 ng/ul	795975	Chrysene-d12	21.562	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cinnamyl cinnamate	264	C18H16O2	000122-69-0	91
2	Cinnamyl cinnamate	264	C18H16O2	000122-69-0	80
3	1H-Indene, 2,3-dihydro-4-propyl-	160	C12H16	092013-16-6	50
4	2,2-Dimethyl-1-(2-vinylphenyl)pr...	188	C13H16O	1000210-99-9	43
5	(Z)-N-(2-Hydroxyethyl)-3-phenyla...	233	C13H15NO3	1000512-63-0	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030223\
Data File : BM038841.D
Acq On : 02 Mar 2023 18:06
Operator : CG/JU
Sample : 01710-08
Misc :
ALS Vial : 13 Sample Multiplier: 1

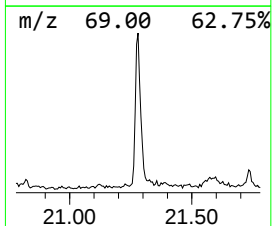
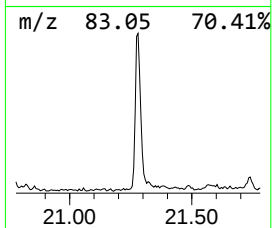
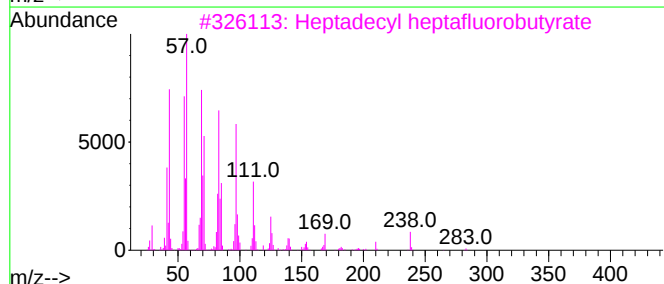
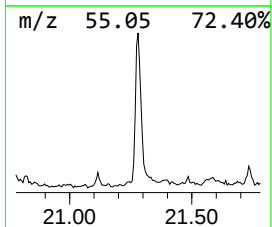
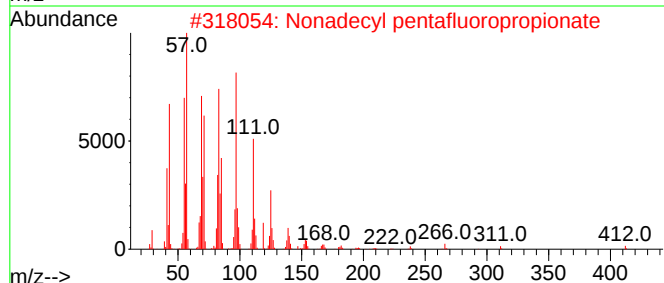
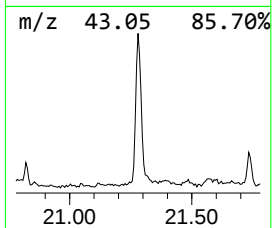
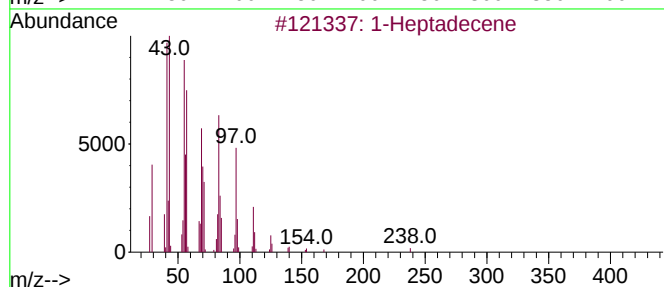
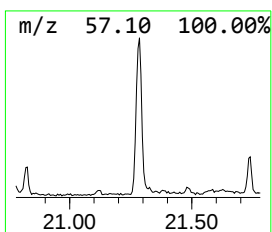
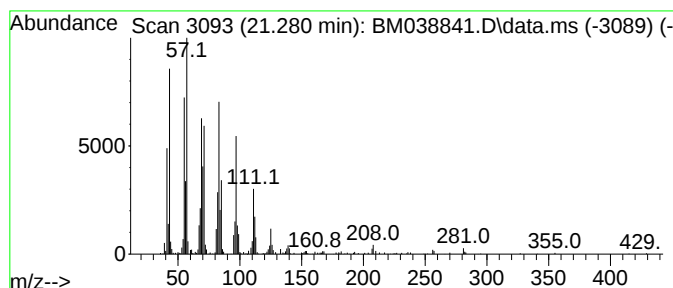
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022823.M
Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.280	2.58 ng/ul	531440	Chrysene-d12		21.562	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1-Heptadecene		238	C17H34	006765-39-5	93
2	Nonadecyl pentafluoropropionate		430	C22H39F5O2	1000351-88-8	91
3	Heptadecyl heptafluorobutyrate		452	C21H35F7O2	959085-66-6	91
4	Ethanol, 2-(tetradecyloxy)-		258	C16H34O2	002136-70-1	90
5	Hexadecyl propyl ether		284	C19H40O	1000406-27-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030223\
Data File : BM038841.D
Acq On : 02 Mar 2023 18:06
Operator : CG/JU
Sample : 01710-08
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DBAA7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022823.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
Butane, 2-metho...	3.116	95.4	ng/ul	6448380	1	8.010	1351590	20.0
Camphene	6.998	2.2	ng/ul	146305	1	8.010	1351590	20.0
Eucalyptol	8.322	18.2	ng/ul	1231410	1	8.010	1351590	20.0
Acetophe none	8.834	2.2	ng/ul	149353	1	8.010	1351590	20.0
endo-Borneol	10.598	3.7	ng/ul	375244	2	10.828	2024110	20.0
Benzophenone	15.998	6.6	ng/ul	973580	3	14.645	2932620	20.0
1,2,3,4,5,6-Hex...	17.098	3.4	ng/ul	685610	4	17.386	4016840	20.0
1H-Inden-1-one,...	18.145	2.0	ng/ul	409002	4	17.386	4016840	20.0
trans-2,3-Epoxy...	18.280	4.1	ng/ul	821717	4	17.386	4016840	20.0
Magnolol	19.921	3.4	ng/ul	701981	5	21.562	4120770	20.0
Benzo[lmn][3,8]...	20.645	3.0	ng/ul	624155	5	21.562	4120770	20.0
Cinnamyl cinnamate	21.115	3.9	ng/ul	795975	5	21.562	4120770	20.0
(DEL) Alkane: S...	21.280	2.6	ng/ul	531440	5	21.562	4120770	20.0