

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030123\
 Data File : BG056781.D
 Acq On : 2 Mar 2023 3:26
 Operator : CG/JU
 Sample : 01649-06
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBZV3

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_G\Methods\SFAM-EPA-BG022323.M
 Title : SVOA CALIBRATION

Signal : TIC: BG056781.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.442	320	324	328	rBV2	4308	6087	1.89%	0.147%
2	7.081	599	603	608	rVB2	7236	10934	3.40%	0.264%
3	7.539	675	681	688	rBV	42346	69525	21.61%	1.677%
4	7.721	705	712	720	rBV	29443	48268	15.00%	1.164%
5	7.862	730	736	737	rBV3	3404	5008	1.56%	0.121%
6	7.939	743	749	757	rVB	37006	59529	18.50%	1.436%
7	8.421	824	831	837	rBV	54542	91637	28.48%	2.211%
8	8.726	879	883	888	rVB2	6913	9487	2.95%	0.229%
9	9.102	941	947	954	rVB	43753	74916	23.29%	1.807%
10	9.243	965	971	976	rBV4	3777	6710	2.09%	0.162%
11	9.590	1022	1030	1037	rBV	39274	71984	22.38%	1.736%
12	9.966	1091	1094	1098	rVB2	1138	1537	0.48%	0.037%
13	10.318	1148	1154	1160	rVB2	33811	57645	17.92%	1.391%
14	10.653	1209	1211	1215	rVB2	979	1098	0.34%	0.026%
15	10.859	1240	1246	1253	rBV	52712	88585	27.54%	2.137%
16	11.023	1270	1274	1277	rVV3	1761	2518	0.78%	0.061%
17	11.065	1277	1281	1287	rVB2	10524	16117	5.01%	0.389%
18	11.253	1305	1313	1322	rVB	77728	138905	43.18%	3.351%
19	11.376	1328	1334	1342	rVB2	30786	53140	16.52%	1.282%
20	12.727	1560	1564	1568	rVB3	1768	2077	0.65%	0.050%
21	12.810	1572	1578	1582	rVB2	691	1319	0.41%	0.032%
22	13.544	1697	1703	1706	rVB2	1861	2802	0.87%	0.068%
23	13.738	1730	1736	1739	rBV	16326	21316	6.63%	0.514%
24	13.767	1739	1741	1746	rVB2	4671	5652	1.76%	0.136%
25	13.820	1746	1750	1753	rVB2	1297	1550	0.48%	0.037%
26	13.896	1759	1763	1767	rBV4	6601	9241	2.87%	0.223%
27	14.349	1835	1840	1844	rVB4	16458	20907	6.50%	0.504%
28	14.408	1844	1850	1857	rBV	103044	154093	47.90%	3.717%
29	14.719	1897	1903	1911	rVB	116469	179744	55.87%	4.336%
30	14.989	1941	1949	1951	rVV	33617	54212	16.85%	1.308%
31	15.025	1951	1955	1961	rVV	143415	216700	67.36%	5.227%
32	15.077	1961	1964	1967	rVB3	2386	3194	0.99%	0.077%
33	15.171	1973	1980	1990	rBV	35201	55414	17.23%	1.337%
34	15.254	1990	1994	1995	rVV2	1617	1431	0.44%	0.035%
35	15.277	1995	1998	2002	rVV3	4470	5401	1.68%	0.130%
36	15.336	2002	2008	2011	rVB3	1075	1754	0.55%	0.042%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
 Title : SVOA CALIBRATION

37	15.759	2076	2080	2085	rVV2	12340	15473	4.81%	0.373%
38	16.006	2114	2122	2128	rVB	161396	233261	72.51%	5.627%
39	16.100	2133	2138	2145	rBV	47217	68004	21.14%	1.640%
40	16.352	2175	2181	2187	rBV	38282	58478	18.18%	1.411%
41	16.429	2191	2194	2197	rVV	1193	1183	0.37%	0.029%
42	16.470	2197	2201	2209	rVB3	2031	2574	0.80%	0.062%
43	16.558	2212	2216	2219	rBV2	1642	1849	0.57%	0.045%
44	16.711	2239	2242	2246	rVB	2131	2409	0.75%	0.058%
45	16.822	2258	2261	2265	rVB2	813	1209	0.38%	0.029%
46	16.916	2273	2277	2280	rVB3	2125	2563	0.80%	0.062%
47	17.351	2345	2351	2355	rBV2	940	1732	0.54%	0.042%
48	17.422	2358	2363	2366	rVV2	7272	9272	2.88%	0.224%
49	17.451	2366	2368	2373	rVB3	3745	4191	1.30%	0.101%
50	17.698	2408	2410	2413	rVV	1207	1557	0.48%	0.038%
51	17.757	2413	2420	2427	rVV	197663	298709	92.85%	7.206%
52	17.851	2430	2436	2443	rVV2	165050	251883	78.30%	6.076%
53	17.945	2443	2452	2455	rVB4	2183	5083	1.58%	0.123%
54	18.391	2524	2528	2532	rBV2	3841	5100	1.59%	0.123%
55	18.497	2542	2546	2551	rVB2	7146	9575	2.98%	0.231%
56	18.597	2558	2563	2564	rBV	21330	26052	8.10%	0.628%
57	18.750	2586	2589	2593	rBV2	2456	2902	0.90%	0.070%
58	18.820	2596	2601	2602	rBV3	832	1084	0.34%	0.026%
59	18.850	2602	2606	2609	rVB2	5731	6405	1.99%	0.155%
60	19.014	2630	2634	2635	rVV2	1051	1033	0.32%	0.025%
61	19.196	2659	2665	2668	rBV5	3732	5107	1.59%	0.123%
62	19.543	2722	2724	2728	rVB2	1739	1676	0.52%	0.040%
63	19.607	2731	2735	2738	rBV3	2315	3456	1.07%	0.083%
64	19.684	2742	2748	2751	rVV6	4229	7675	2.39%	0.185%
65	19.748	2754	2759	2763	rBV4	2469	4930	1.53%	0.119%
66	19.819	2769	2771	2772	rBV	2318	1572	0.49%	0.038%
67	19.848	2772	2776	2777	rBV2	2682	3484	1.08%	0.084%
68	19.936	2789	2791	2793	rBV3	1246	1143	0.36%	0.028%
69	19.983	2794	2799	2801	rBV4	2356	3703	1.15%	0.089%
70	20.113	2815	2821	2830	rVB	218716	315793	98.16%	7.618%
71	20.207	2835	2837	2841	rBV4	1628	1799	0.56%	0.043%
72	20.330	2856	2858	2861	rBV3	1120	1294	0.40%	0.031%
73	20.448	2876	2878	2879	rBV	1796	1117	0.35%	0.027%
74	20.465	2879	2881	2885	rVV3	4880	6038	1.88%	0.146%
75	20.536	2888	2893	2897	rVV2	25143	34049	10.58%	0.821%
76	20.853	2944	2947	2950	rVB3	5679	5693	1.77%	0.137%

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77	20.924	2957	2959	2962	rVB4	2642	1985	0.62%	0.048%
78	21.065	2972	2983	2987	rBV2	21601	43128	13.41%	1.040%
79	21.288	3018	3021	3026	rVB5	8316	11174	3.47%	0.270%
80	21.352	3031	3032	3040	rVB6	3775	5651	1.76%	0.136%
81	21.458	3046	3050	3056	rVB	42132	57568	17.89%	1.389%
82	21.546	3060	3065	3068	rBV4	5820	8856	2.75%	0.214%
83	21.640	3075	3081	3091	rBV3	33509	61698	19.18%	1.488%
84	22.063	3146	3153	3162	rVB2	187712	318899	99.13%	7.693%
85	22.492	3222	3226	3231	rVB3	9694	16515	5.13%	0.398%
86	22.604	3243	3245	3247	rBV2	1764	1547	0.48%	0.037%
87	22.792	3276	3277	3280	rBV3	2171	2010	0.62%	0.048%
88	22.886	3288	3293	3299	rBV6	4823	11524	3.58%	0.278%
89	23.297	3361	3363	3364	rVB	2668	1153	0.36%	0.028%
90	23.620	3411	3418	3419	rBV5	8669	13556	4.21%	0.327%
91	23.773	3442	3444	3447	rBV4	2777	2621	0.81%	0.063%
92	23.944	3471	3473	3478	rVB5	3996	6364	1.98%	0.154%
93	25.260	3695	3697	3700	rBV3	2147	2583	0.80%	0.062%
94	25.342	3700	3711	3725	rVV2	93889	289349	89.94%	6.980%
95	25.589	3741	3753	3763	rBV2	106631	321704	100.00%	7.760%
96	26.505	3907	3909	3912	rVB3	1573	1006	0.31%	0.024%
97	28.479	4244	4245	4248	rVB2	2380	1302	0.40%	0.031%
98	29.713	4453	4455	4458	rVB3	2496	2056	0.64%	0.050%
99	30.424	4574	4576	4578	rBV3	1568	1353	0.42%	0.033%
100	30.459	4581	4582	4584	rBV2	2405	1350	0.42%	0.033%

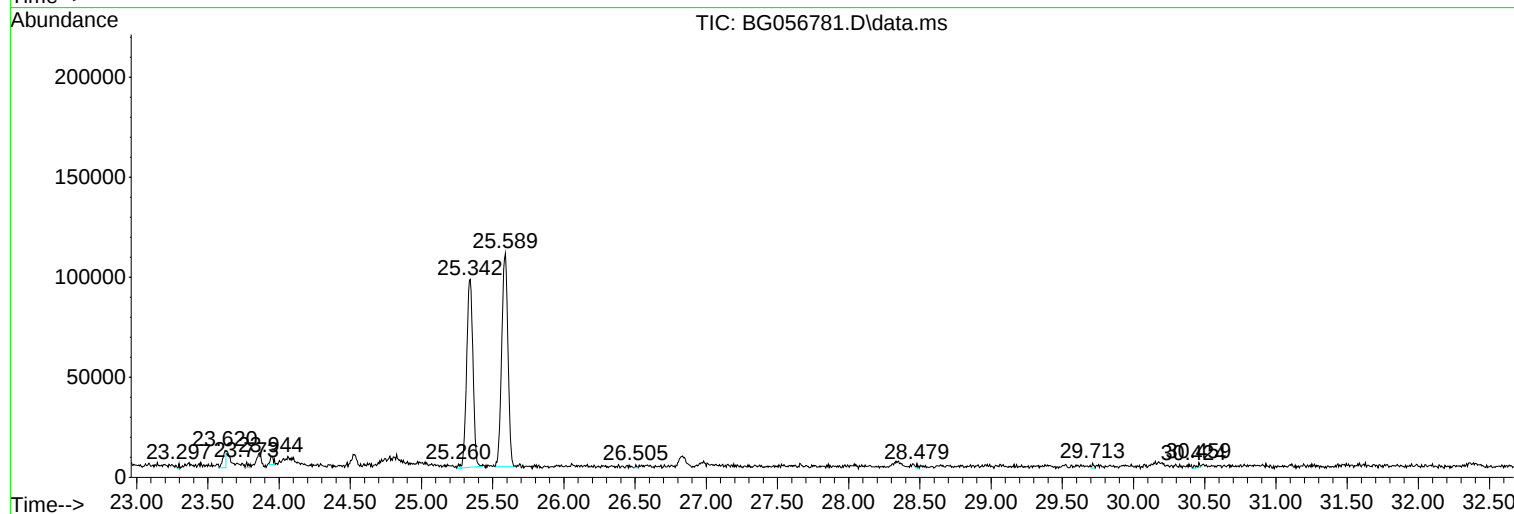
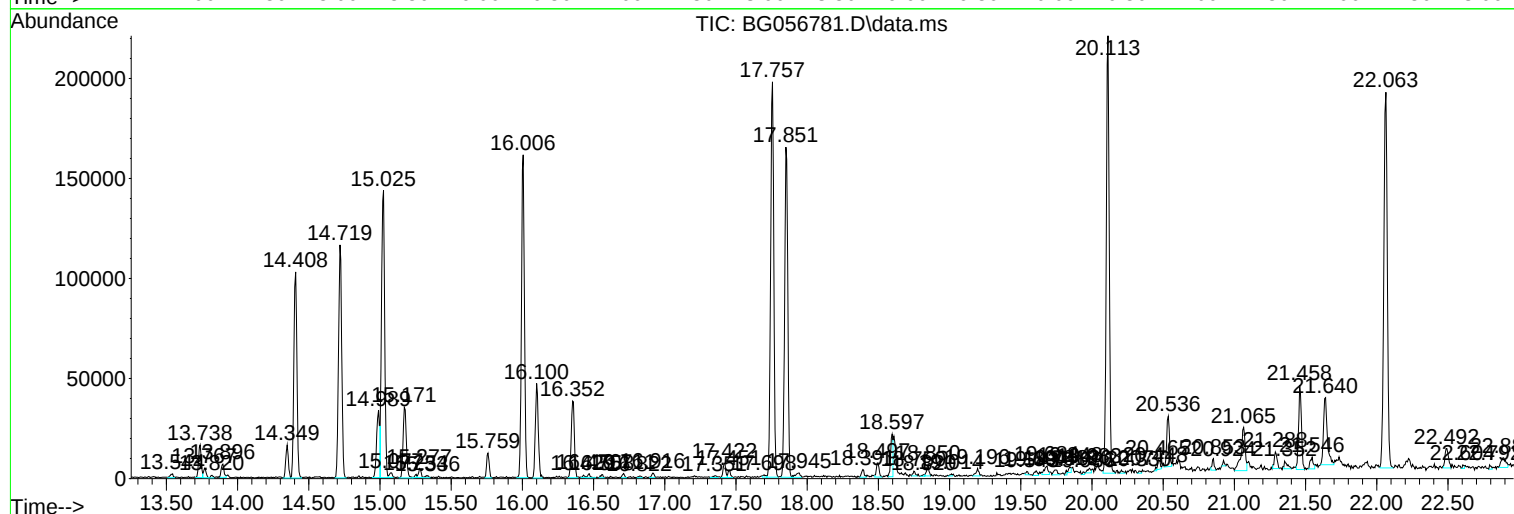
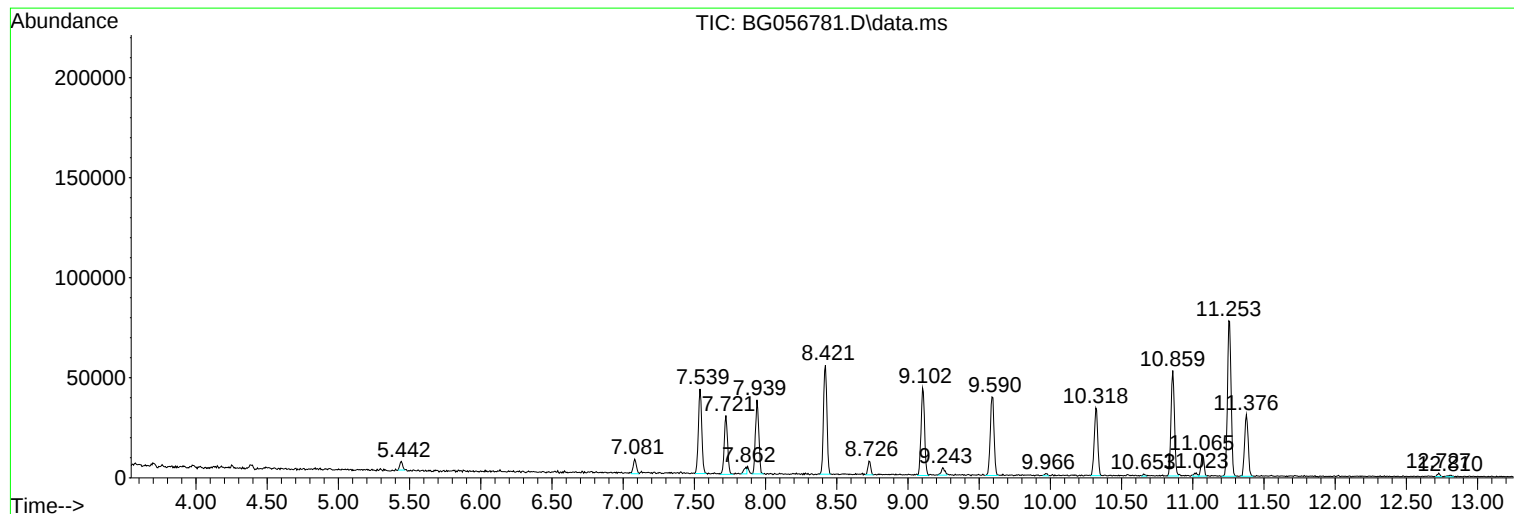
Sum of corrected areas: 4145499

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

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



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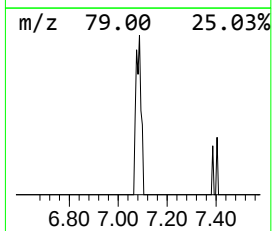
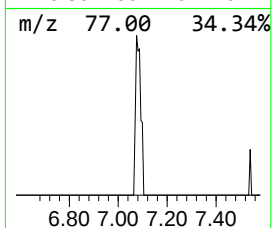
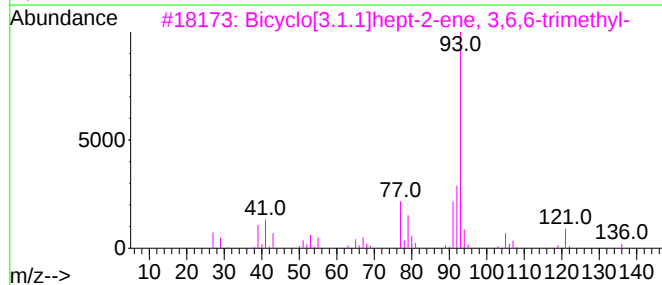
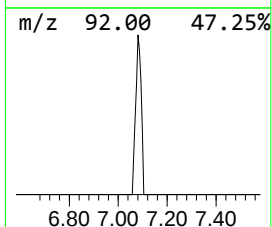
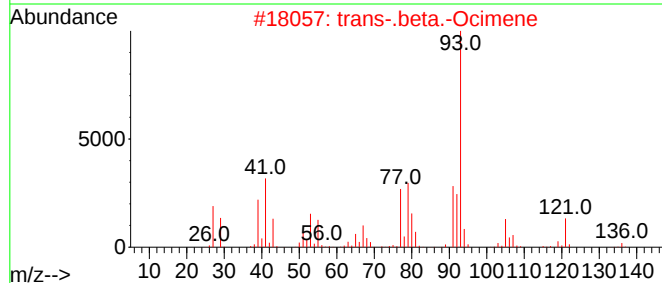
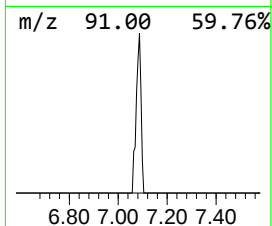
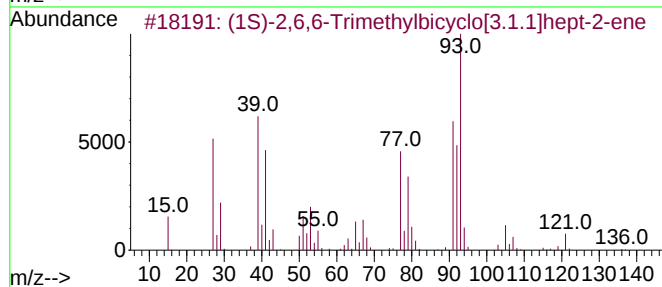
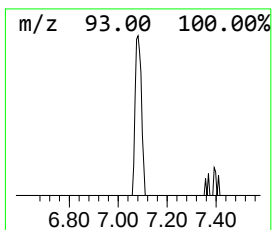
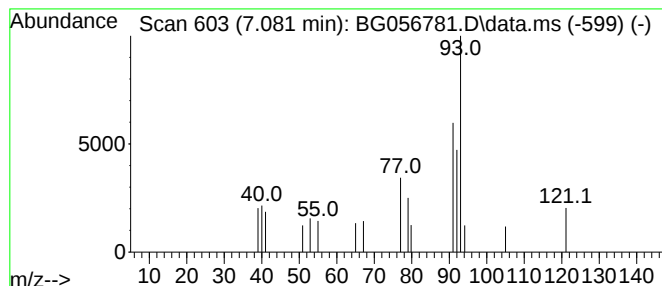
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (1S)-2,6,6-Trimethylbicyclo... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.081	2.39 ng/ul	10934	1,4-Dichlorobenzene-d4	8.421

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1S)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-26-4	83		
2	trans-.beta.-Ocimene	136	C10H16	003779-61-1	72		
3	Bicyclo[3.1.1]hept-2-ene, 3,6,6-trimethyl-	136	C10H16	004889-83-2	64		
4	3,5-Methanocyclopentapyrazole, 3,6-dimethyl-	164	C10H16N2	087143-58-6	50		
5	1-Hexen-3-yne, 5,5-dimethyl-	108	C8H12	004911-58-4	38		



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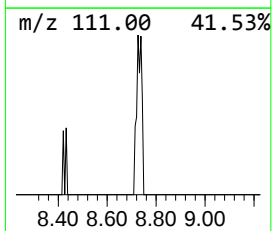
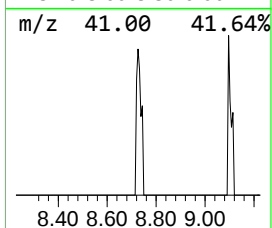
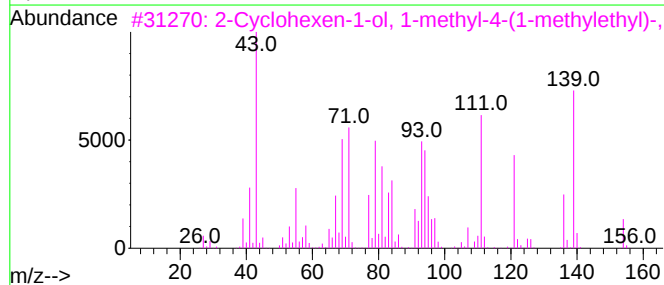
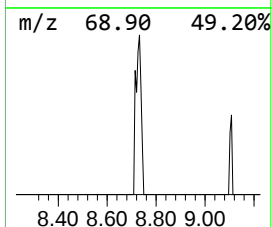
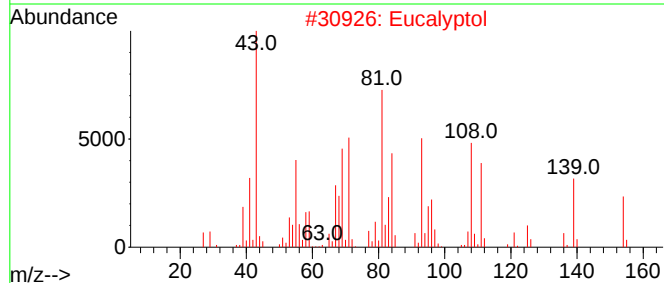
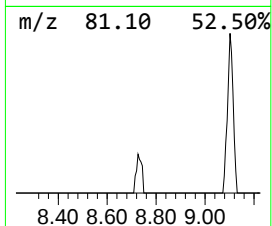
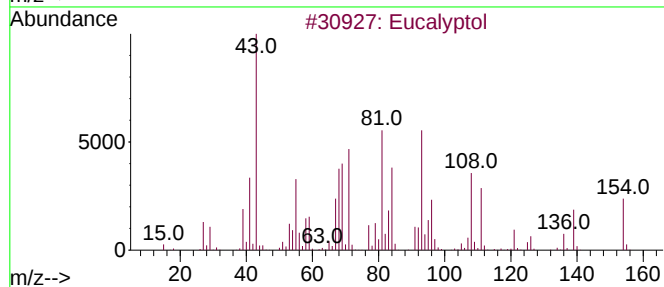
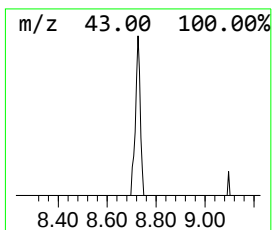
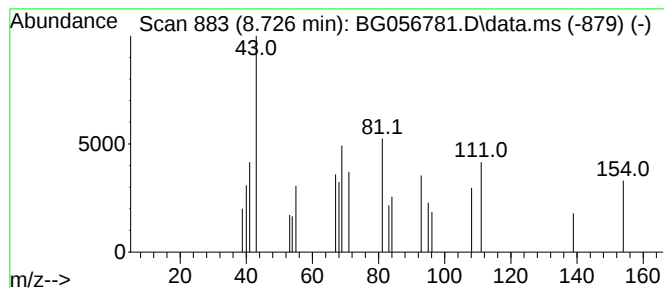
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Eucalyptol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.726	2.07 ng/ul	9487	1,4-Dichlorobenzene-d4	8.421

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eucalyptol	154	C10H18O	000470-82-6	70
2			Eucalyptol	154	C10H18O	000470-82-6	59
3			2-Cyclohexen-1-ol, 1-methyl-4-(1...	154	C10H18O	029803-82-5	33
4			5-Hepten-2-one, 6-methyl-	126	C8H14O	000110-93-0	32
5			2-Heptanone, 3-propylidene-	154	C10H18O	032064-70-3	30



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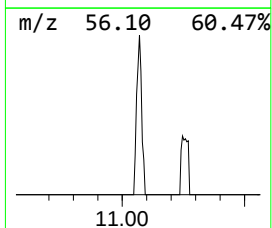
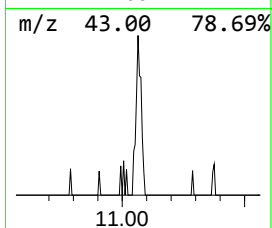
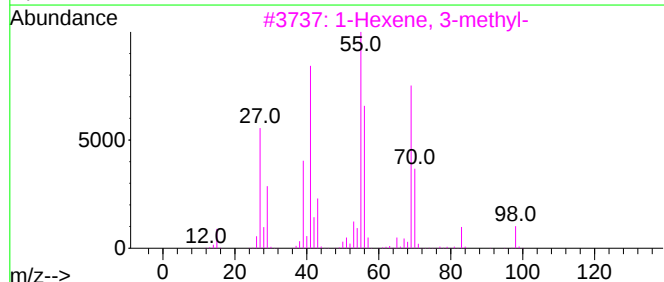
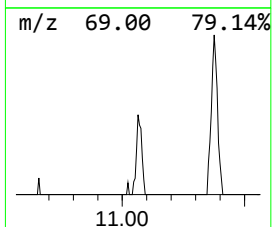
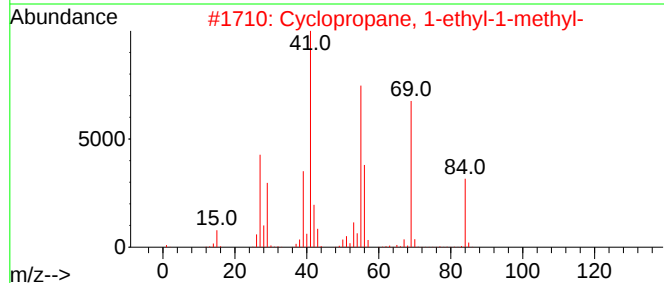
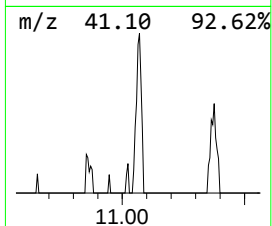
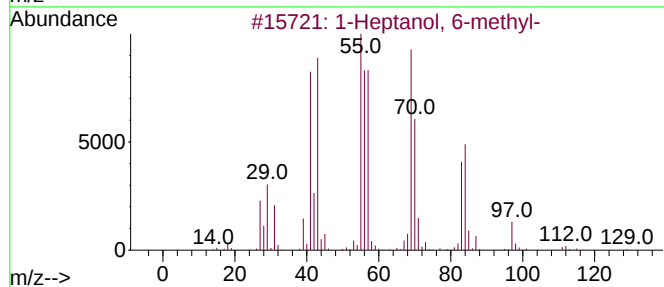
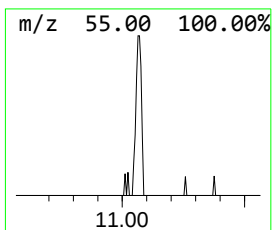
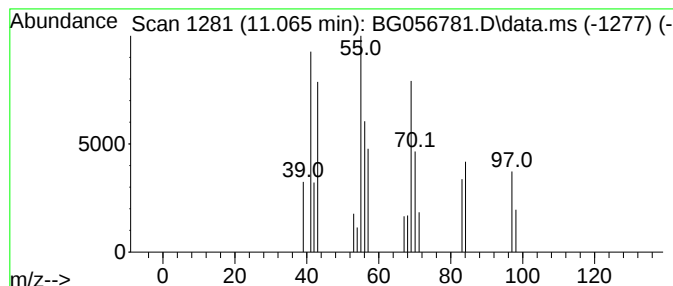
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Heptanol, 6-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.065	2.32 ng/ul	16117	Naphthalene-d8	11.253

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptanol, 6-methyl-			130	C8H18O	001653-40-3	64
2	Cyclopropane, 1-ethyl-1-methyl-			84	C6H12	053778-43-1	52
3	1-Hexene, 3-methyl-			98	C7H14	003404-61-3	50
4	Pentane, 3-methylene-			84	C6H12	000760-21-4	46
5	1-Nonanol			144	C9H20O	000143-08-8	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030123\
Data File : BG056781.D
Acq On : 2 Mar 2023 3:26
Operator : CG/JU
Sample : 01649-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZV3

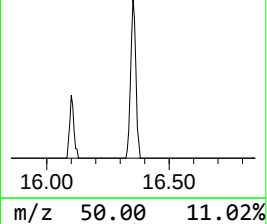
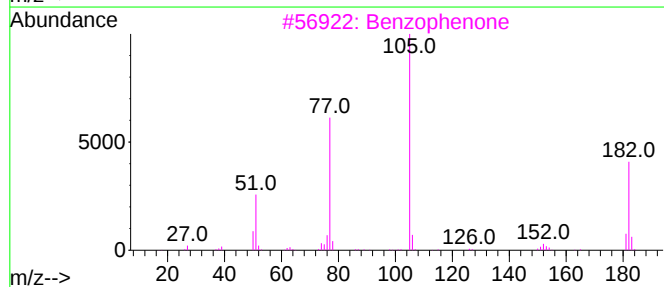
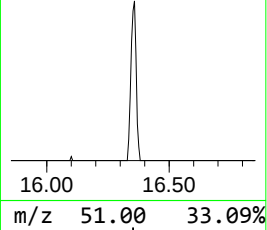
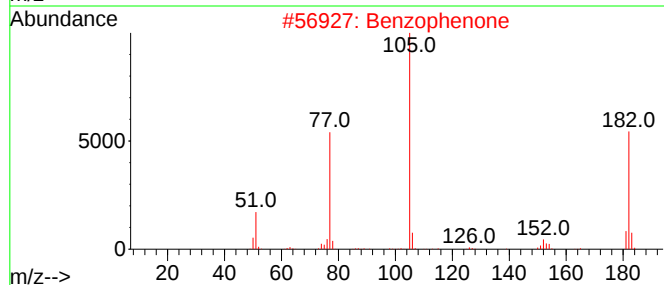
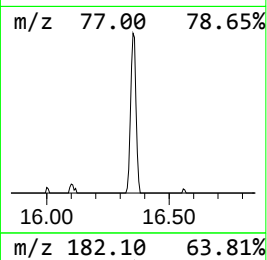
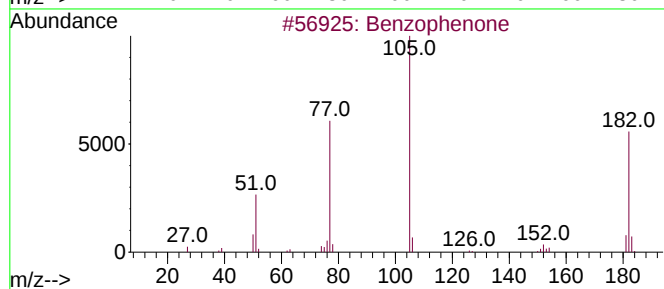
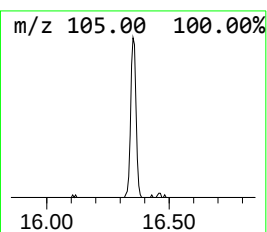
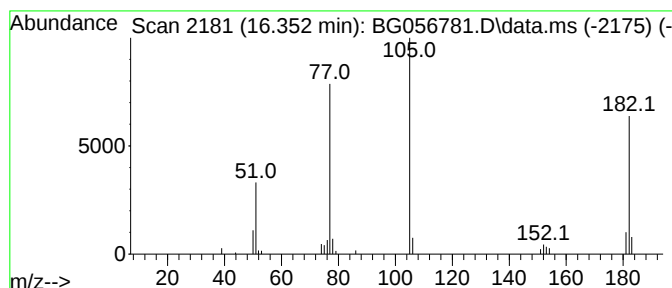
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzophenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.352	5.40 ng/ul	58478	Acenaphthene-d10	15.025

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030123\
Data File : BG056781.D
Acq On : 2 Mar 2023 3:26
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Sample : 01649-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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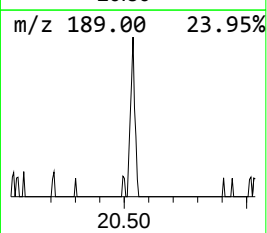
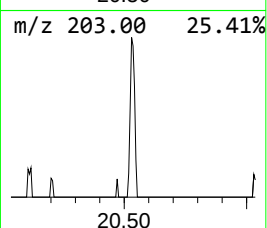
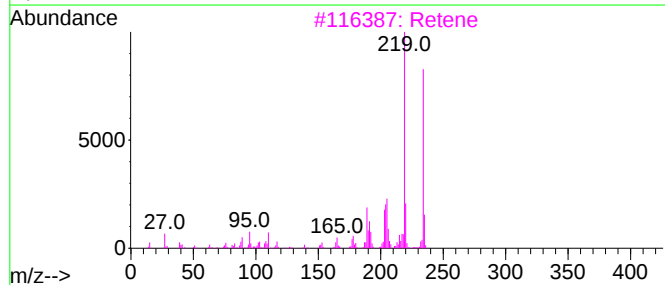
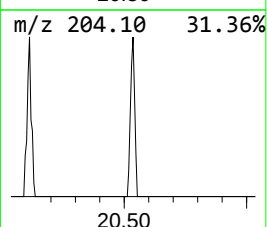
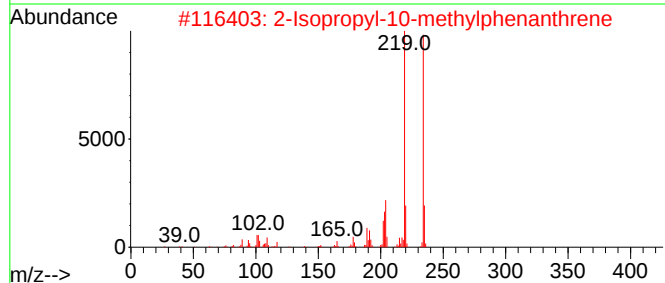
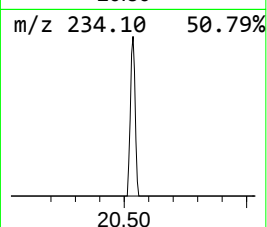
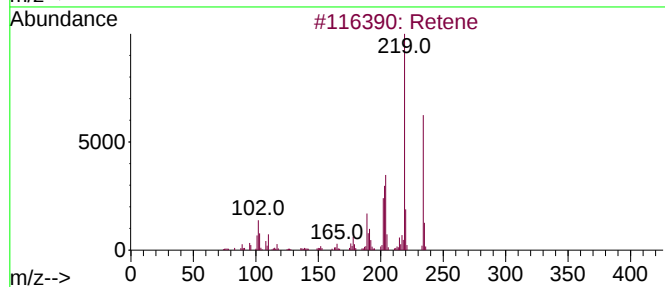
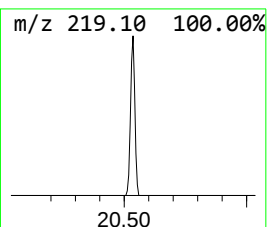
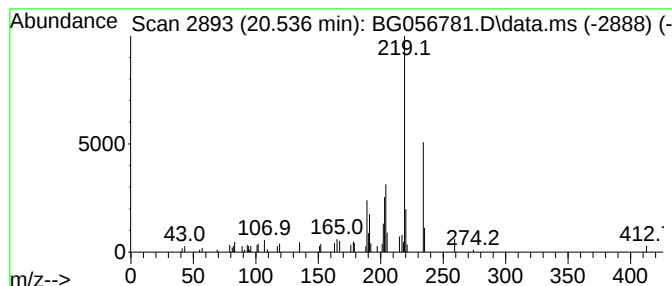
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Retene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.536	2.14 ng/ul	34049	Chrysene-d12	22.063

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Retene			234	C18H18	000483-65-8	93
2	2-Isopropyl-10-methylphenanthrene			234	C18H18	066552-97-4	91
3	Retene			234	C18H18	000483-65-8	90
4	Retene			234	C18H18	000483-65-8	90
5	Retene			234	C18H18	000483-65-8	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030123\
Data File : BG056781.D
Acq On : 2 Mar 2023 3:26
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Sample : 01649-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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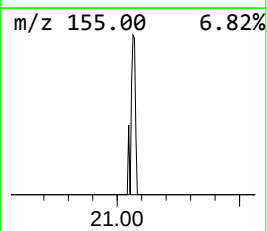
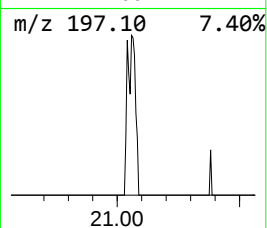
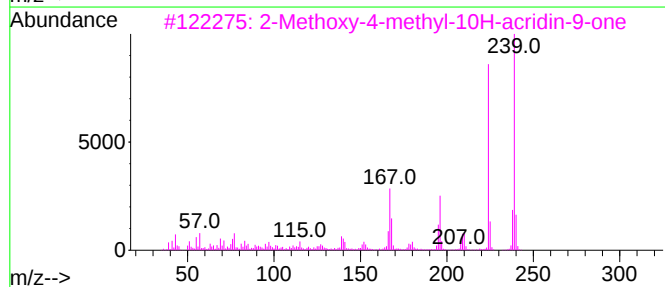
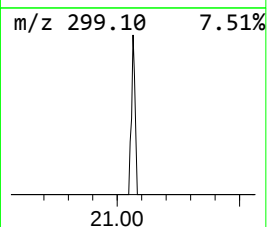
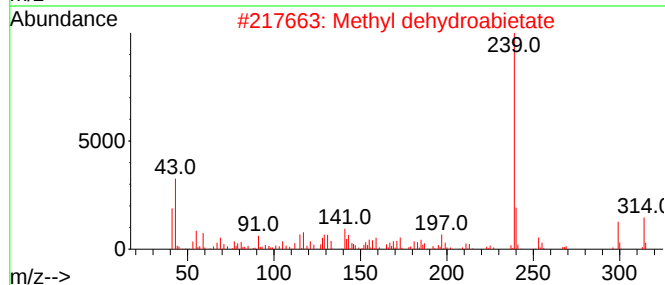
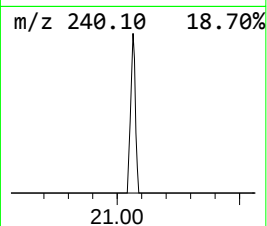
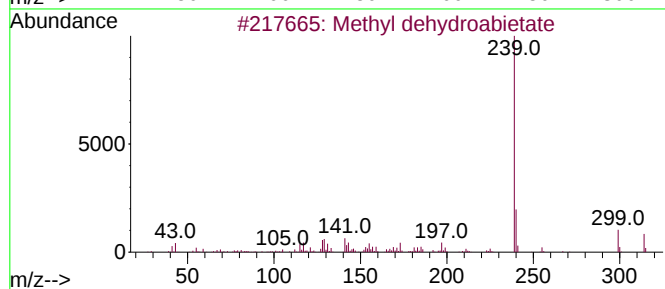
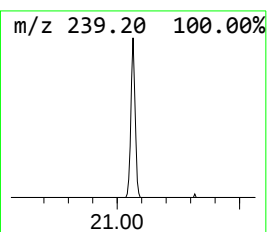
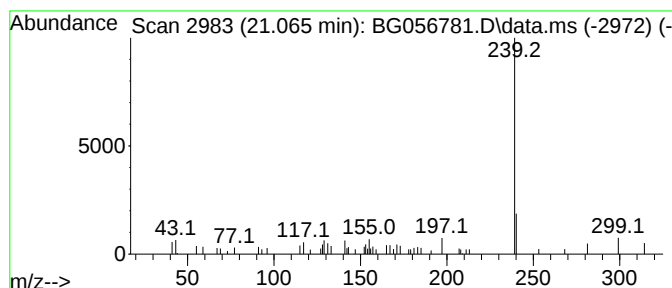
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Methyl dehydroabietate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.065	2.70 ng/ul	43128	Chrysene-d12	22.063

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl dehydroabietate	314	C21H30O2	001235-74-1	87
2			Methyl dehydroabietate	314	C21H30O2	001235-74-1	87
3			2-Methoxy-4-methyl-10H-acridin-9...	239	C15H13NO2	1000210-25-6	59
4			2-Methyl-3-phenylsulfanyl-1H-indole	239	C15H13NS	1000322-52-7	53
5			Indole, 3-imidazolo[2,1-b]thiazo...	239	C13H9N3S	292855-05-1	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030123\
Data File : BG056781.D
Acq On : 2 Mar 2023 3:26
Operator : CG/JU
Sample : 01649-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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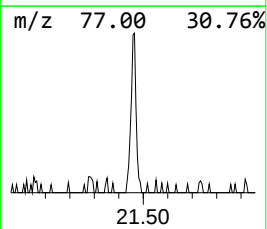
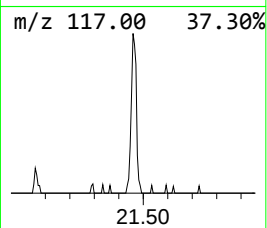
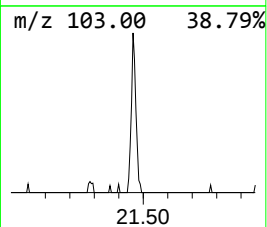
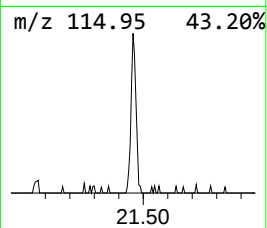
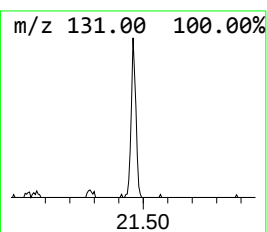
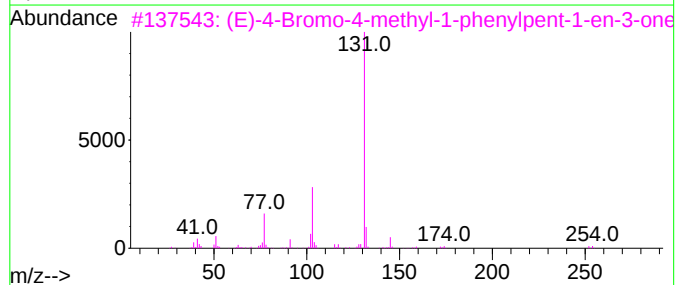
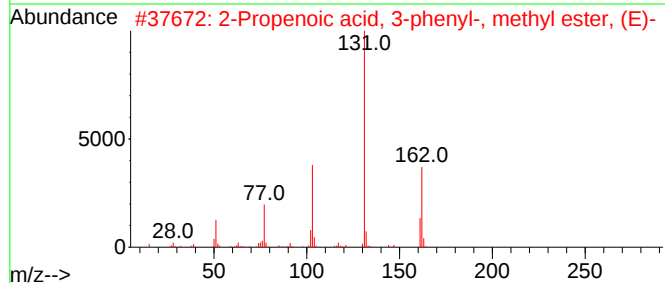
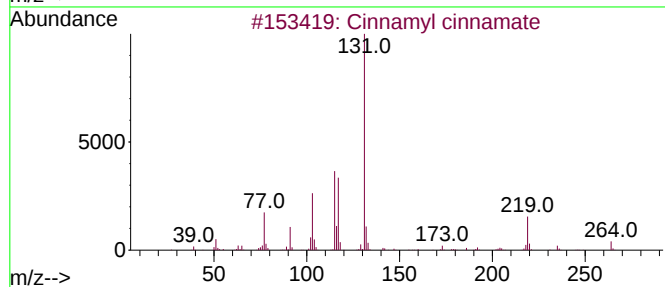
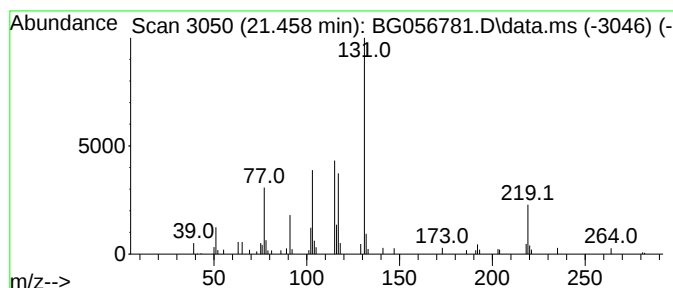
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Cinnamyl cinnamate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.458	3.61 ng/ul	57568	Chrysene-d12	22.063

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cinnamyl cinnamate	264	C18H16O2	000122-69-0	70
2			2-Propenoic acid, 3-phenyl-, met...	162	C10H10O2	001754-62-7	47
3			(E)-4-Bromo-4-methyl-1-phenylpen...	252	C12H13BrO	1000443-02-8	47
4			2-Nitrophenyl cinnamamide	268	C15H12N2O3	111273-37-1	46
5			p-Vinylbenzohydrazide	162	C9H10N2O	1000244-74-9	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030123\
Data File : BG056781.D
Acq On : 2 Mar 2023 3:26
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Sample : 01649-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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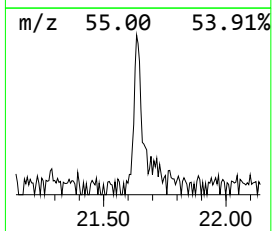
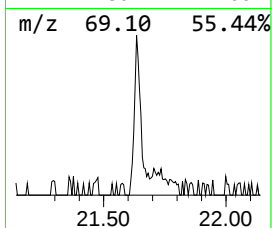
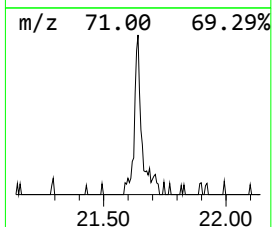
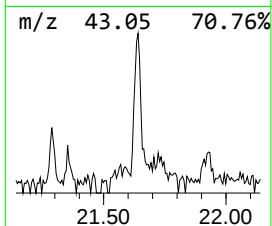
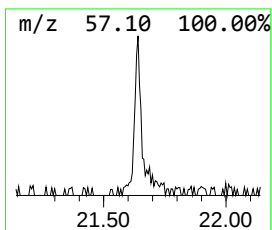
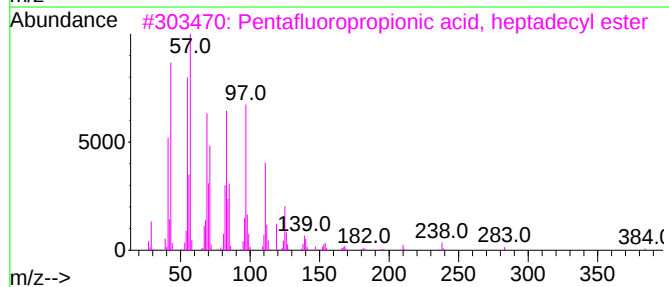
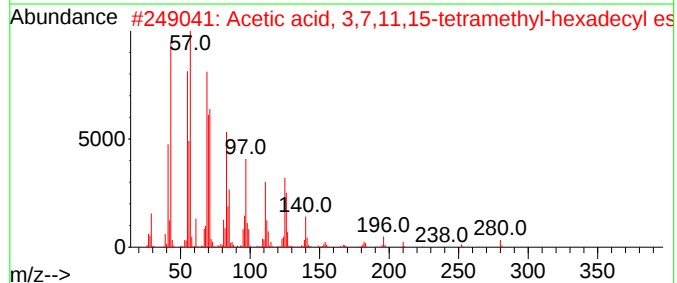
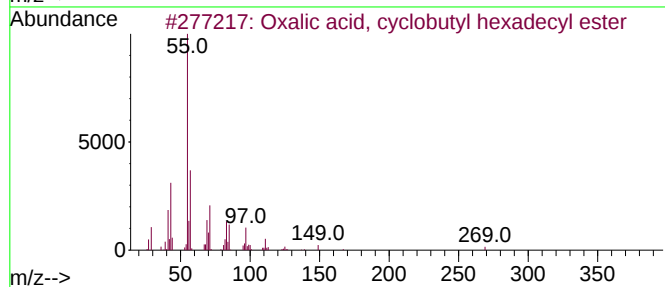
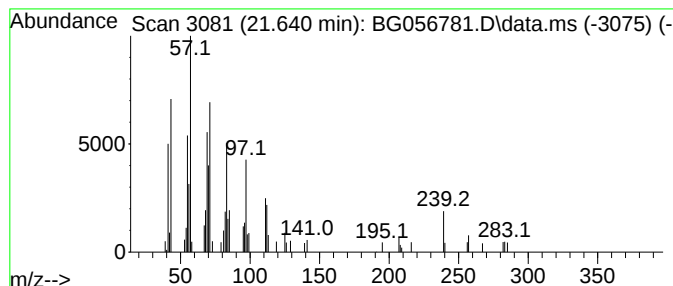
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Oxalic acid, cyclobutyl hex... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.640	3.87 ng/ul	61698	Chrysene-d12	22.063

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, cyclobutyl hexadecy...	368	C22H40O4	1000309-70-6	64
2			Acetic acid, 3,7,11,15-tetrameth...	340	C22H44O2	1000193-63-0	64
3			Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	58
4			1-Hexacosene	364	C26H52	018835-33-1	58
5			3-Eicosene, (E)-	280	C20H40	074685-33-9	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030123\
Data File : BG056781.D
Acq On : 2 Mar 2023 3:26
Operator : CG/JU
Sample : 01649-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZV3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.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
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Eucalyptol	8.726	2.1	ng/ul	9487	1	8.421	91637	20.0
1-Heptanol, 6-m...	11.065	2.3	ng/ul	16117	2	11.253	138905	20.0
Benzophenone	16.352	5.4	ng/ul	58478	3	15.025	216700	20.0
Retene	20.536	2.1	ng/ul	34049	5	22.063	318899	20.0
Methyl dehydroa...	21.065	2.7	ng/ul	43128	5	22.063	318899	20.0
Cinnamyl cinnamate	21.458	3.6	ng/ul	57568	5	22.063	318899	20.0
Oxalic acid, cy...	21.640	3.9	ng/ul	61698	5	22.063	318899	20.0