

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
 Data File : BG056824.D
 Acq On : 4 Mar 2023 5:32
 Operator : CG/JU
 Sample : 01711-13
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBAC9

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: BG056824.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.437	320	323	331	rVB2	5694	9123	2.68%	0.178%
2	7.083	598	603	608	rVB2	13350	19983	5.86%	0.391%
3	7.394	653	656	662	rVB3	3927	5875	1.72%	0.115%
4	7.541	675	681	689	rVB	42967	74502	21.86%	1.456%
5	7.717	706	711	719	rVB	29713	49886	14.64%	0.975%
6	7.852	729	734	739	rVV4	12509	19963	5.86%	0.390%
7	7.940	742	749	756	rBV2	38030	65763	19.30%	1.285%
8	8.416	824	830	837	rVB	56335	93421	27.41%	1.826%
9	8.546	849	852	856	rBV2	1944	2536	0.74%	0.050%
10	8.634	863	867	873	rVB4	3203	4739	1.39%	0.093%
11	9.104	941	947	955	rVB2	47976	82340	24.16%	1.609%
12	9.245	965	971	980	rVB3	6223	11778	3.46%	0.230%
13	9.591	1023	1030	1038	rVB	44931	79225	23.25%	1.549%
14	10.320	1148	1154	1160	rBV3	35530	62943	18.47%	1.230%
15	10.860	1239	1246	1252	rBV	58313	99448	29.18%	1.944%
16	11.254	1306	1313	1321	rVB	86966	150749	44.24%	2.947%
17	11.378	1329	1334	1339	rBV2	2584	4103	1.20%	0.080%
18	12.024	1439	1444	1447	rVB2	1180	1806	0.53%	0.035%
19	13.786	1742	1744	1747	rVV	1076	1351	0.40%	0.026%
20	13.816	1747	1749	1754	rVV3	1937	2807	0.82%	0.055%
21	13.892	1757	1762	1767	rVV2	22911	31175	9.15%	0.609%
22	14.092	1793	1796	1800	rBV3	1673	1468	0.43%	0.029%
23	14.180	1806	1811	1816	rVB3	18094	24779	7.27%	0.484%
24	14.345	1834	1839	1844	rBV3	23881	33273	9.76%	0.650%
25	14.403	1844	1849	1860	rVB	98301	149953	44.00%	2.931%
26	14.562	1870	1876	1881	rVB2	46008	65118	19.11%	1.273%
27	14.615	1881	1885	1888	rBV4	2551	3089	0.91%	0.060%
28	14.721	1896	1903	1909	rBV2	123256	196621	57.70%	3.843%
29	14.785	1911	1914	1916	rBV3	1516	1689	0.50%	0.033%
30	14.850	1920	1925	1928	rBV2	6969	8619	2.53%	0.168%
31	14.973	1942	1946	1948	rVV3	3698	4954	1.45%	0.097%
32	15.020	1948	1954	1959	rVV2	151224	243364	71.42%	4.757%
33	15.067	1959	1962	1967	rVV3	23758	35473	10.41%	0.693%
34	15.126	1967	1972	1975	rVV5	8519	12700	3.73%	0.248%
35	15.173	1975	1980	1986	rVV2	33032	52040	15.27%	1.017%
36	15.244	1988	1992	1997	rVB5	4425	8125	2.38%	0.159%

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Smoothing : OFF

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Stop Thrs : 0

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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

37	15.314	1997	2004	2008	rBV4	11580	24713	7.25%	0.483%
38	15.373	2010	2014	2022	rVB5	16578	34158	10.02%	0.668%
39	15.761	2077	2080	2083	rVB3	1249	1379	0.40%	0.027%
40	15.896	2097	2103	2104	rBV4	1911	2568	0.75%	0.050%
41	15.937	2108	2110	2115	rVV2	1732	2688	0.79%	0.053%
42	16.002	2115	2121	2128	rVV	171819	259880	76.26%	5.080%
43	16.101	2131	2138	2143	rBV2	39139	57301	16.81%	1.120%
44	16.225	2157	2159	2165	rVB2	2979	3856	1.13%	0.075%
45	16.354	2176	2181	2186	rBV	19031	26745	7.85%	0.523%
46	16.466	2195	2200	2205	rVV4	8403	13238	3.88%	0.259%
47	17.241	2330	2332	2337	rBV	1095	1869	0.55%	0.037%
48	17.453	2363	2368	2372	rVV5	4438	6128	1.80%	0.120%
49	17.752	2410	2419	2427	rBV	209349	311270	91.34%	6.084%
50	17.852	2430	2436	2443	rBV	180586	265348	77.87%	5.187%
51	18.340	2513	2519	2520	rVV4	1743	2444	0.72%	0.048%
52	18.487	2540	2544	2545	rBV4	3960	4943	1.45%	0.097%
53	18.593	2558	2562	2569	rBV2	31476	51761	15.19%	1.012%
54	18.892	2609	2613	2619	rVB2	56278	82878	24.32%	1.620%
55	19.068	2638	2643	2648	rVB2	85855	122989	36.09%	2.404%
56	19.151	2654	2657	2661	rVB3	3967	3729	1.09%	0.073%
57	19.215	2666	2668	2671	rBV3	1363	1592	0.47%	0.031%
58	19.403	2698	2700	2703	rVB4	1795	1495	0.44%	0.029%
59	19.439	2703	2706	2710	rBV3	1895	2168	0.64%	0.042%
60	19.497	2713	2716	2720	rBV2	1956	2296	0.67%	0.045%
61	19.609	2731	2735	2737	rBV4	1703	2200	0.65%	0.043%
62	19.674	2741	2746	2750	rVV5	15025	21730	6.38%	0.425%
63	19.727	2753	2755	2756	rBV2	1808	1788	0.52%	0.035%
64	19.844	2770	2775	2785	rBV	46245	68895	20.22%	1.347%
65	19.962	2794	2795	2798	rBV3	1928	1491	0.44%	0.029%
66	20.067	2811	2813	2815	rBV3	2243	2257	0.66%	0.044%
67	20.108	2815	2820	2828	rVB	213791	323509	94.93%	6.324%
68	20.314	2853	2855	2856	rBV2	1883	1544	0.45%	0.030%
69	20.355	2859	2862	2865	rBV4	2016	2562	0.75%	0.050%
70	20.473	2879	2882	2885	rBV4	3299	4499	1.32%	0.088%
71	20.526	2889	2891	2894	rBV4	3381	3444	1.01%	0.067%
72	20.596	2900	2903	2904	rBV3	4429	4891	1.44%	0.096%
73	20.878	2949	2951	2954	rVB5	4300	2747	0.81%	0.054%
74	20.984	2966	2969	2974	rVB4	7564	10365	3.04%	0.203%
75	21.060	2979	2982	2986	rVV2	10022	11862	3.48%	0.232%
76	21.096	2986	2988	2993	rVV5	4680	6749	1.98%	0.132%

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 Title : SVOA CALIBRATION

77	21.278	3017	3019	3020	rBV	3024	2387	0.70%	0.047%
78	21.378	3032	3036	3038	rBV5	3654	4852	1.42%	0.095%
79	21.460	3045	3050	3055	rVB2	35498	51253	15.04%	1.002%
80	21.636	3075	3080	3094	rVB2	74116	131017	38.45%	2.561%
81	21.930	3125	3130	3135	rBV6	11813	21094	6.19%	0.412%
82	22.065	3146	3153	3162	rVB	187336	340774	100.00%	6.661%
83	22.224	3176	3180	3181	rBV3	3759	4618	1.36%	0.090%
84	22.435	3214	3216	3217	rBV2	2622	1800	0.53%	0.035%
85	22.488	3222	3225	3230	rVB5	7485	12237	3.59%	0.239%
86	22.911	3290	3297	3313	rVB4	37420	106287	31.19%	2.078%
87	23.405	3377	3381	3388	rBV6	5703	9850	2.89%	0.193%
88	23.992	3479	3481	3482	rBV2	2812	1645	0.48%	0.032%
89	24.027	3483	3487	3495	rVB8	10149	21166	6.21%	0.414%
90	24.527	3565	3572	3580	rVB2	43416	91440	26.83%	1.787%
91	24.603	3580	3585	3591	rBV7	4606	11474	3.37%	0.224%
92	25.343	3700	3711	3721	rVB2	102563	296181	86.91%	5.789%
93	25.590	3743	3753	3766	rVB2	112033	338560	99.35%	6.618%
94	26.154	3843	3849	3859	rVB2	10465	33141	9.73%	0.648%
95	26.254	3865	3866	3870	rBV4	2444	2251	0.66%	0.044%
96	26.830	3953	3964	3976	rBV3	42515	140746	41.30%	2.751%
97	26.959	3982	3986	3987	rBV4	10771	13595	3.99%	0.266%
98	28.910	4317	4318	4320	rBV2	3241	2109	0.62%	0.041%
99	29.580	4430	4432	4437	rBV3	2351	3285	0.96%	0.064%
100	31.959	4835	4837	4838	rBV2	2594	1449	0.43%	0.028%

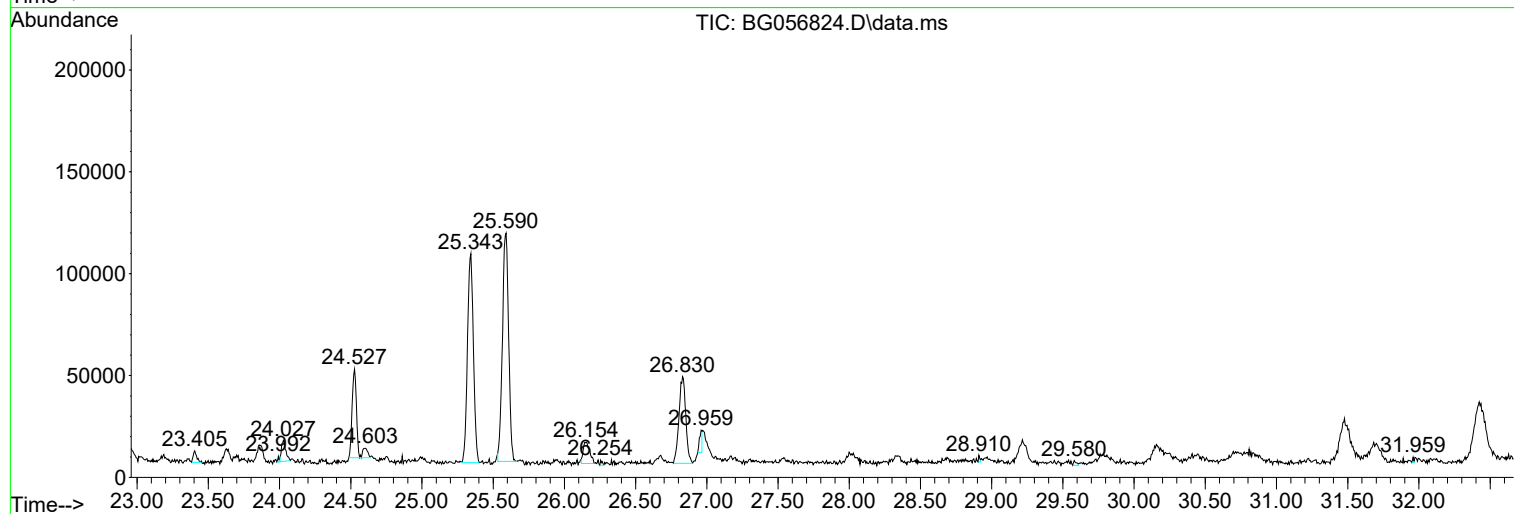
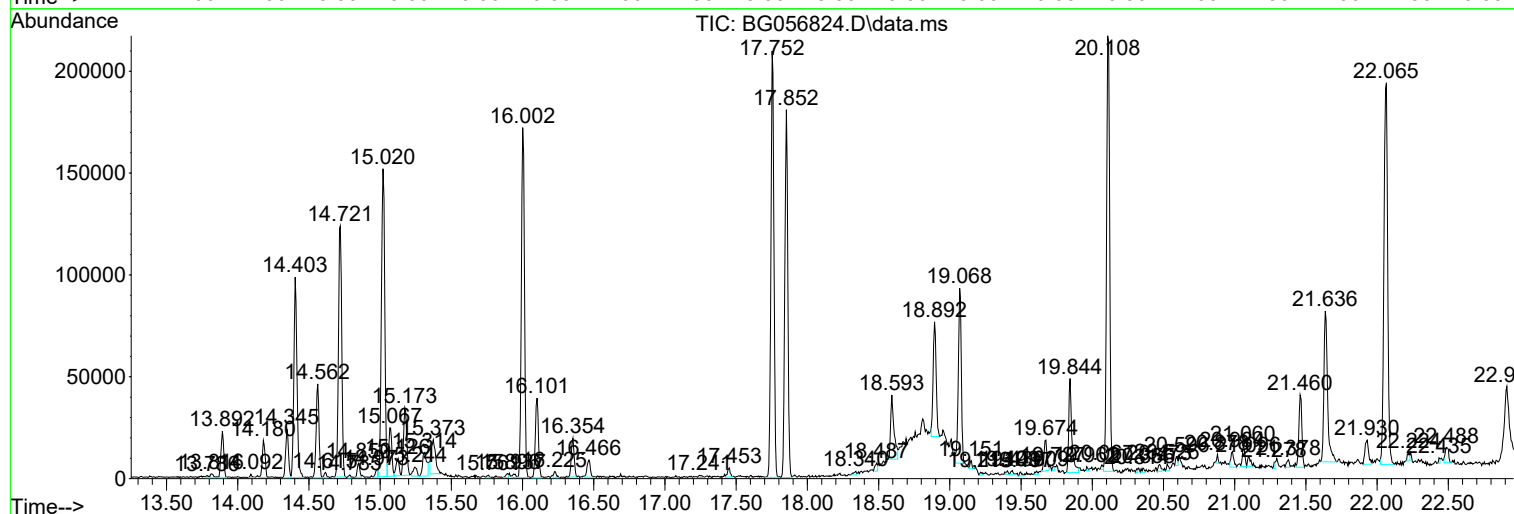
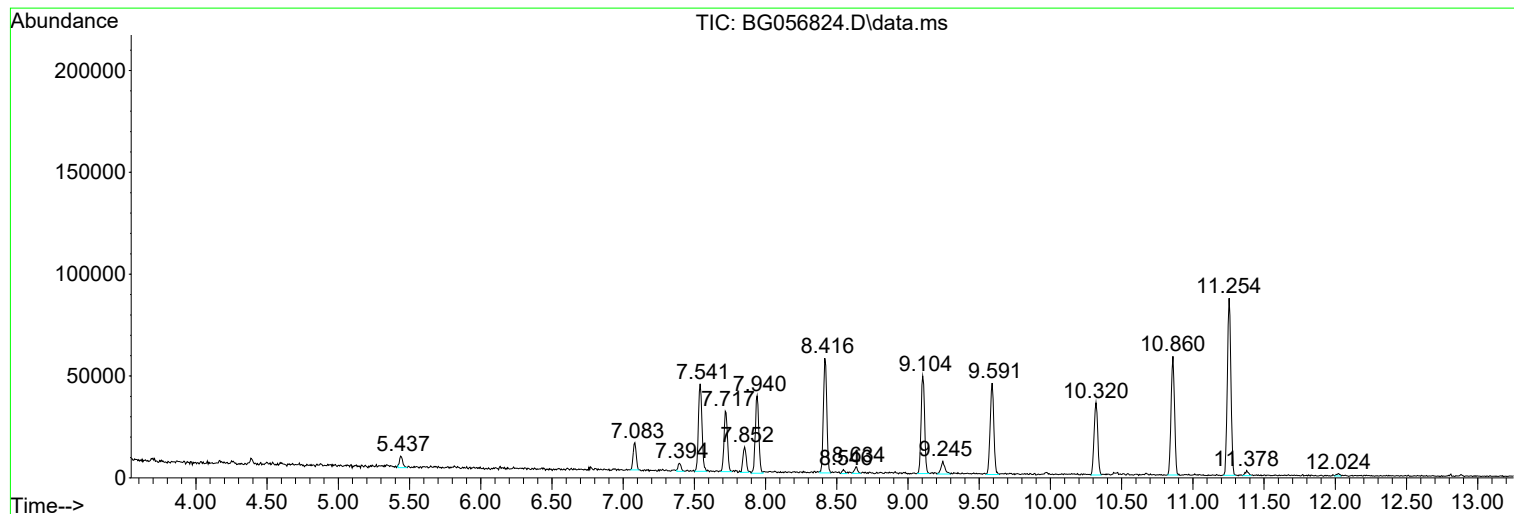
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

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



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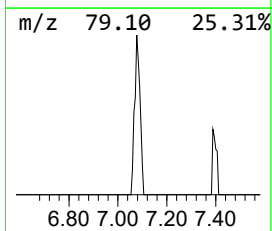
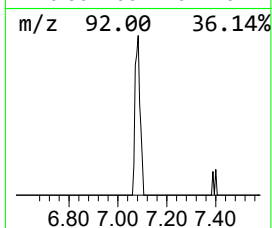
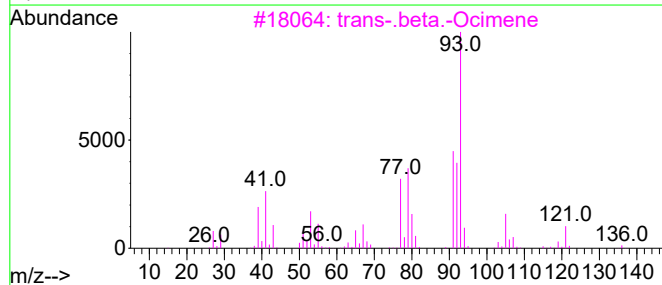
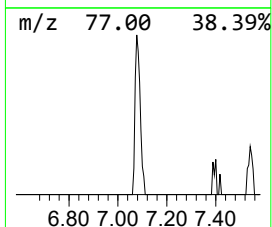
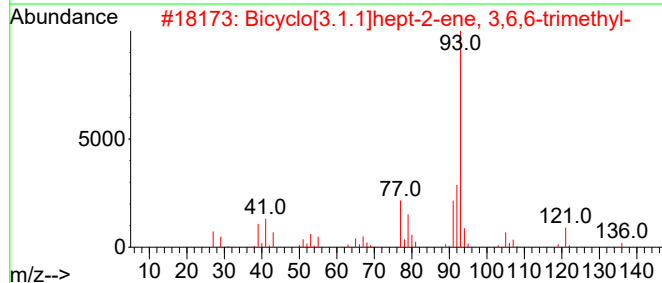
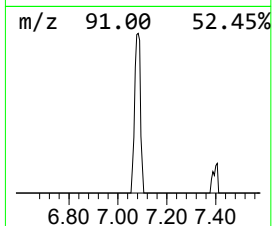
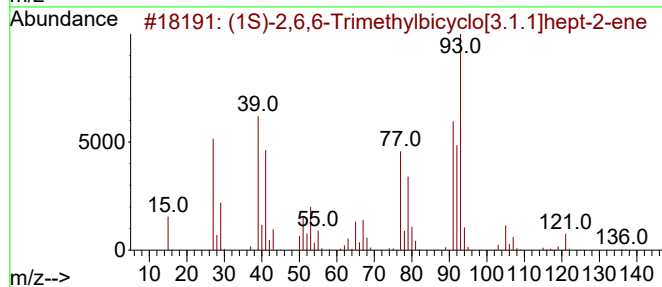
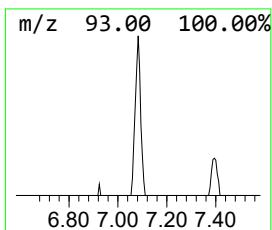
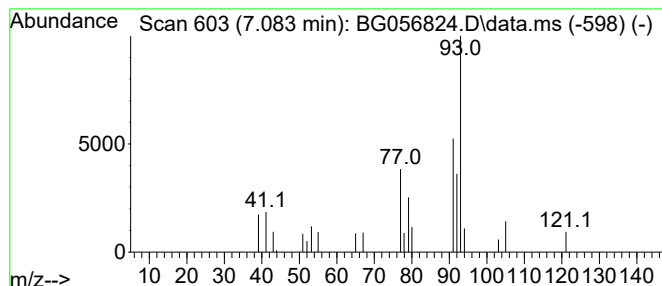
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (1S)-2,6,6-Trimethylbicyclo... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.083	4.28 ng/ul	19983	1,4-Dichlorobenzene-d4	8.422

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	Bicyclo[3.1.1]hept-2-ene, 3,6,6-trimethyl-	136	C10H16	004889-83-2	78		
3	trans-.beta.-Ocimene	136	C10H16	003779-61-1	72		
4	trans-.beta.-Ocimene	136	C10H16	003779-61-1	64		
5	3,5-Methanocyclopentapyrazole, 3,4,5-trimethyl-	164	C10H16N2	087143-58-6	59		



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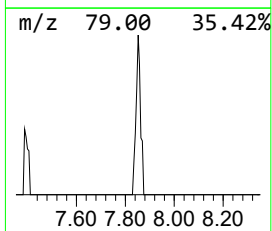
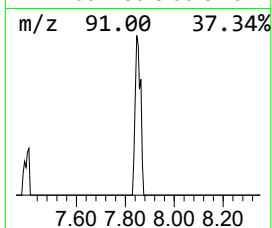
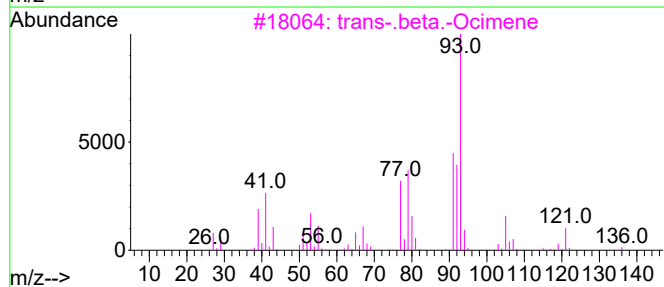
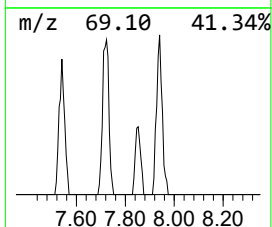
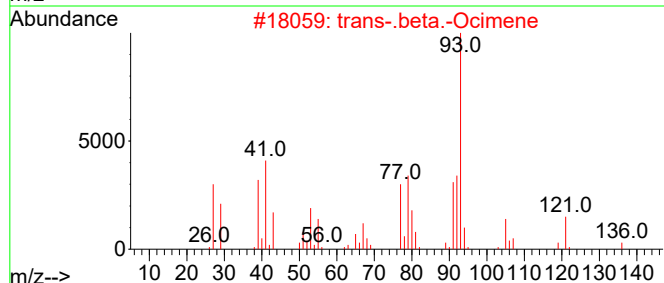
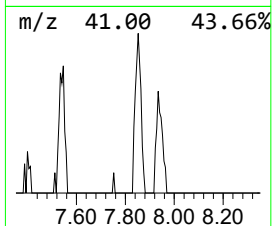
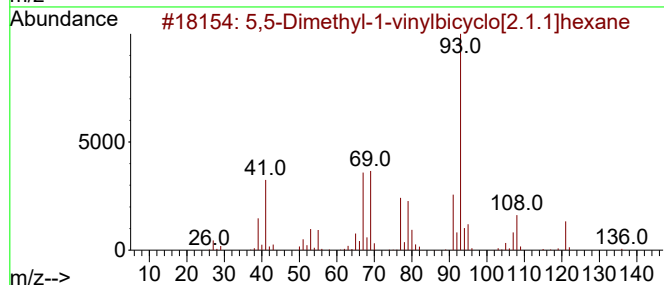
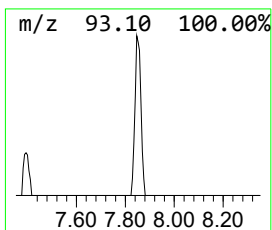
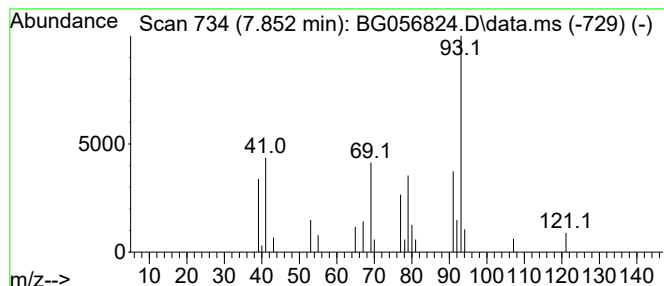
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 5,5-Dimethyl-1-vinylbicyclo... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.852	4.27 ng/ul	19963	1,4-Dichlorobenzene-d4	8.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5,5-Dimethyl-1-vinylbicyclo[2.1....	136	C10H16	016626-39-4	72
2			trans-.beta.-Ocimene	136	C10H16	003779-61-1	64
3			trans-.beta.-Ocimene	136	C10H16	003779-61-1	56
4			3,5-Methanocyclopentapyrazole, 3...	164	C10H16N2	087143-58-6	45
5			2-Pyridinemethanamine, N-methyl-	122	C7H10N2	021035-59-6	40



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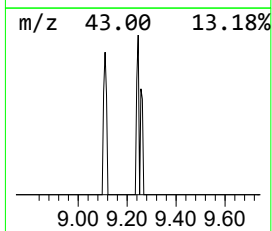
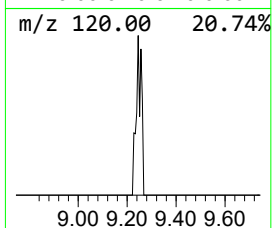
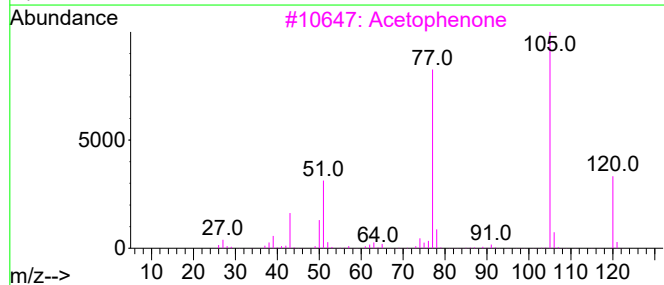
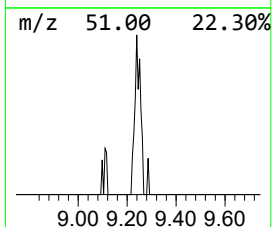
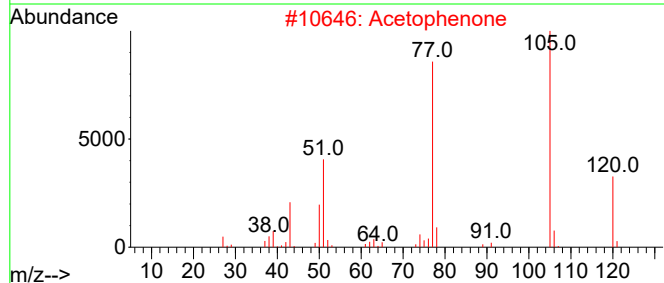
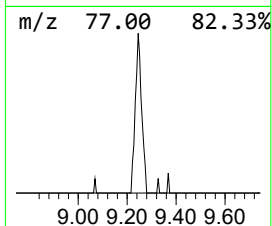
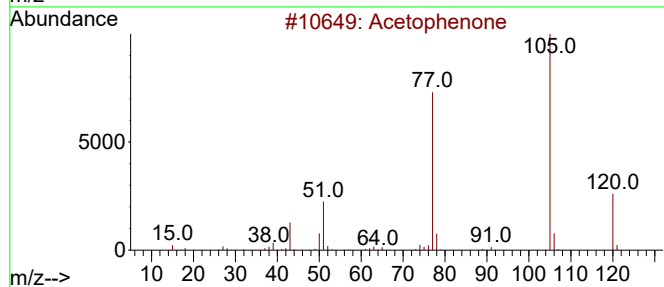
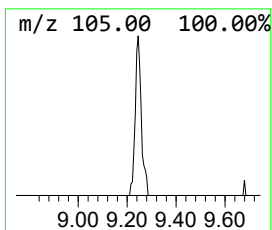
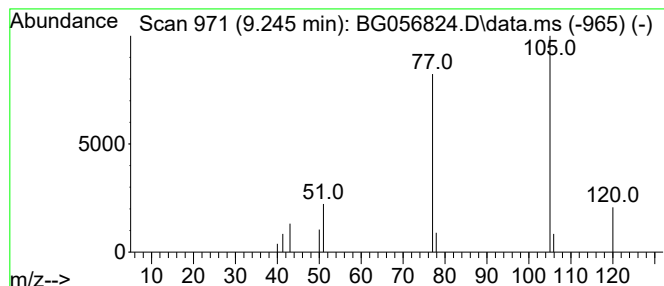
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Acetophe none Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.245	2.52 ng/ul	11778	1,4-Dichlorobenzene-d4	8.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetophenone	120	C8H8O	000098-86-2	90
2			Acetophenone	120	C8H8O	000098-86-2	90
3			Acetophenone	120	C8H8O	000098-86-2	90
4			Acetophenone	120	C8H8O	000098-86-2	83
5			Acetophenone	120	C8H8O	000098-86-2	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
Data File : BG056824.D
Acq On : 4 Mar 2023 5:32
Operator : CG/JU
Sample : 01711-13
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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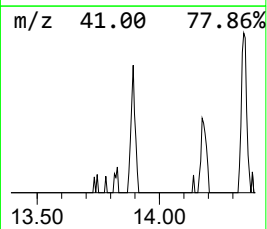
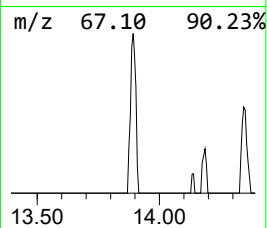
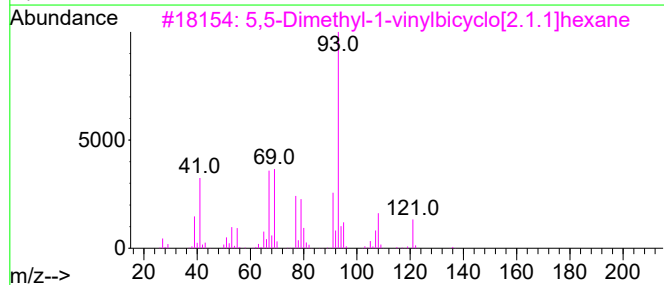
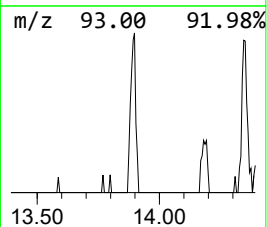
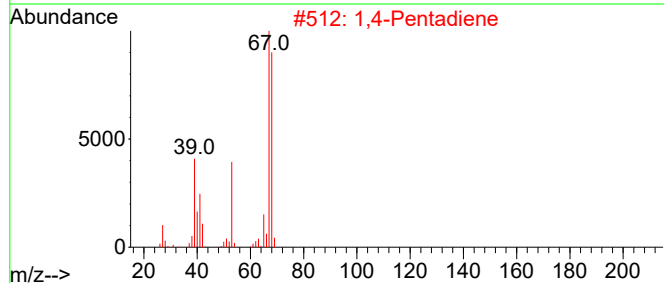
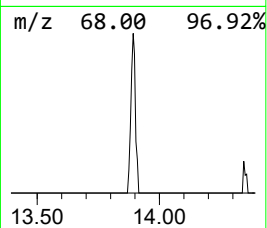
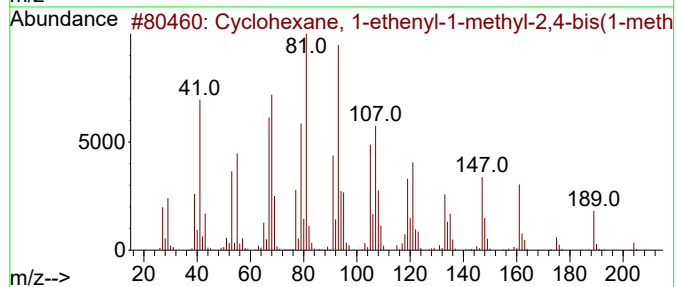
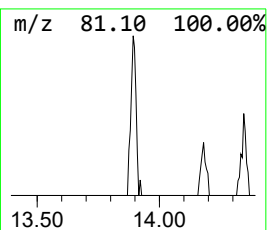
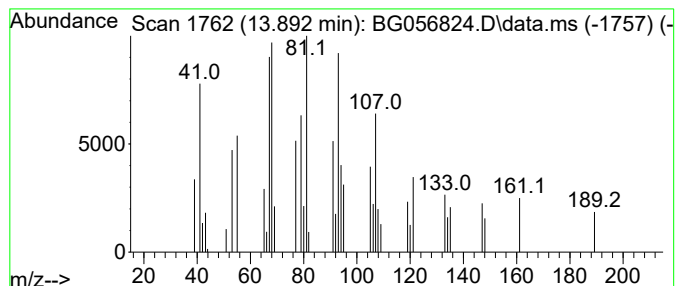
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-01 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.892	2.56 ng/ul	31175	Acenaphthene-d10	15.020

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-ethenyl-1-methyl-...	204	C15H24	000515-13-9	46
2			1,4-Pentadiene	68	C5H8	000591-93-5	43
3			5,5-Dimethyl-1-vinylbicyclo[2.1....	136	C10H16	016626-39-4	43
4			1,3-Pentadiene	68	C5H8	000504-60-9	38
5			D-Limonene	136	C10H16	005989-27-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
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Operator : CG/JU
Sample : 01711-13
Misc :
ALS Vial : 19 Sample Multiplier: 1

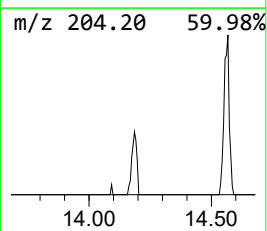
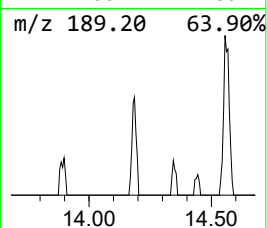
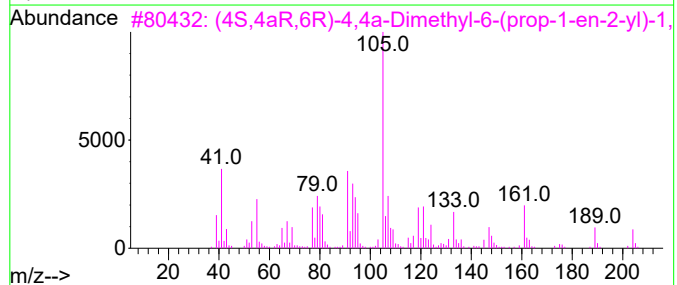
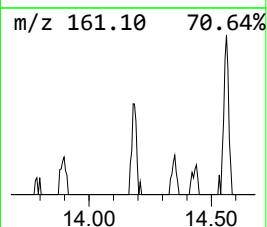
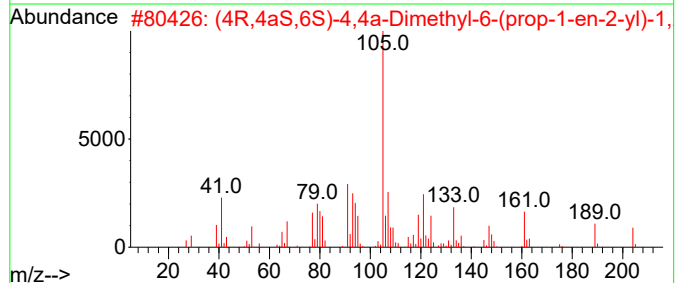
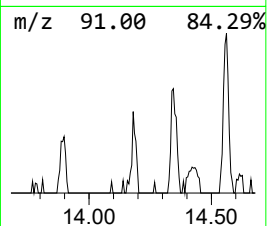
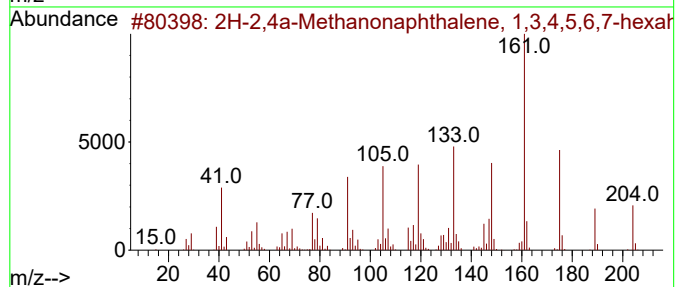
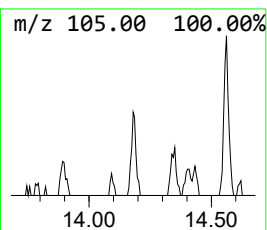
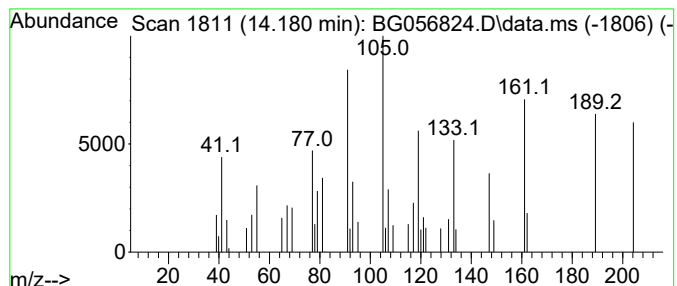
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 2H-2,4a-Methanonaphthalene,... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.180	2.04 ng/ul	24779	Acenaphthene-d10	15.020	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-2,4a-Methanonaphthalene, 1,3,...	204	C15H24	001135-66-6	76
2	(4R,4aS,6S)-4,4a-Dimethyl-6-(pro...	204	C15H24	823810-22-6	76
3	(4S,4aR,6R)-4,4a-Dimethyl-6-(pro...	204	C15H24	054868-40-5	58
4	(-)-Aristolene	204	C15H24	006831-16-9	58
5	(4S,4aR,6R)-4,4a-Dimethyl-6-(pro...	204	C15H24	054868-40-5	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
Data File : BG056824.D
Acq On : 4 Mar 2023 5:32
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Misc :
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ClientSampleId :
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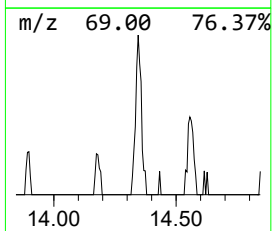
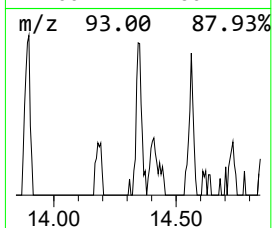
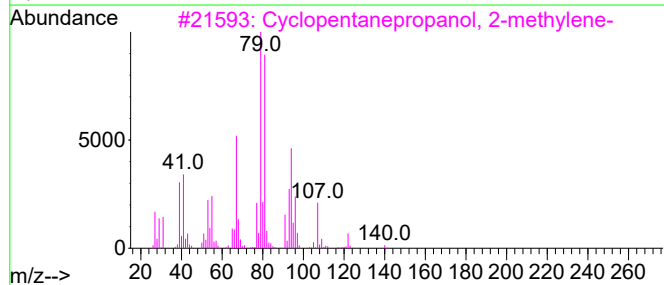
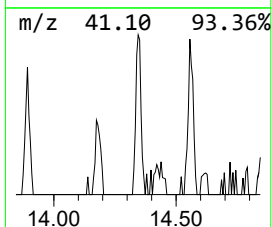
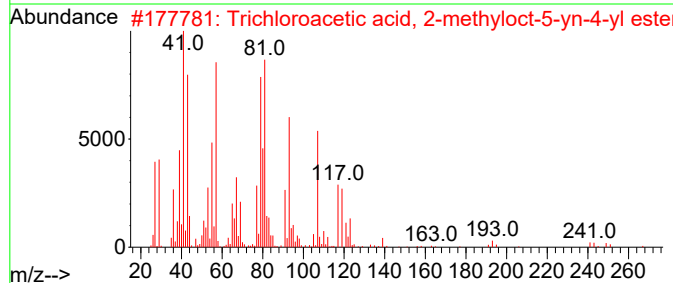
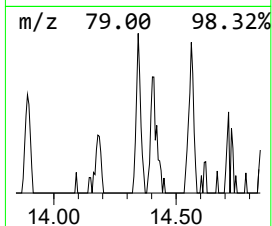
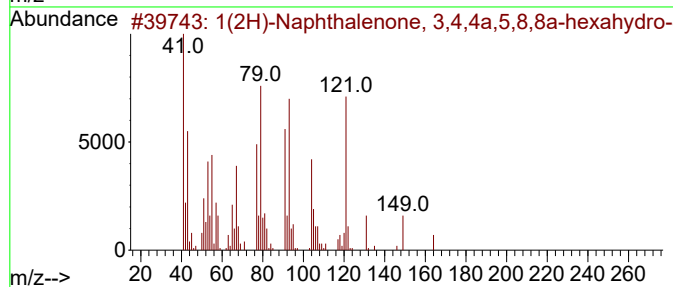
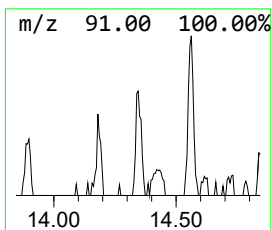
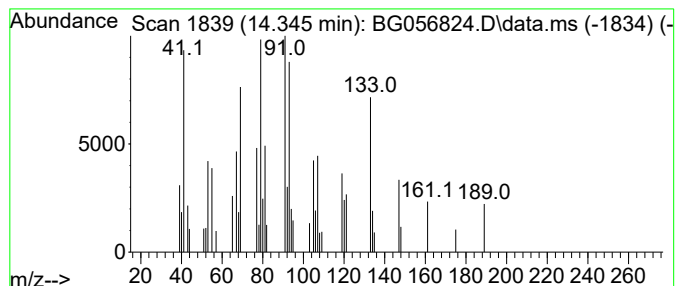
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-02 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.345	2.73 ng/ul	33273	Acenaphthene-d10	15.020

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1(2H)-Naphthalenone, 3,4,4a,5,8,...	164	C11H16O	021841-29-2	43		
2	Trichloroacetic acid, 2-methyloc...	284	C11H15Cl3O2	1000299-25-8	43		
3	Cyclopentanepropanol, 2-methylene-	140	C9H16O	053544-48-2	38		
4	2,3-Diazabicyclo[2.2.1]hept-2-en...	164	C10H16N2	1000221-84-7	35		
5	(2R,4R)-p-Mentha-[1(7),8]-diene,...	168	C10H16O2	1000292-74-1	35		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
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Operator : CG/JU
Sample : 01711-13
Misc :
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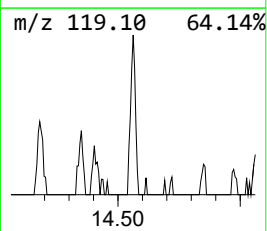
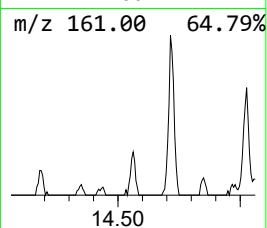
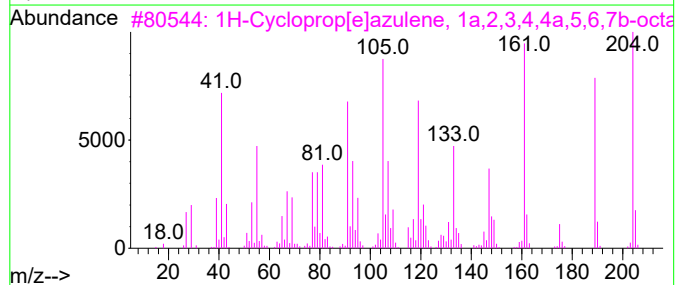
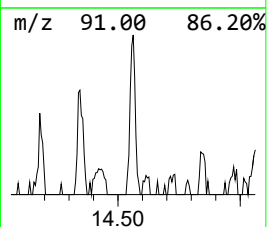
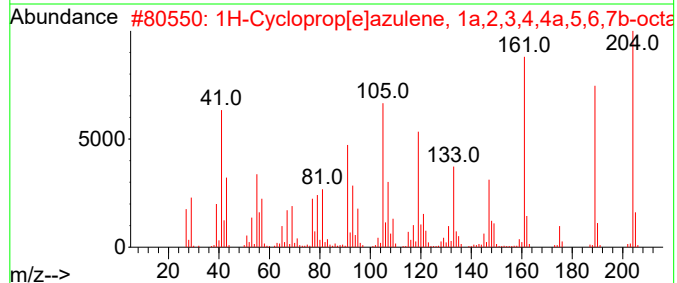
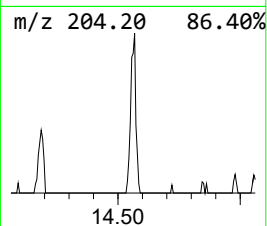
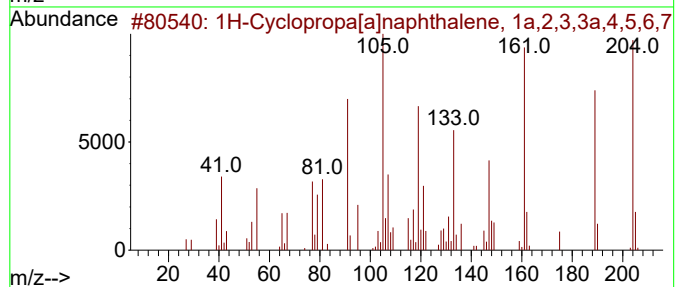
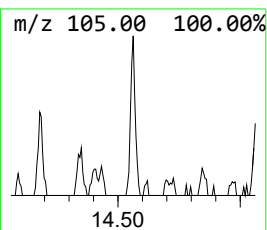
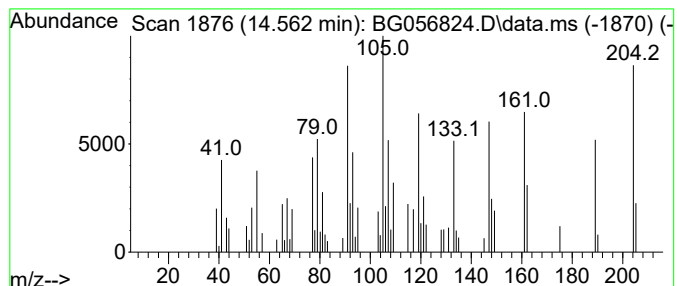
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1H-Cyclopropa[a]naphthalene... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.562	5.35 ng/ul	65118	Acenaphthene-d10	15.020	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Cyclopropa[a]naphthalene, 1a,...	204	C15H24	000489-29-2	95
2	1H-Cycloprop[e]azulene, 1a,2,3,4...	204	C15H24	000489-40-7	95
3	1H-Cycloprop[e]azulene, 1a,2,3,4...	204	C15H24	000489-40-7	94
4	Naphthalene, decahydro-4a-methyl...	204	C15H24	017066-67-0	89
5	8,8,9,9-Tetramethyl-3,4,5,6,7,8-...	204	C15H24	026783-22-2	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
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Acq On : 4 Mar 2023 5:32
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Sample : 01711-13
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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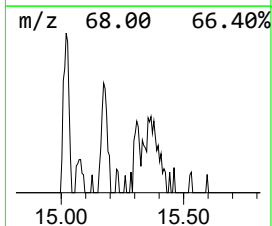
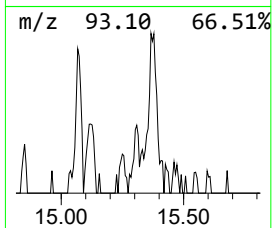
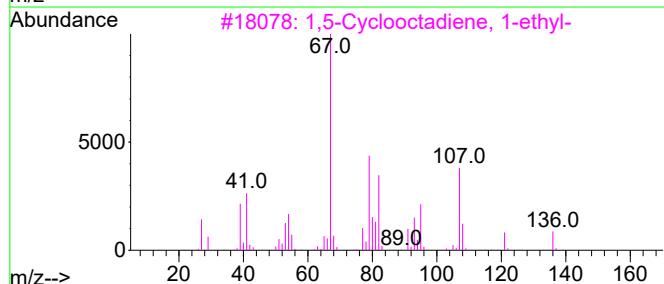
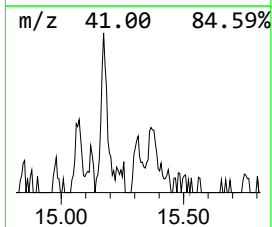
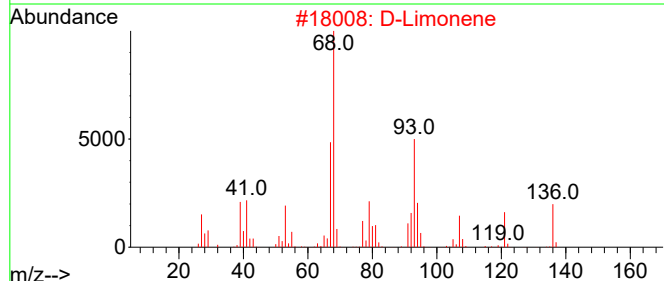
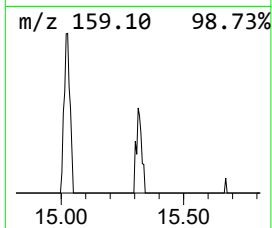
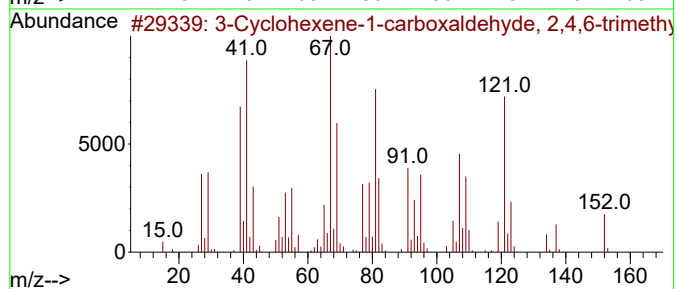
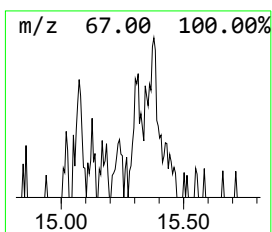
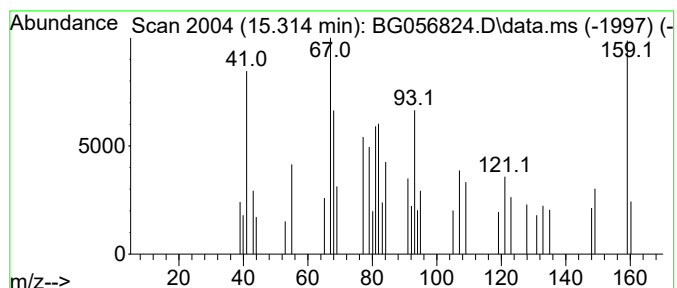
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-03 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.314	2.03 ng/ul	24713	Acenaphthene-d10	15.020

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Cyclohexene-1-carboxaldehyde, ...	152	C10H16O	001423-46-7	14
2			D-Limonene	136	C10H16	005989-27-5	14
3			1,5-Cyclooctadiene, 1-ethyl-	136	C10H16	016539-10-9	14
4			9-Oxabicyclo[3.3.1]non-6-en-2-ol...	140	C8H12O2	029946-76-7	10
5			Fragranyl isovaalerate	238	C15H26O2	051193-03-4	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
Data File : BG056824.D
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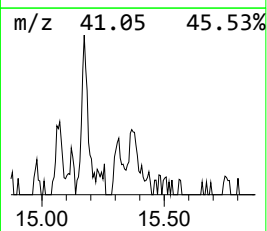
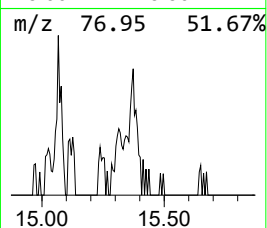
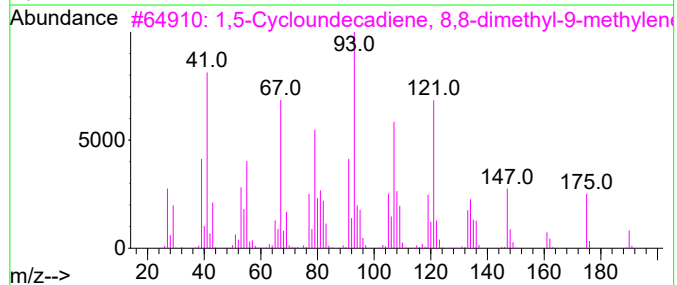
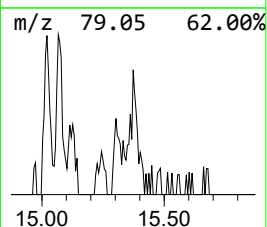
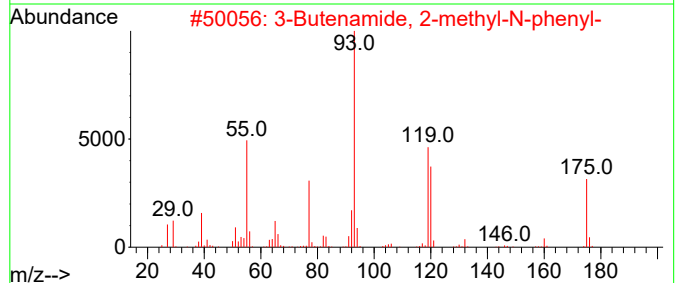
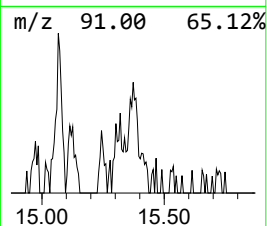
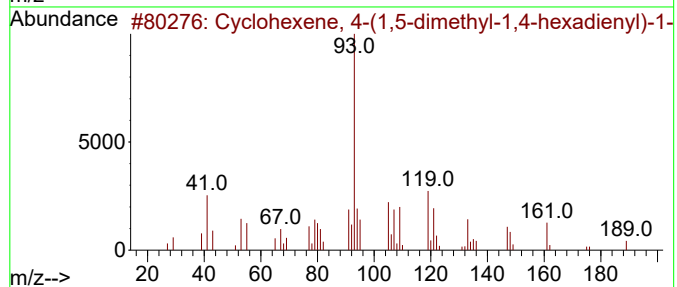
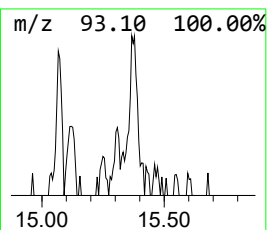
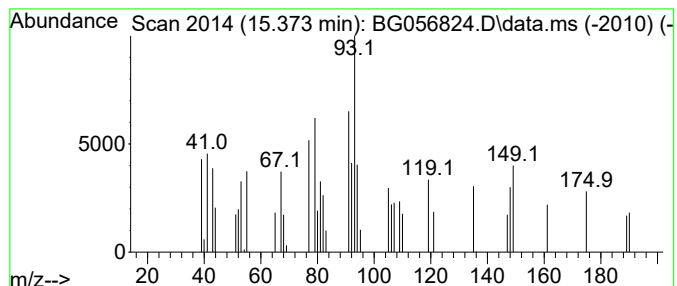
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-04 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.373	2.81 ng/ul	34158	Acenaphthene-d10	15.020

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene, 4-(1,5-dimethyl-1,4...	204	C15H24	017627-44-0	27
2			3-Butenamide, 2-methyl-N-phenyl-	175	C11H13NO	080188-12-1	22
3			1,5-Cycloundecadiene, 8,8-dimeth...	190	C14H22	062338-54-9	16
4			5-Hydroxy-4-methyl-5-phenyl-pent...	281	C18H19NO2	081912-11-0	14
5			Cycloheptanol, 1-methyl-2-methyl...	140	C9H16O	1000163-97-5	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
Data File : BG056824.D
Acq On : 4 Mar 2023 5:32
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Misc :
ALS Vial : 19 Sample Multiplier: 1

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ClientSampleId :
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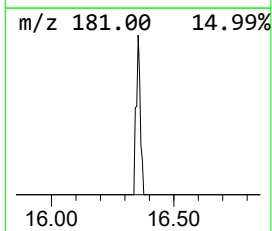
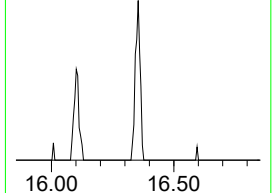
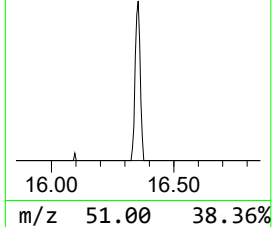
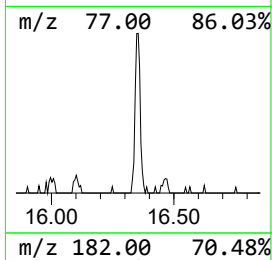
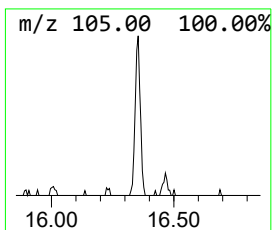
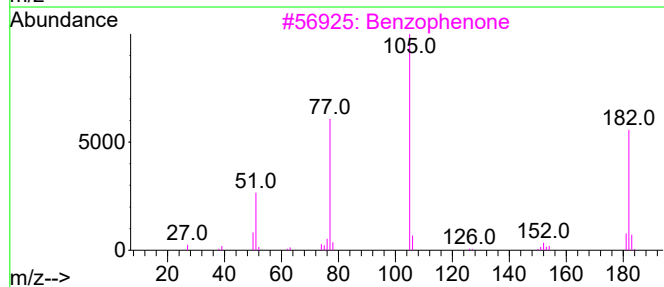
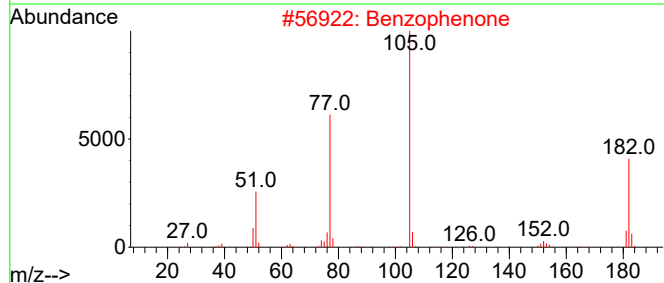
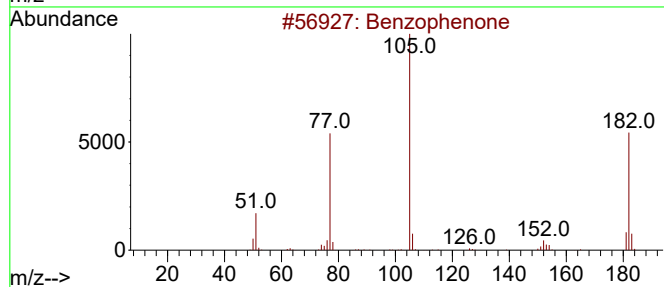
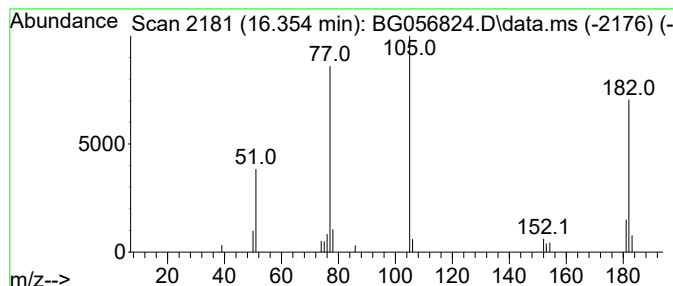
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzophenone Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.354	2.20 ng/ul	26745	Acenaphthene-d10	15.020

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	91
2			Benzophenone	182	C13H10O	000119-61-9	83
3			Benzophenone	182	C13H10O	000119-61-9	83
4			Benzophenone	182	C13H10O	000119-61-9	78
5			1,2,4,5-Tetroxane, 3,3,6,6-tetra...	396	C26H20O4	016204-36-7	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
Data File : BG056824.D
Acq On : 4 Mar 2023 5:32
Operator : CG/JU
Sample : 01711-13
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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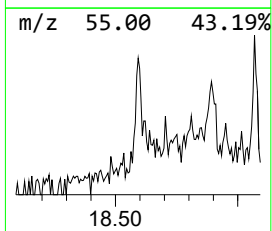
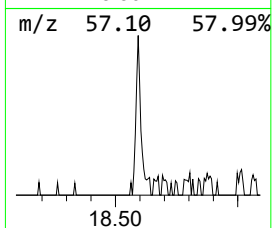
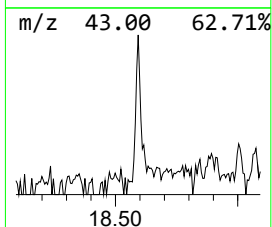
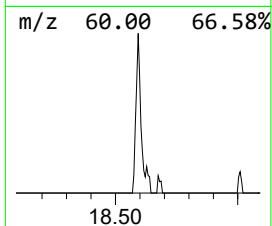
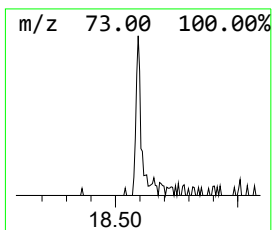
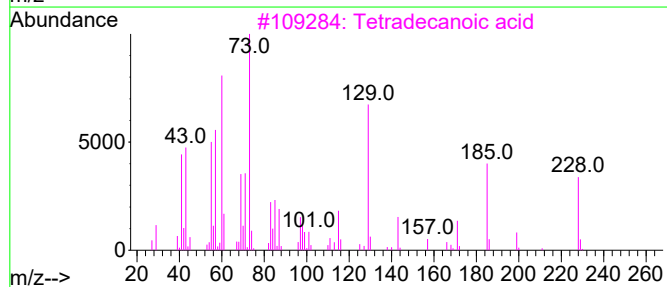
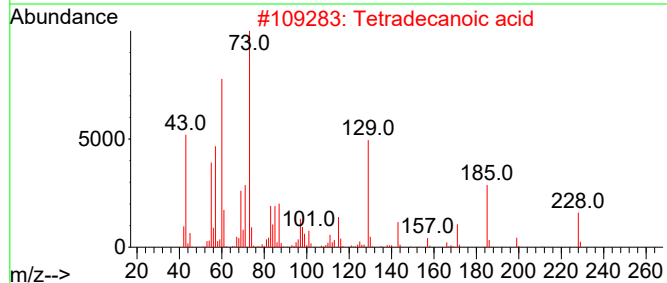
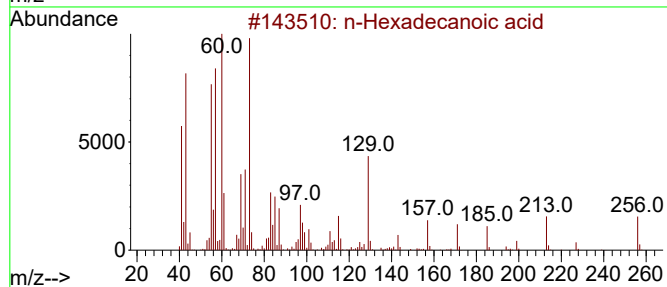
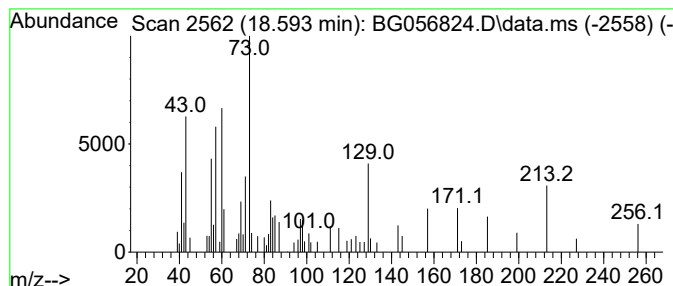
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.593	3.33 ng/ul	51761	Phenanthrene-d10	17.752

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	53
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	52
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	47
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
Data File : BG056824.D
Acq On : 4 Mar 2023 5:32
Operator : CG/JU
Sample : 01711-13
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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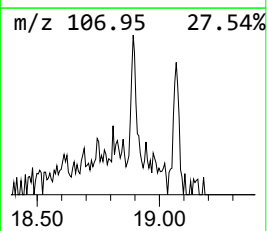
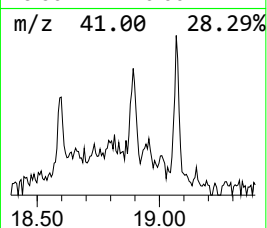
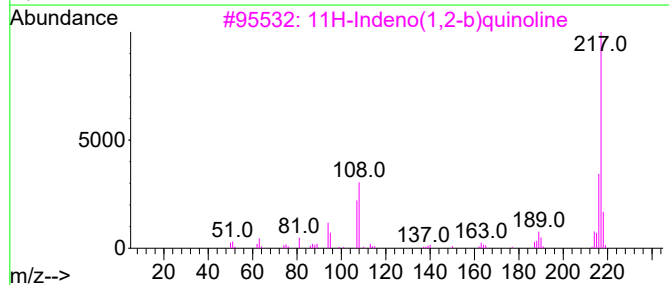
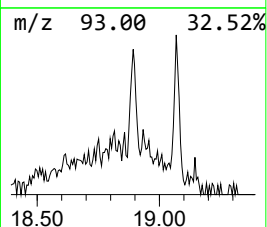
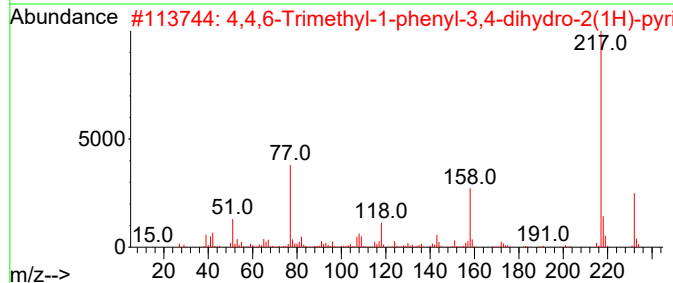
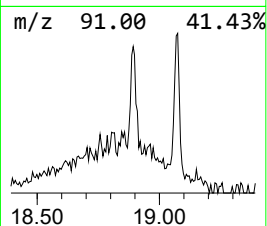
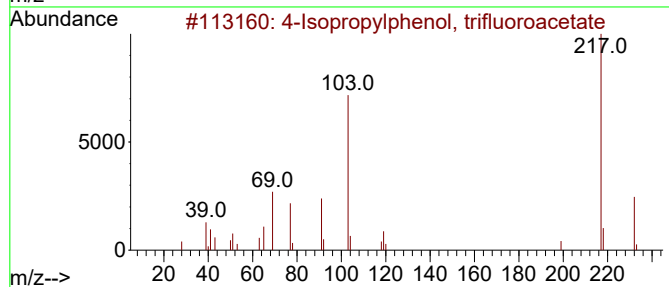
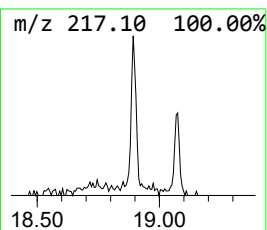
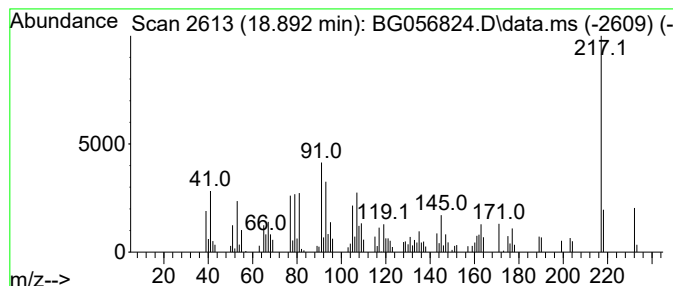
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 unknown-05 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.892	5.33 ng/ul	82878	Phenanthrene-d10	17.752

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Isopropylphenol, trifluoroacetate	232	C11H11F3O2	1000365-67-8	49
2			4,4,6-Trimethyl-1-phenyl-3,4-dih...	232	C13H16N2S	016325-43-2	49
3			11H-Indeno(1,2-b)quinoline	217	C16H11N	000243-51-6	47
4			1H-Pyrrolo[3,4-c]quinoline-1,3(2...	217	C11H11N3O2	298688-31-0	46
5			Benzo(b)carbazole	217	C16H11N	000243-28-7	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
Data File : BG056824.D
Acq On : 4 Mar 2023 5:32
Operator : CG/JU
Sample : 01711-13
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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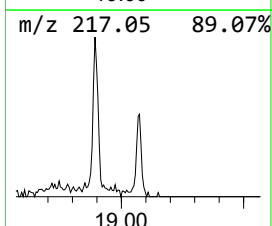
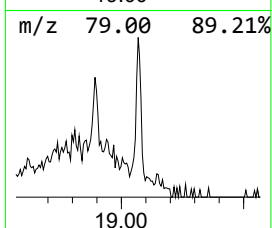
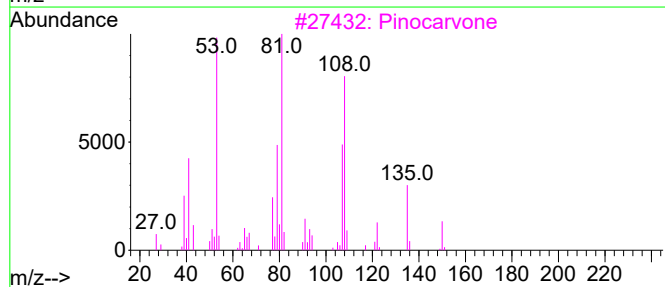
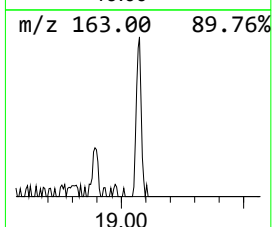
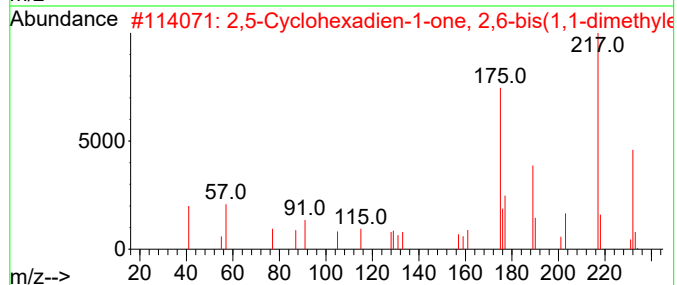
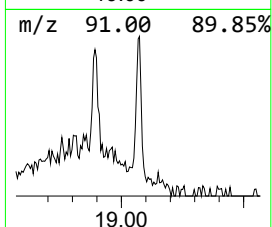
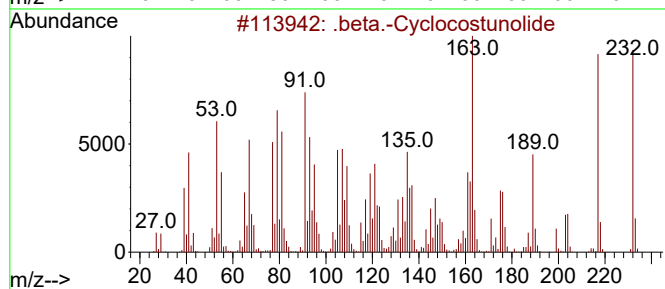
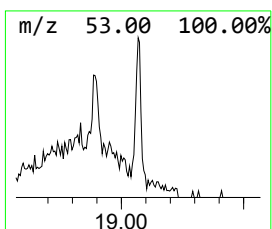
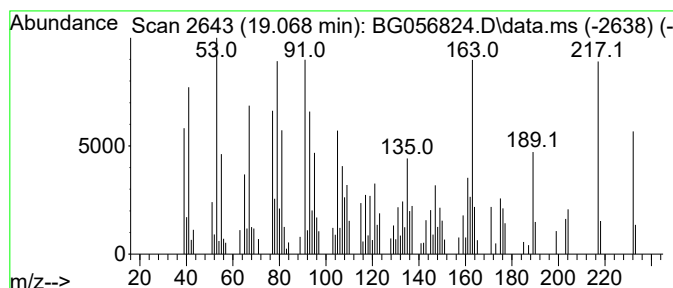
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 .beta.-Cyclocostunolide Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.069	7.90 ng/ul	122989	Phenanthrene-d10	17.752

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			.beta.-Cyclocostunolide	232	C15H20O2	002221-82-1	92
2			2,5-Cyclohexadien-1-one, 2,6-bis...	232	C16H24O	006738-27-8	45
3			Pinocarvone	150	C10H14O	030460-92-5	38
4			Hexadecahydro-benzo[de]anthracene	232	C17H28	1010371-48-1	35
5			1-Butyne, 3-methyl-	68	C5H8	000598-23-2	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
Data File : BG056824.D
Acq On : 4 Mar 2023 5:32
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Sample : 01711-13
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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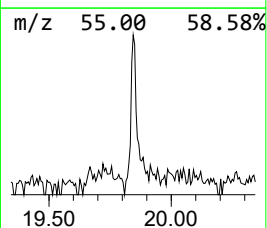
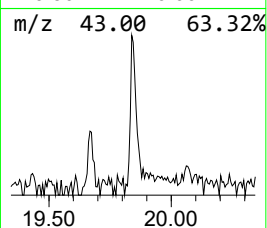
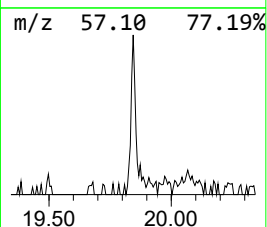
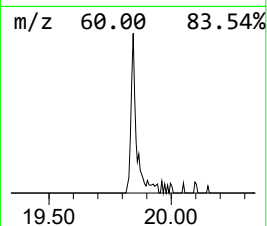
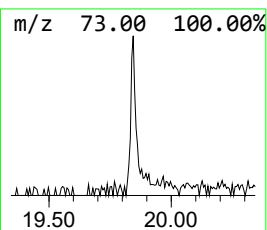
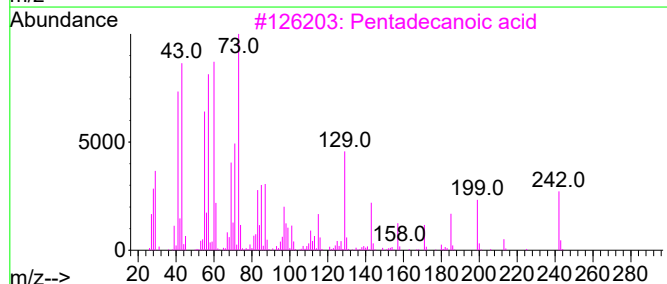
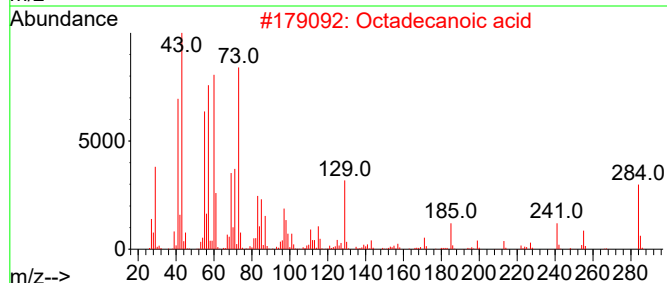
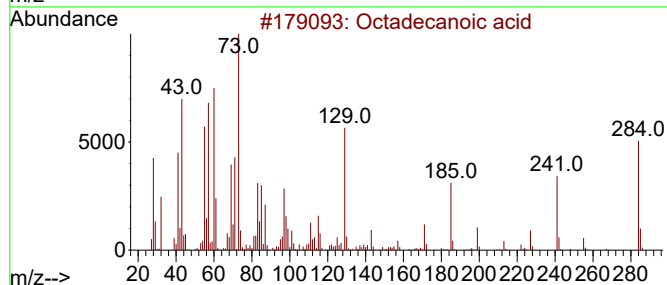
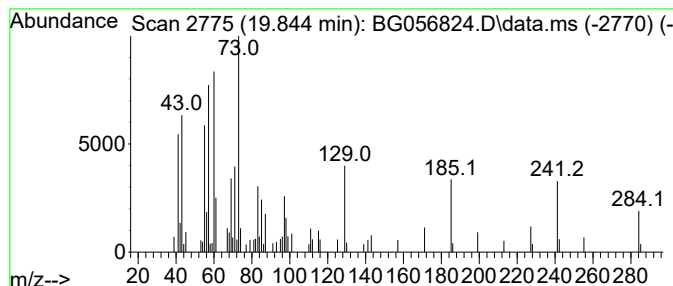
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.844	4.43 ng/ul	68895	Phenanthrene-d10	17.752

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	81
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	76
4			Octadecanoic acid	284	C18H36O2	000057-11-4	74
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
Data File : BG056824.D
Acq On : 4 Mar 2023 5:32
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Sample : 01711-13
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBAC9

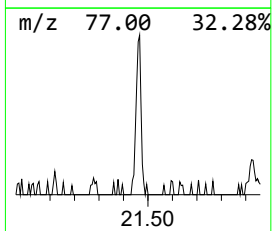
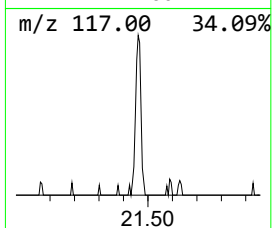
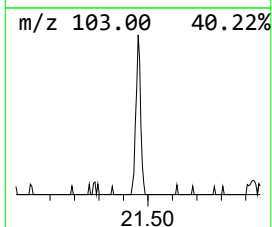
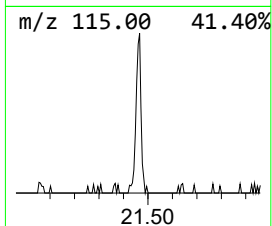
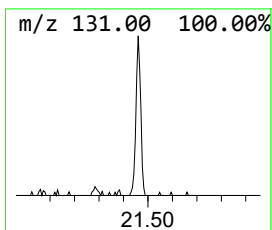
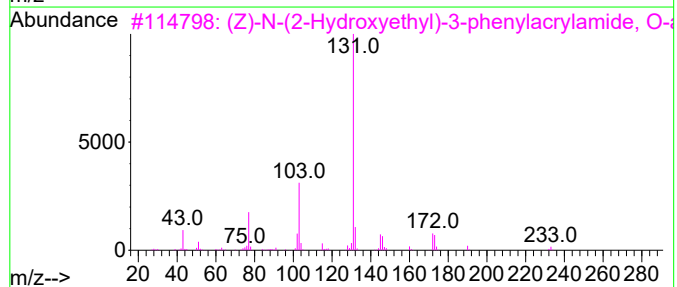
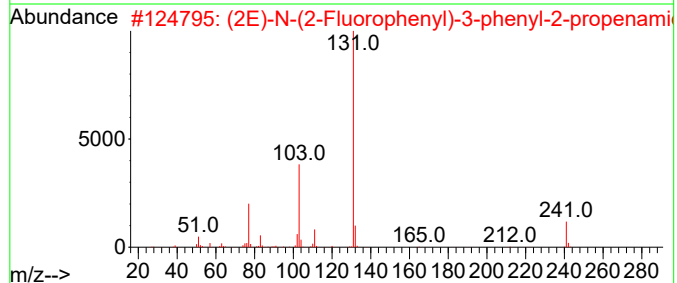
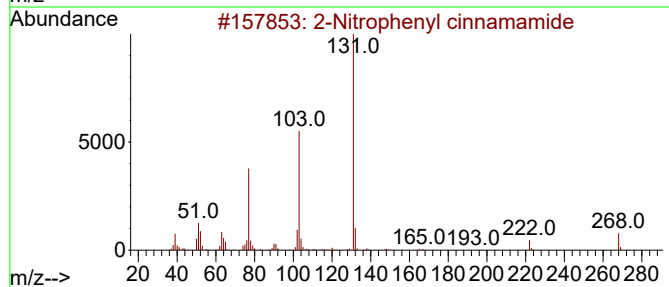
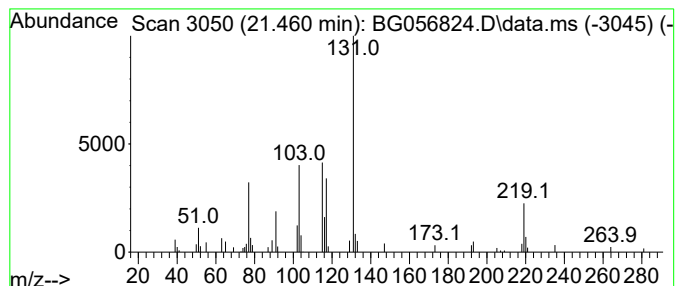
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 unknown-06 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.460	3.01 ng/ul	51253	Chrysene-d12	22.065

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Nitrophenyl cinnamamide	268	C15H12N2O3	111273-37-1	49
2			(2E)-N-(2-Fluorophenyl)-3-phenyl...	241	C15H12FNO	025893-50-9	47
3			(Z)-N-(2-Hydroxyethyl)-3-phenyla...	233	C13H15NO3	1000512-63-0	47
4			(2E)-N-(4-Methylphenyl)-3-phenyl...	237	C16H15NO	134430-88-9	47
5			N-[(2Z)-3-Phenylprop-2-enoyl]gly...	233	C13H15NO3	1000498-00-6	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
Data File : BG056824.D
Acq On : 4 Mar 2023 5:32
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Sample : 01711-13
Misc :
ALS Vial : 19 Sample Multiplier: 1

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ClientSampleId :
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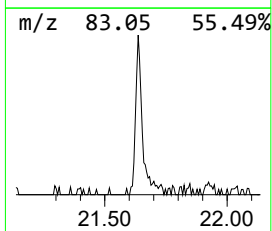
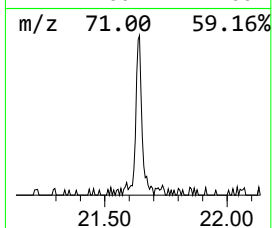
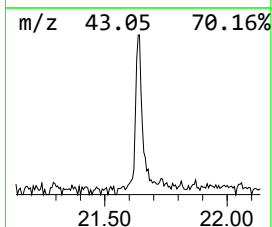
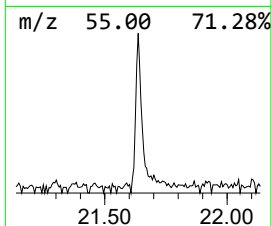
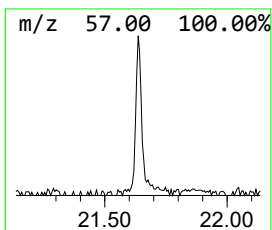
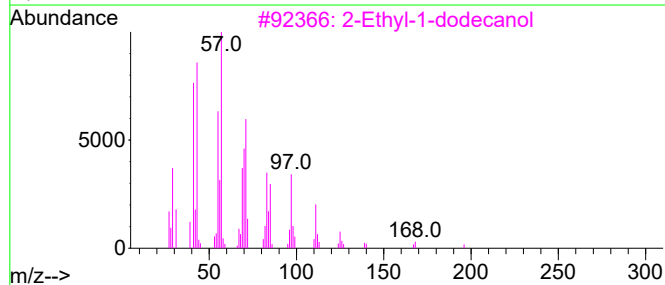
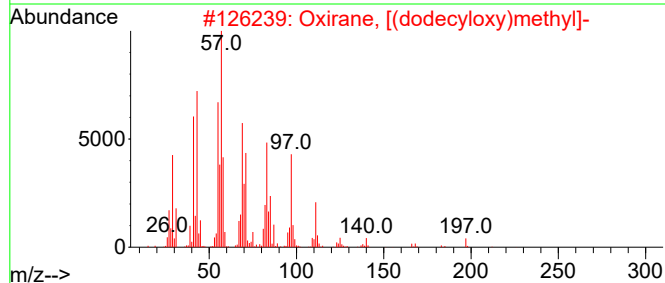
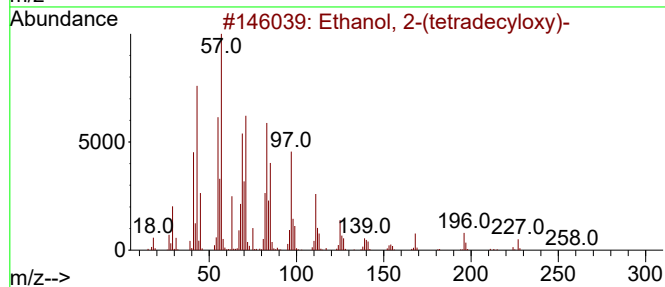
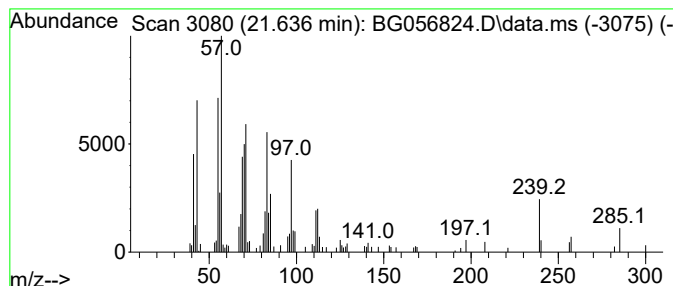
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Ethanol, 2-(tetradecyloxy)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.636	7.69 ng/ul	131017	Chrysene-d12	22.065

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	62
2			Oxirane, [(dodecyloxy)methyl]-	242	C15H30O2	002461-18-9	58
3			2-Ethyl-1-dodecanol	214	C14H30O	019780-33-7	58
4			Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	52
5			Oxalic acid, isobutyl pentadecyl...	356	C21H40O4	1000309-38-0	52



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Data File : BG056824.D
Acq On : 4 Mar 2023 5:32
Operator : CG/JU
Sample : 01711-13
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DBAC9

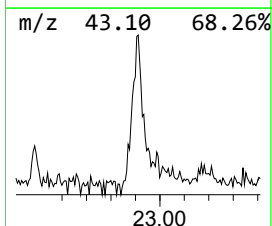
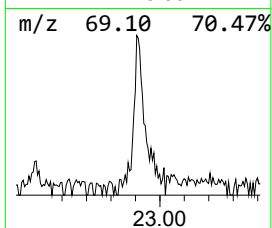
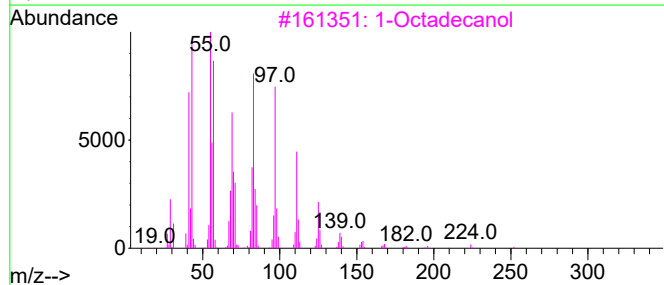
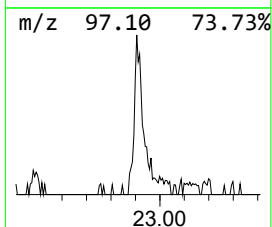
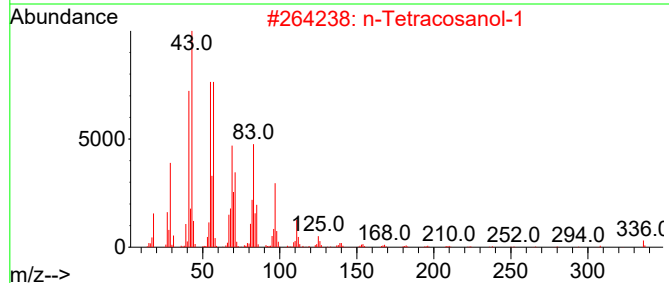
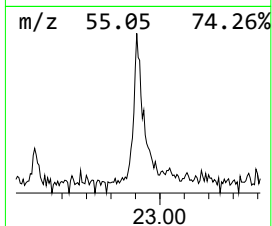
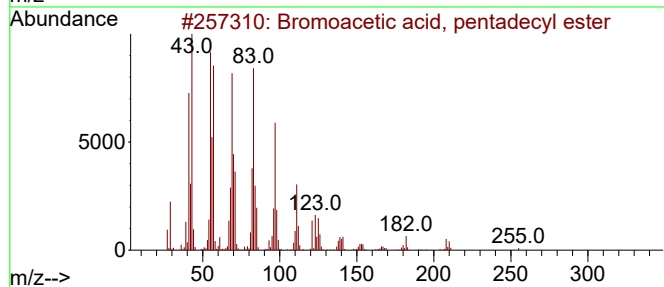
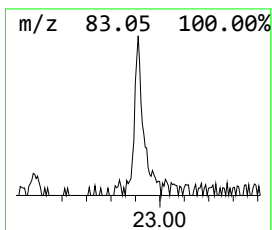
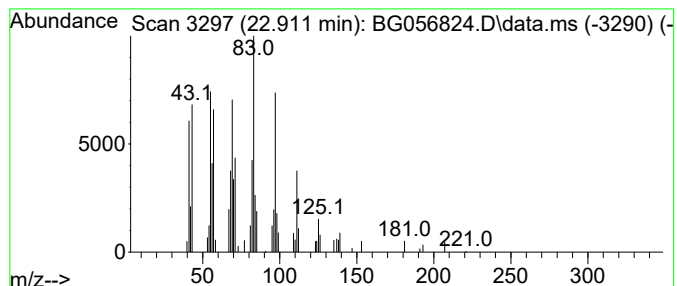
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Bromoacetic acid, pentadecy... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.911	6.24 ng/ul	106287	Chrysene-d12	22.065

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bromoacetic acid, pentadecyl ester	348	C17H33BrO2	131143-01-6	91
2			n-Tetracosanol-1	354	C24H50O	000506-51-4	91
3			1-Octadecanol	270	C18H38O	000112-92-5	91
4			n-Heptadecanol-1	256	C17H36O	001454-85-9	90
5			1-Eicosanol	298	C20H42O	000629-96-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
Data File : BG056824.D
Acq On : 4 Mar 2023 5:32
Operator : CG/JU
Sample : 01711-13
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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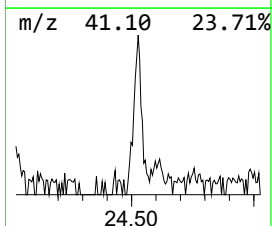
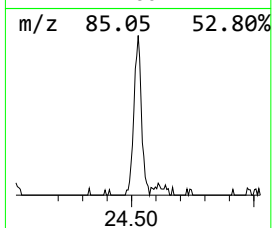
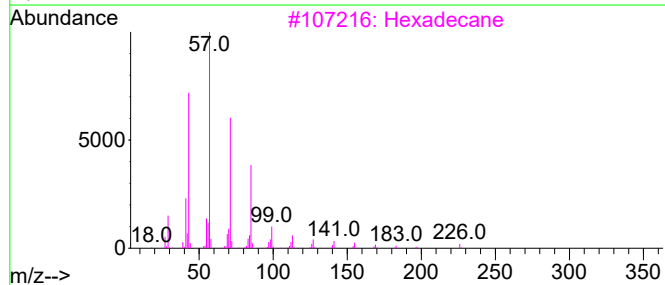
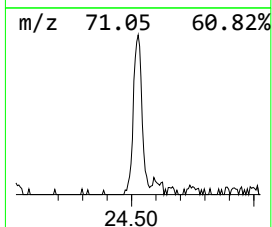
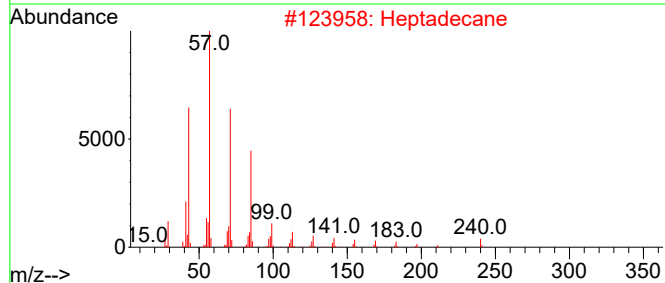
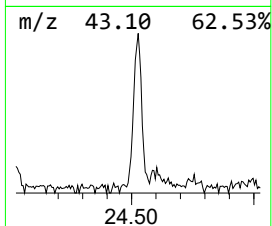
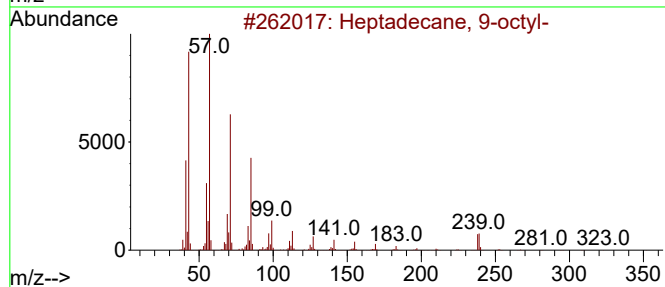
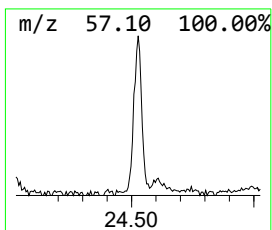
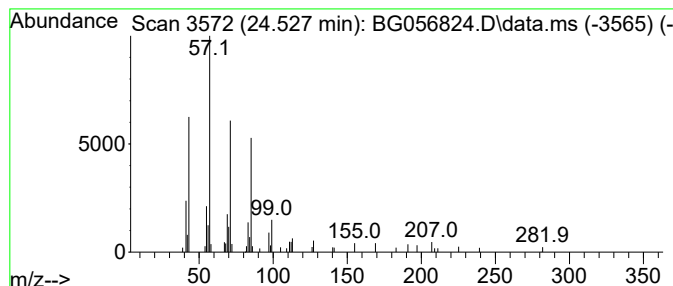
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.527	5.40 ng/ul	91440	Perylene-d12	25.584

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
2			Heptadecane	240	C17H36	000629-78-7	86
3			Hexadecane	226	C16H34	000544-76-3	86
4			Heneicosane	296	C21H44	000629-94-7	80
5			Tetradecane	198	C14H30	000629-59-4	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
Data File : BG056824.D
Acq On : 4 Mar 2023 5:32
Operator : CG/JU
Sample : 01711-13
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBAC9

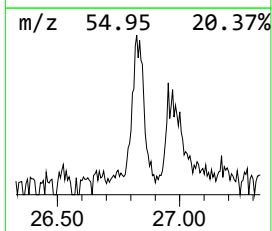
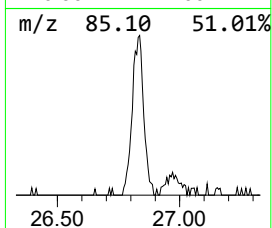
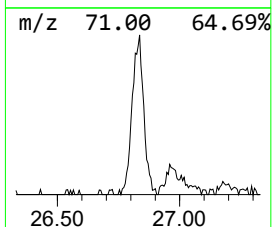
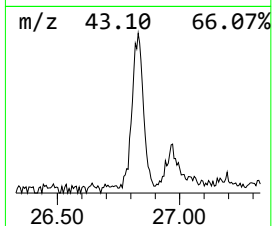
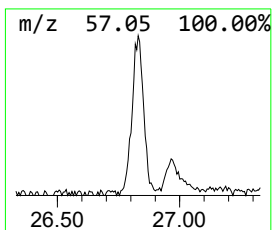
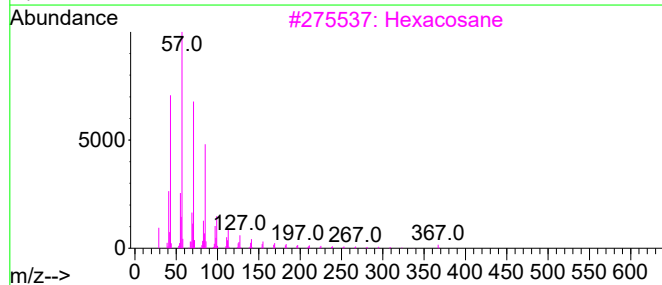
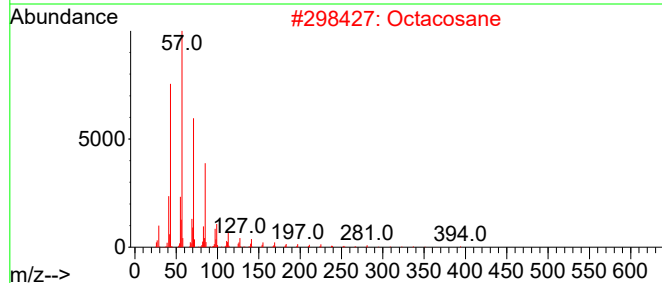
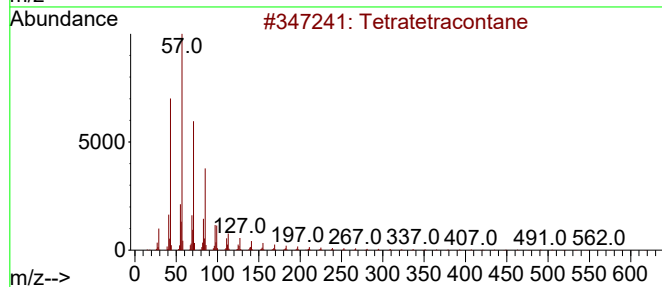
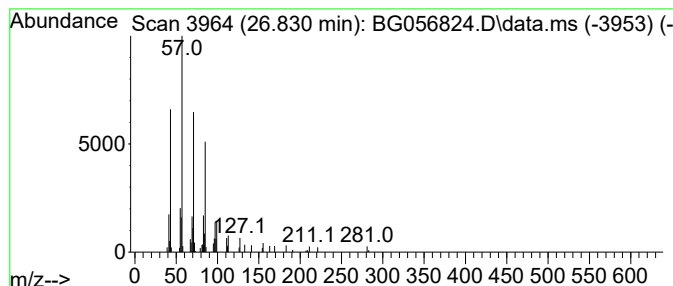
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.830	8.31 ng/ul	140746	Perylene-d12	25.584

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetratetracontane	619	C44H90	007098-22-8	91
2		Octacosane	394	C28H58	000630-02-4	91
3		Hexacosane	366	C26H54	000630-01-3	90
4		Eicosane	282	C20H42	000112-95-8	87
5		Tetratetracontane	619	C44H90	007098-22-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
 Data File : BG056824.D
 Acq On : 4 Mar 2023 5:32
 Operator : CG/JU
 Sample : 01711-13
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBAC9

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
(1S)-2,6,6-Trim...	7.083	4.3	ng/ul	19983	1	8.422	93421	20.0
5,5-Dimethyl-1-...	7.852	4.3	ng/ul	19963	1	8.422	93421	20.0
Acetophe none	9.245	2.5	ng/ul	11778	1	8.422	93421	20.0
unknown-01	13.892	2.6	ng/ul	31175	3	15.020	243364	20.0
2H-2,4a-Methano...	14.180	2.0	ng/ul	24779	3	15.020	243364	20.0
unknown-02	14.345	2.7	ng/ul	33273	3	15.020	243364	20.0
1H-Cyclopropa[a...	14.562	5.3	ng/ul	65118	3	15.020	243364	20.0
unknown-03	15.314	2.0	ng/ul	24713	3	15.020	243364	20.0
unknown-04	15.373	2.8	ng/ul	34158	3	15.020	243364	20.0
Benzophenone	16.354	2.2	ng/ul	26745	3	15.020	243364	20.0
n-Hexadecanoic ...	18.593	3.3	ng/ul	51761	4	17.752	311270	20.0
unknown-05	18.892	5.3	ng/ul	82878	4	17.752	311270	20.0
.beta.-Cyclocos...	19.069	7.9	ng/ul	122989	4	17.752	311270	20.0
Octadecanoic acid	19.844	4.4	ng/ul	68895	4	17.752	311270	20.0
unknown-06	21.460	3.0	ng/ul	51253	5	22.065	340774	20.0
Ethanol, 2-(tet...	21.636	7.7	ng/ul	131017	5	22.065	340774	20.0
Bromoacetic aci...	22.911	6.2	ng/ul	106287	5	22.065	340774	20.0
(DEL) Alkane: S...	24.527	5.4	ng/ul	91440	6	25.584	338560	20.0
(DEL) Alkane: S...	26.830	8.3	ng/ul	140746	6	25.584	338560	20.0