

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
 Data File : BG057229.D
 Acq On : 28 Apr 2023 22:55
 Operator : CG/JU
 Sample : 02417-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW8W7

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-BG042823.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG057229.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	6.774	547	551	556	rVV2	3805	6147	1.86%	0.111%
2	7.438	661	664	670	rVB	2887	3834	1.16%	0.069%
3	7.514	670	677	681	rBV	40772	77021	23.29%	1.391%
4	7.679	700	705	713	rVB	27806	46310	14.00%	0.836%
5	7.749	713	717	722	rVB	4347	6529	1.97%	0.118%
6	7.902	738	743	750	rVV2	33574	56615	17.12%	1.022%
7	7.967	750	754	758	rVV3	4007	5535	1.67%	0.100%
8	8.014	758	762	767	rVB2	10134	15293	4.62%	0.276%
9	8.325	810	815	818	rBV3	4010	6119	1.85%	0.110%
10	8.378	818	824	830	rVV	50770	81601	24.67%	1.474%
11	8.489	837	843	849	rBV2	13102	19197	5.80%	0.347%
12	8.889	902	911	924	rVB6	15071	39125	11.83%	0.707%
13	9.006	926	931	936	rBV	29113	45308	13.70%	0.818%
14	9.077	936	943	949	rBV2	44840	75630	22.87%	1.366%
15	9.183	955	961	963	rBV2	5893	8770	2.65%	0.158%
16	9.329	981	986	991	rBV2	13833	21983	6.65%	0.397%
17	9.382	991	995	1000	rVB	12258	17512	5.29%	0.316%
18	9.482	1006	1012	1018	rVB	22970	35570	10.75%	0.642%
19	9.553	1018	1024	1028	rBV	36621	70537	21.33%	1.274%
20	9.741	1049	1056	1063	rVB4	6204	13190	3.99%	0.238%
21	9.823	1065	1070	1074	rBV	10215	15753	4.76%	0.284%
22	10.017	1097	1103	1108	rVV	29041	45727	13.83%	0.826%
23	10.076	1108	1113	1119	rVB	67270	110131	33.30%	1.989%
24	10.281	1137	1148	1158	rBV3	34835	79187	23.94%	1.430%
25	10.381	1158	1165	1173	rVV2	36598	77308	23.37%	1.396%
26	10.534	1182	1191	1194	rVV3	2983	6187	1.87%	0.112%
27	10.598	1197	1202	1207	rVV2	29566	51009	15.42%	0.921%
28	10.657	1207	1212	1217	rVV3	11364	19680	5.95%	0.355%
29	10.745	1222	1227	1228	rBV2	5092	6988	2.11%	0.126%
30	10.775	1230	1232	1235	rVV2	6802	6898	2.09%	0.125%
31	10.827	1235	1241	1248	rVV2	48747	85124	25.74%	1.537%
32	10.886	1248	1251	1259	rVB	10179	16639	5.03%	0.300%
33	11.074	1275	1283	1289	rBV	4820	9121	2.76%	0.165%
34	11.145	1291	1295	1297	rBV	4357	5922	1.79%	0.107%
35	11.215	1301	1307	1313	rBV2	77354	165332	49.99%	2.986%
36	11.339	1322	1328	1334	rBV2	18177	33412	10.10%	0.603%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
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37	11.409	1334	1340	1346	rVB	12163	19784	5.98%	0.357%
38	11.703	1383	1390	1396	rBV	48504	70904	21.44%	1.280%
39	11.820	1405	1410	1414	rBV3	3388	5714	1.73%	0.103%
40	12.108	1453	1459	1463	rVB4	6200	9786	2.96%	0.177%

41	12.161	1463	1468	1469	rBV	2848	3890	1.18%	0.070%
42	12.290	1483	1490	1496	rBV7	6350	14410	4.36%	0.260%
43	12.402	1502	1509	1519	rVB5	4849	11132	3.37%	0.201%
44	12.502	1519	1526	1533	rBV	13879	22965	6.94%	0.415%
45	12.566	1533	1537	1543	rVB3	3933	5323	1.61%	0.096%

46	12.842	1579	1584	1590	rVB	12374	17561	5.31%	0.317%
47	12.977	1603	1607	1612	rVB3	8691	13075	3.95%	0.236%
48	13.060	1614	1621	1627	rBV4	11826	18137	5.48%	0.328%
49	13.218	1644	1648	1654	rVB	8378	12928	3.91%	0.233%
50	13.348	1663	1670	1678	rBV2	33136	52953	16.01%	0.956%

51	13.471	1685	1691	1697	rVV5	5619	11243	3.40%	0.203%
52	13.829	1746	1752	1754	rBV3	2579	4148	1.25%	0.075%
53	13.923	1759	1768	1771	rBV5	6183	13843	4.19%	0.250%
54	14.017	1778	1784	1790	rBV4	2301	3730	1.13%	0.067%
55	14.129	1798	1803	1805	rBV3	2971	4641	1.40%	0.084%

56	14.164	1805	1809	1814	rVB4	4636	7740	2.34%	0.140%
57	14.329	1829	1837	1841	rBV	69622	109053	32.97%	1.969%
58	14.376	1841	1845	1852	rVV	109060	156006	47.17%	2.817%
59	14.546	1870	1874	1878	rBV2	3626	4430	1.34%	0.080%
60	14.687	1890	1898	1905	rBV	112598	176515	53.37%	3.188%

61	14.904	1929	1935	1940	rBV	16649	23466	7.10%	0.424%
62	14.987	1943	1949	1954	rVV	136369	212841	64.35%	3.844%
63	15.034	1954	1957	1965	rVB3	20465	28762	8.70%	0.519%
64	15.180	1973	1982	1991	rBV2	30196	55096	16.66%	0.995%
65	15.298	1998	2002	2004	rBV5	4239	5588	1.69%	0.101%

66	15.380	2011	2016	2018	rBV4	5788	7663	2.32%	0.138%
67	15.738	2072	2077	2082	rBV	75319	98388	29.75%	1.777%
68	15.973	2111	2117	2122	rBV	156951	230216	69.61%	4.157%
69	16.020	2122	2125	2130	rVB2	9655	12704	3.84%	0.229%
70	16.085	2130	2136	2145	rVB2	40420	63437	19.18%	1.146%

71	16.326	2169	2177	2181	rBV	21130	32013	9.68%	0.578%
72	16.396	2186	2189	2194	rVV2	6027	8856	2.68%	0.160%
73	16.661	2232	2234	2239	rVV5	3815	5287	1.60%	0.095%
74	17.007	2289	2293	2301	rVB4	3757	7306	2.21%	0.132%
75	17.377	2350	2356	2360	rBV3	2933	5157	1.56%	0.093%

76	17.454	2361	2369	2379	rBV	108972	159335	48.18%	2.877%
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77	17.724	2409	2415	2420	rBV	176733	271659	82.14%	4.906%
78	17.771	2420	2423	2427	rVV	34594	46186	13.96%	0.834%
79	17.824	2427	2432	2443	rVV2	159292	251073	75.91%	4.534%
80	17.918	2443	2448	2451	rVV3	3460	5613	1.70%	0.101%
81	18.270	2504	2508	2513	rVV2	22647	30915	9.35%	0.558%
82	18.388	2524	2528	2533	rVB4	3366	5460	1.65%	0.099%
83	18.564	2554	2558	2567	rBV4	20301	40133	12.13%	0.725%
84	18.640	2568	2571	2577	rVB	7749	11420	3.45%	0.206%
85	18.793	2590	2597	2600	rBV6	14641	25590	7.74%	0.462%
86	18.922	2616	2619	2623	rBV3	4863	6316	1.91%	0.114%
87	19.075	2639	2645	2651	rBV4	13953	21823	6.60%	0.394%
88	19.128	2651	2654	2662	rVB8	5557	10988	3.32%	0.198%
89	19.369	2692	2695	2700	rVB6	2835	4362	1.32%	0.079%
90	19.422	2700	2704	2708	rBV5	2712	5138	1.55%	0.093%
91	19.610	2734	2736	2738	rBV3	4154	4221	1.28%	0.076%
92	19.757	2756	2761	2766	rVB	76932	106837	32.30%	1.929%
93	20.091	2812	2818	2821	rBV	210473	326844	98.82%	5.902%
94	21.020	2972	2976	2981	rBV	149068	177191	53.58%	3.200%
95	21.595	3070	3074	3085	rVB6	22905	49084	14.84%	0.886%
96	21.865	3115	3120	3126	rBV	51077	80447	24.32%	1.453%
97	22.036	3141	3149	3154	rBV3	162976	330731	100.00%	5.972%
98	24.444	3548	3559	3565	rBV4	39773	140500	42.48%	2.537%
99	25.302	3697	3705	3714	rBV2	86292	241322	72.97%	4.358%
100	25.543	3738	3746	3755	rVB2	79653	231441	69.98%	4.179%

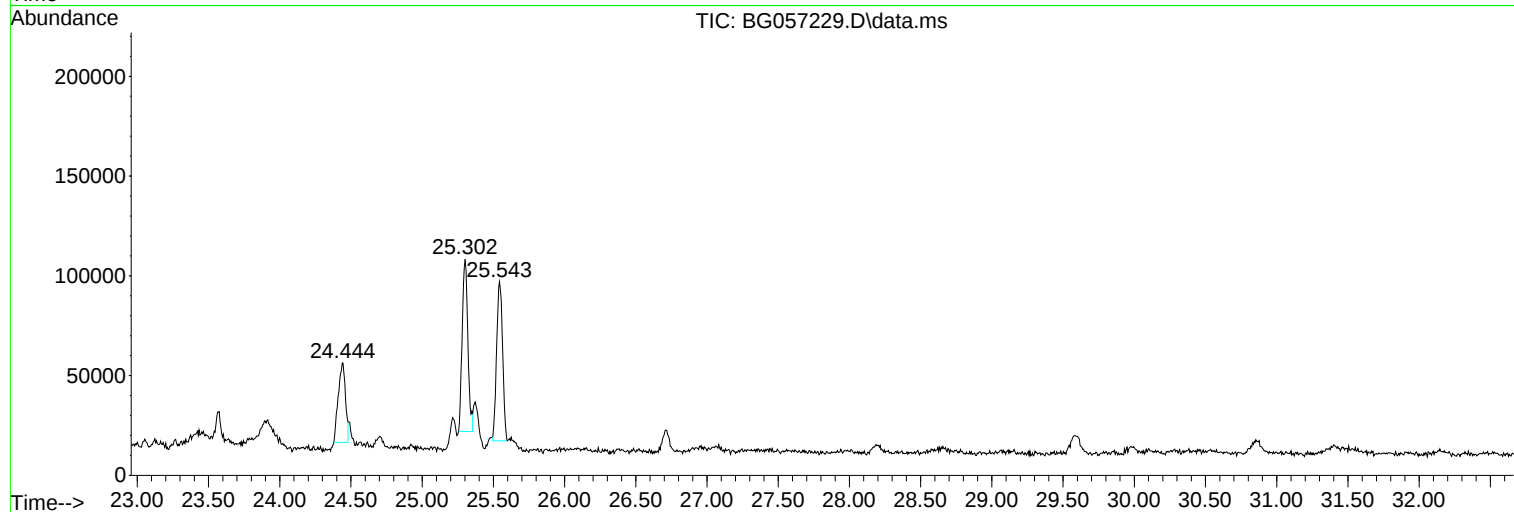
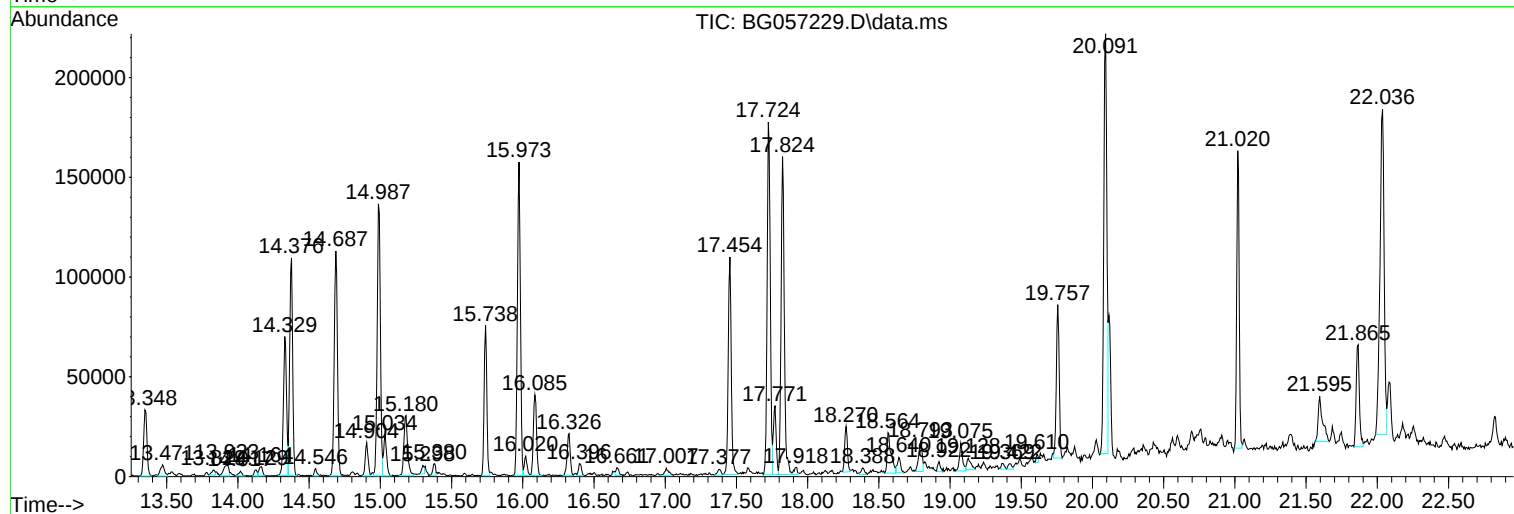
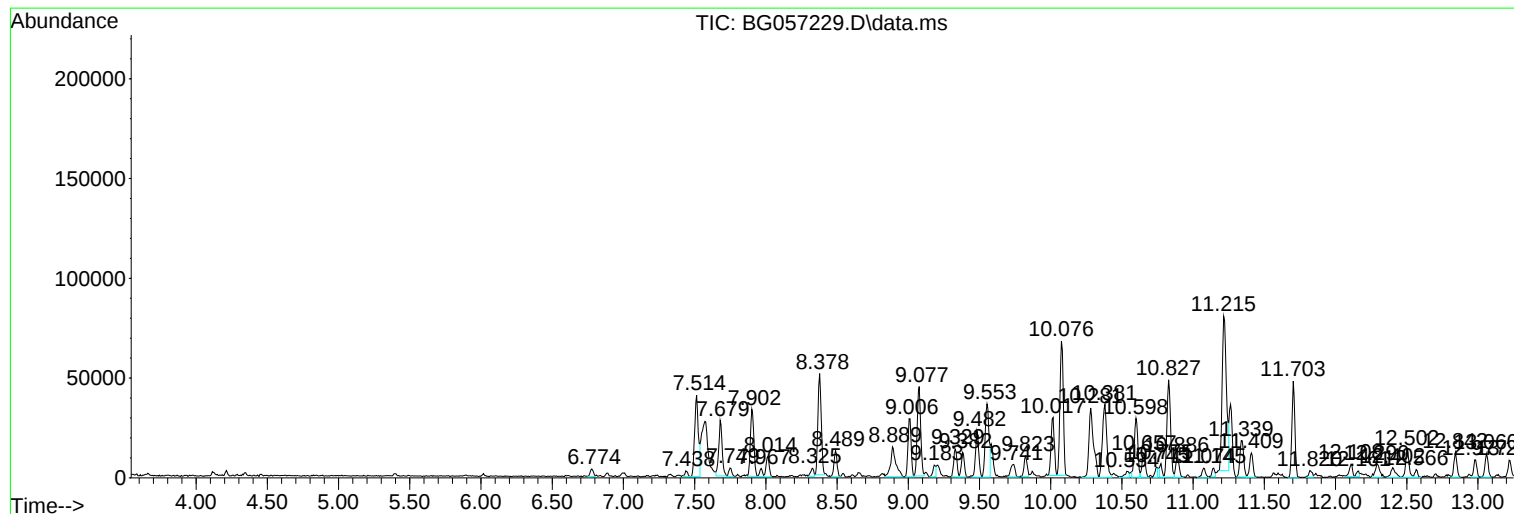
Sum of corrected areas: 5537564

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION

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



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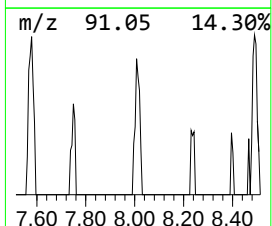
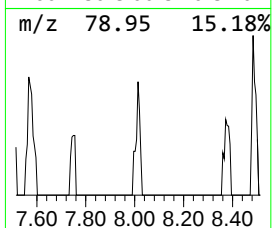
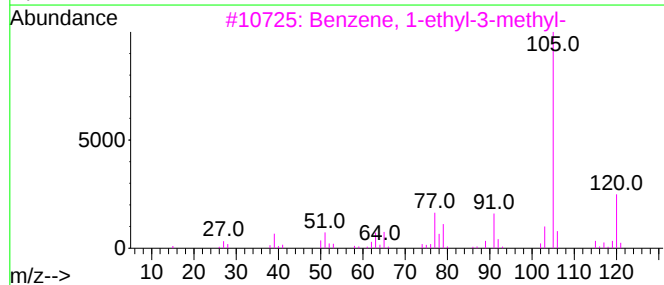
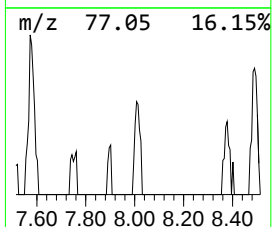
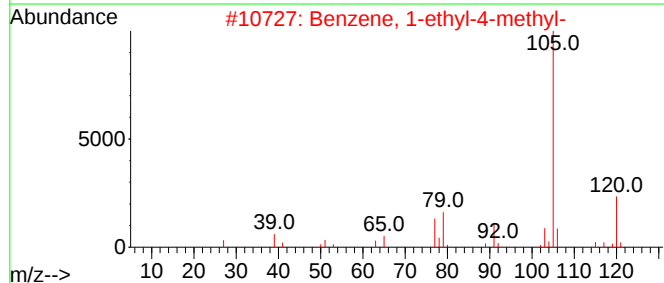
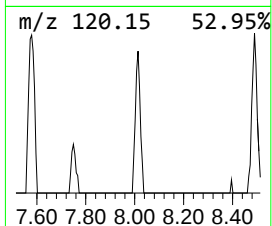
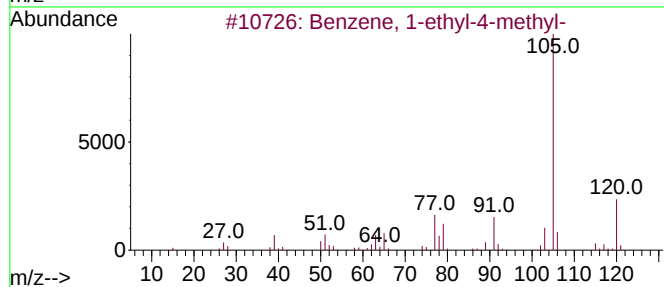
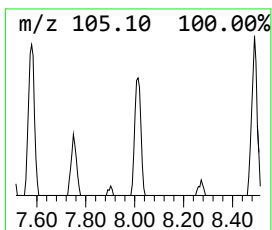
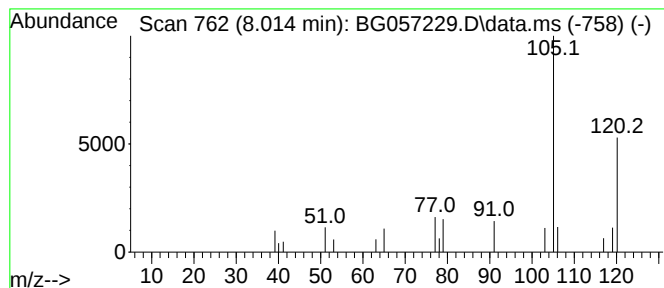
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, 1-ethyl-4-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.014	3.75 ng/ul	15293	1,4-Dichlorobenzene-d4	8.378

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	87
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	86
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	83



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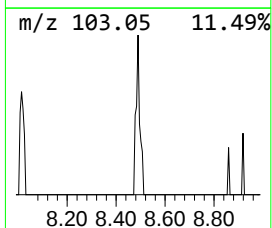
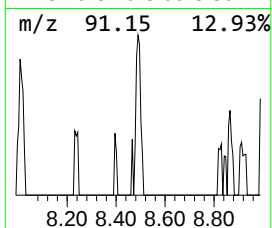
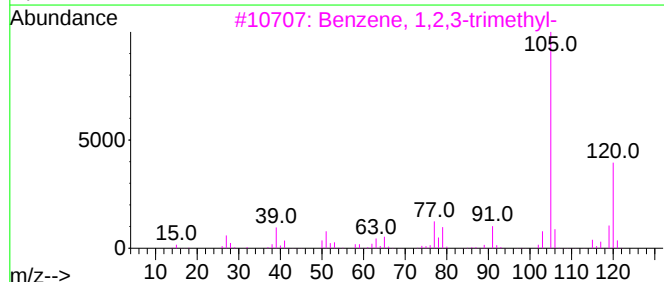
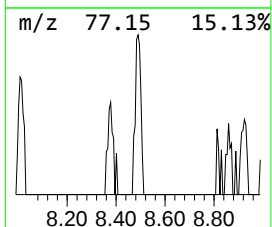
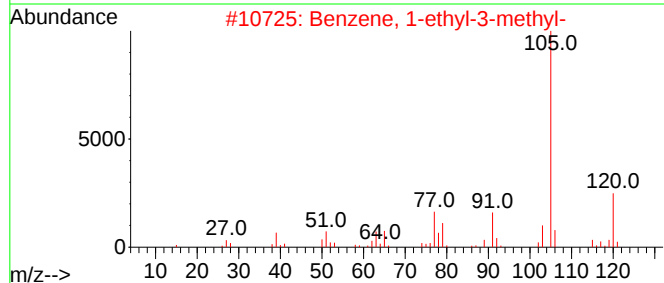
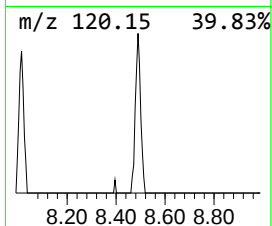
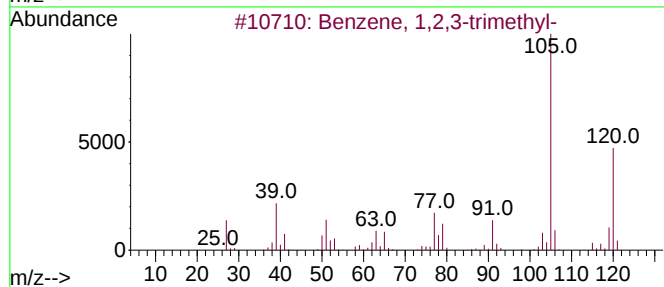
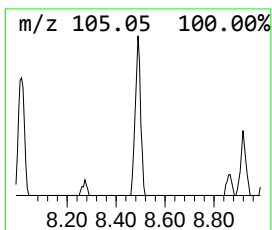
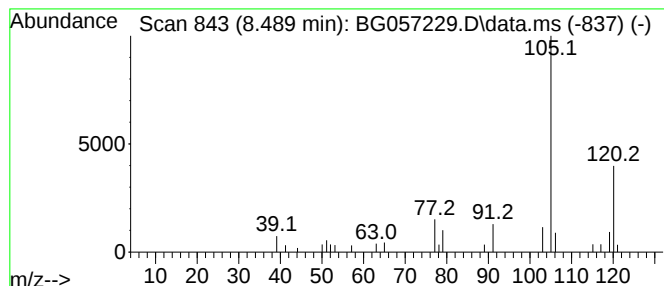
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,2,3-trimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.489	4.71 ng/ul	19197	1,4-Dichlorobenzene-d4	8.378

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Mesitylene	120	C9H12	000108-67-8	90



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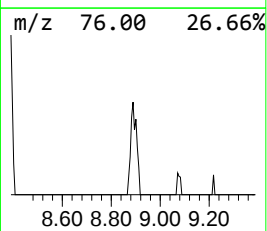
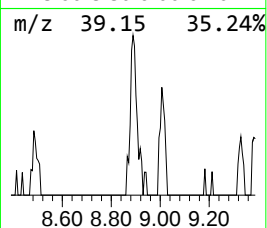
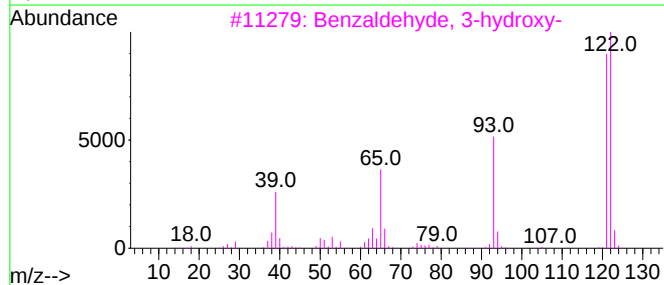
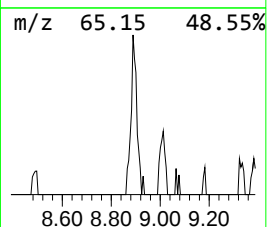
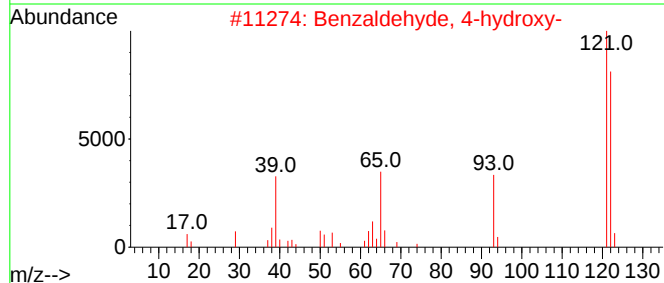
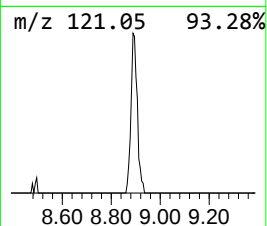
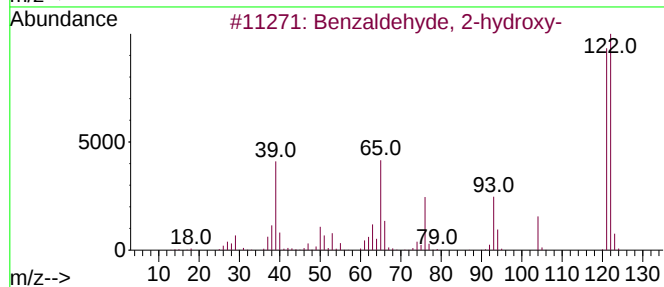
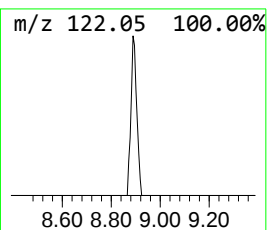
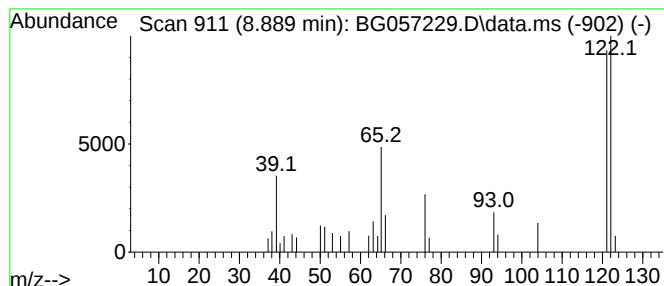
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzaldehyde, 2-hydroxy- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.889	9.59 ng/ul	39125	1,4-Dichlorobenzene-d4	8.378

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehyde, 2-hydroxy-	122	C7H6O2	000090-02-8	92
2			Benzaldehyde, 4-hydroxy-	122	C7H6O2	000123-08-0	91
3			Benzaldehyde, 3-hydroxy-	122	C7H6O2	000100-83-4	90
4			Benzaldehyde, 4-hydroxy-	122	C7H6O2	000123-08-0	90
5			Benzaldehyde, 3-hydroxy-	122	C7H6O2	000100-83-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
Operator : CG/JU
Sample : 02417-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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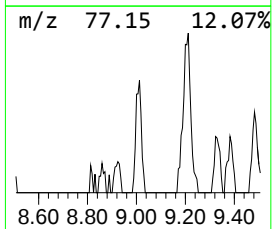
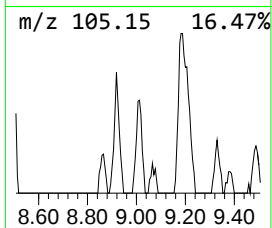
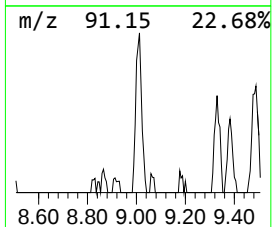
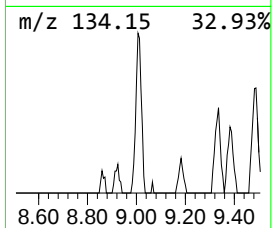
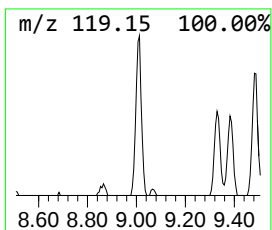
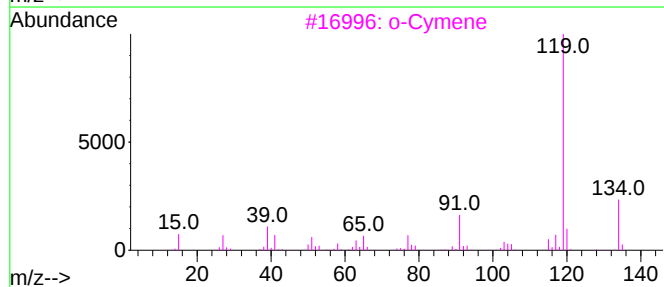
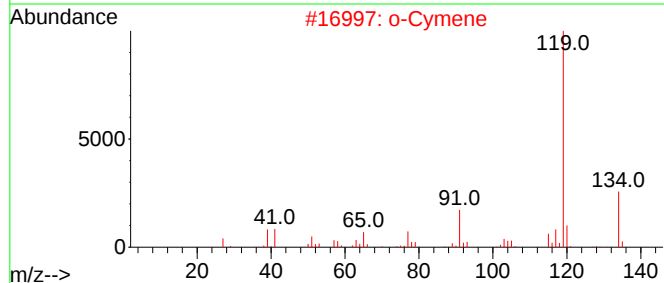
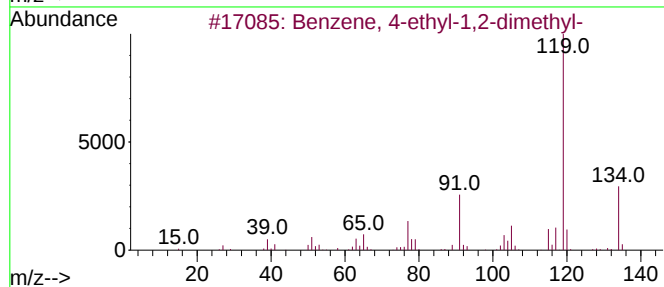
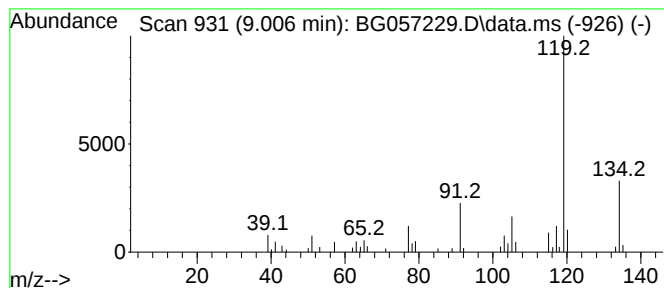
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.006	11.10 ng/ul	45308	1,4-Dichlorobenzene-d4	8.378

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	97
2			o-Cymene	134	C10H14	000527-84-4	94
3			o-Cymene	134	C10H14	000527-84-4	94
4			o-Cymene	134	C10H14	000527-84-4	94
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
Operator : CG/JU
Sample : 02417-01
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Instrument :
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ClientSampleId :
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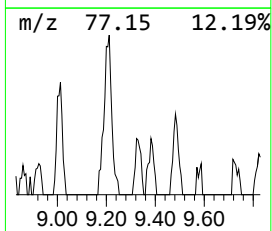
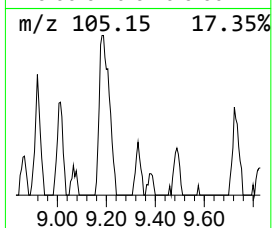
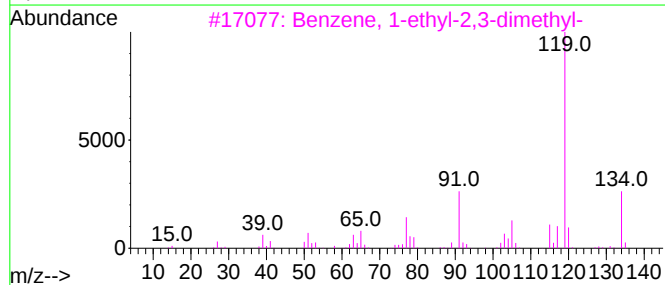
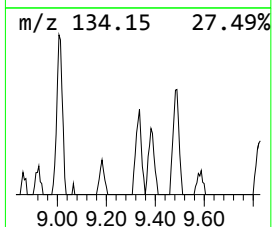
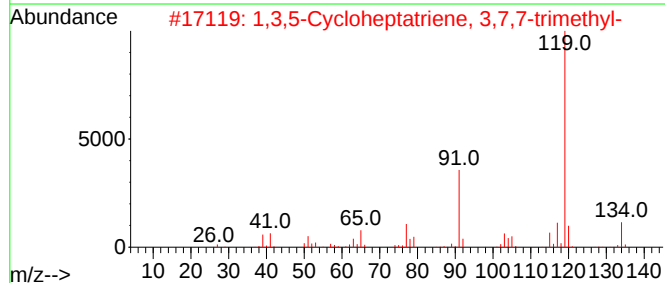
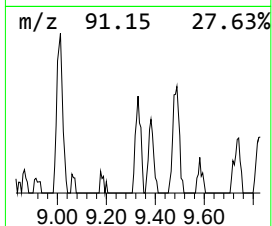
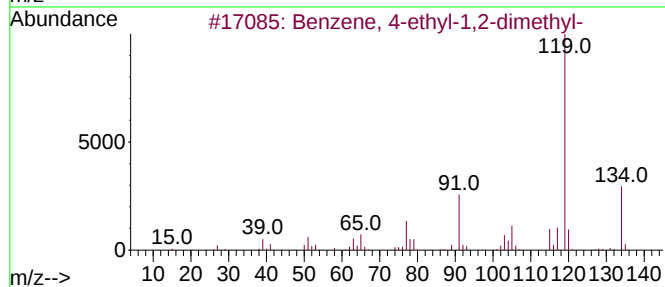
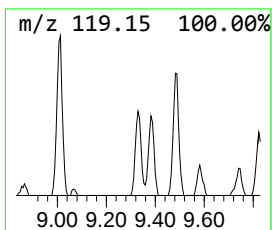
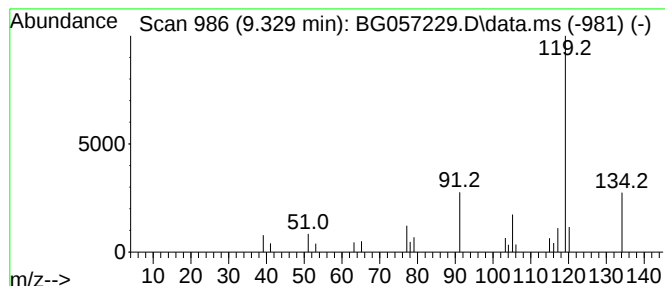
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,3,5-Cycloheptatriene, 3,7... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.329	5.39 ng/ul	21983	1,4-Dichlorobenzene-d4	8.378

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	91
2			1,3,5-Cycloheptatriene, 3,7,7-tr...	134	C10H14	003479-89-8	87
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	86
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	86
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
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Sample : 02417-01
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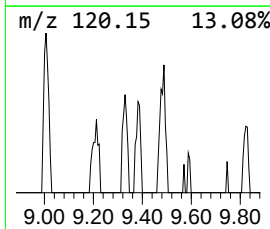
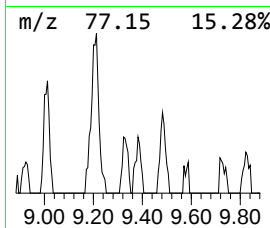
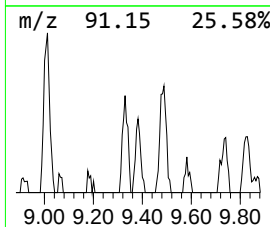
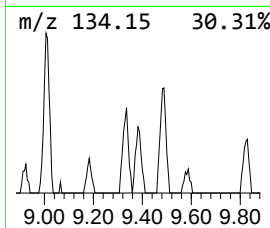
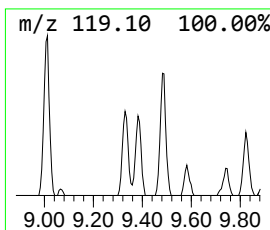
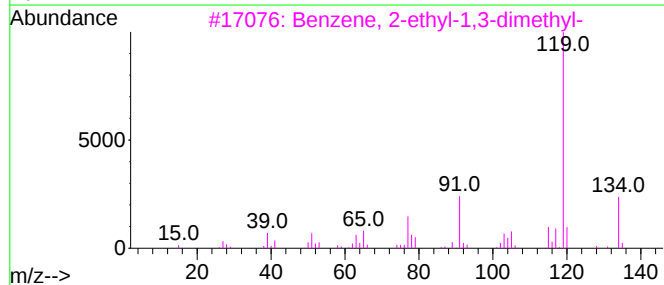
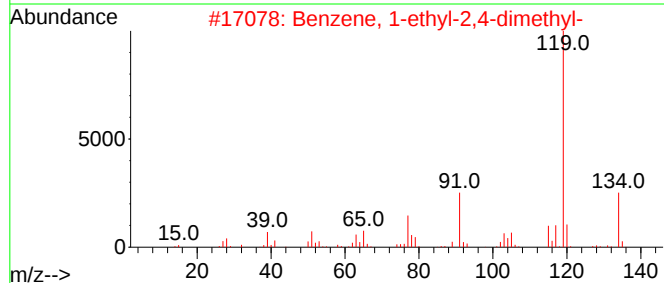
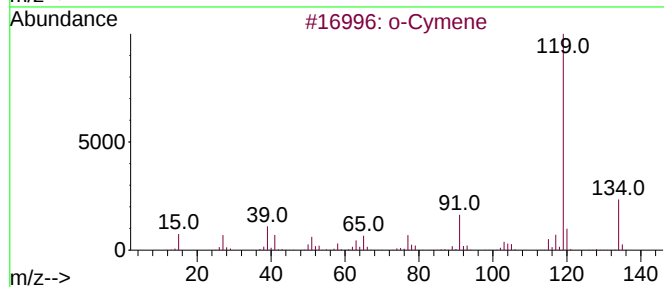
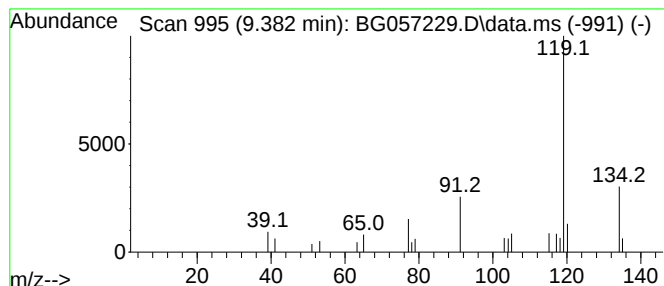
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 o-Cymene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.382	4.29 ng/ul	17512	1,4-Dichlorobenzene-d4	8.378

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	91
2			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	91
3			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	91
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
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Sample : 02417-01
Misc :
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Instrument :
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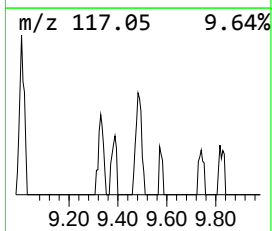
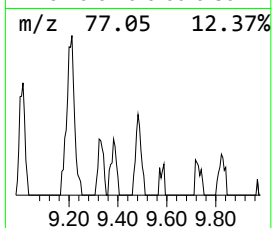
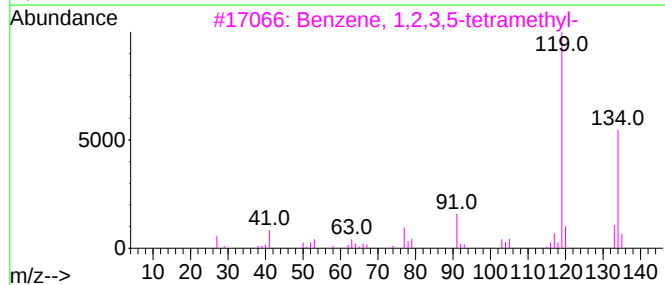
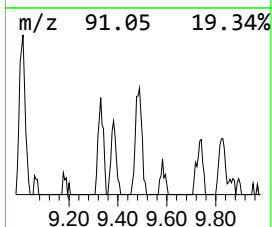
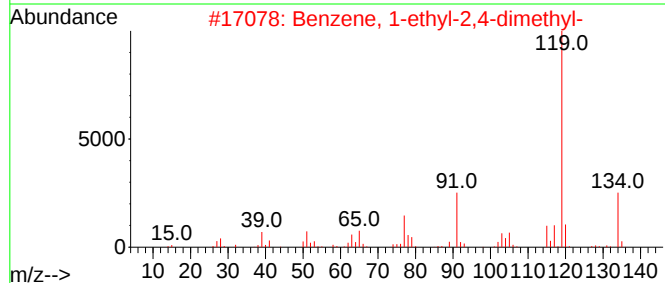
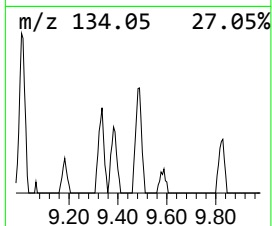
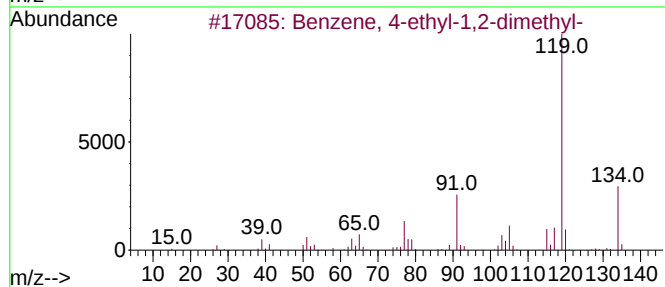
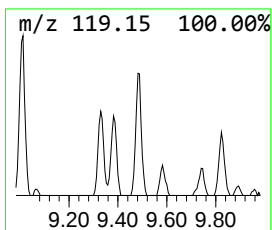
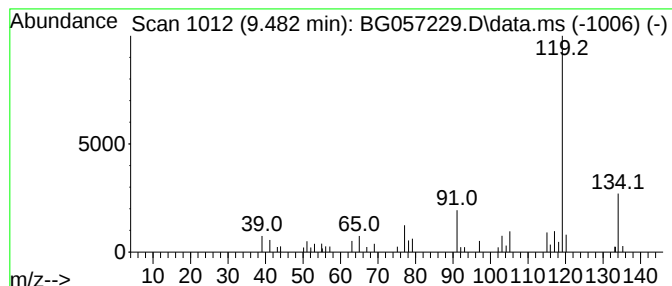
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.482	8.72 ng/ul	35570	1,4-Dichlorobenzene-d4	8.378

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
2			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	93
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
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Misc :
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Instrument :
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ClientSampleId :
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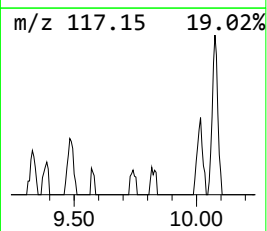
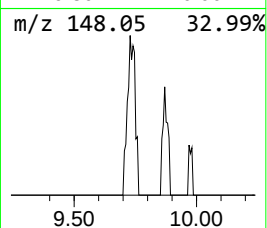
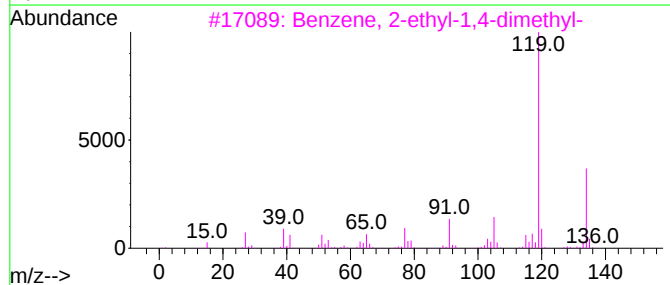
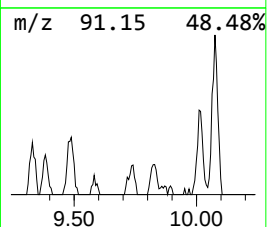
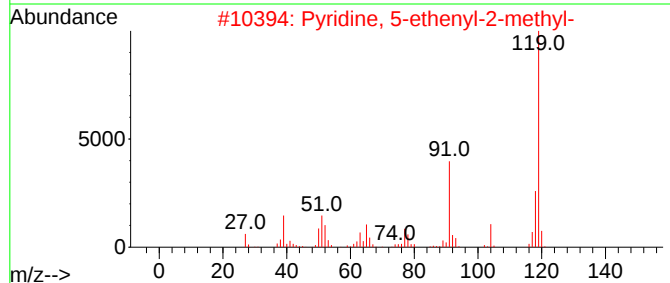
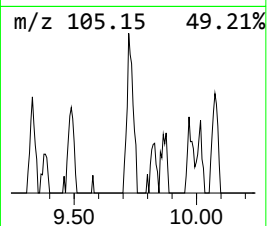
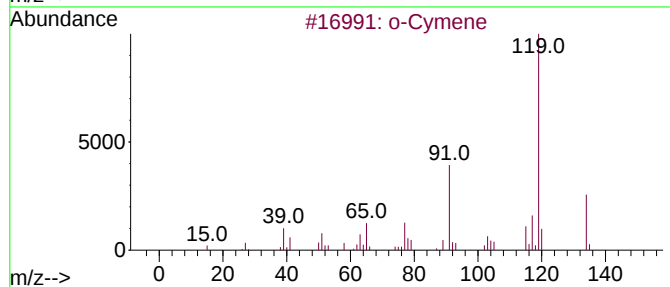
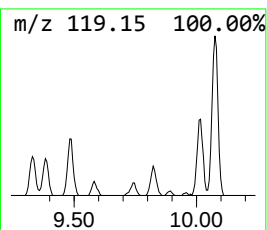
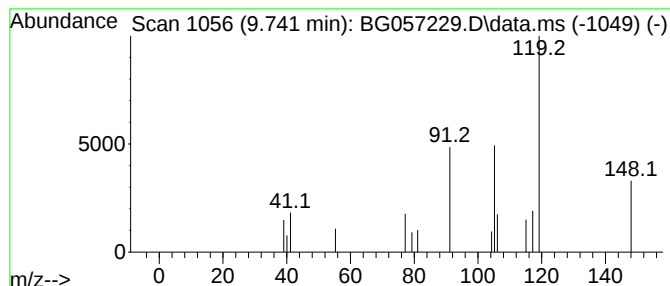
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-01 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.741	3.23 ng/ul	13190	1,4-Dichlorobenzene-d4	8.378

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	50
2			Pyridine, 5-ethenyl-2-methyl-	119	C8H9N	000140-76-1	9
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	9
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	9
5			o-Cymene	134	C10H14	000527-84-4	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
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ClientSampleId :
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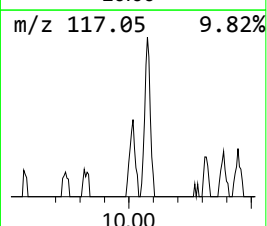
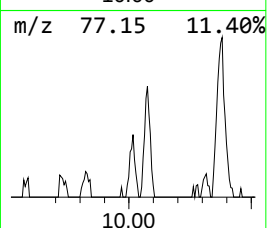
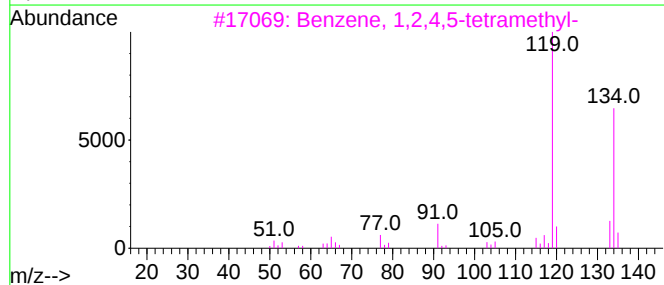
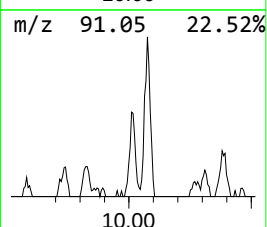
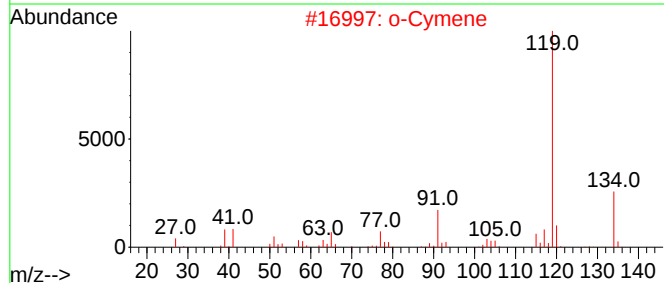
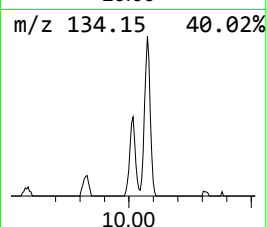
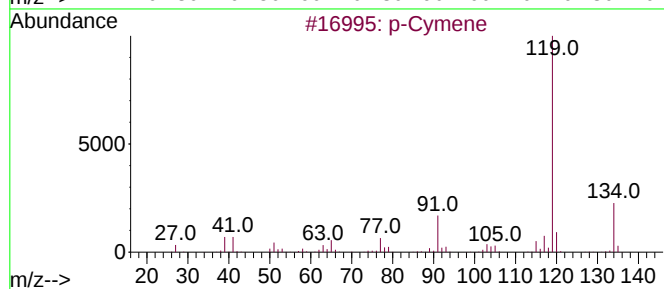
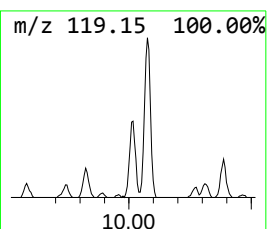
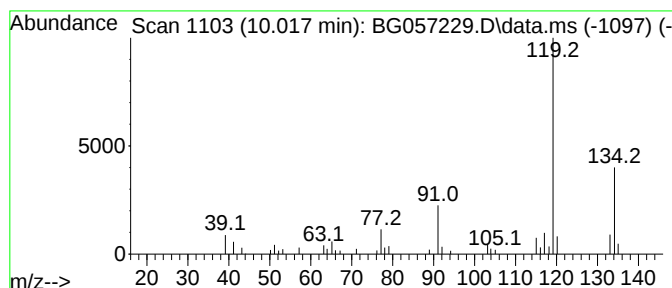
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 p-Cymene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.017	5.53 ng/ul	45727	Naphthalene-d8	11.215

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Cymene	134	C10H14	000099-87-6	94
2			o-Cymene	134	C10H14	000527-84-4	94
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
4			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
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ALS Vial : 19 Sample Multiplier: 1

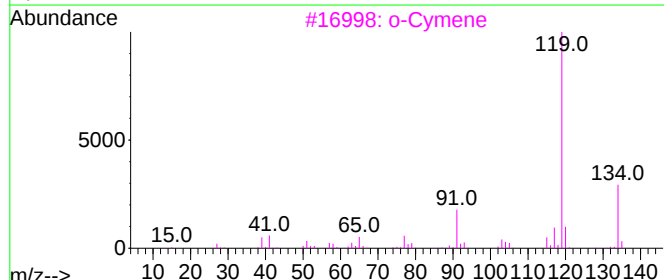
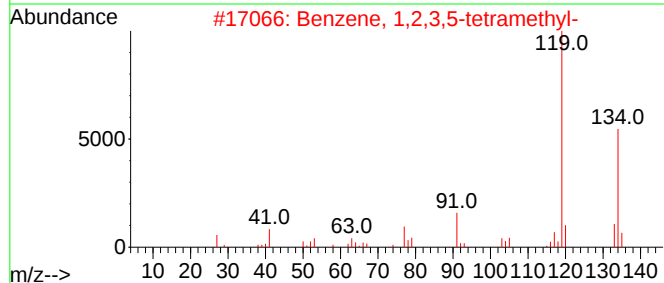
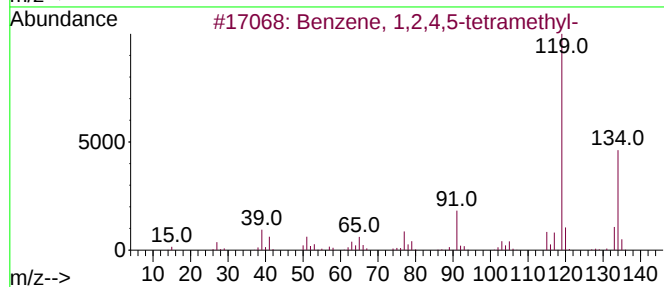
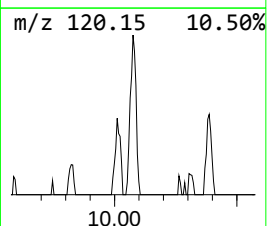
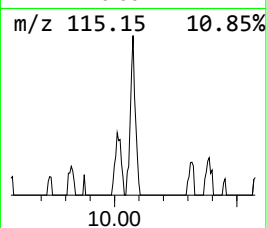
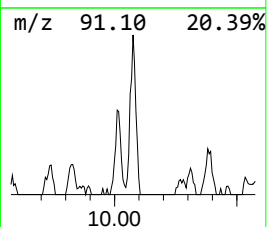
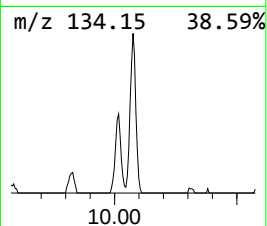
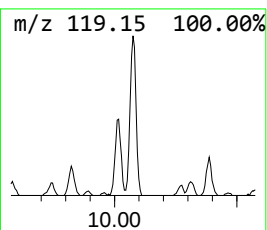
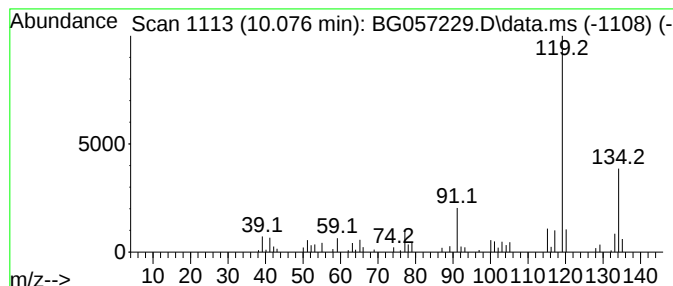
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ClientSampleId :
EW8W7

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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.076	13.32 ng/ul	110131	Naphthalene-d8	11.215	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
3	o-Cymene	134	C10H14	000527-84-4	93
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	93
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
Operator : CG/JU
Sample : 02417-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

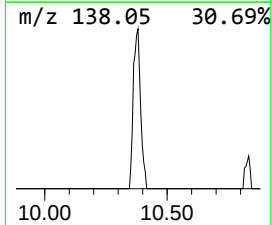
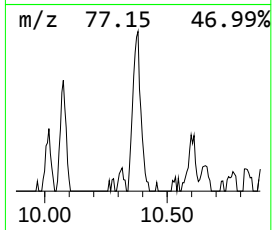
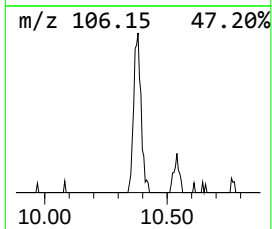
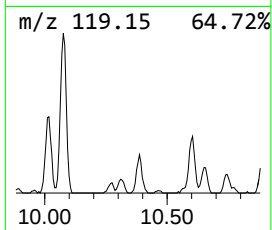
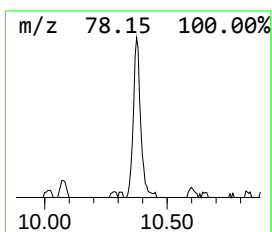
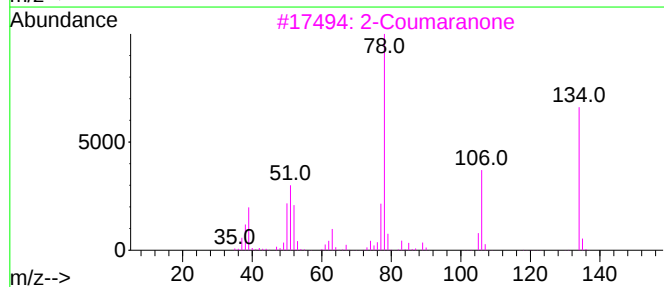
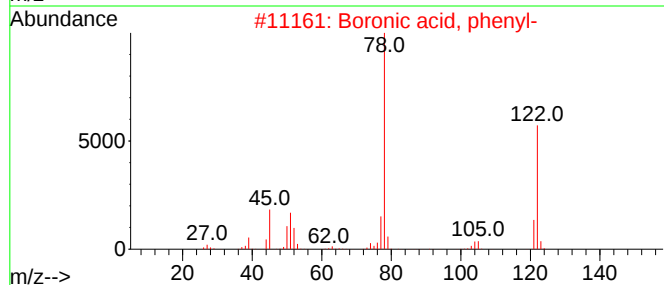
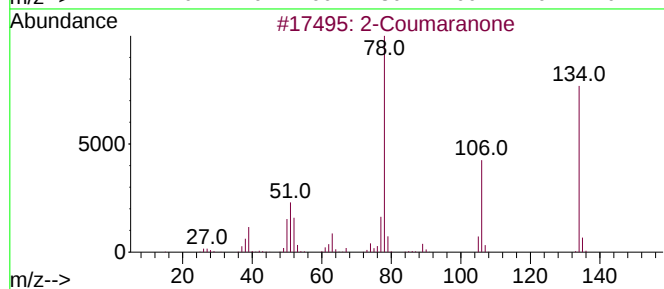
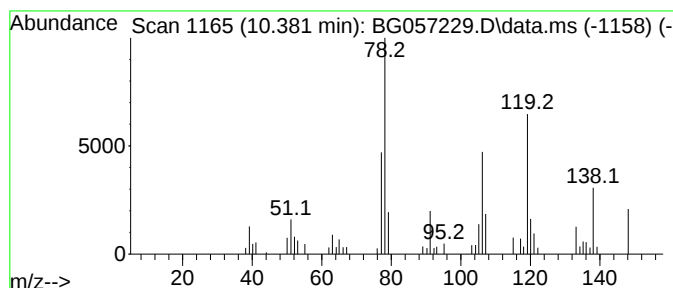
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 2-Coumaranone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.381	9.35 ng/ul	77308	Naphthalene-d8	11.215
Hit#	of 5	Tentative ID	MW MolForm	CAS# Qual
1		2-Coumaranone	134 C8H6O2	000553-86-6 58
2		Boronic acid, phenyl-	122 C6H7BO2	000098-80-6 50
3		2-Coumaranone	134 C8H6O2	000553-86-6 50
4		Salicyl alcohol	124 C7H8O2	000090-01-7 45
5		Tricyclo[4.2.2.0(2,5)]dec-7-en-3...	148 C10H12O	035237-74-2 42



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
Operator : CG/JU
Sample : 02417-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

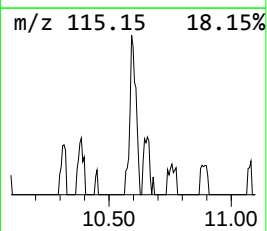
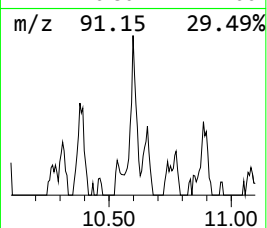
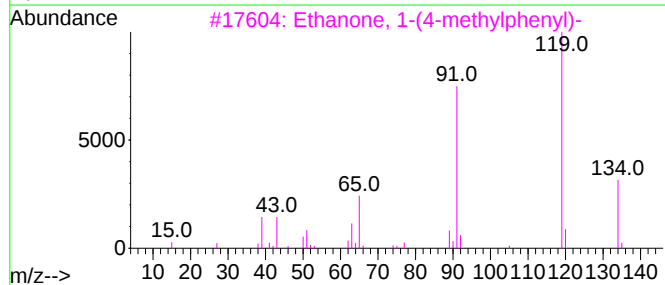
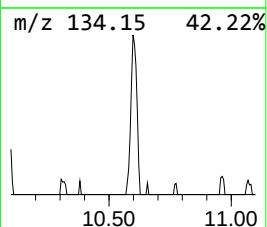
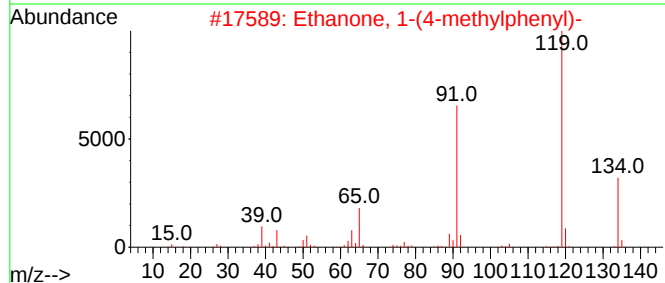
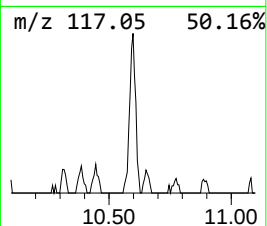
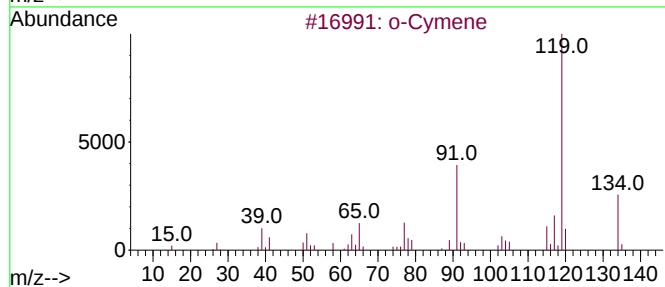
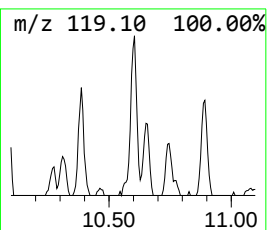
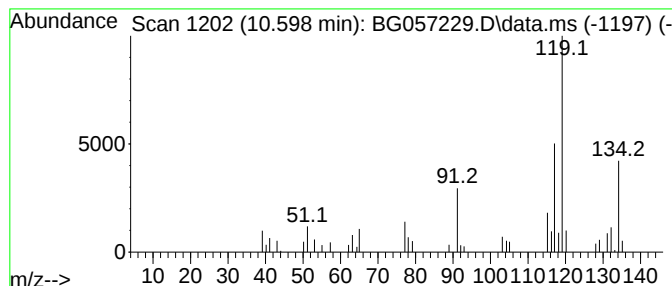
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 Ethanone, 1-(4-methylphenyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
10.598	6.17 ng/ul	51009	Naphthalene-d8		11.215	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	o-Cymene		134	C10H14	000527-84-4	70
2	Ethanone, 1-(4-methylphenyl)-		134	C9H10O	000122-00-9	64
3	Ethanone, 1-(4-methylphenyl)-		134	C9H10O	000122-00-9	64
4	Benzene, 1,2,4,5-tetramethyl-		134	C10H14	000095-93-2	64
5	o-Cymene		134	C10H14	000527-84-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
Operator : CG/JU
Sample : 02417-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

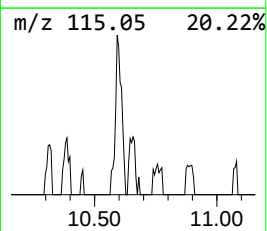
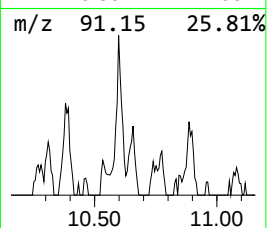
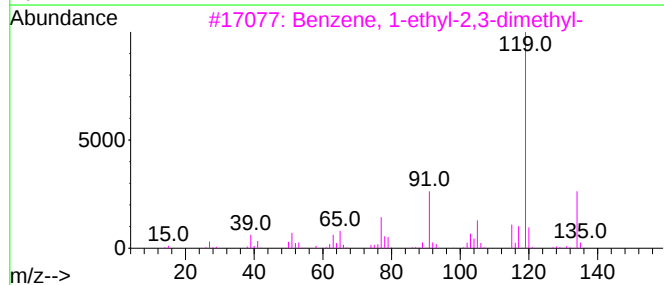
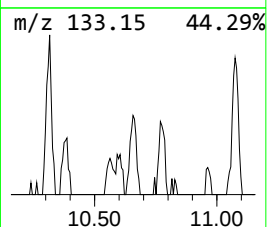
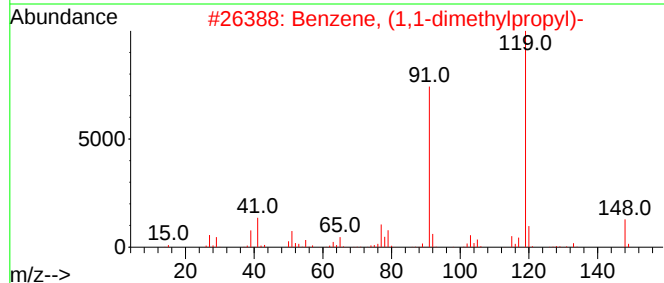
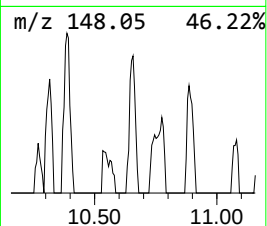
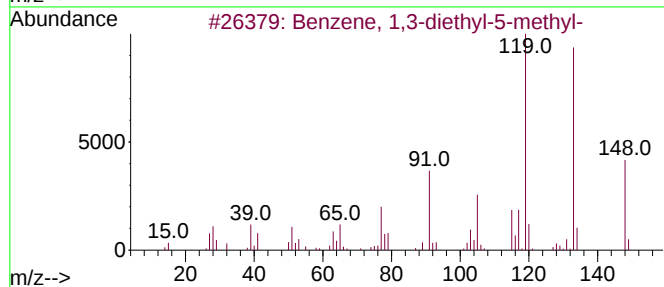
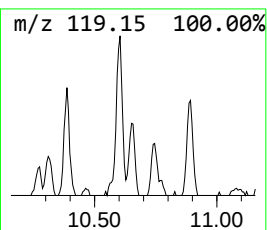
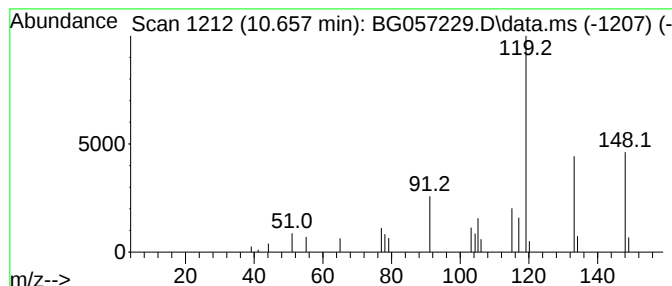
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Benzene, 1,3-diethyl-5-methyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.657	2.38 ng/ul	19680	Naphthalene-d8	11.215	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	53
2	Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	50
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	32
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	32
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
Operator : CG/JU
Sample : 02417-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

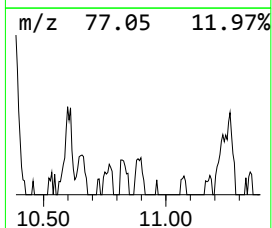
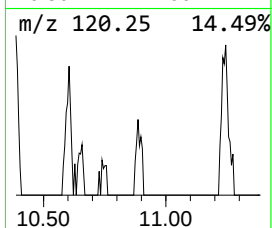
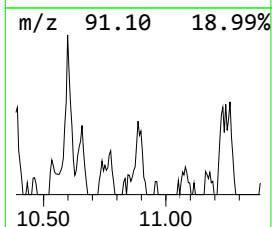
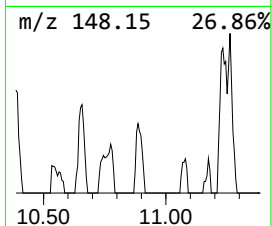
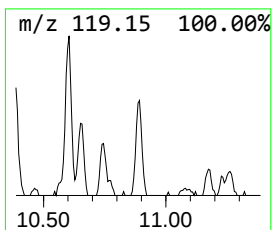
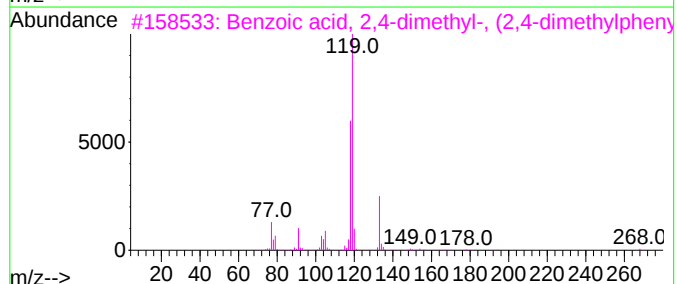
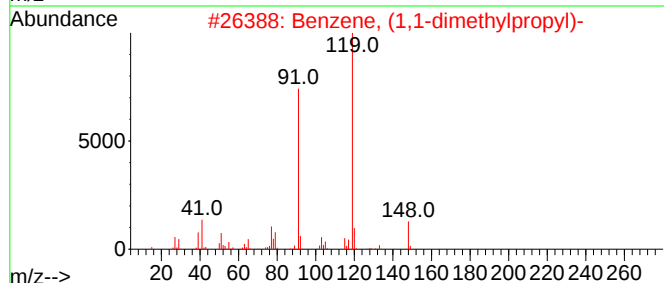
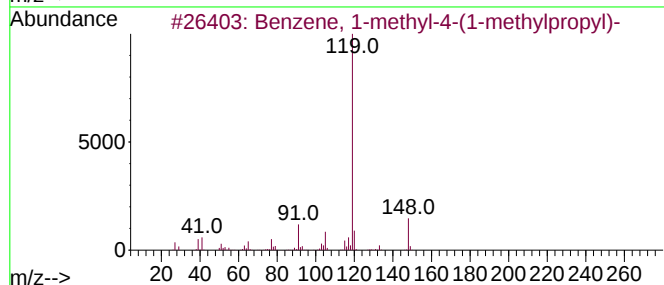
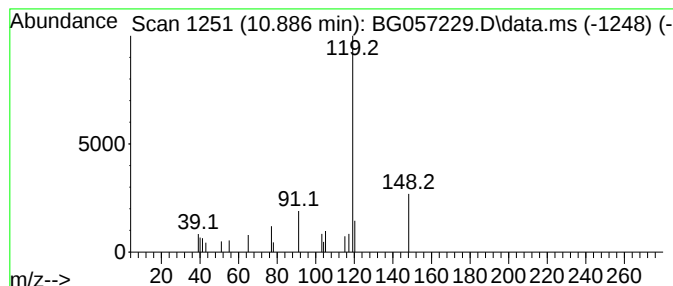
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 14 Benzene, 1-methyl-4-(1-meth... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.886	2.01 ng/ul	16639	Naphthalene-d8	11.215	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	86
2	Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	80
3	Benzoic acid, 2,4-dimethyl-, (2,...	268	C18H20O2	055000-43-6	72
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	72
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
Operator : CG/JU
Sample : 02417-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW8W7

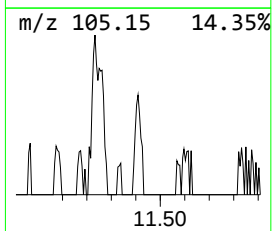
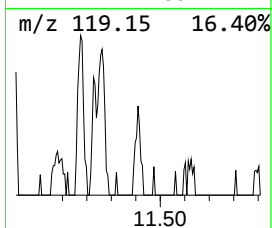
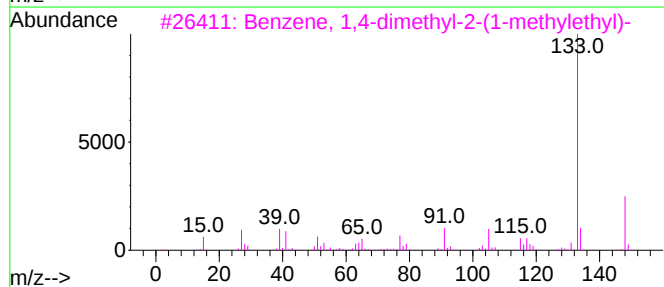
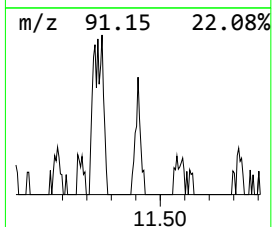
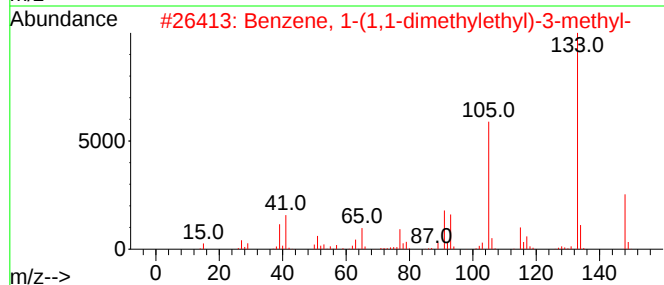
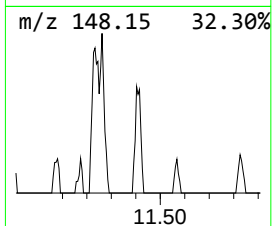
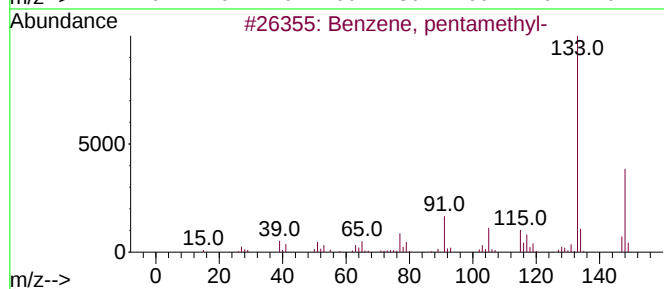
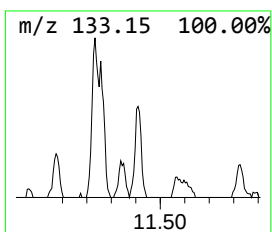
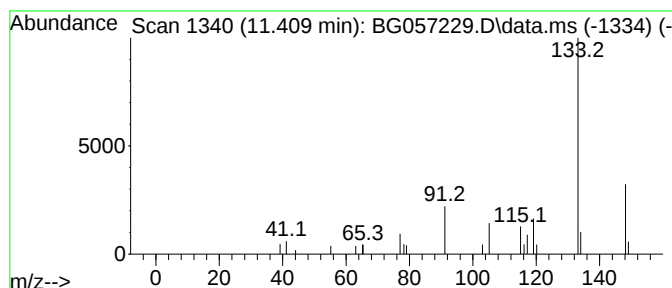
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Benzene, pentamethyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.409	2.39 ng/ul	19784	Naphthalene-d8	11.215

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, pentamethyl-	148	C11H16	000700-12-9	87
2			Benzene, 1-(1,1-dimethylethyl)-3...	148	C11H16	001075-38-3	87
3			Benzene, 1,4-dimethyl-2-(1-methy...	148	C11H16	004132-72-3	86
4			Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	86
5			3,4-Dimethylcumene	148	C11H16	1000370-34-1	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
Operator : CG/JU
Sample : 02417-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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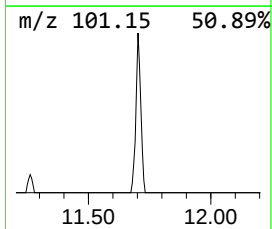
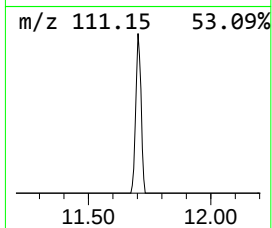
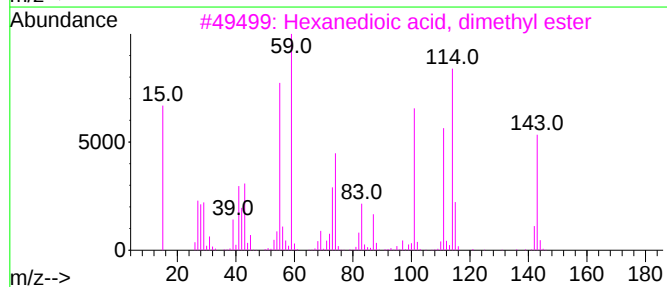
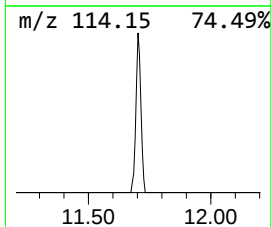
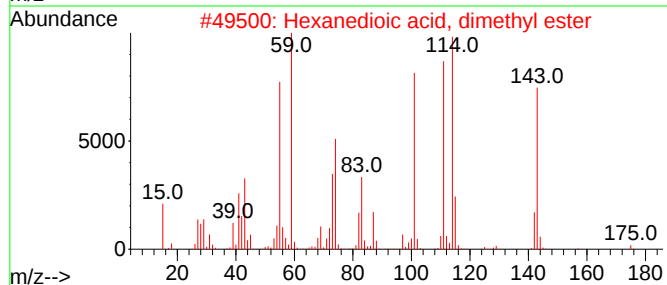
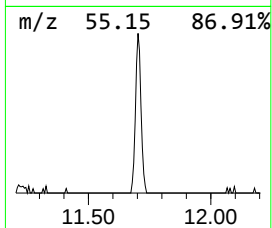
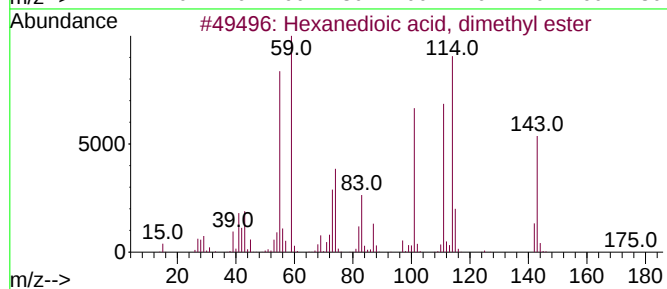
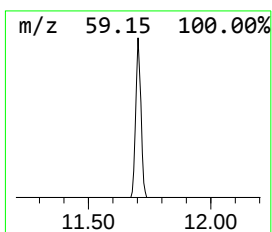
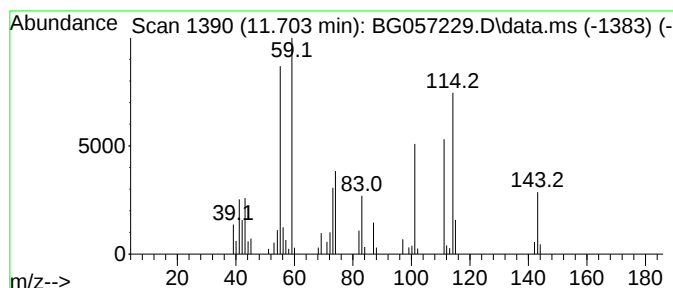
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Hexanedioic acid, dimethyl ... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.703	8.58 ng/ul	70904	Naphthalene-d8	11.215

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexanedioic acid, dimethyl ester	174	C8H14O4	000627-93-0	91
2		Hexanedioic acid, dimethyl ester	174	C8H14O4	000627-93-0	91
3		Hexanedioic acid, dimethyl ester	174	C8H14O4	000627-93-0	78
4		Hexanedioic acid, dimethyl ester	174	C8H14O4	000627-93-0	64
5		Pentanedioic acid, 2-methyl-, di...	174	C8H14O4	014035-94-0	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
Operator : CG/JU
Sample : 02417-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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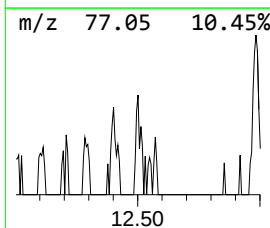
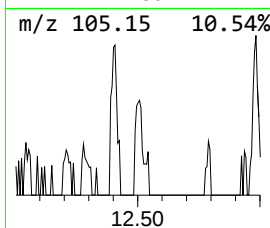
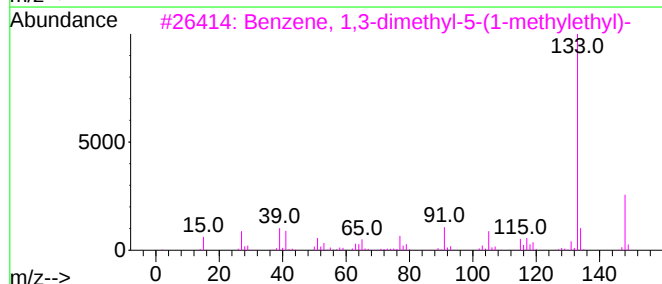
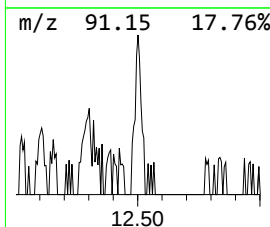
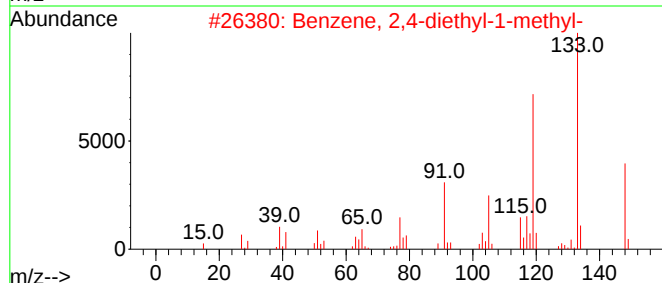
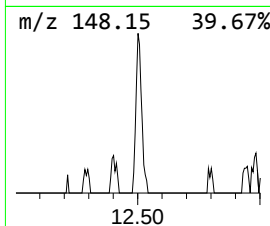
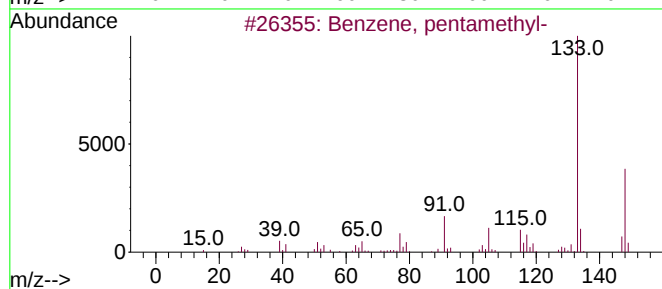
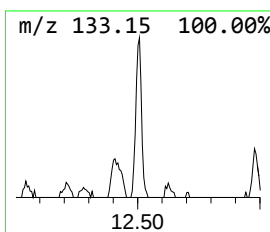
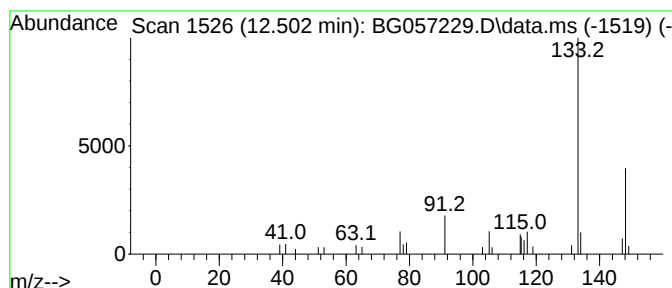
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Benzene, 2,4-diethyl-1-methyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.502	2.78 ng/ul	22965	Naphthalene-d8	11.215

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, pentamethyl-	148	C11H16	000700-12-9	94
2			Benzene, 2,4-diethyl-1-methyl-	148	C11H16	001758-85-6	91
3			Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	91
4			Benzene, pentamethyl-	148	C11H16	000700-12-9	91
5			2-tert-Butyltoluene	148	C11H16	001074-92-6	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
Operator : CG/JU
Sample : 02417-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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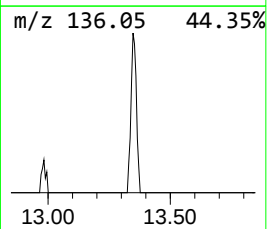
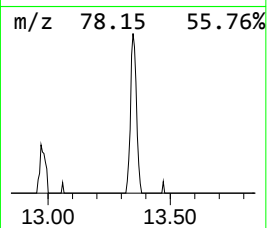
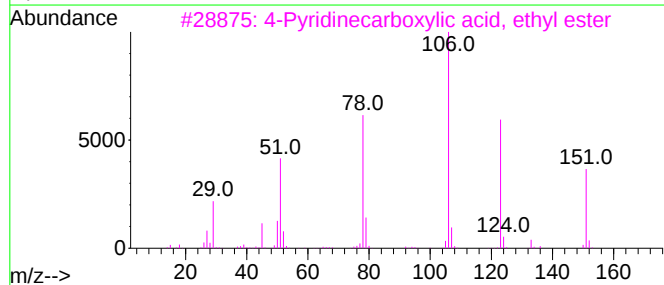
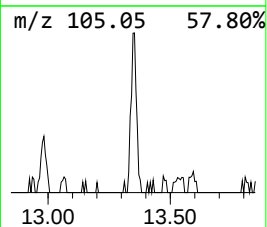
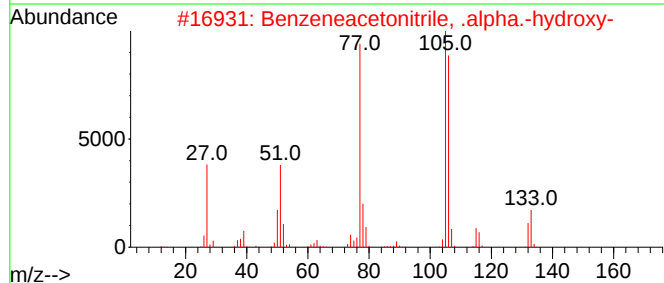
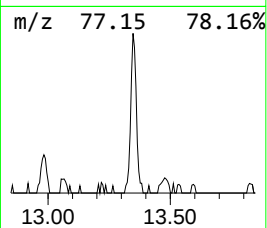
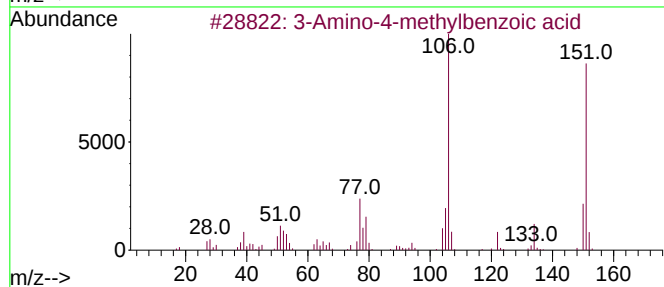
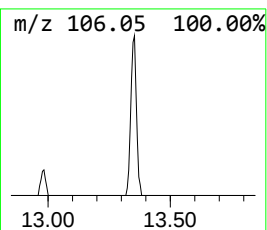
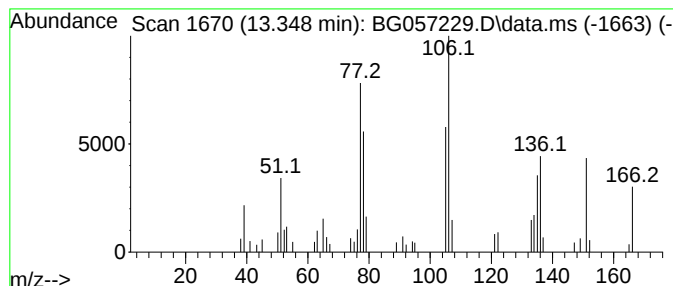
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 3-Amino-4-methylbenzoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.348	4.98 ng/ul	52953	Acenaphthene-d10	14.986

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Amino-4-methylbenzoic acid	151	C8H9NO2	002458-12-0	50
2			Benzeneacetonitrile, .alpha.-hyd...	133	C8H7NO	000532-28-5	43
3			4-Pyridinecarboxylic acid, ethyl...	151	C8H9NO2	001570-45-2	38
4			N-Phenylglycine	151	C8H9NO2	000103-01-5	35
5			4-Pyridinecarboxylic acid, ethyl...	151	C8H9NO2	001570-45-2	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
Operator : CG/JU
Sample : 02417-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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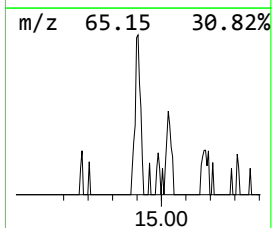
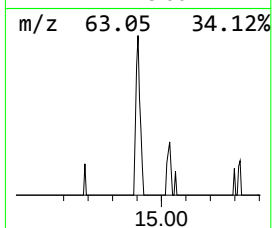
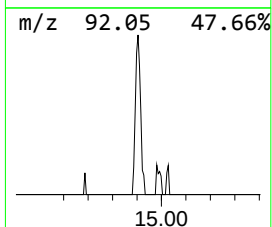
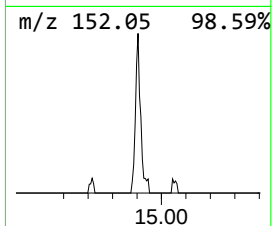
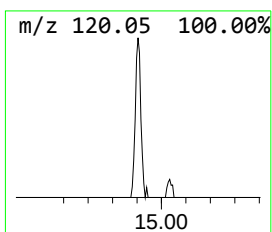
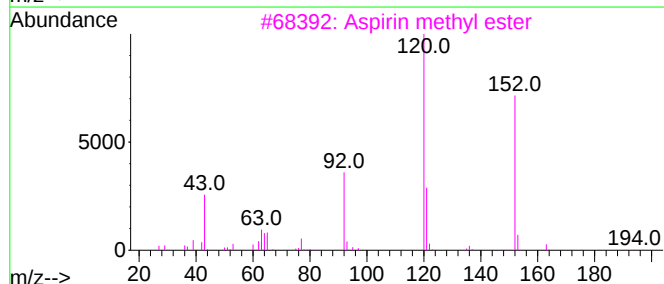
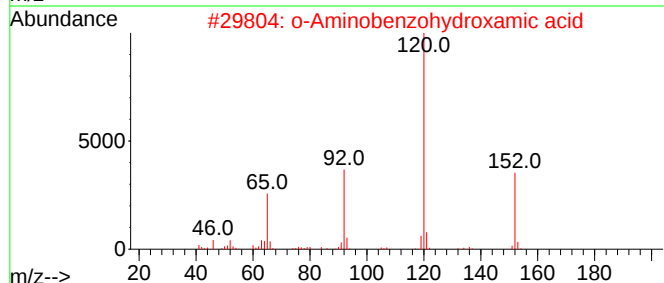
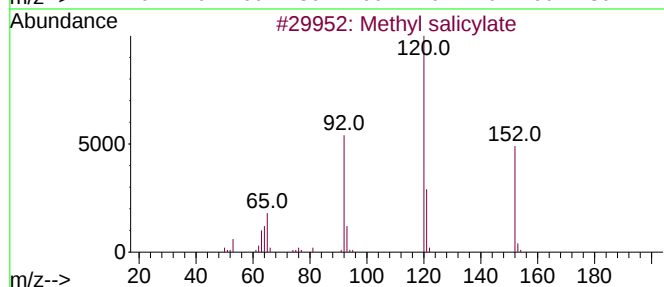
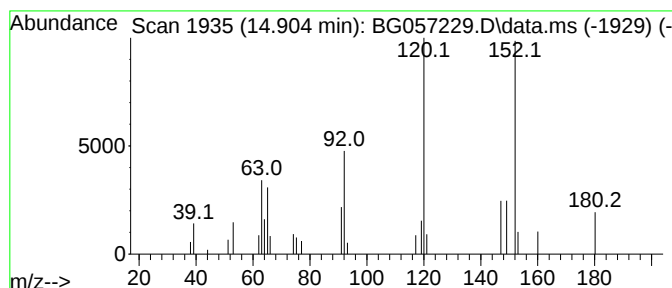
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 Methyl salicylate Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.904	2.21 ng/ul	23466	Acenaphthene-d10	14.986

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl salicylate	152	C8H8O3	000119-36-8	58
2			o-Aminobenzohydroxamic acid	152	C7H8N2O2	005623-04-1	50
3			Aspirin methyl ester	194	C10H10O4	000580-02-9	42
4			2-Hydroxy-3-cyanopyridine	120	C6H4N2O	020577-27-9	35
5			1-Propanone, 1-(4-aminophenyl)-	149	C9H11NO	000070-69-9	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
Operator : CG/JU
Sample : 02417-01
Misc :
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Instrument :
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ClientSampleId :
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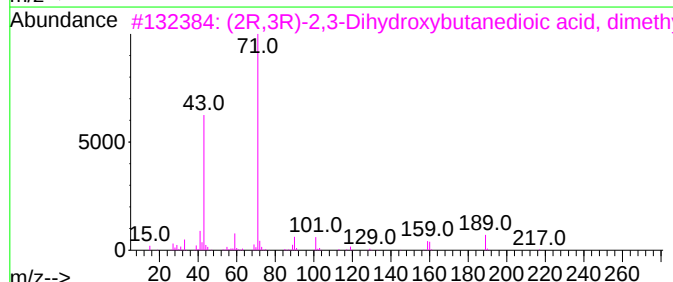
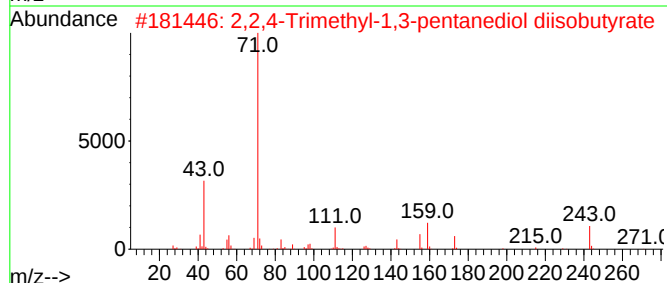
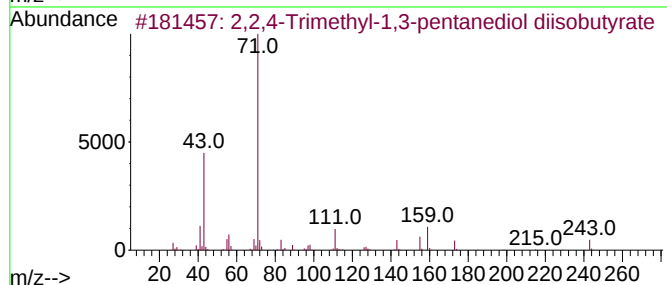
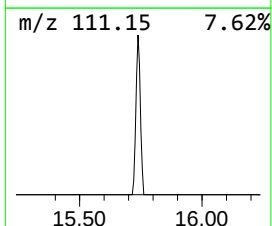
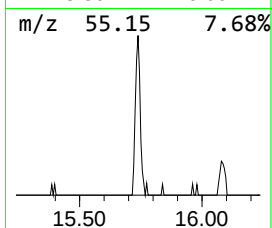
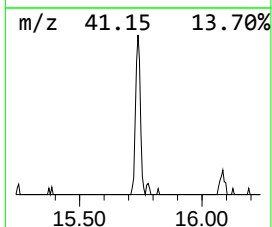
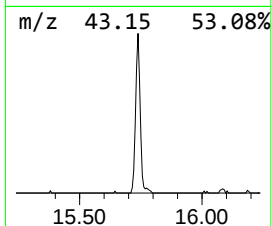
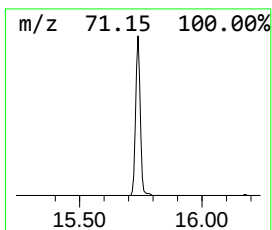
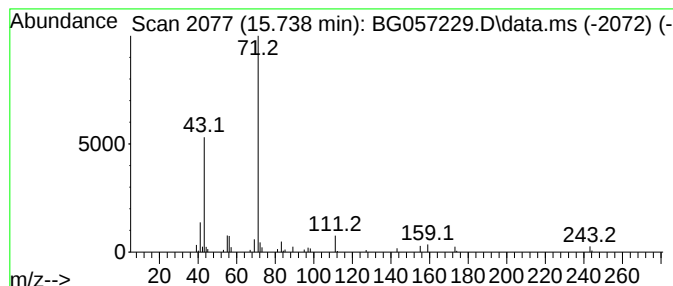
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 2,2,4-Trimethyl-1,3-pentane... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.738	9.25 ng/ul	98388	Acenaphthene-d10	14.986

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2,4-Trimethyl-1,3-pentanediol ...	286	C16H30O4	006846-50-0	83	
2	2,2,4-Trimethyl-1,3-pentanediol ...	286	C16H30O4	006846-50-0	83	
3	(2R,3R)-2,3-Dihydroxybutanedioic...	248	C10H16O7	1000462-99-8	64	
4	Oxalic acid, dineopentyl ester	230	C12H22O4	1010309-72-7	59	
5	Oxalic acid, neopentyl octyl ester	272	C15H28O4	1000309-73-2	45	



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
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Sample : 02417-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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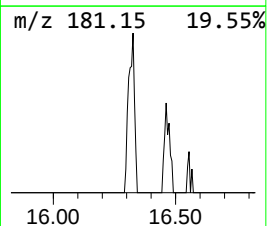
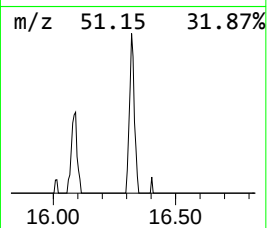
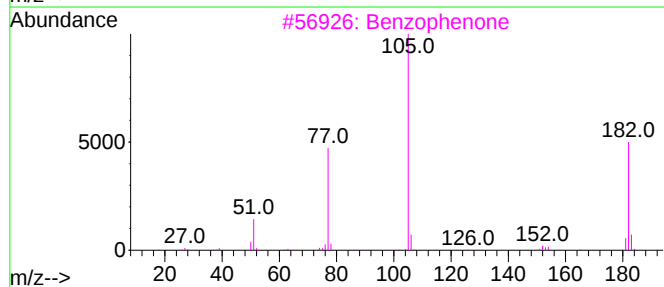
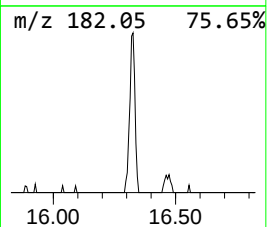
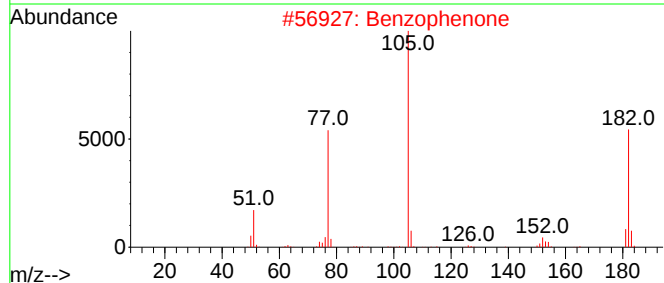
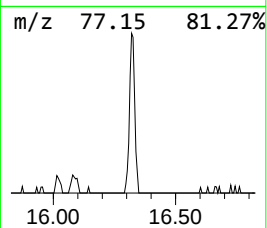
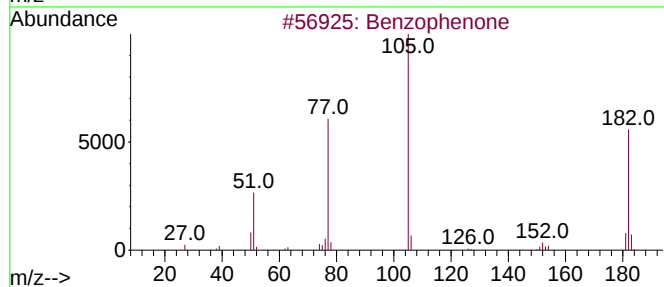
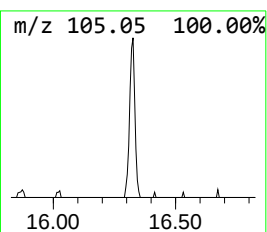
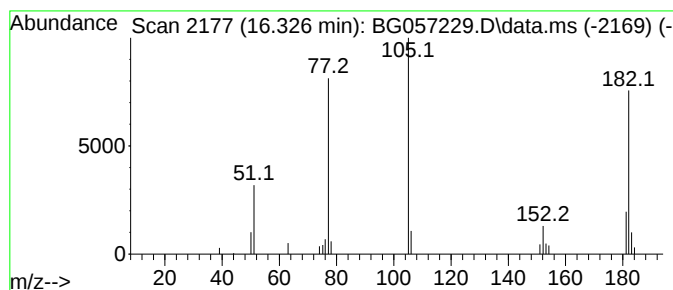
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 Benzophenone Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.326	3.01 ng/ul	32013	Acenaphthene-d10	14.986

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	90
2			Benzophenone	182	C13H10O	000119-61-9	87
3			Benzophenone	182	C13H10O	000119-61-9	86
4			Benzophenone	182	C13H10O	000119-61-9	64
5			Azobenzene, (E)-	182	C12H10N2	017082-12-1	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
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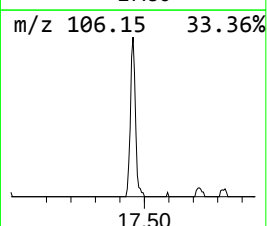
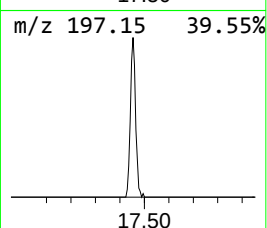
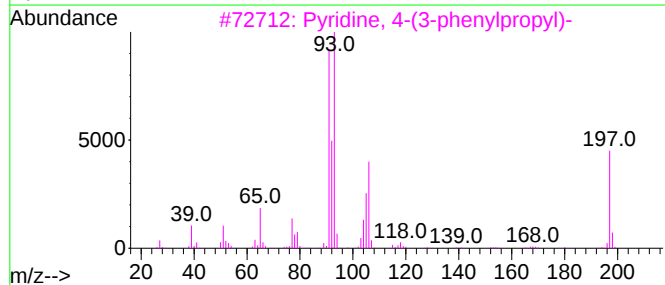
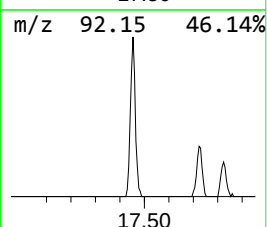
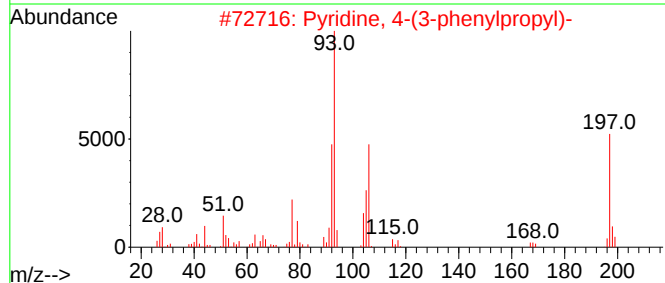
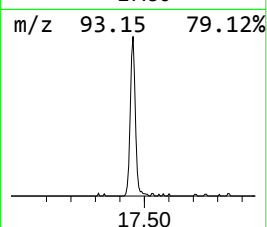
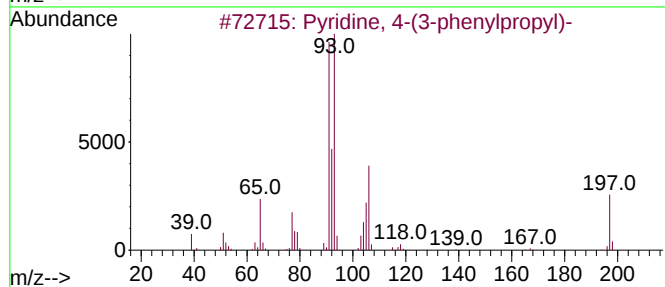
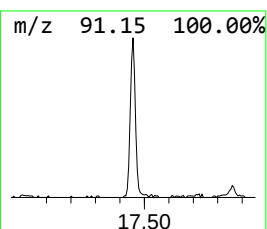
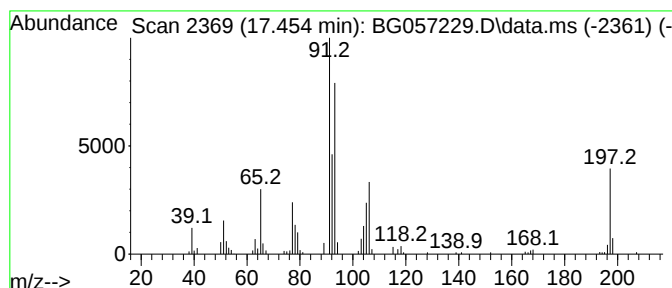
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 22 Pyridine, 4-(3-phenylpropyl)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.454	11.73 ng/ul	159335	Phenanthrene-d10		17.724	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Pyridine, 4-(3-phenylpropyl)-		197	C14H15N	002057-49-0	97
2	Pyridine, 4-(3-phenylpropyl)-		197	C14H15N	002057-49-0	96
3	Pyridine, 4-(3-phenylpropyl)-		197	C14H15N	002057-49-0	95
4	Tricyclo[4.1.0.0(2,4)]heptane, 5...		202	C13H14S	071055-11-3	38
5	Pyridine, 2-methyl-		93	C6H7N	000109-06-8	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
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ALS Vial : 19 Sample Multiplier: 1

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ClientSampleId :
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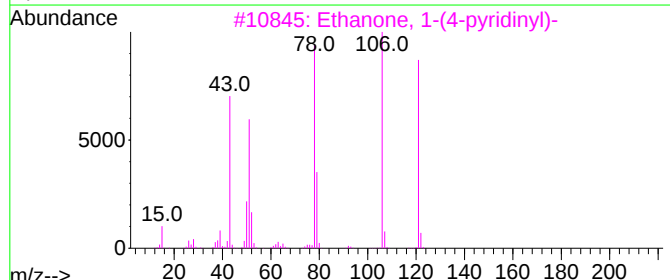
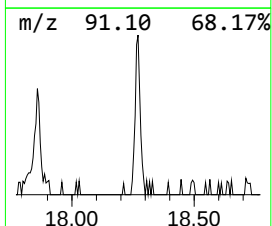
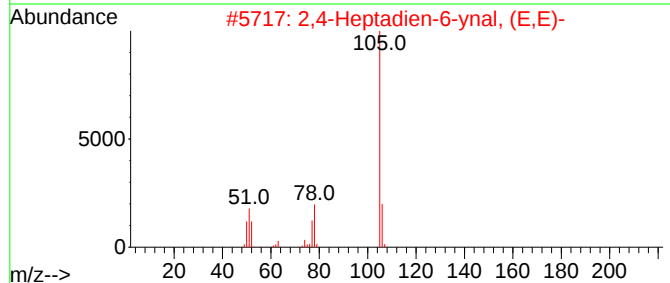
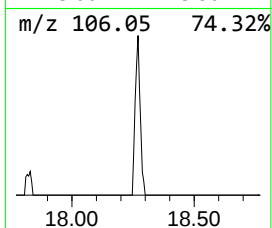
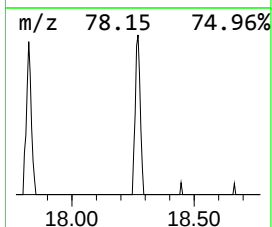
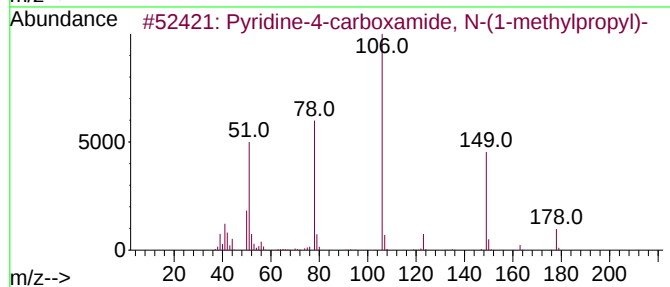
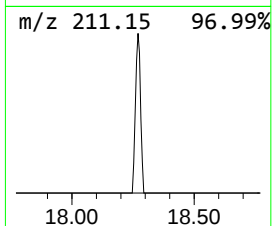
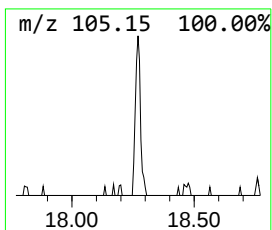
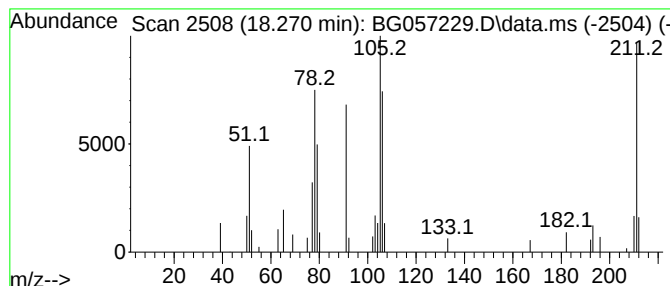
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 Pyridine-4-carboxamide, N-(... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.270	2.28 ng/ul	30915	Phenanthrene-d10	17.724

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyridine-4-carboxamide, N-(1-met...	178	C10H14N2O	311793-05-2	59
2			2,4-Heptadien-6-ynal, (E,E)-	106	C7H6O	007200-04-6	59
3			Ethanone, 1-(4-pyridinyl)-	121	C7H7NO	001122-54-9	59
4			Ethanone, 1-(3-pyridinyl)-	121	C7H7NO	000350-03-8	59
5			Nicorandil	211	C8H9N3O4	065141-46-0	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
Operator : CG/JU
Sample : 02417-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW8W7

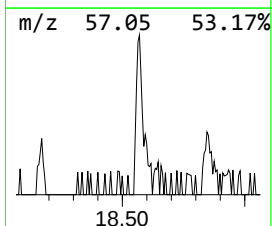
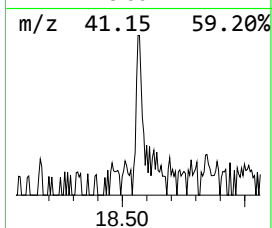
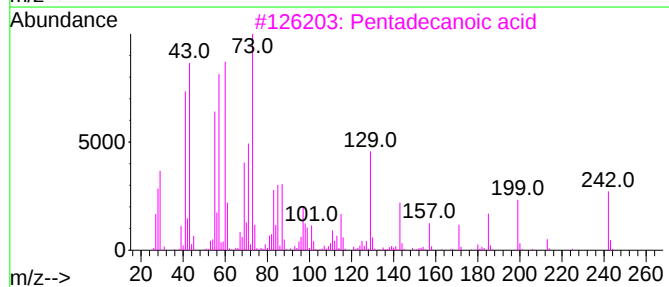
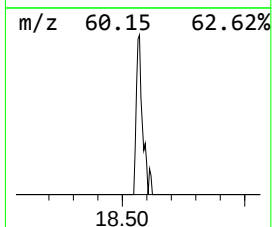
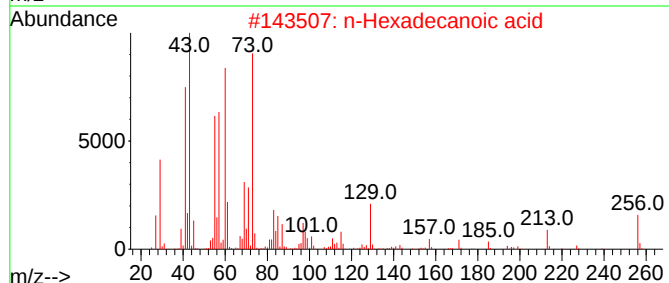
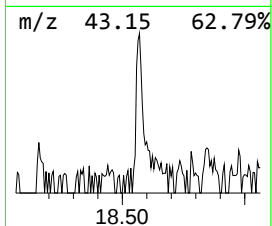
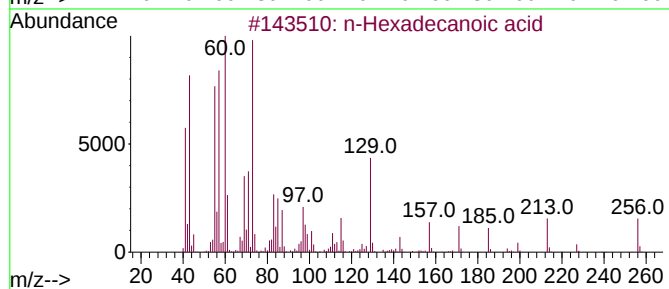
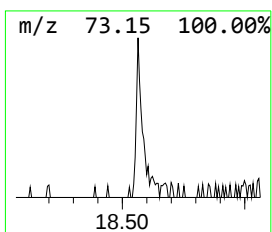
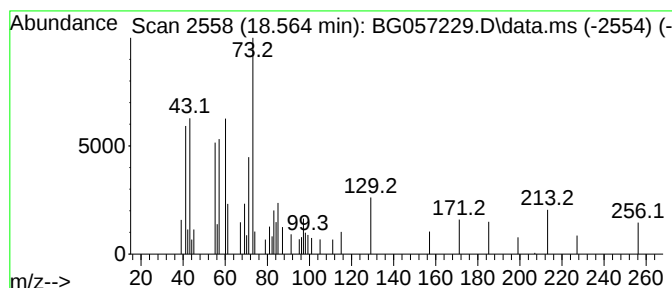
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 n-Hexadecanoic acid Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.564	2.95 ng/ul	40133	Phenanthrene-d10	17.724

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	89
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	74
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	53
5		Decanoic acid, silver(1+) salt	278	C10H19AgO2	013126-67-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
Operator : CG/JU
Sample : 02417-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

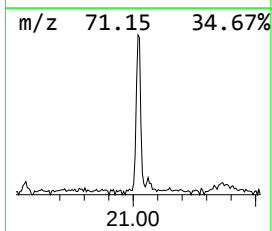
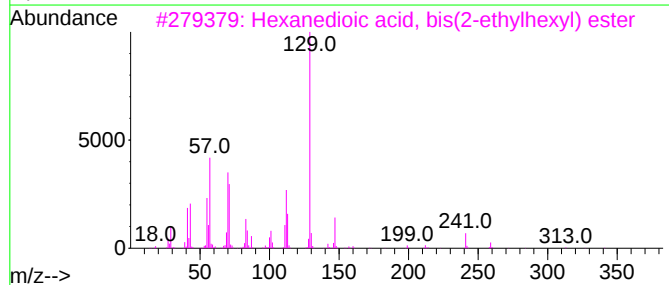
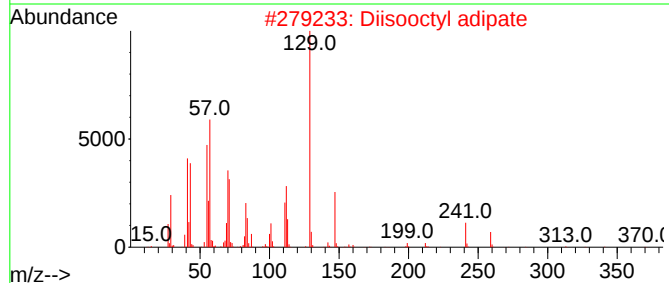
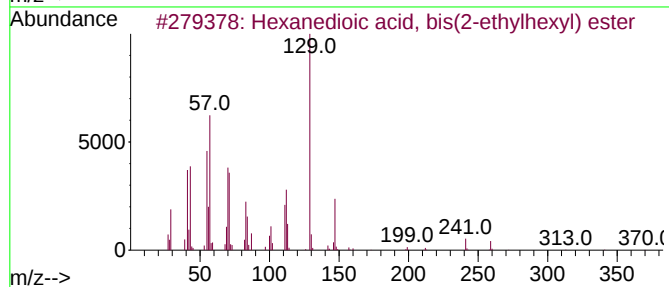
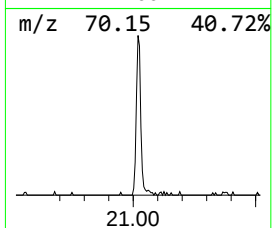
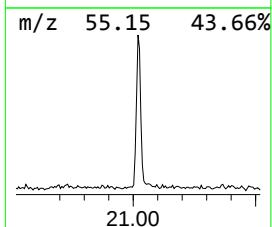
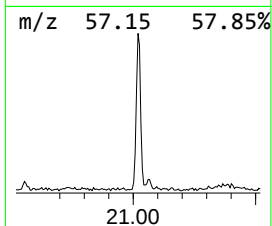
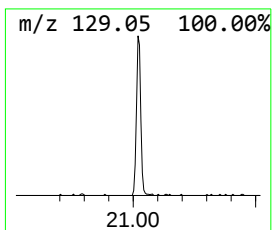
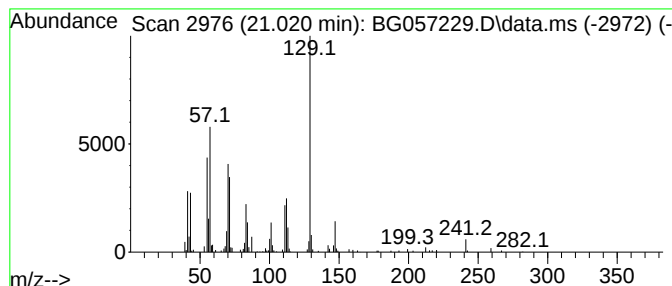
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 25 Hexanedioic acid, bis(2-eth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.020	10.72 ng/ul	177191	Chrysene-d12	22.036	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	94
2	Diisooctyl adipate	370	C22H42O4	001330-86-5	90
3	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	83
4	Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	83
5	Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
Operator : CG/JU
Sample : 02417-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW8W7

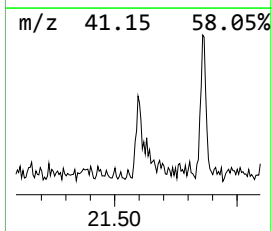
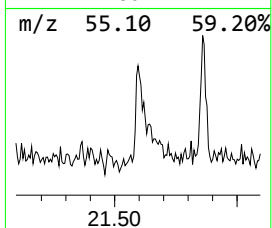
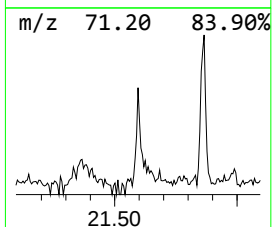
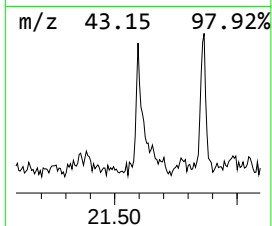
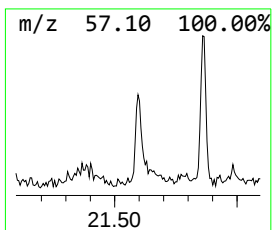
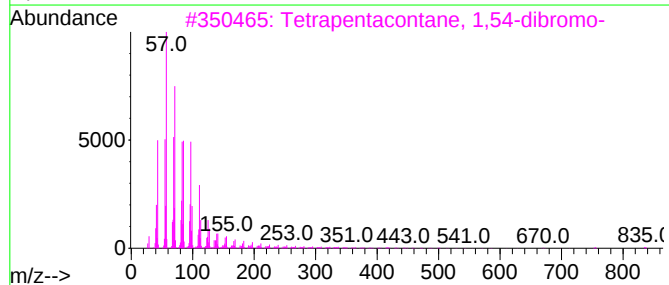
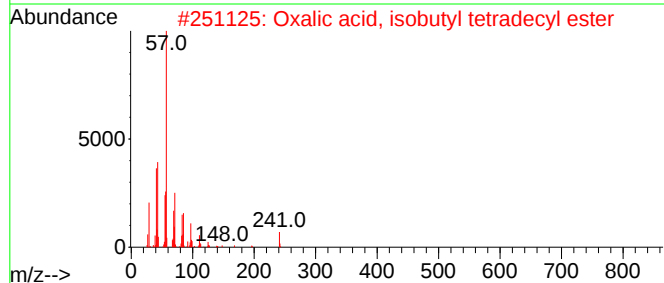
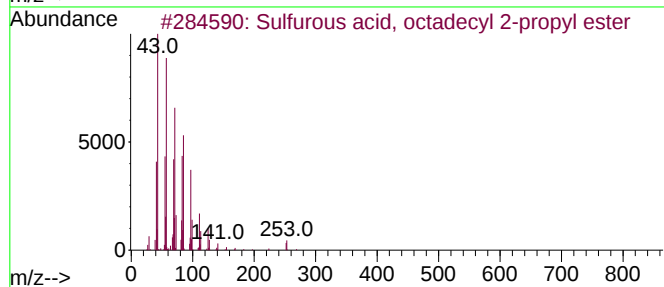
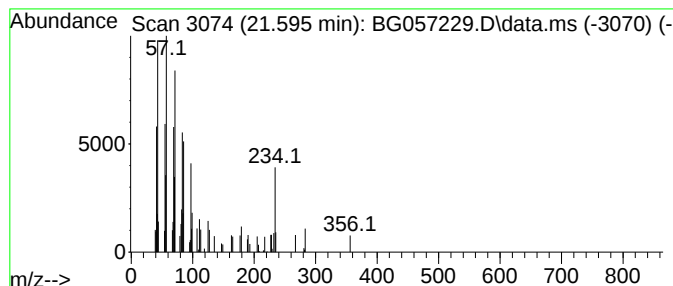
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 Sulfurous acid, octadecyl 2... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.595	2.97 ng/ul	49084	Chrysene-d12	22.036

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	58
2			Oxalic acid, isobutyl tetradecyl...	342	C20H38O4	1000309-37-9	58
3			Tetrapentacontane, 1,54-dibromo-	915	C54H108Br2	1000156-09-4	58
4			Ethanol, 2-(octadecyloxy)-	314	C20H42O2	002136-72-3	52
5			Carbonic acid, octadecyl vinyl e...	340	C21H40O3	1000382-54-4	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057229.D
Acq On : 28 Apr 2023 22:55
Operator : CG/JU
Sample : 02417-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW8W7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-ethy...	8.014	3.8	ng/ul	15293	1	8.378	81601	20.0
Benzene, 1,2,3-...	8.489	4.7	ng/ul	19197	1	8.378	81601	20.0
Benzaldehyde, 2...	8.889	9.6	ng/ul	39125	1	8.378	81601	20.0
Benzene, 4-ethy...	9.006	11.1	ng/ul	45308	1	8.378	81601	20.0
1,3,5-Cyclohept...	9.329	5.4	ng/ul	21983	1	8.378	81601	20.0
o-Cymene	9.382	4.3	ng/ul	17512	1	8.378	81601	20.0
Benzene, 1-ethy...	9.482	8.7	ng/ul	35570	1	8.378	81601	20.0
unknown-01	9.741	3.2	ng/ul	13190	1	8.378	81601	20.0
p-Cymene	10.017	5.5	ng/ul	45727	2	11.215	165332	20.0
Benzene, 1,2,4,...	10.076	13.3	ng/ul	110131	2	11.215	165332	20.0
2-Coumaranone	10.381	9.3	ng/ul	77308	2	11.215	165332	20.0
Ethanone, 1-(4-...	10.598	6.2	ng/ul	51009	2	11.215	165332	20.0
Benzene, 1,3-di...	10.657	2.4	ng/ul	19680	2	11.215	165332	20.0
Benzene, 1-meth...	10.886	2.0	ng/ul	16639	2	11.215	165332	20.0
Benzene, pentam...	11.409	2.4	ng/ul	19784	2	11.215	165332	20.0
Hexanedioic aci...	11.703	8.6	ng/ul	70904	2	11.215	165332	20.0
Benzene, 2,4-di...	12.502	2.8	ng/ul	22965	2	11.215	165332	20.0
3-Amino-4-methy...	13.348	5.0	ng/ul	52953	3	14.986	212841	20.0
Methyl salicylate	14.904	2.2	ng/ul	23466	3	14.986	212841	20.0
2,2,4-Trimethyl...	15.738	9.3	ng/ul	98388	3	14.986	212841	20.0
Benzophenone	16.326	3.0	ng/ul	32013	3	14.986	212841	20.0
Pyridine, 4-(3-...	17.454	11.7	ng/ul	159335	4	17.724	271659	20.0
Pyridine-4-carb...	18.270	2.3	ng/ul	30915	4	17.724	271659	20.0
n-Hexadecanoic ...	18.564	3.0	ng/ul	40133	4	17.724	271659	20.0
Hexanedioic aci...	21.020	10.7	ng/ul	177191	5	22.036	330731	20.0
Sulfurous acid,...	21.595	3.0	ng/ul	49084	5	22.036	330731	20.0