

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
 Data File : BG051349.D
 Acq On : 6 Dec 2021 14:04
 Operator : CG/JU
 Sample : M4942-03 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGKQ1

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-BG112321.M
 Title : SVOA CALIBRATION

Signal : TIC: BG051349.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.656	364	369	376	rBV	901489	1570242	3.23%	0.446%
2	5.797	385	393	409	rBV	2361561	5434611	11.18%	1.542%
3	6.167	449	456	465	rVB2	1370551	3115524	6.41%	0.884%
4	6.649	526	538	543	rBV2	326456	866112	1.78%	0.246%
5	7.143	618	622	628	rVV2	831772	1370419	2.82%	0.389%
6	7.260	636	642	647	rBV	1129171	1911319	3.93%	0.542%
7	7.319	647	652	657	rVB2	1077391	1846446	3.80%	0.524%
8	7.395	657	665	672	rVB2	1898394	3666819	7.55%	1.040%
9	7.566	687	694	699	rBV	484537	986617	2.03%	0.280%
10	7.789	725	732	735	rBV	2934000	4642288	9.55%	1.317%
11	7.836	735	740	746	rVB	2105720	3745170	7.71%	1.063%
12	8.077	771	781	786	rBV4	313130	897065	1.85%	0.255%
13	8.142	786	792	797	rVB	952832	1592728	3.28%	0.452%
14	8.236	797	808	814	rBV4	723351	1656079	3.41%	0.470%
15	8.306	814	820	824	rBV2	758199	1421153	2.92%	0.403%
16	8.694	880	886	890	rVV3	605330	1378755	2.84%	0.391%
17	8.753	890	896	902	rVB2	778431	1976093	4.07%	0.561%
18	8.829	902	909	920	rBV2	1662037	3227350	6.64%	0.916%
19	8.935	920	927	931	rBV	725595	1290277	2.66%	0.366%
20	9.293	977	988	997	rBV2	819524	1821169	3.75%	0.517%
21	9.411	997	1008	1013	rBV	3820625	7179236	14.77%	2.037%
22	9.540	1018	1030	1038	rBV	463988	1692320	3.48%	0.480%
23	9.640	1041	1047	1059	rVV5	407700	1240603	2.55%	0.352%
24	9.822	1072	1078	1084	rVV2	537524	1051656	2.16%	0.298%
25	9.887	1084	1089	1093	rVV2	491729	890542	1.83%	0.253%
26	10.110	1123	1127	1135	rVV4	416297	864332	1.78%	0.245%
27	10.186	1135	1140	1144	rVV3	400087	824995	1.70%	0.234%
28	10.251	1144	1151	1157	rVB2	456737	1144862	2.36%	0.325%
29	10.404	1171	1177	1182	rVV2	804210	1442089	2.97%	0.409%
30	10.956	1265	1271	1277	rBV	2399996	3977722	8.19%	1.129%
31	11.097	1284	1295	1299	rBV	4333459	9756808	20.08%	2.769%
32	11.138	1299	1302	1308	rVB	712250	1066707	2.20%	0.303%
33	11.996	1440	1448	1453	rBV	777341	1418759	2.92%	0.403%
34	12.395	1509	1516	1524	rBV	2587476	3971577	8.17%	1.127%
35	12.677	1557	1564	1570	rVB	2998357	5270112	10.84%	1.495%
36	12.889	1592	1600	1607	rVB	1544110	2824072	5.81%	0.801%

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Integrator: RTE

Smoothing : OFF

Filtering: 5

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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-BG112321.M
 Title : SVOA CALIBRATION

37	13.277	1660	1666	1671	rBV	969356	1600396	3.29%	0.454%
38	13.330	1671	1675	1680	rVB	697686	959958	1.98%	0.272%
39	13.623	1717	1725	1729	rBV	3491793	5557703	11.44%	1.577%
40	13.858	1760	1765	1777	rVB	433559	823373	1.69%	0.234%
41	14.005	1782	1790	1796	rVB2	795849	2044739	4.21%	0.580%
42	14.152	1808	1815	1817	rBV3	1064428	2106869	4.34%	0.598%
43	14.270	1832	1835	1839	rVB2	943561	1214336	2.50%	0.345%
44	14.387	1850	1855	1859	rBV	630774	914367	1.88%	0.259%
45	14.693	1901	1907	1911	rVB	3065014	4288659	8.82%	1.217%
46	14.787	1915	1923	1927	rBV6	354723	917054	1.89%	0.260%
47	15.028	1959	1964	1971	rBV2	418242	985254	2.03%	0.280%
48	15.221	1993	1997	2006	rVV4	789183	2347760	4.83%	0.666%
49	15.298	2006	2010	2018	rVB3	978630	1602487	3.30%	0.455%
50	15.439	2029	2034	2039	rBV	434012	931695	1.92%	0.264%
51	15.598	2055	2061	2063	rBV4	743494	1469837	3.02%	0.417%
52	15.639	2063	2068	2072	rVB	3018395	4422118	9.10%	1.255%
53	15.744	2081	2086	2090	rBV2	598401	964230	1.98%	0.274%
54	16.038	2132	2136	2140	rBV	827207	1229004	2.53%	0.349%
55	16.079	2140	2143	2148	rVB3	736393	1244668	2.56%	0.353%
56	16.502	2210	2215	2217	rBV	2845897	4245508	8.74%	1.205%
57	16.632	2232	2237	2241	rBV2	1017012	1648636	3.39%	0.468%
58	16.778	2256	2262	2265	rBV2	448510	927186	1.91%	0.263%
59	17.296	2345	2350	2354	rBV	2392094	3133445	6.45%	0.889%
60	17.343	2354	2358	2362	rVB	1050216	1464898	3.01%	0.416%
61	17.448	2372	2376	2381	rVB	690273	937490	1.93%	0.266%
62	17.830	2438	2441	2449	rVB	791616	1176307	2.42%	0.334%
63	18.036	2471	2476	2481	rVB	2283329	3083819	6.35%	0.875%
64	18.500	2545	2555	2565	rBV3	2819989	7896591	16.25%	2.241%
65	18.723	2589	2593	2597	rVB	1611094	1940208	3.99%	0.551%
66	18.999	2637	2640	2648	rVB2	878520	1288326	2.65%	0.366%
67	19.346	2694	2699	2708	rBV2	2661972	3625527	7.46%	1.029%
68	19.546	2729	2733	2738	rBV2	1238223	1603174	3.30%	0.455%
69	19.634	2741	2748	2754	rVB2	1356692	3154922	6.49%	0.895%
70	19.781	2770	2773	2776	rVB	1185995	1339026	2.76%	0.380%
71	19.828	2778	2781	2785	rVB	779560	946953	1.95%	0.269%
72	19.916	2793	2796	2799	rBV	1181129	1346453	2.77%	0.382%
73	19.951	2799	2802	2809	rVB2	1438720	2404864	4.95%	0.682%
74	20.227	2845	2849	2852	rBV	1323579	1681376	3.46%	0.477%
75	20.445	2882	2886	2891	rBV	2614623	2990356	6.15%	0.849%
76	20.909	2950	2965	2968	rBV3	12079481	48596940	100.00%	13.790%

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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

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77	21.214	3013	3017	3022	rVB2	2365710	3301074	6.79%	0.937%
78	21.367	3040	3043	3051	rVB	860011	1237260	2.55%	0.351%
79	21.473	3056	3061	3066	rVB2	4198039	6138432	12.63%	1.742%
80	21.767	3099	3111	3116	rVB3	12011275	47436710	97.61%	13.460%
81	21.873	3124	3129	3133	rBV2	1257172	2108511	4.34%	0.598%
82	22.037	3152	3157	3163	rBV2	2959428	4981175	10.25%	1.413%
83	22.090	3164	3166	3176	rVB2	687537	1067154	2.20%	0.303%
84	22.454	3226	3228	3232	rVB	753969	963881	1.98%	0.274%
85	22.548	3233	3244	3250	rVB5	4436612	11106178	22.85%	3.151%
86	22.677	3260	3266	3277	rVB2	2692273	5112544	10.52%	1.451%
87	22.918	3303	3307	3311	rBV	603280	1139087	2.34%	0.323%
88	23.018	3317	3324	3337	rVB	2753157	7714361	15.87%	2.189%
89	23.171	3345	3350	3358	rBV5	901741	1962199	4.04%	0.557%
90	23.400	3382	3389	3398	rVB2	2010320	4241367	8.73%	1.203%
91	23.541	3409	3413	3424	rVB	589095	1360756	2.80%	0.386%
92	23.900	3468	3474	3485	rBV	984116	2341631	4.82%	0.664%
93	24.252	3527	3534	3541	rBV	1262358	2755960	5.67%	0.782%
94	24.669	3598	3605	3609	rBV	1325595	3591011	7.39%	1.019%
95	24.934	3643	3650	3666	rVB3	1612214	4993307	10.27%	1.417%
96	25.245	3696	3703	3711	rVB3	1028333	2711963	5.58%	0.770%
97	26.438	3898	3906	3918	rVB2	740858	2382286	4.90%	0.676%
98	26.949	3982	3993	4005	rVB	1136673	4108192	8.45%	1.166%
99	27.155	4014	4028	4040	rBV4	1522351	6399004	13.17%	1.816%
100	27.395	4060	4069	4087	rVB7	723123	3176320	6.54%	0.901%

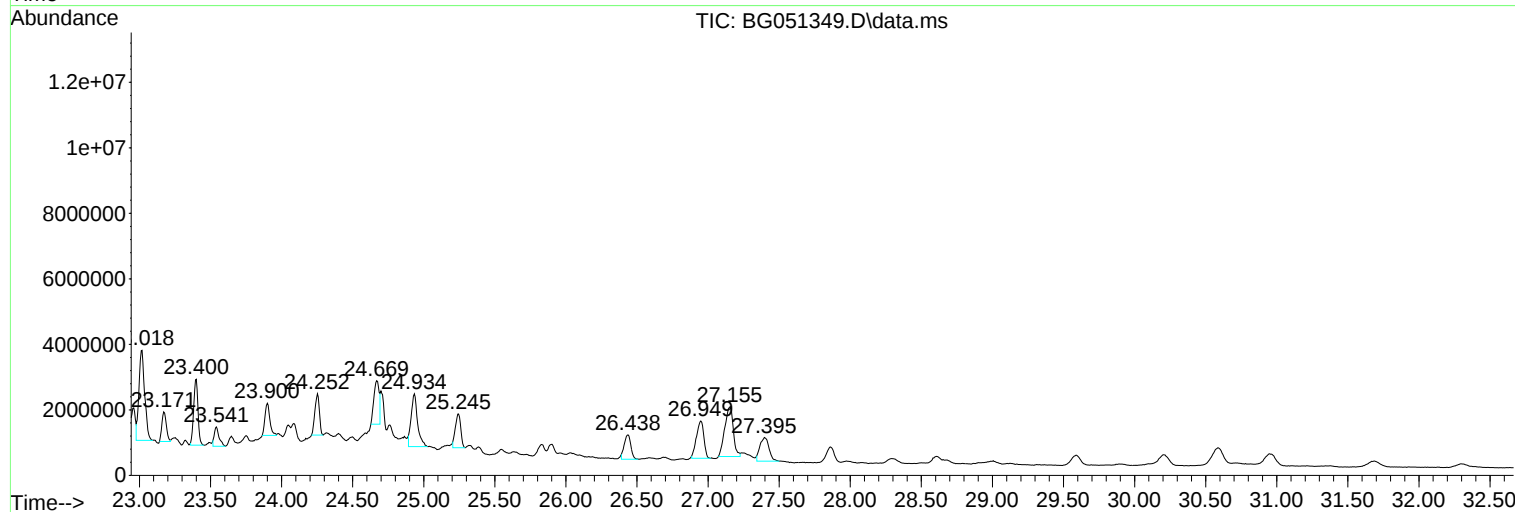
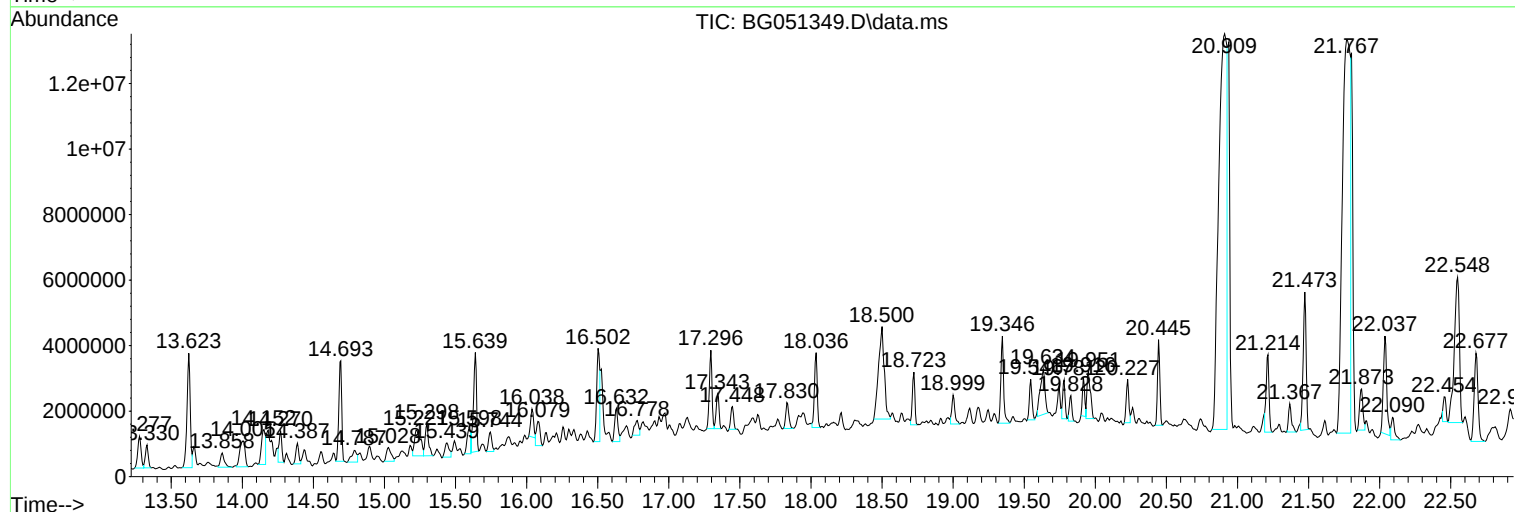
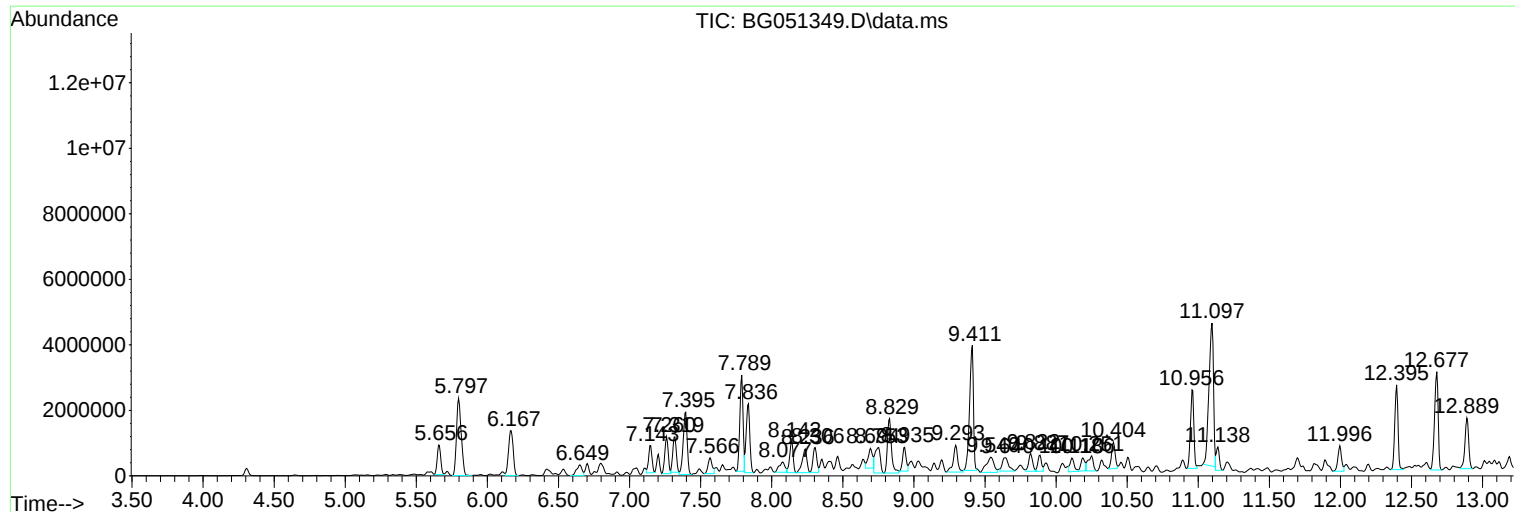
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Instrument :
BNA_G
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

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



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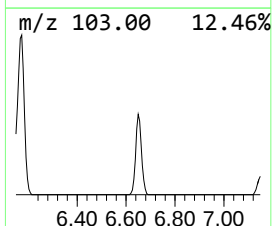
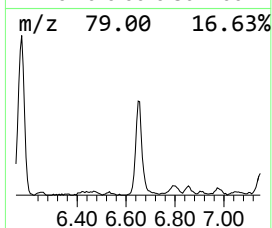
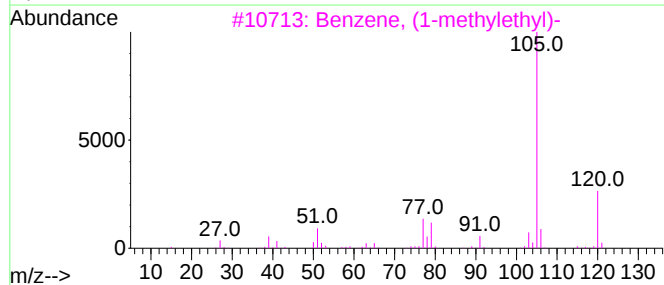
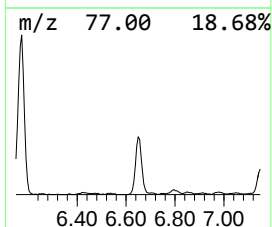
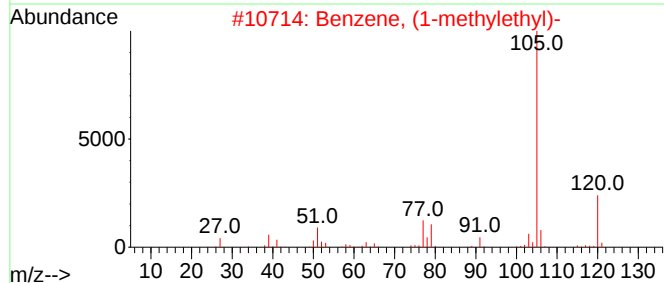
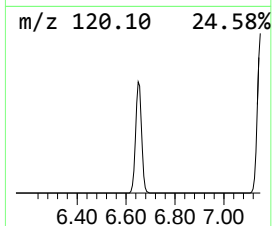
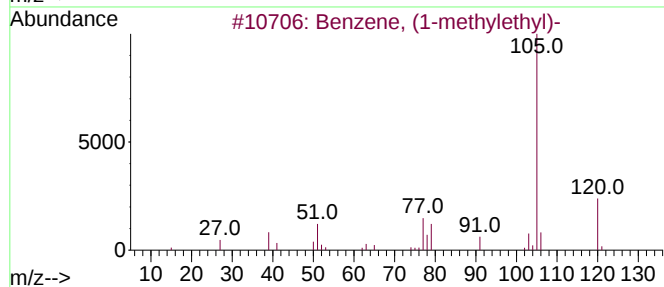
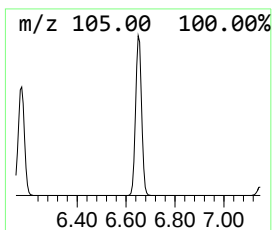
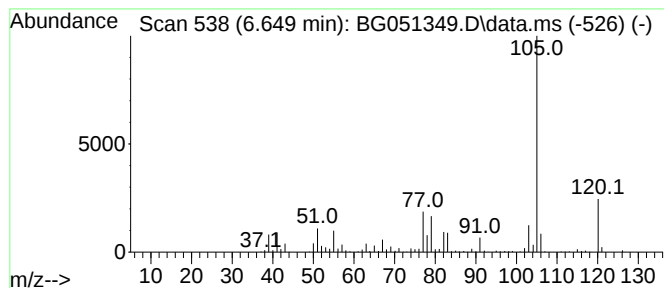
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, (1-methylethyl)- Concentration Rank 64

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.649	10.46 ng/ul	866112	1,4-Dichlorobenzene-d4	8.194

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	93
2			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	90
3			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	90
4			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	87
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	87



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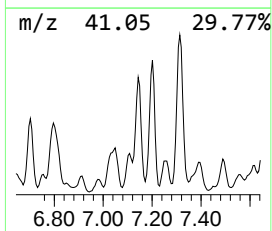
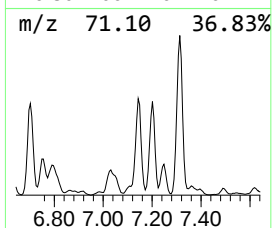
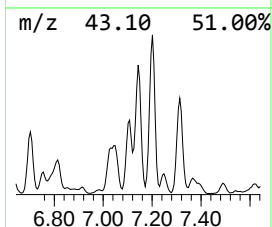
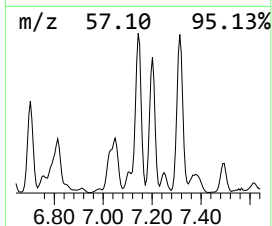
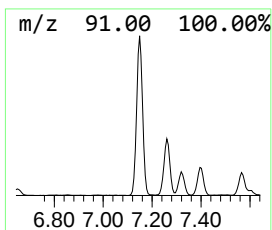
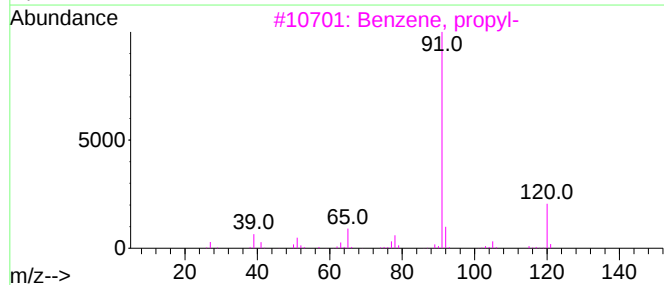
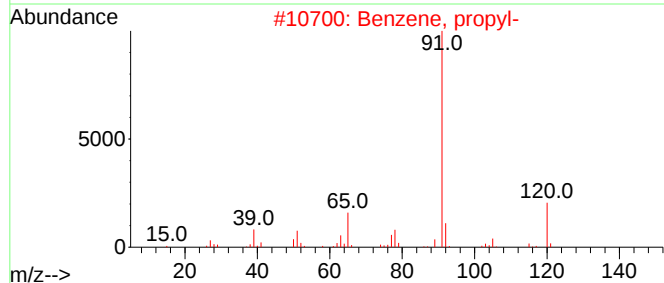
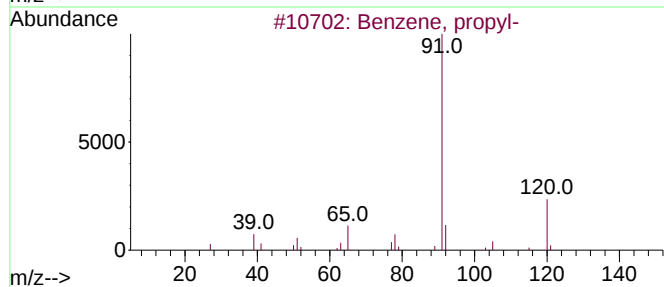
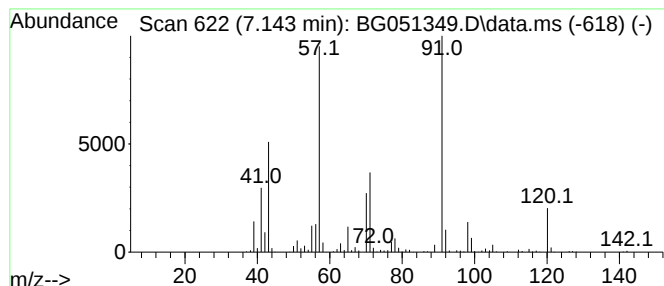
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, propyl- Concentration Rank 55

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.143	16.55 ng/ul	1370420	1,4-Dichlorobenzene-d4	8.194

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	83
2			Benzene, propyl-	120	C9H12	000103-65-1	83
3			Benzene, propyl-	120	C9H12	000103-65-1	46
4			Benzene, propyl-	120	C9H12	000103-65-1	41
5			N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	35



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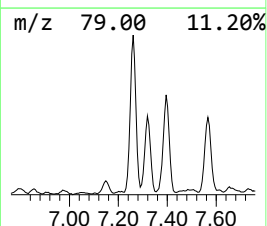
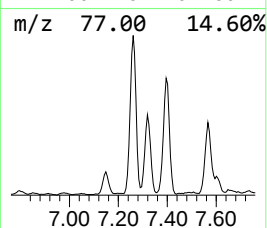
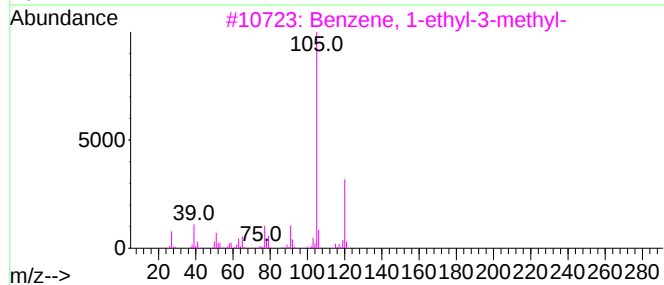
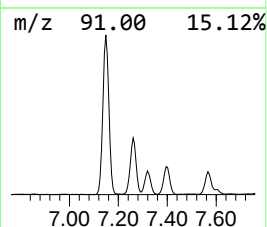
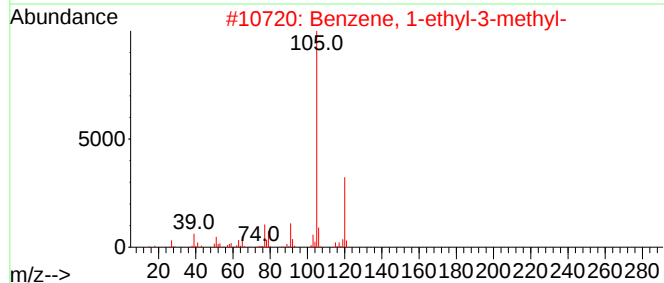
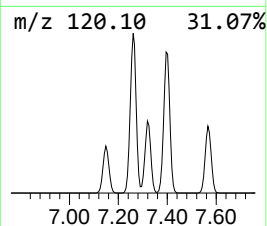
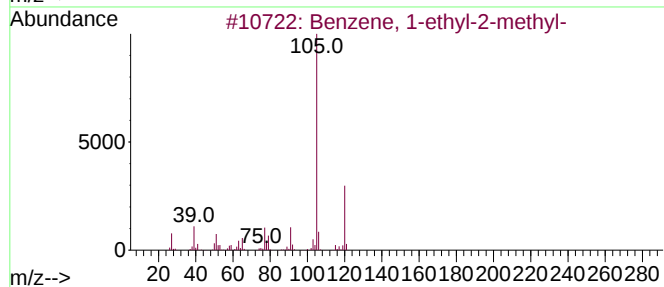
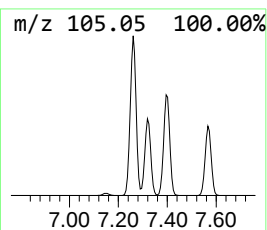
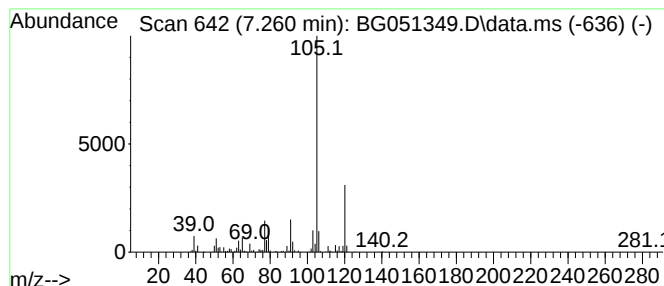
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1-ethyl-2-methyl- Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.260	23.08 ng/ul	1911320	1,4-Dichlorobenzene-d4	8.194

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

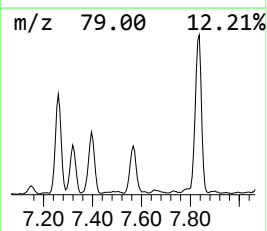
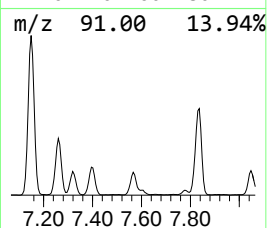
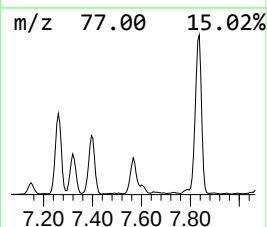
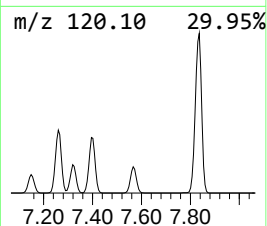
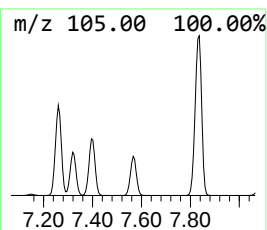
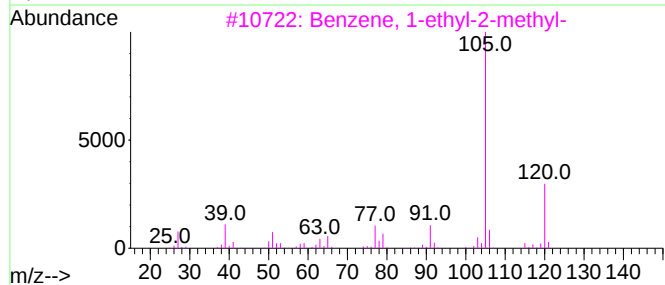
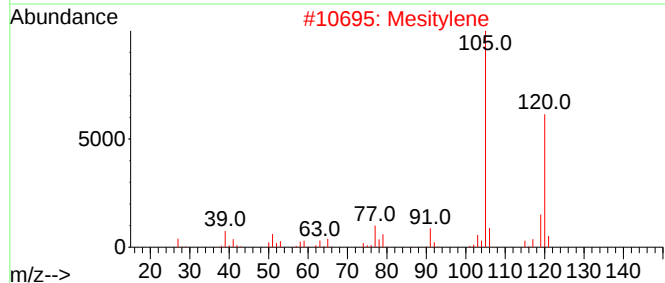
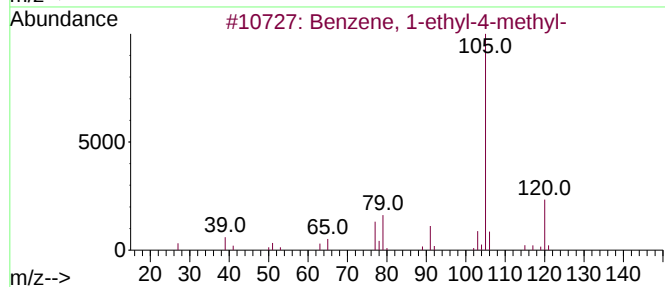
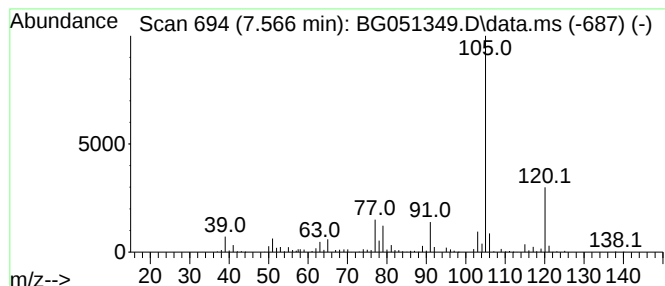
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1-ethyl-4-methyl- Concentration Rank 60

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.566	11.92 ng/ul	986617	1,4-Dichlorobenzene-d4	8.194

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
2			Mesitylene	120	C9H12	000108-67-8	91
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
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Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

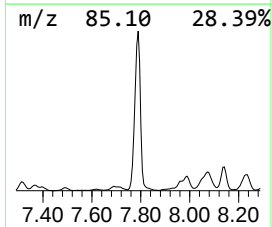
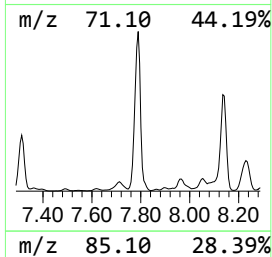
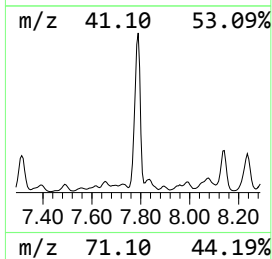
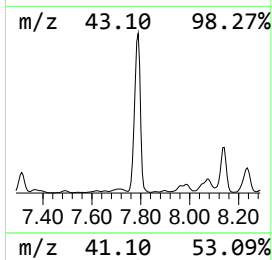
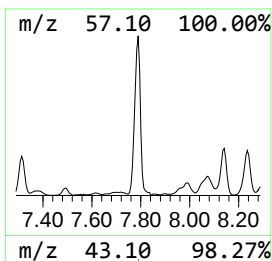
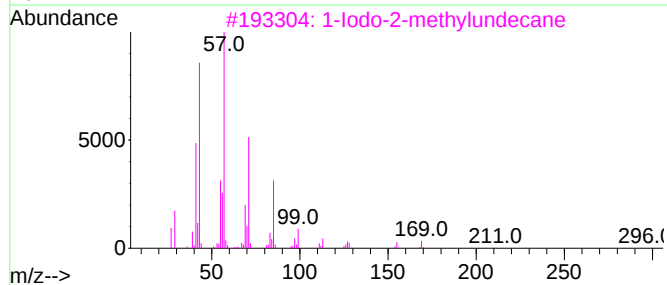
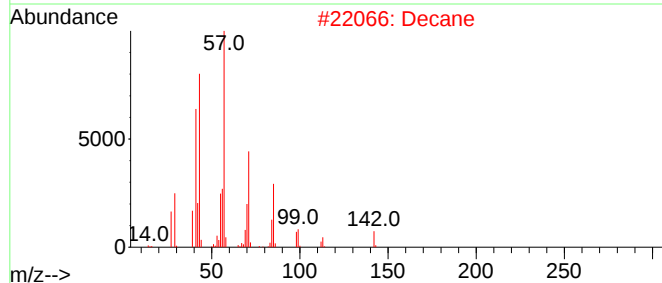
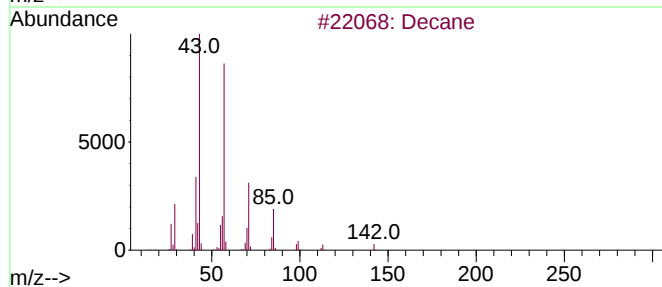
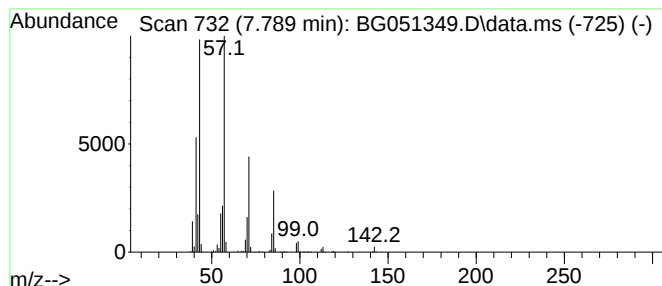
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD			R.T.
7.789	56.06 ng/ul	4642290	1,4-Dichlorobenzene-d4			8.194
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Decane		142	C10H22	000124-18-5	90
2	Decane		142	C10H22	000124-18-5	90
3	1-Iodo-2-methylundecane		296	C12H25I	073105-67-6	83
4	Tridecane		184	C13H28	000629-50-5	72
5	Decane		142	C10H22	000124-18-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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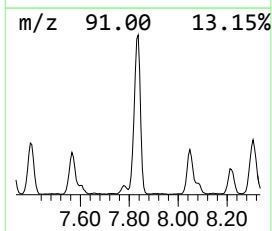
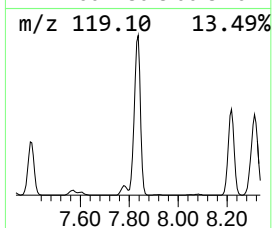
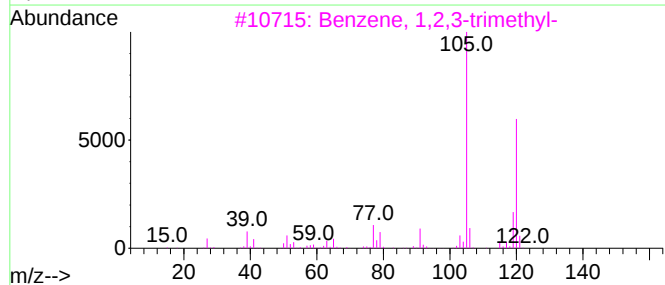
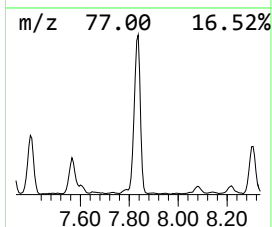
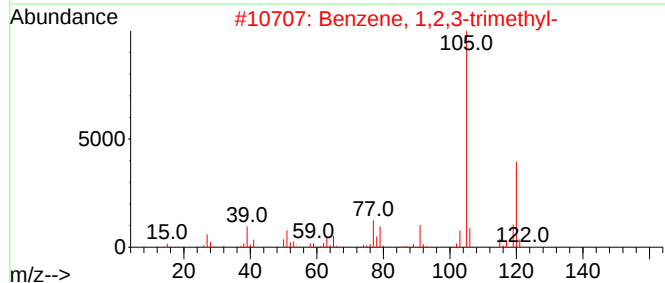
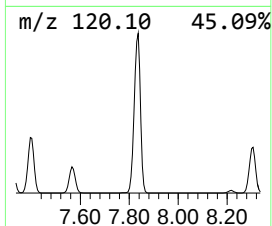
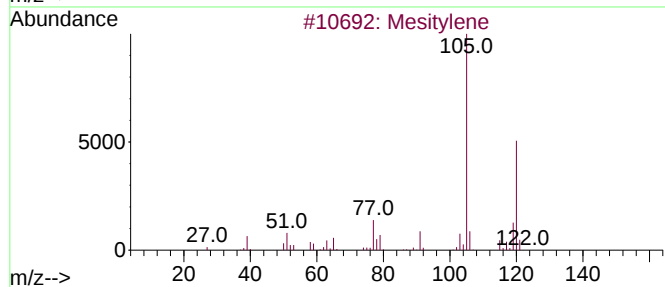
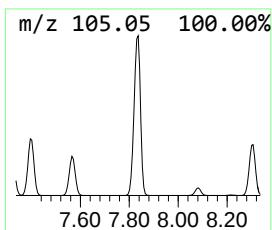
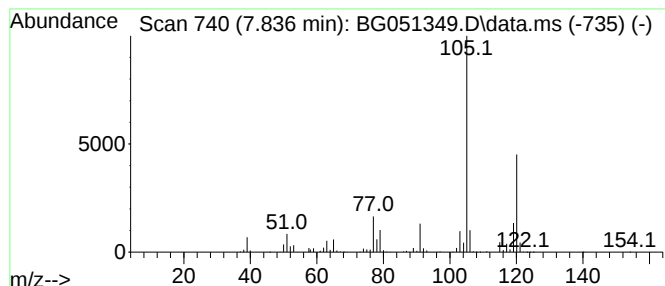
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Mesitylene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.836	45.23 ng/ul	3745170	1,4-Dichlorobenzene-d4	8.194

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5			Mesitylene	120	C9H12	000108-67-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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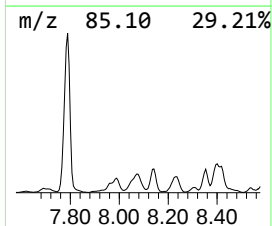
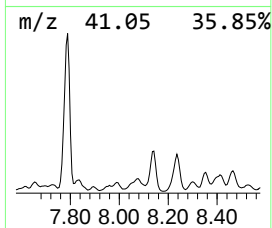
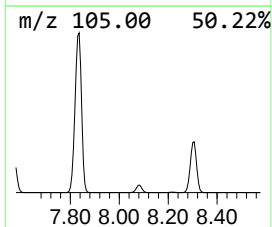
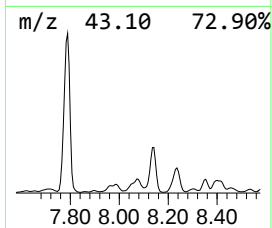
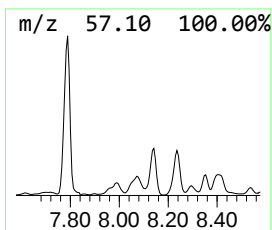
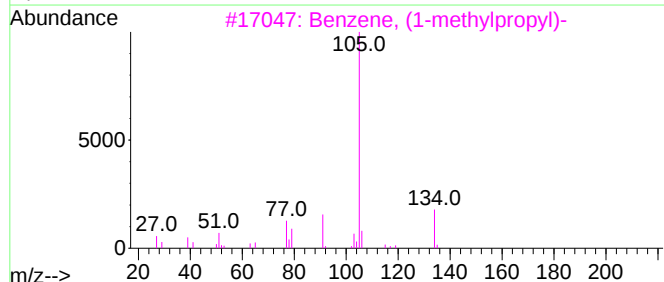
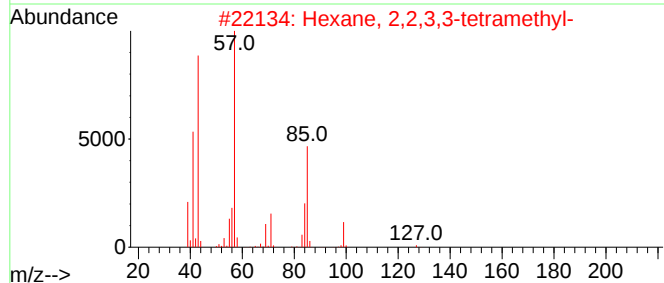
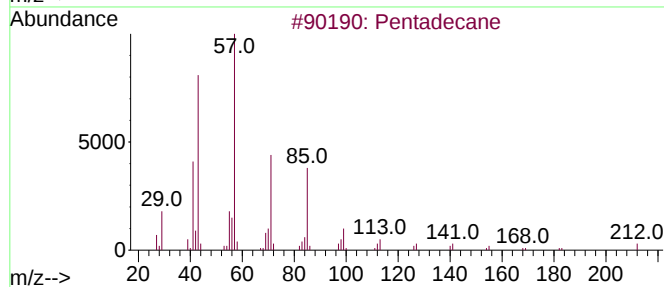
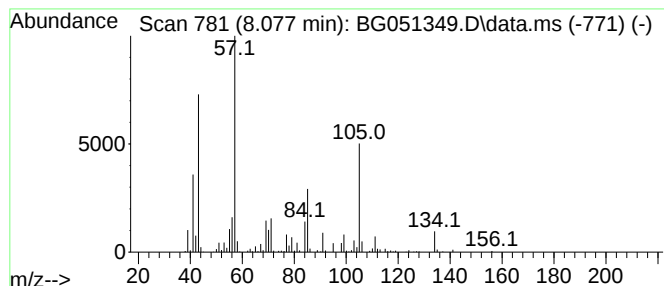
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 62

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.077	10.83 ng/ul	897065	1,4-Dichlorobenzene-d4	8.194

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane	212	C15H32	000629-62-9	43
2			Hexane, 2,2,3,3-tetramethyl-	142	C10H22	013475-81-5	35
3			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	35
4			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	30
5			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
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Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

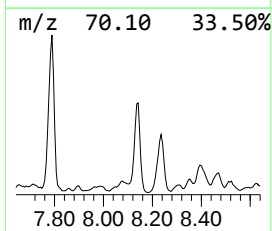
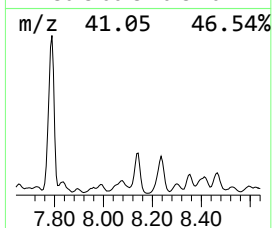
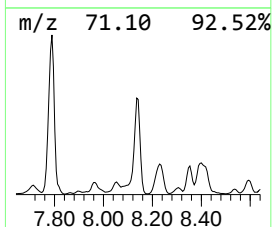
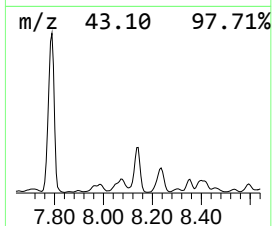
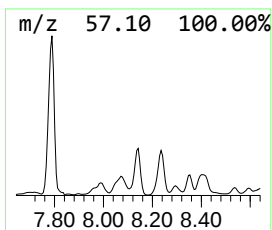
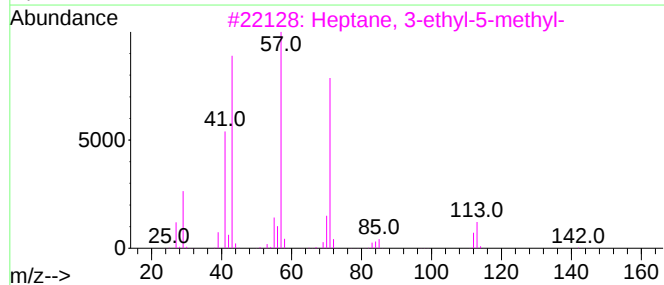
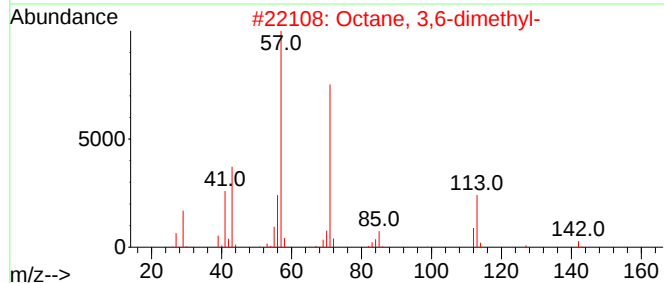
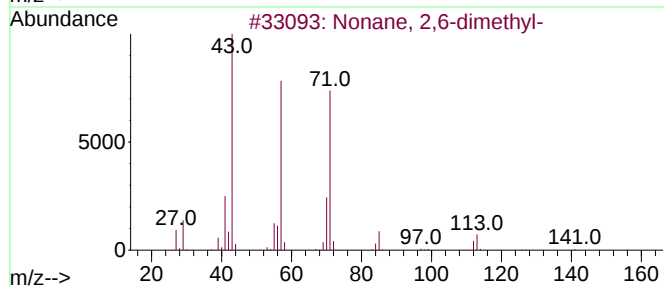
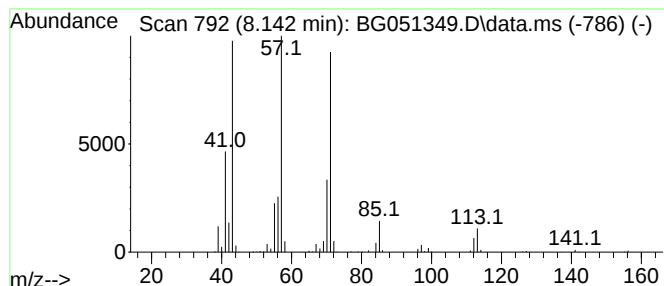
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.142	19.23 ng/ul	1592730	1,4-Dichlorobenzene-d4	8.194

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	86
2		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	78
3		Heptane, 3-ethyl-5-methyl-	142	C10H22	052896-90-9	64
4		Sulfurous acid, octyl 2-propyl e...	236	C11H24O3S	1000309-11-9	59
5		Heptane, 2,5,5-trimethyl-	142	C10H22	001189-99-7	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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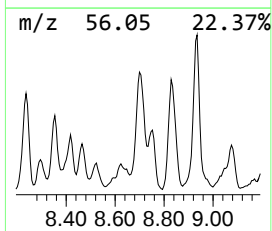
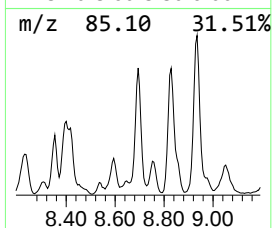
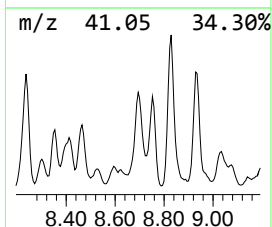
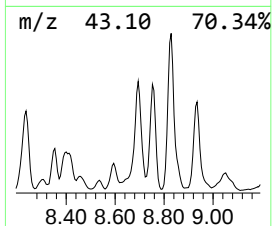
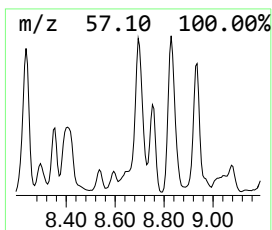
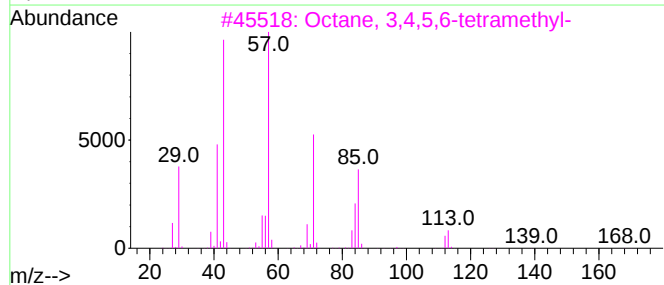
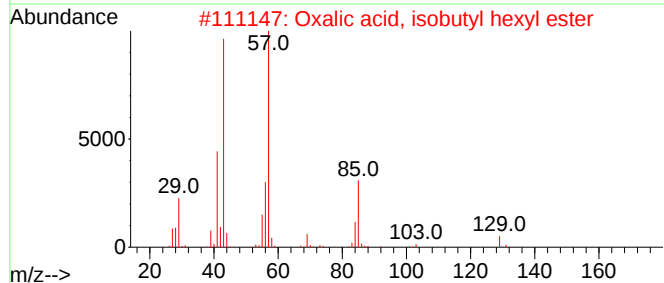
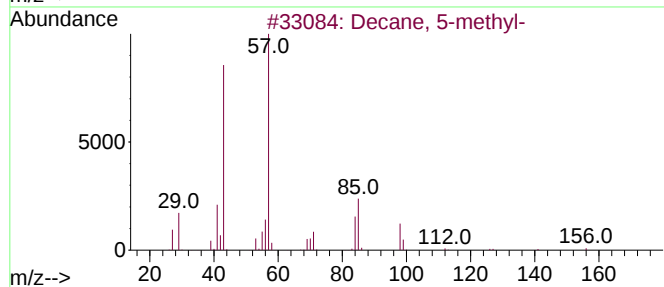
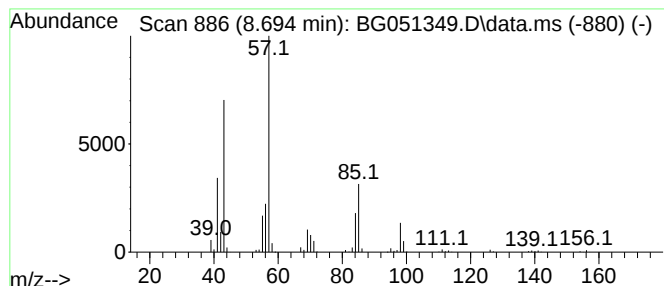
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 54

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.694	16.65 ng/ul	1378760	1,4-Dichlorobenzene-d4	8.194

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane, 5-methyl-	156	C11H24	013151-35-4	81
2		Oxalic acid, isobutyl hexyl ester	230	C12H22O4	1000309-37-1	59
3		Octane, 3,4,5,6-tetramethyl-	170	C12H26	062185-21-1	53
4		Undecane, 5-methyl-	170	C12H26	001632-70-8	53
5		Sulfurous acid, butyl decyl ester	278	C14H30O3S	1000309-17-7	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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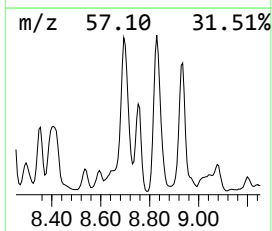
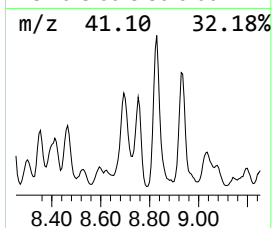
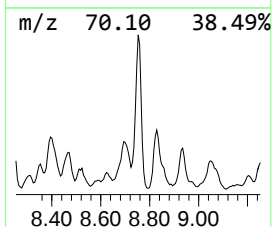
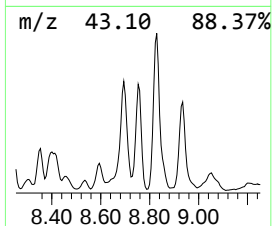
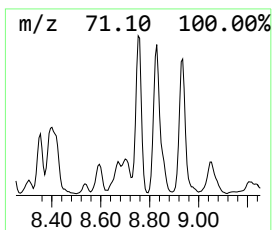
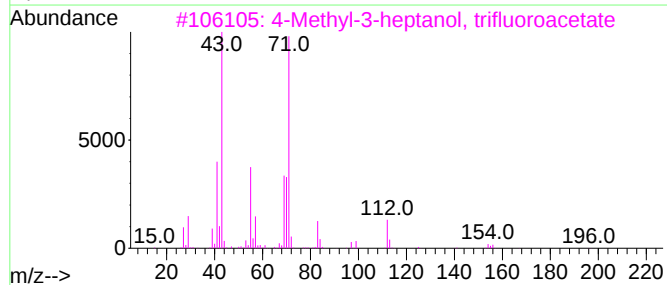
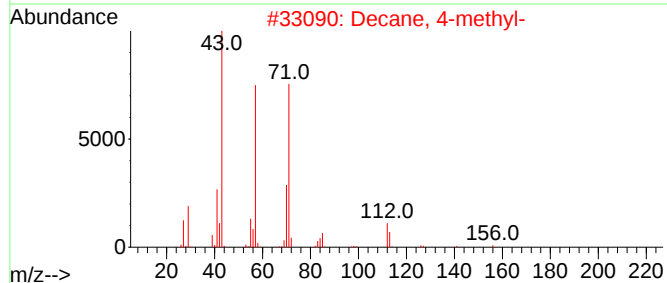
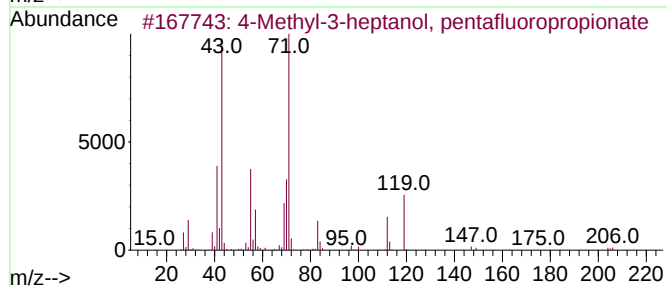
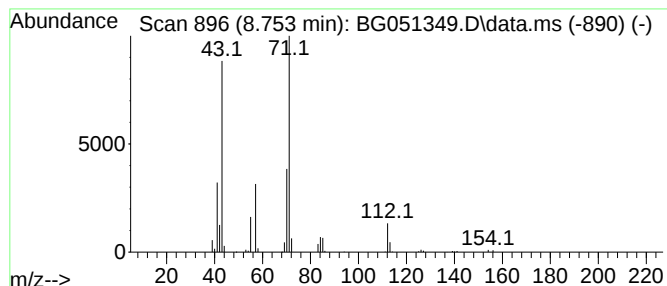
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 4-Methyl-3-heptanol, pentafl... Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD			R.T.
8.753	23.86 ng/ul	1976090	1,4-Dichlorobenzene-d4			8.194

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Methyl-3-heptanol, pentafluoro...	276	C11H17F5O2	1000365-51-7	83	
2	Decane, 4-methyl-	156	C11H24	002847-72-5	80	
3	4-Methyl-3-heptanol, trifluoroac...	226	C10H17F3O2	1000365-51-8	74	
4	Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	72	
5	2,4,4-Trimethyl-1-hexene	126	C9H18	051174-12-0	64	



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

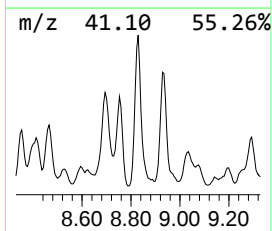
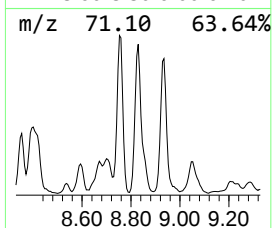
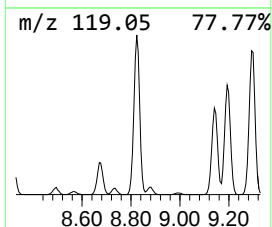
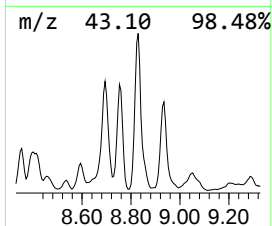
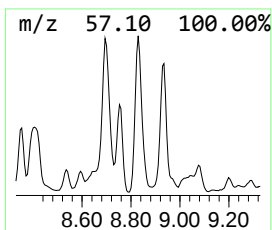
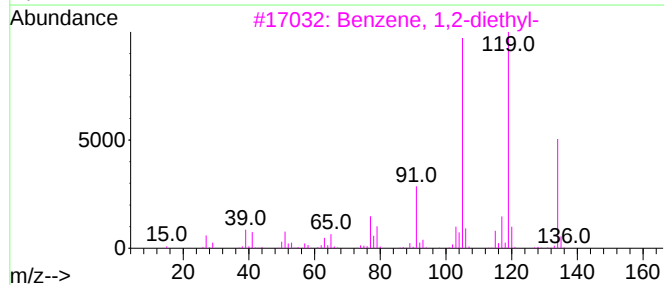
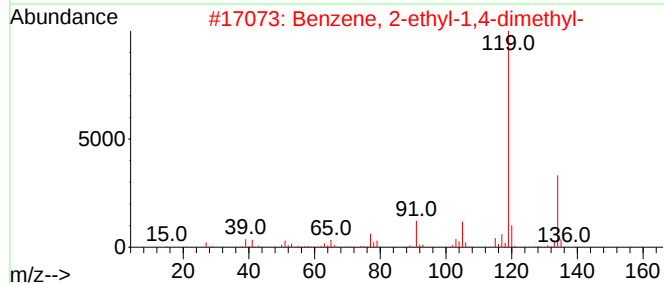
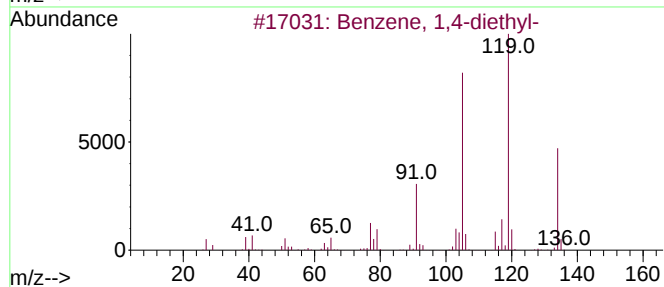
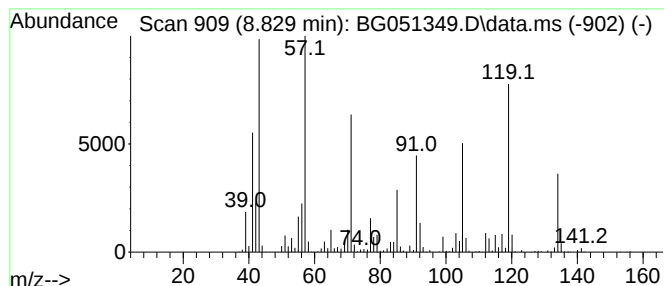
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Benzene, 1,4-diethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.829	38.98 ng/ul	3227350	1,4-Dichlorobenzene-d4	8.194

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	89
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	83
3			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	60
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	59
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

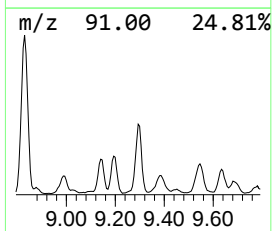
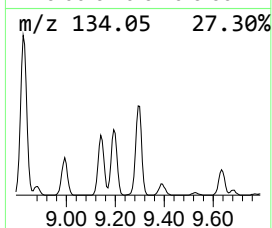
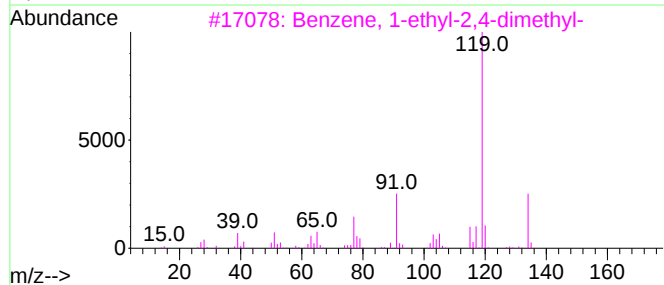
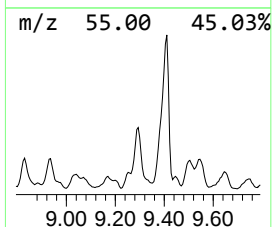
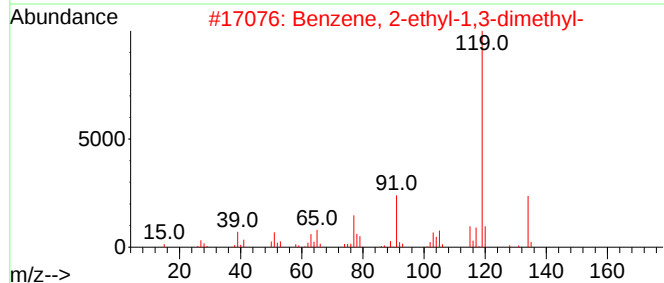
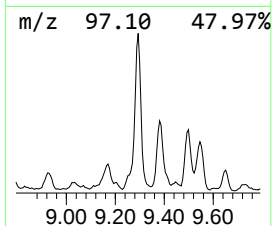
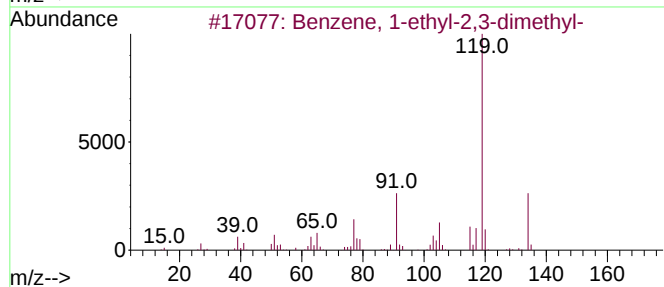
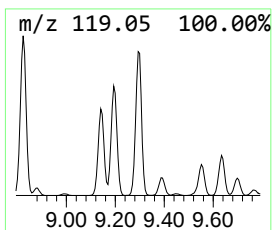
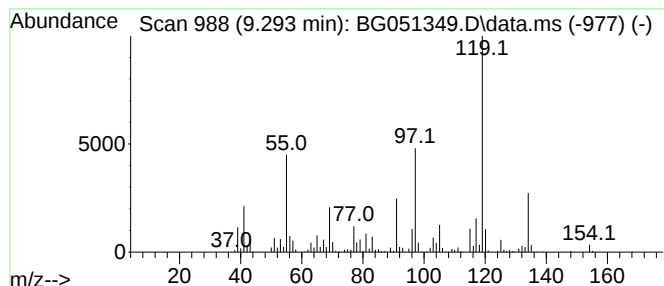
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.293	21.99 ng/ul	1821170	1,4-Dichlorobenzene-d4	8.194

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93
5			p-Cymene	134	C10H14	000099-87-6	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

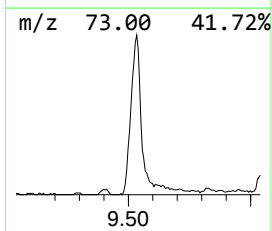
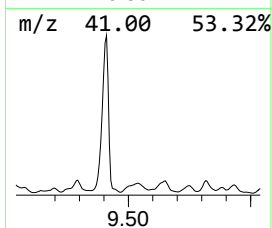
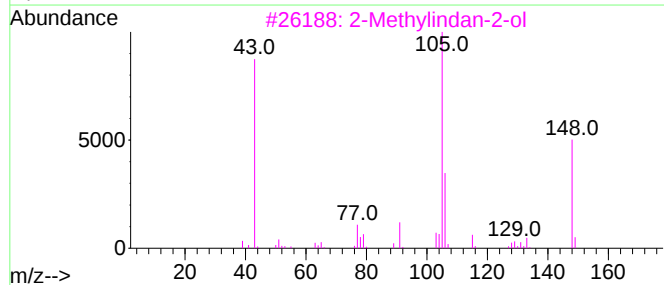
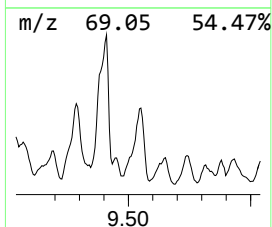
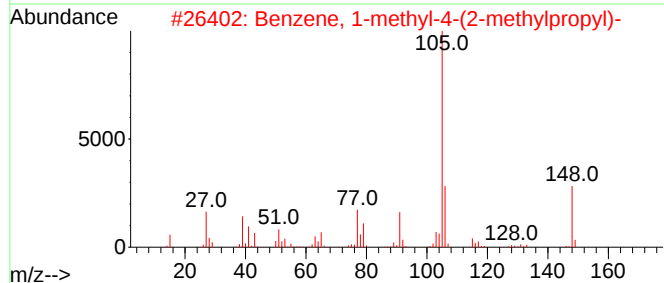
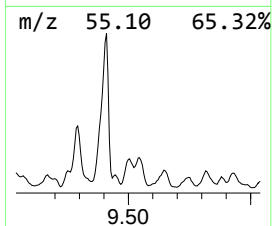
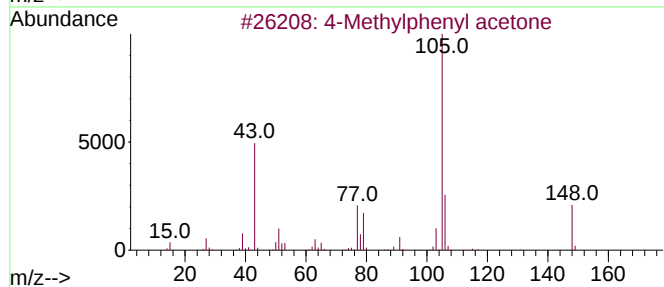
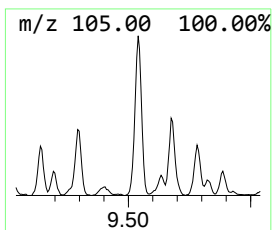
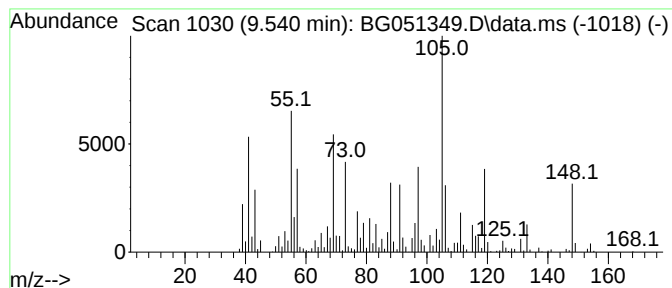
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 unknown-01 Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.540	20.44 ng/ul	1692320	1,4-Dichlorobenzene-d4	8.194

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		4-Methylphenyl acetone	148	C10H12O	002096-86-8	38
2		Benzene, 1-methyl-4-(2-methylpro...	148	C11H16	005161-04-6	30
3		2-Methylindan-2-ol	148	C10H12O	033223-84-6	22
4		3-Buten-2-ol, 4-phenyl-, (E)-	148	C10H12O	1000163-70-9	18
5		2-Butanone, 4-phenyl-	148	C10H12O	002550-26-7	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

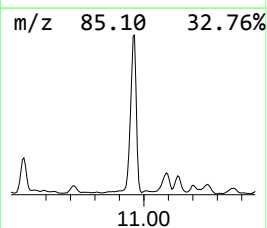
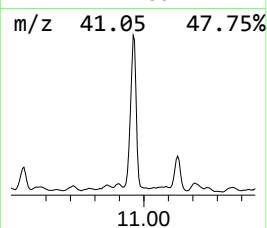
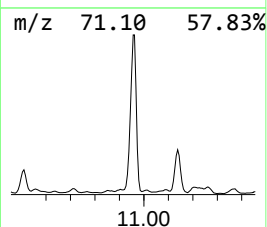
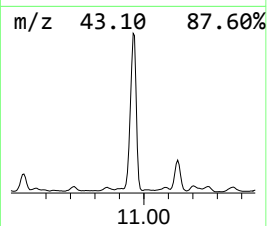
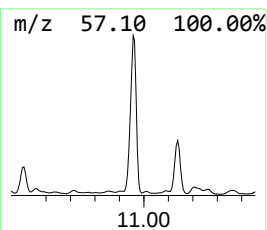
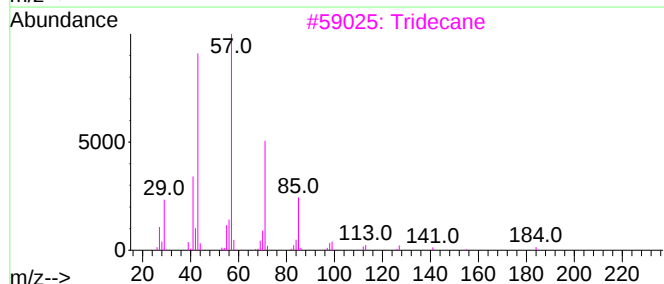
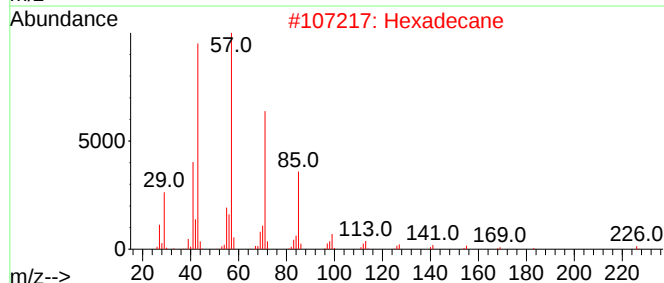
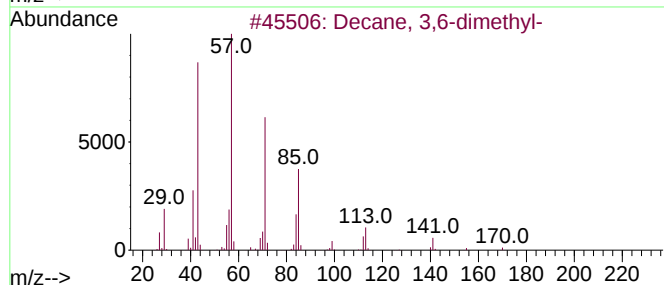
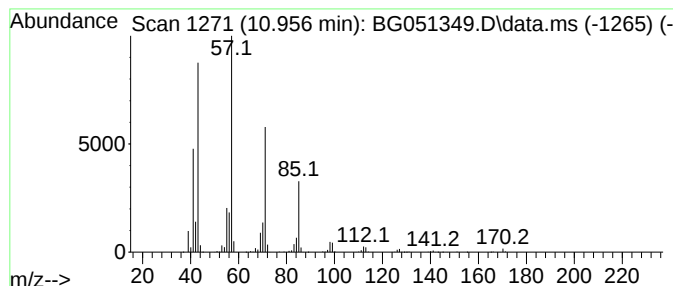
Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.956	20.00 ng/ul	3977720	Naphthalene-d8	11.026	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	87
2	Hexadecane	226	C16H34	000544-76-3	83
3	Tridecane	184	C13H28	000629-50-5	80
4	Undecane, 5-methyl-	170	C12H26	001632-70-8	80
5	Carbonic acid, dodecyl vinyl ester	256	C15H28O3	1000382-54-8	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

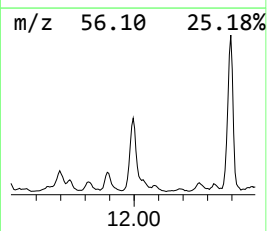
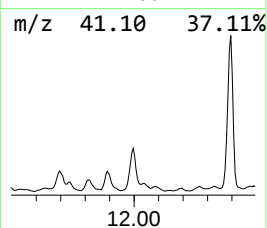
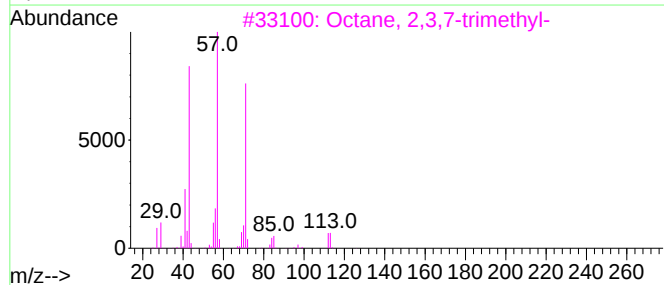
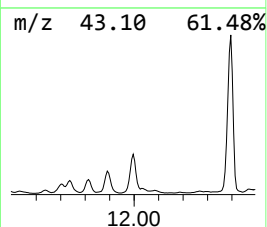
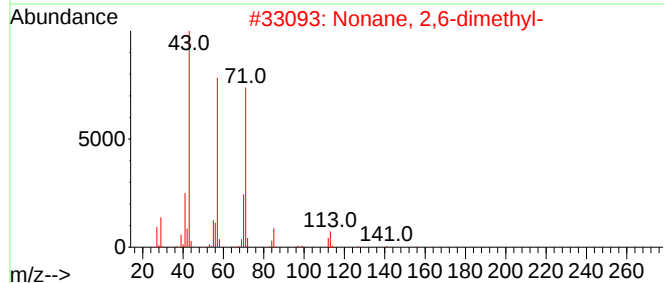
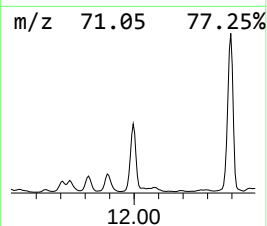
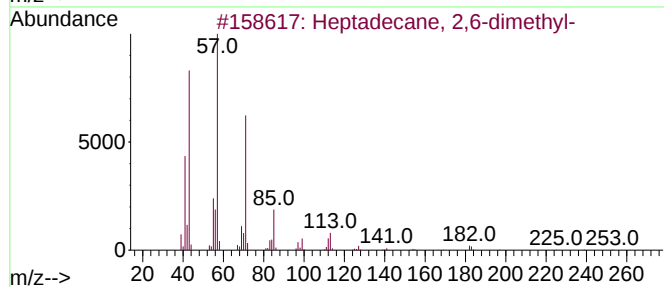
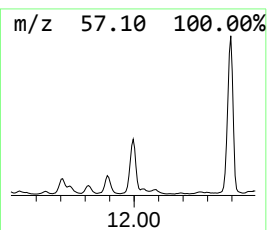
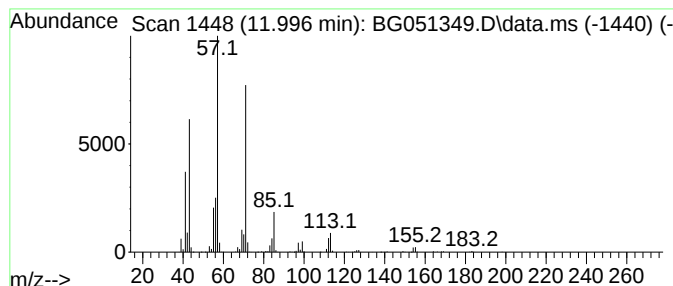
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 69

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.996	7.13 ng/ul	1418760	Naphthalene-d8	11.026

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	83
2		Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	80
3		Octane, 2,3,7-trimethyl-	156	C11H24	062016-34-6	72
4		Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	72
5		Oxalic acid, bis(2-ethylhexyl) e...	314	C18H34O4	1000309-39-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

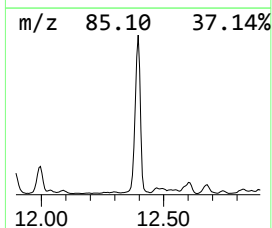
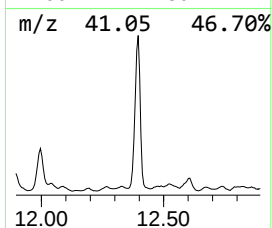
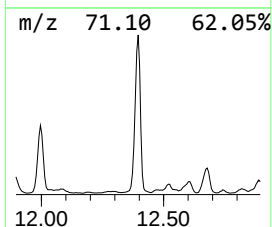
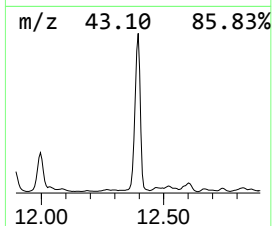
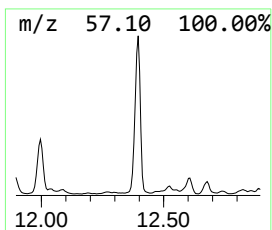
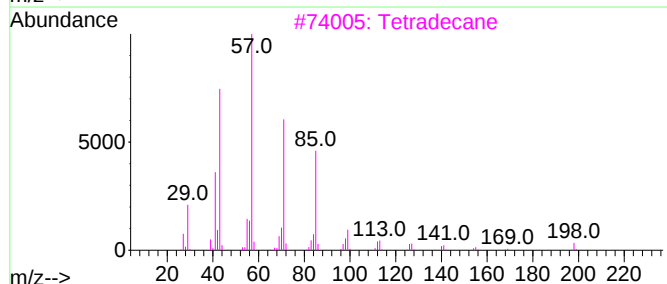
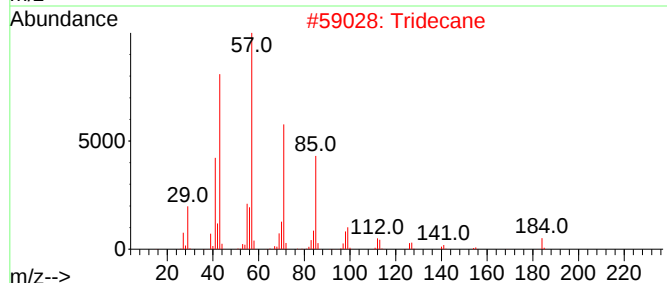
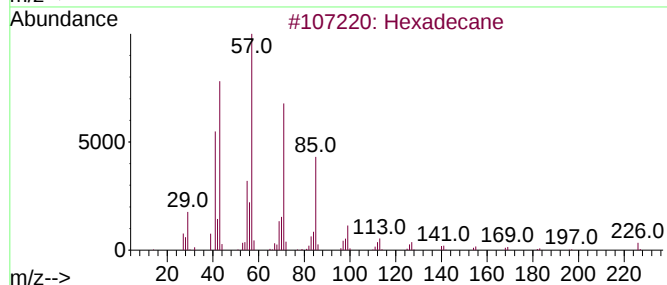
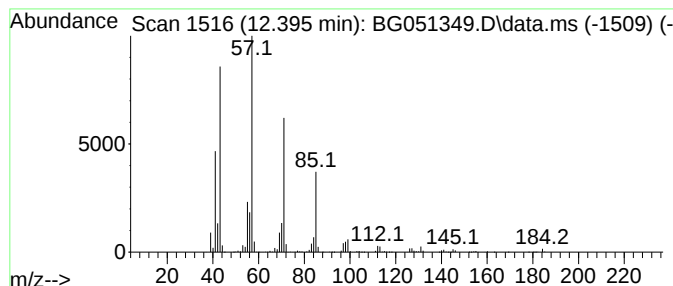
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.395	19.97 ng/ul	3971580	Naphthalene-d8	11.026

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	90
2		Tridecane	184	C13H28	000629-50-5	86
3		Tetradecane	198	C14H30	000629-59-4	86
4		Hexadecane	226	C16H34	000544-76-3	86
5		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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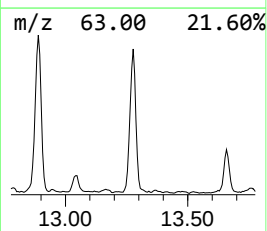
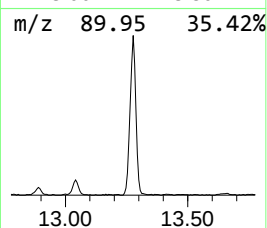
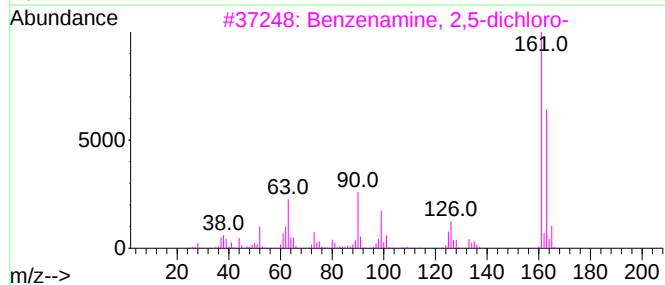
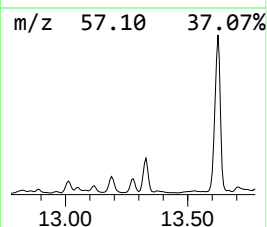
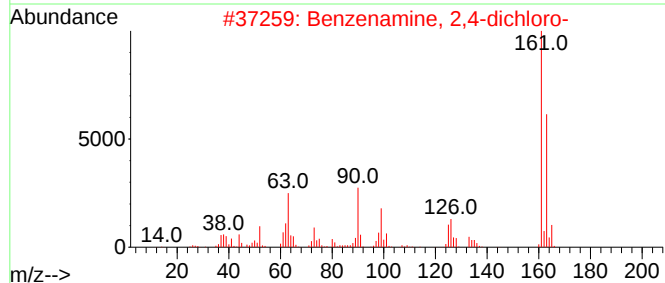
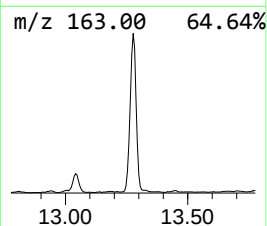
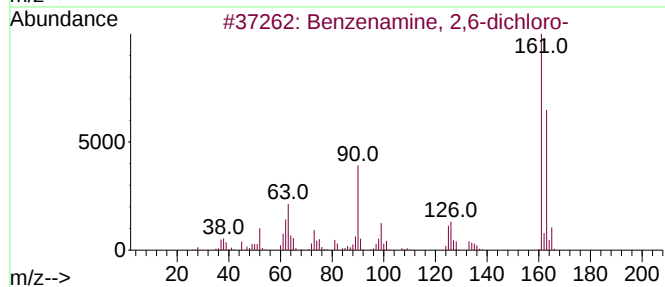
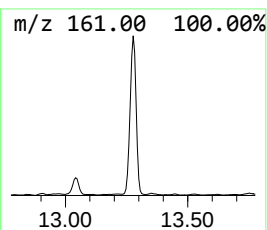
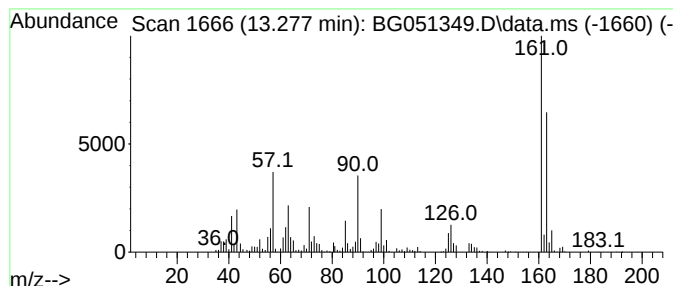
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 Benzenamine, 2,6-dichloro- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.277	34.90 ng/ul	1600400	Acenaphthene-d10	14.828

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenamine, 2,6-dichloro-	161	C6H5Cl2N	000608-31-1	97
2			Benzenamine, 2,4-dichloro-	161	C6H5Cl2N	000554-00-7	96
3			Benzenamine, 2,5-dichloro-	161	C6H5Cl2N	000095-82-9	96
4			Benzenamine, 2,3-dichloro-	161	C6H5Cl2N	000608-27-5	96
5			Benzenamine, 2,4-dichloro-	161	C6H5Cl2N	000554-00-7	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

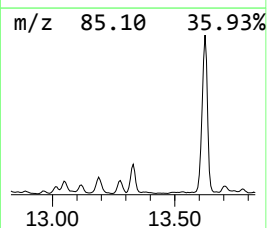
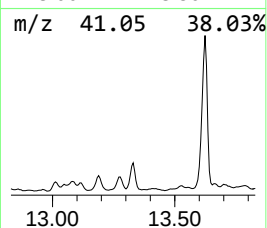
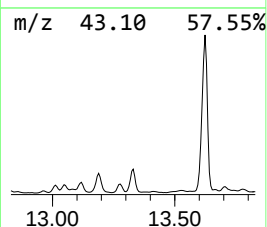
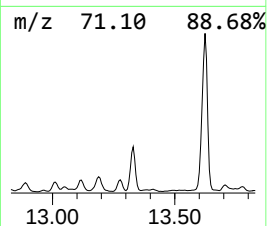
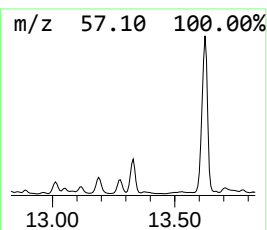
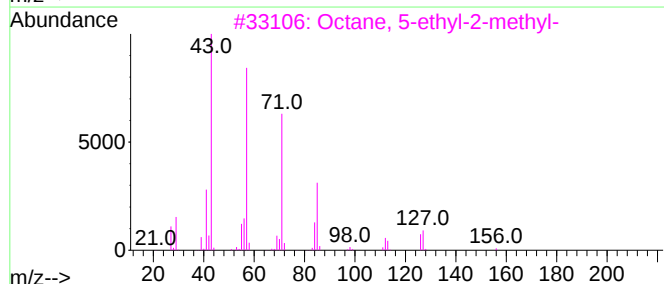
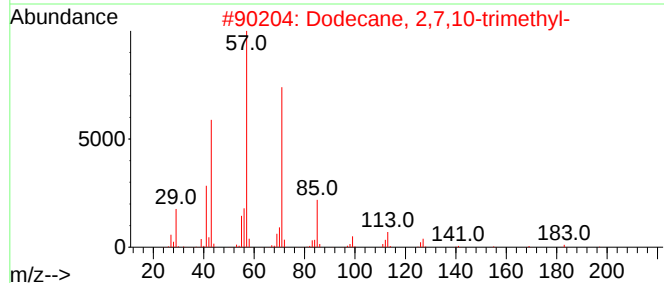
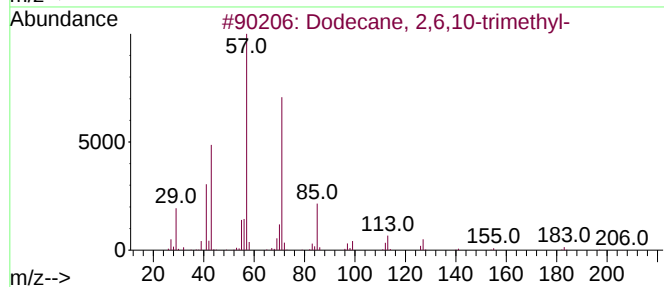
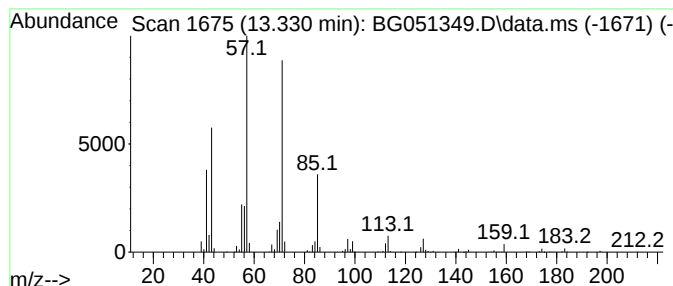
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.330	20.94 ng/ul	959958	Acenaphthene-d10		14.828	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Dodecane, 2,6,10-trimethyl-		212	C15H32	003891-98-3	90
2	Dodecane, 2,7,10-trimethyl-		212	C15H32	074645-98-0	90
3	Octane, 5-ethyl-2-methyl-		156	C11H24	062016-18-6	64
4	Nonane, 2,6-dimethyl-		156	C11H24	017302-28-2	58
5	Octane, 2-bromo-		192	C8H17Br	000557-35-7	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

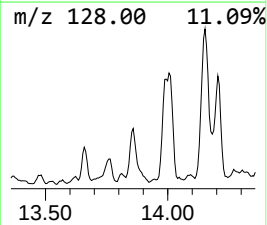
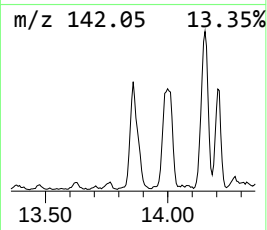
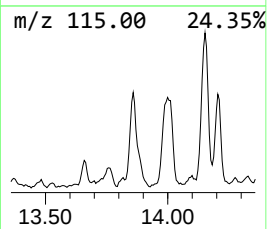
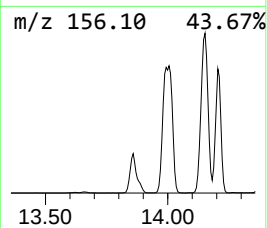
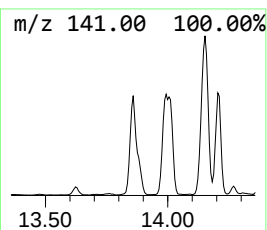
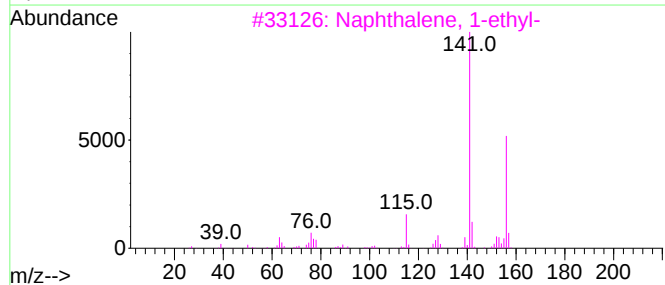
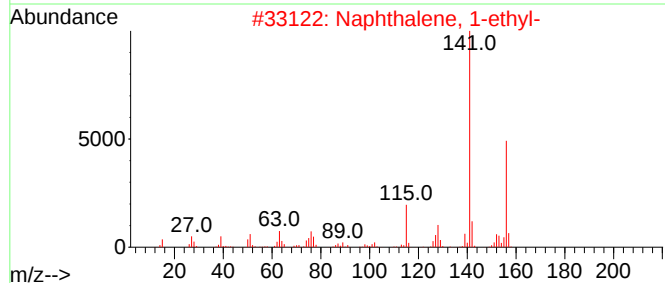
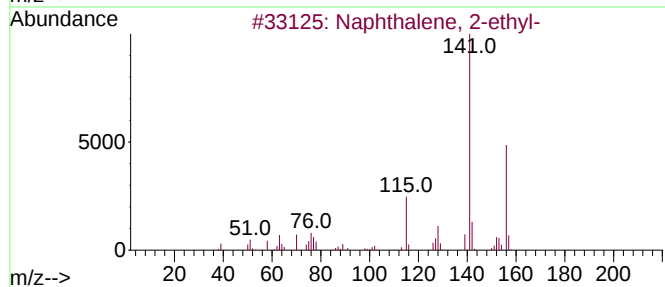
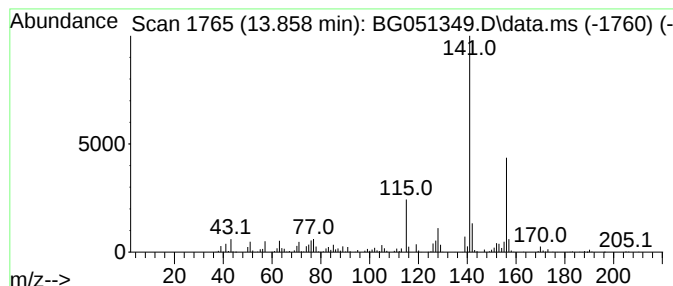
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Quant Title : SVOA CALIBRATION

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

Peak Number 28 Naphthalene, 2-ethyl- Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.858	17.96 ng/ul	823373	Acenaphthene-d10	14.828

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	96
2		Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	95
3		Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	94
4		Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	94
5		Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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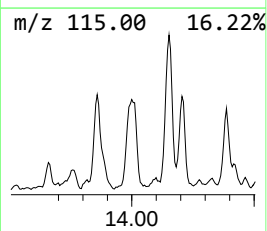
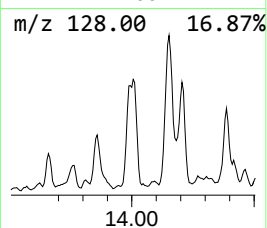
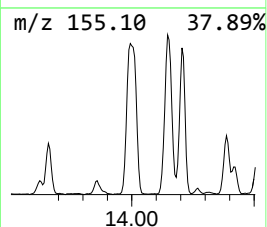
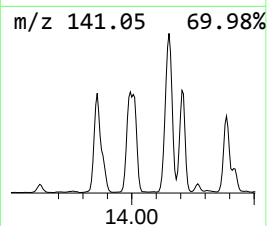
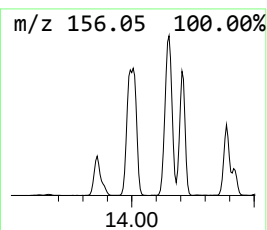
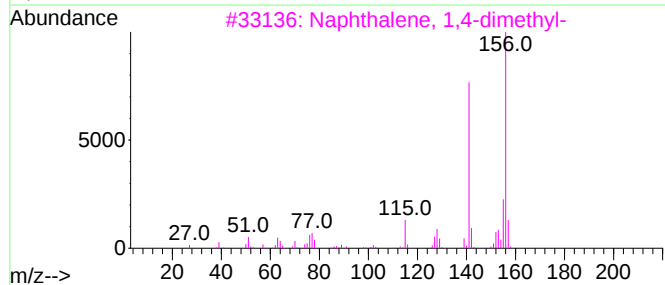
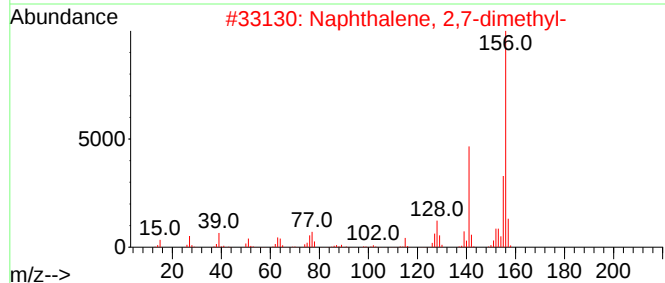
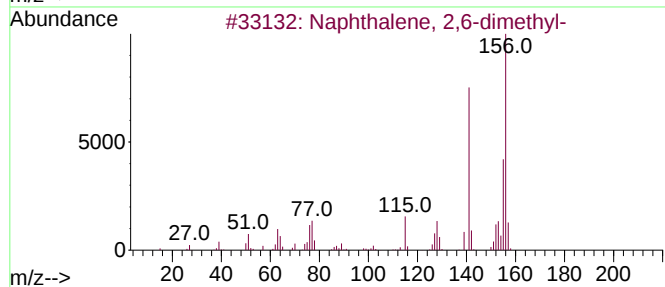
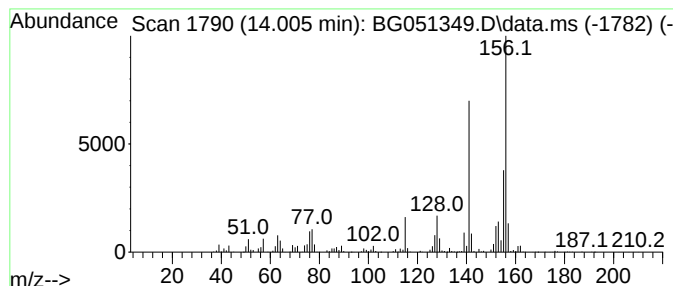
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Quant Title : SVOA CALIBRATION

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

Peak Number 29 Naphthalene, 2,6-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.005	44.59 ng/ul	2044740	Acenaphthene-d10	14.828

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
5			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

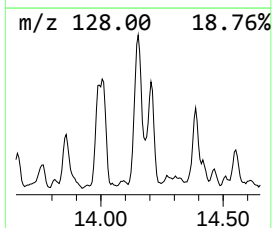
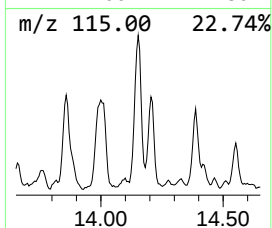
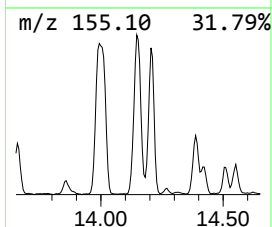
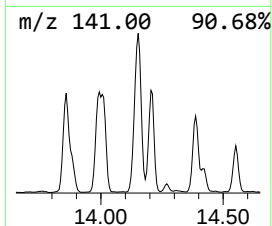
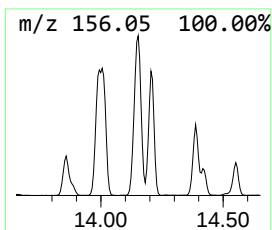
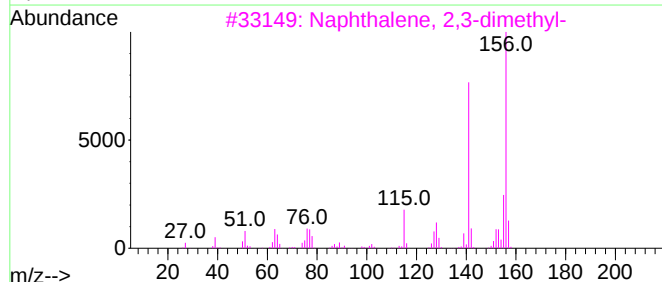
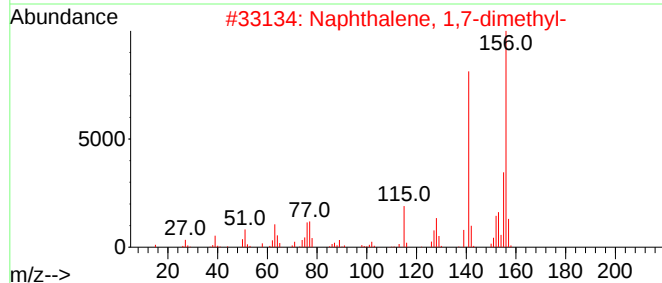
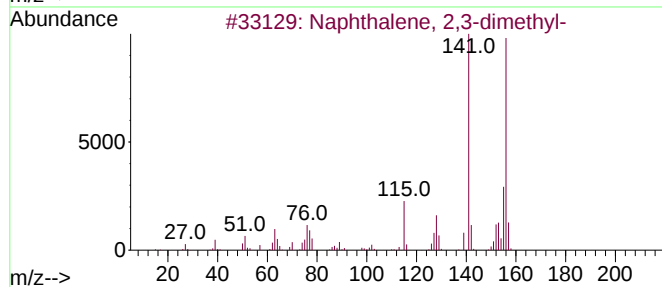
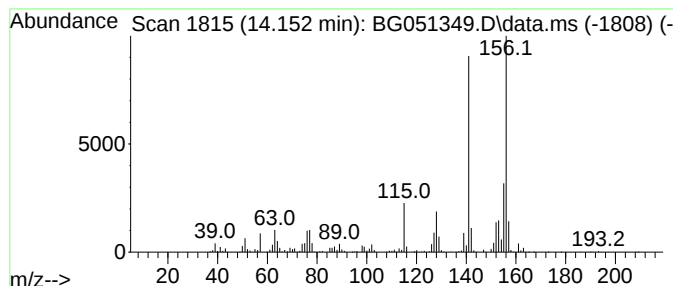
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Quant Title : SVOA CALIBRATION

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

Peak Number 30 Naphthalene, 2,3-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.152	45.95 ng/ul	2106870	Acenaphthene-d10	14.828

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
3			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
5			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

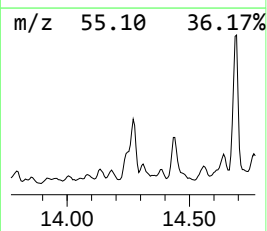
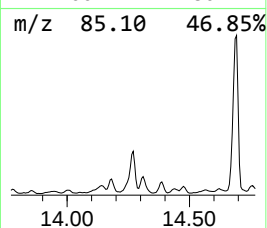
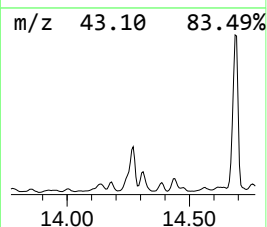
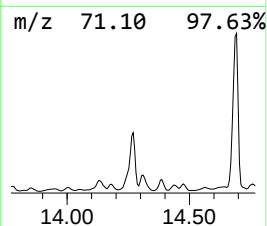
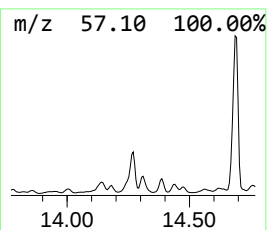
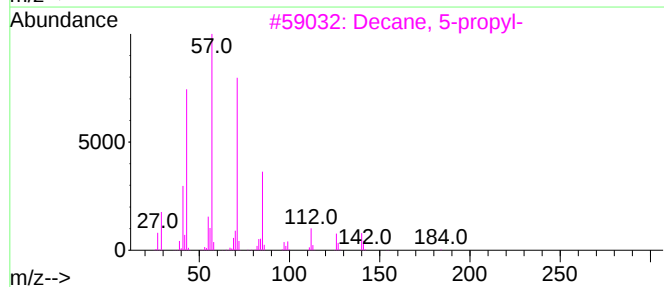
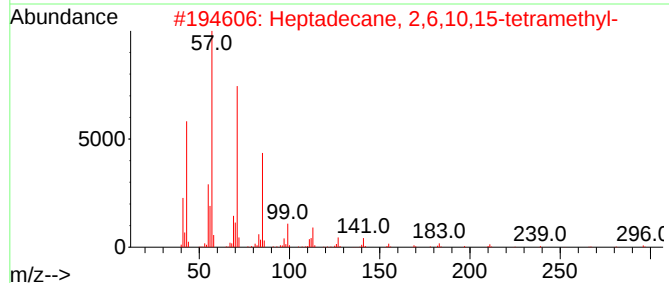
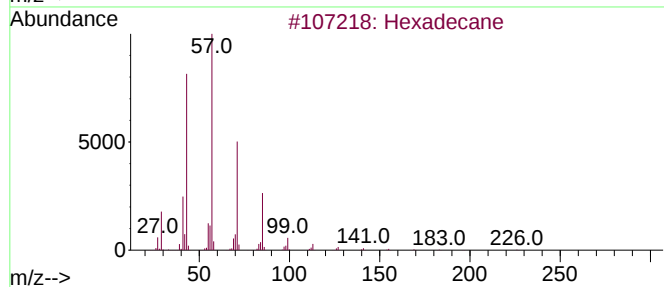
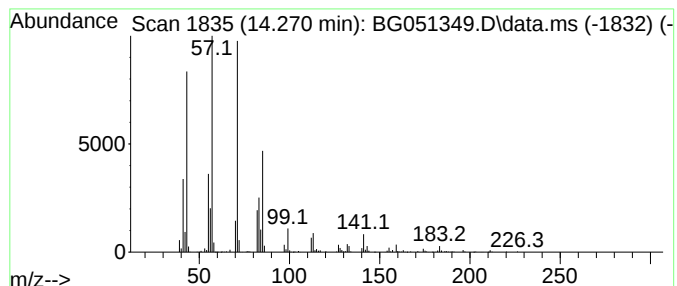
Instrument :
BNA_G
ClientSampleId :
BGKQ1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 31 (DEL) Alkane: Straight-Chai... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.270	26.48 ng/ul	1214340	Acenaphthene-d10	14.828	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	70
2	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	64
3	Decane, 5-propyl-	184	C13H28	017312-62-8	62
4	Dodecane, 4-methyl-	184	C13H28	006117-97-1	58
5	1-Iodo-2-methylnonane	268	C10H21I	1000101-47-9	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

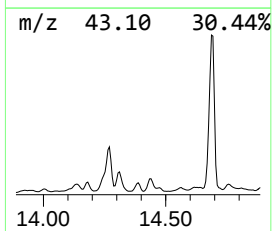
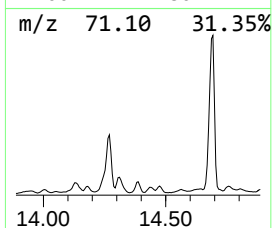
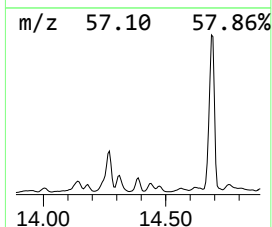
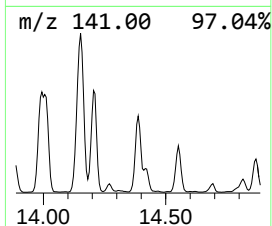
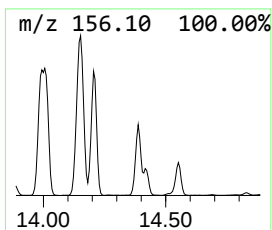
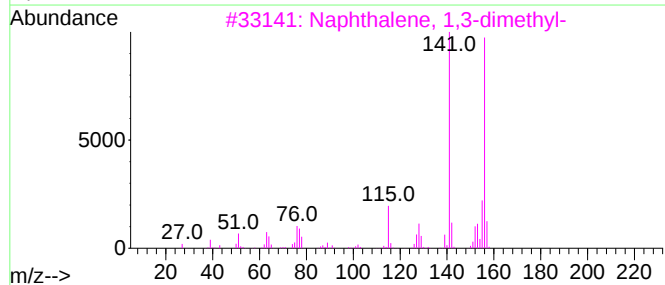
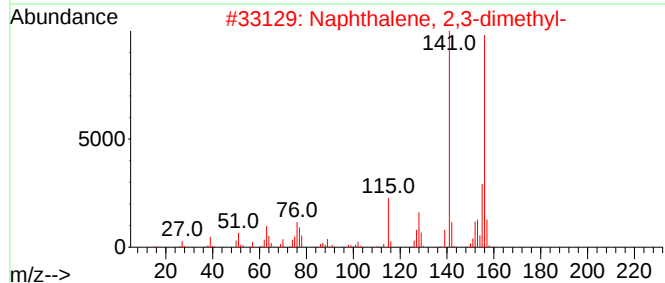
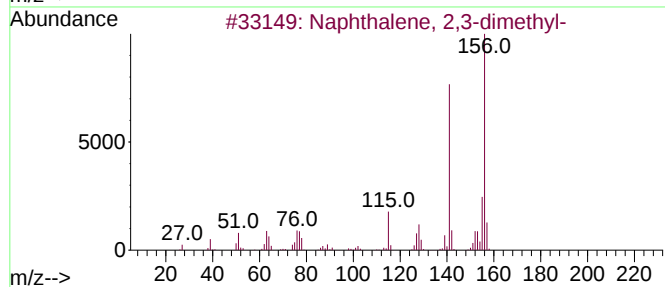
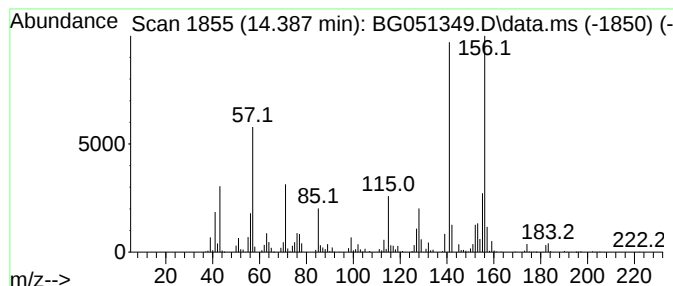
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BNA_G
ClientSampleId :
BGKQ1

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Quant Title : SVOA CALIBRATION

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

Peak Number 32 Naphthalene, 1,3-dimethyl- Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.387	19.94 ng/ul	914367	Acenaphthene-d10	14.828	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
3	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
5	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

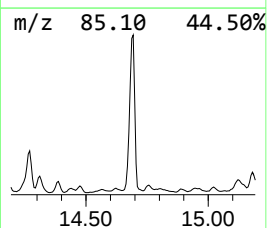
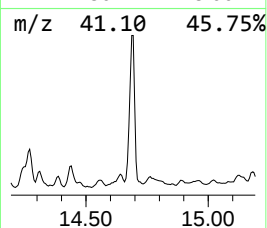
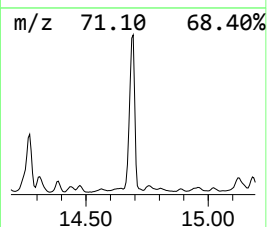
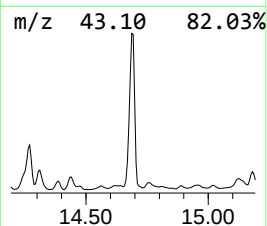
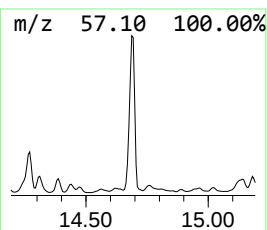
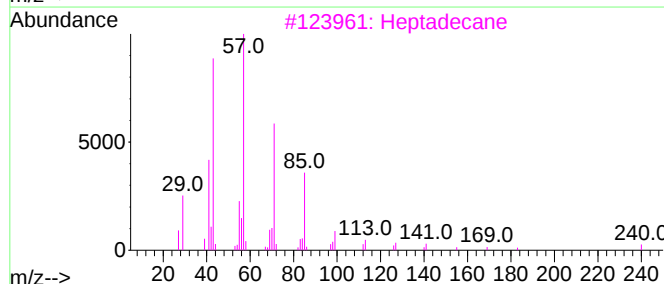
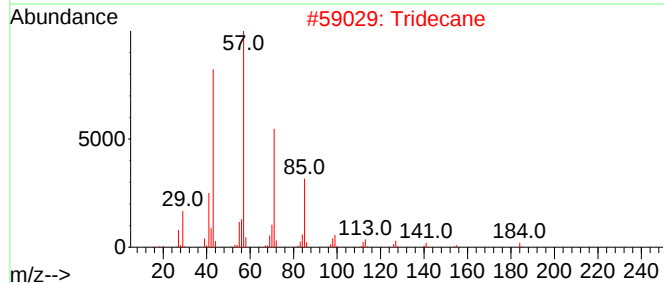
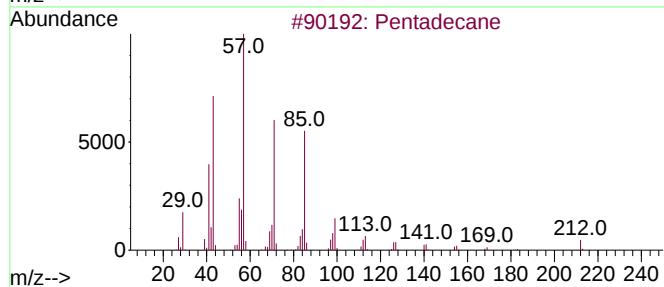
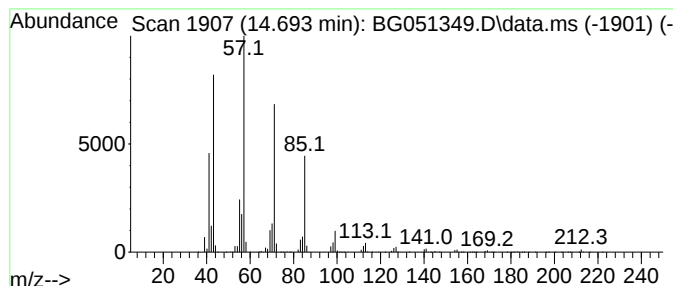
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Quant Title : SVOA CALIBRATION

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

Peak Number 33 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.693	93.53 ng/ul	4288660	Acenaphthene-d10	14.828

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane	212	C15H32	000629-62-9	90
2		Tridecane	184	C13H28	000629-50-5	90
3		Heptadecane	240	C17H36	000629-78-7	90
4		Tetradecane	198	C14H30	000629-59-4	90
5		Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

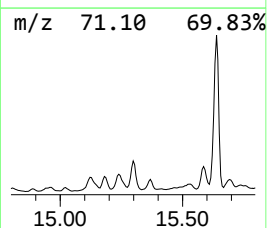
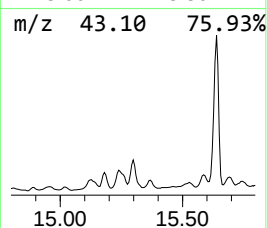
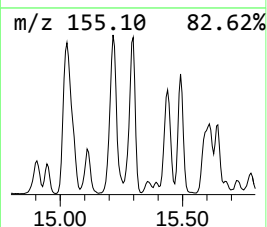
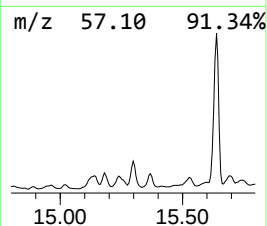
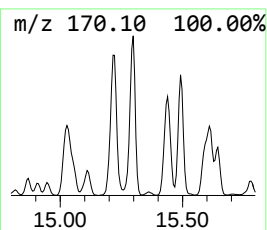
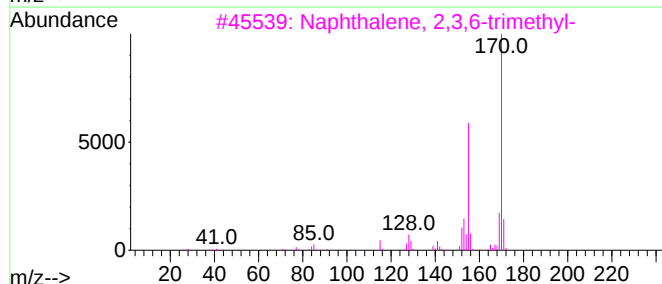
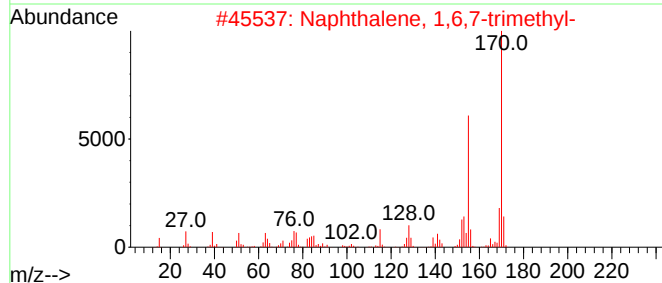
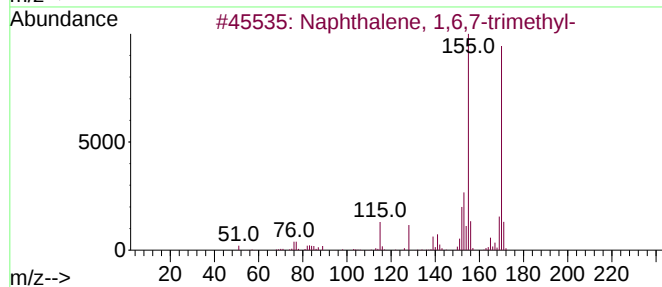
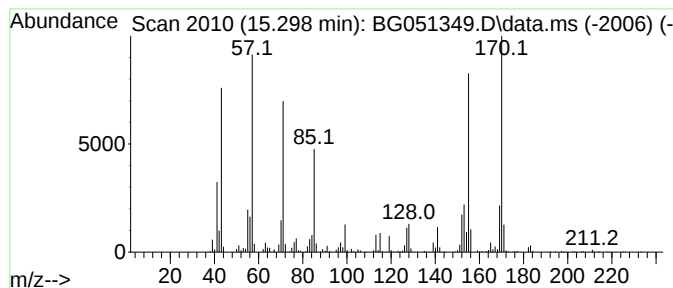
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Quant Title : SVOA CALIBRATION

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

Peak Number 34 Naphthalene, 2,3,6-trimethyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.298	34.95 ng/ul	1602490	Acenaphthene-d10	14.828

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	93
2			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	91
3			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	89
4			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	89
5			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

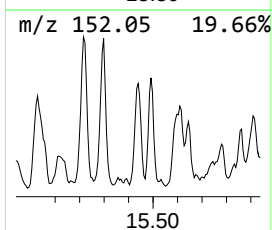
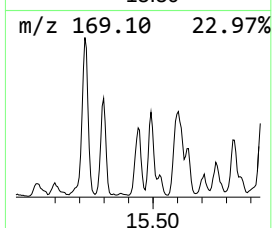
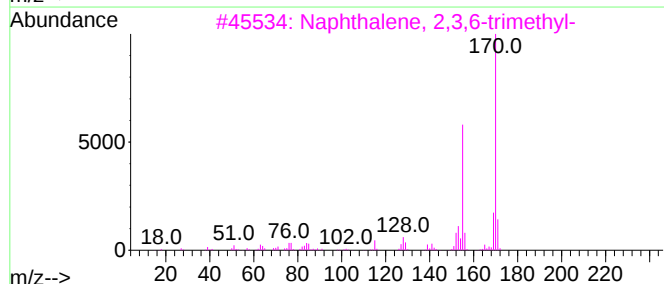
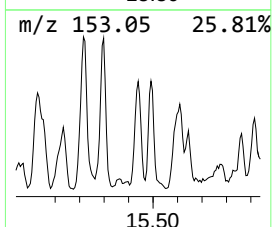
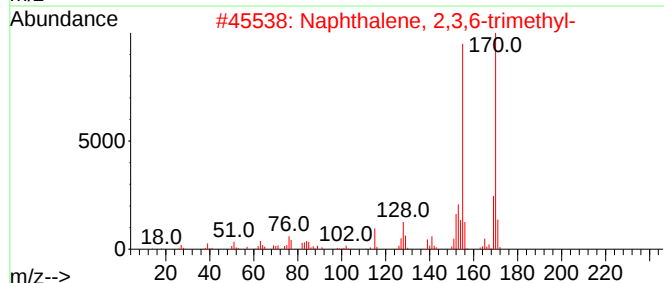
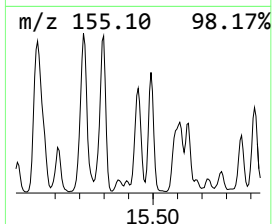
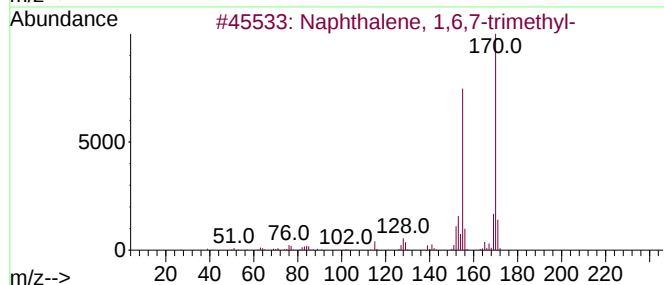
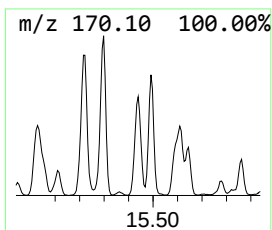
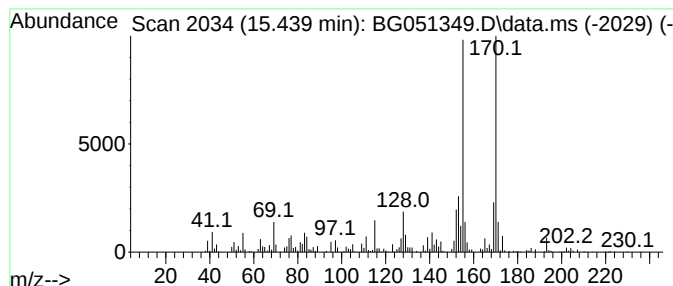
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Quant Title : SVOA CALIBRATION

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

Peak Number 35 Naphthalene, 1,6,7-trimethyl- Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.439	20.32 ng/ul	931695	Acenaphthene-d10	14.828

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	97
2			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
3			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
4			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
5			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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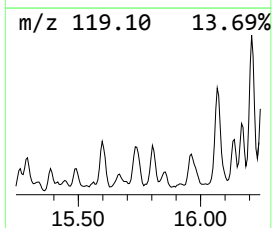
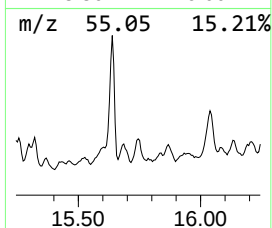
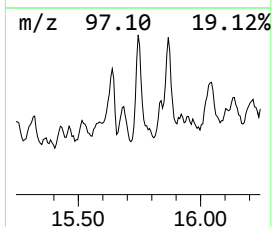
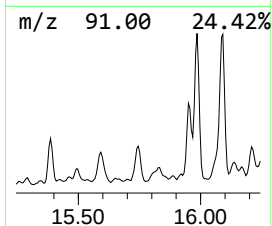
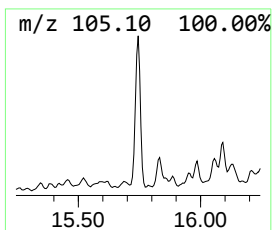
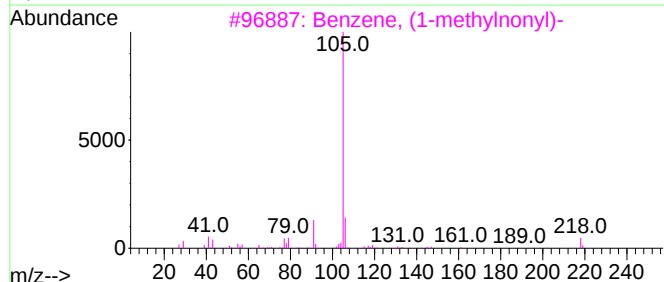
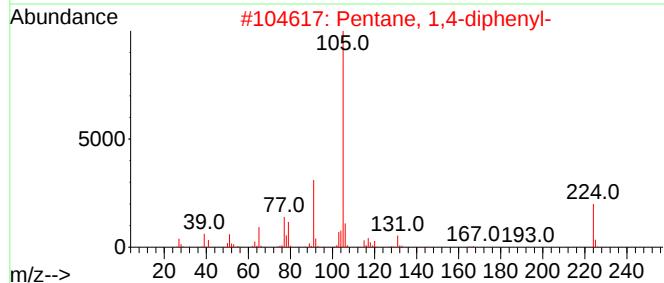
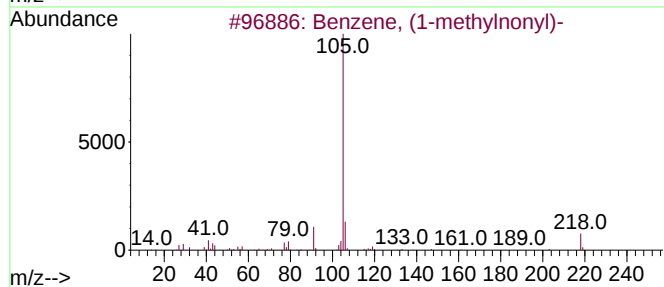
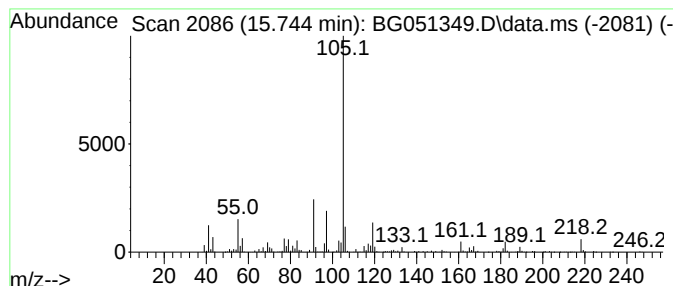
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Quant Title : SVOA CALIBRATION

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

Peak Number 36 unknown-02 Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.744	21.03 ng/ul	964230	Acenaphthene-d10	14.828

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methylnonyl)-	218	C16H26	004537-13-7	43
2			Pentane, 1,4-diphenyl-	224	C17H20	006443-80-7	41
3			Benzene, (1-methylnonyl)-	218	C16H26	004537-13-7	38
4			1-Ethanone, 1-[4-[(4-methylpheny...	240	C16H16O2	1000338-24-3	35
5			(.+/-)-2-Phenylpropanoic Acid, ...	220	C14H20O2	1000453-22-3	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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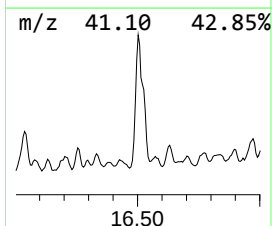
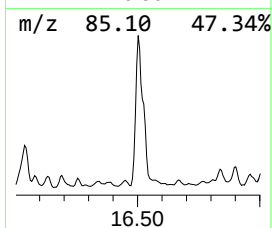
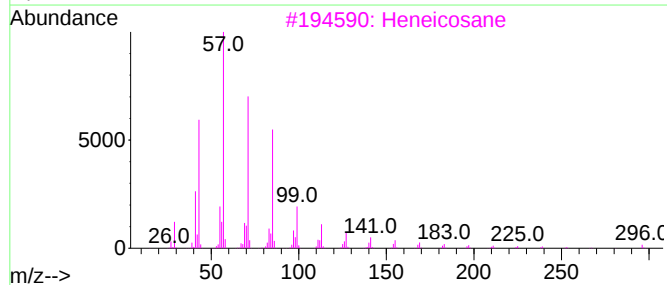
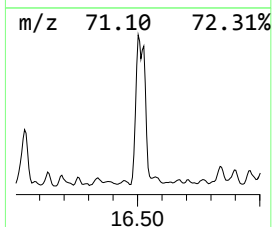
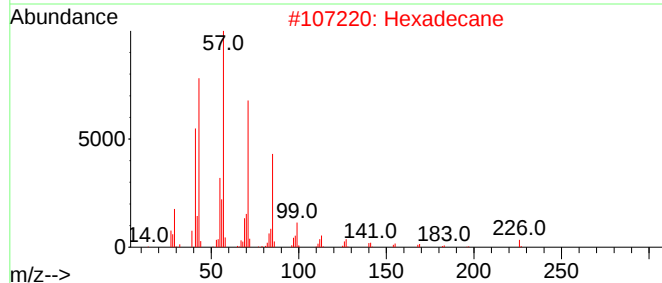
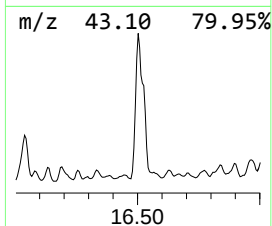
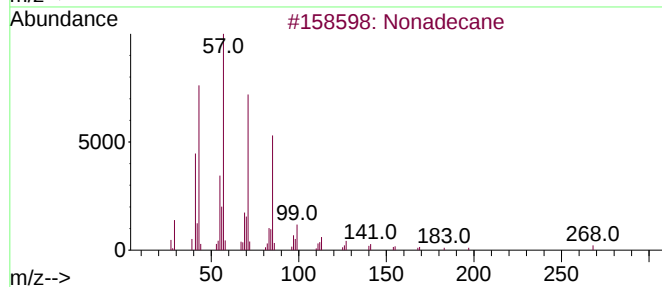
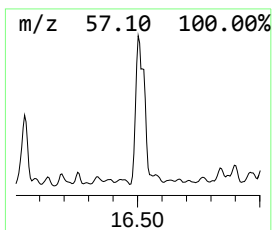
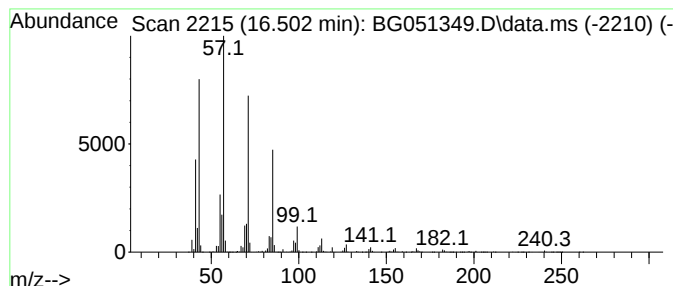
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Quant Title : SVOA CALIBRATION

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

Peak Number 37 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.502	90.57 ng/ul	4245510	Phenanthrene-d10	17.583

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	91
2		Hexadecane	226	C16H34	000544-76-3	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Nonadecane	268	C19H40	000629-92-5	91
5		Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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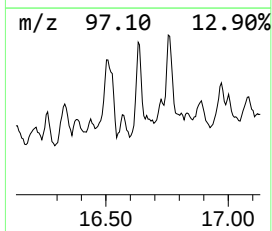
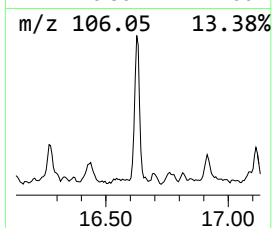
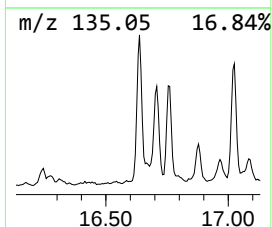
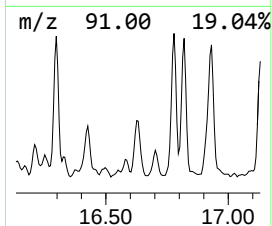
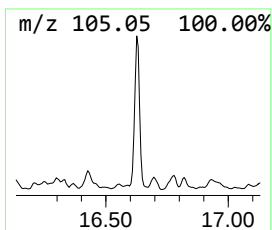
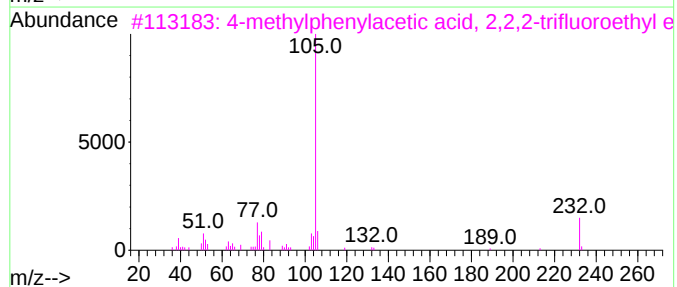
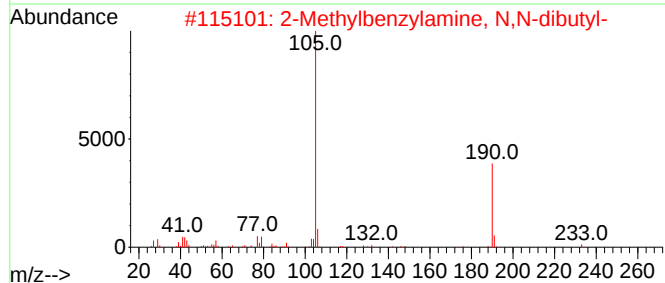
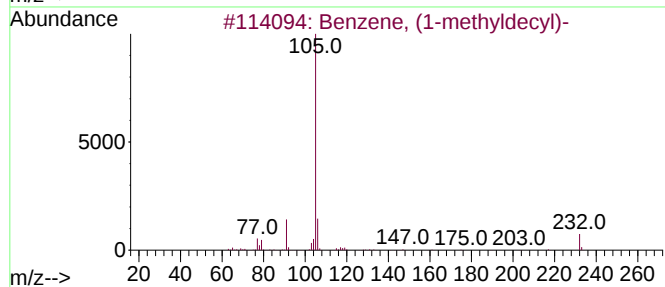
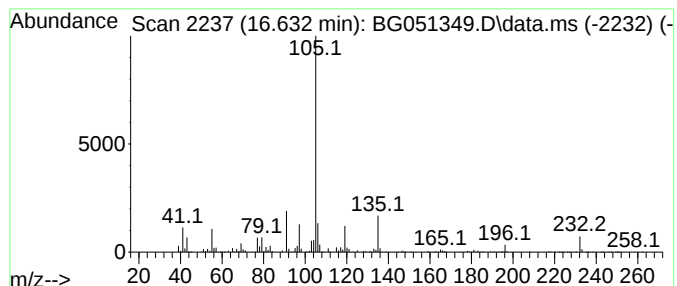
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Quant Title : SVOA CALIBRATION

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

Peak Number 38 Benzene, (1-methyldecyl)- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.632	35.17 ng/ul	1648640	Phenanthrene-d10	17.583

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyldecyl)-	232	C17H28	004536-88-3	90
2			2-Methylbenzylamine, N,N-dibutyl-	233	C16H27N	1000310-39-6	38
3			4-methylphenylacetic acid, 2,2,2...	232	C11H11F3O2	1000376-55-2	38
4			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	35
5			Ethyl meta-tolylacetate	178	C11H14O2	040061-55-0	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

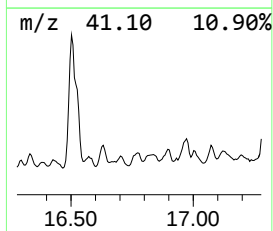
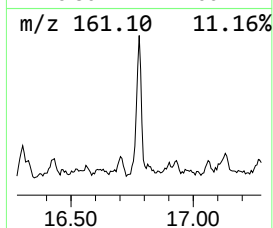
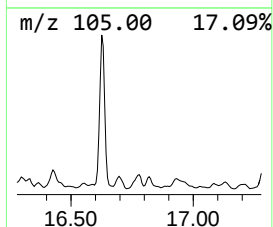
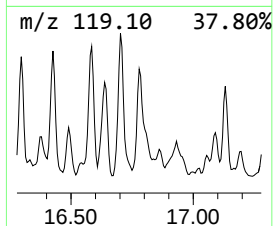
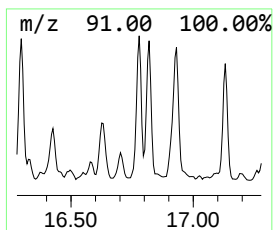
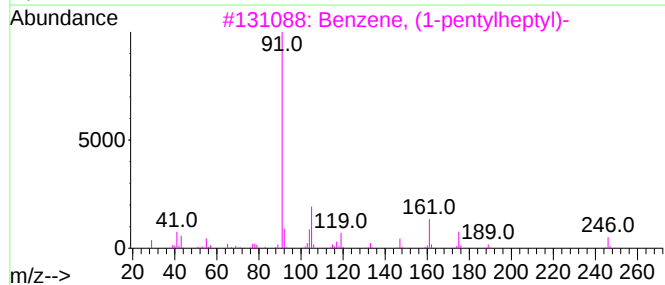
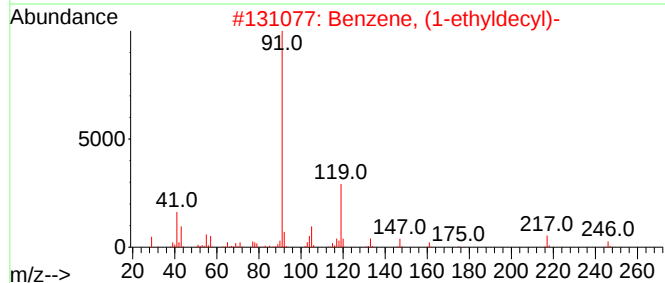
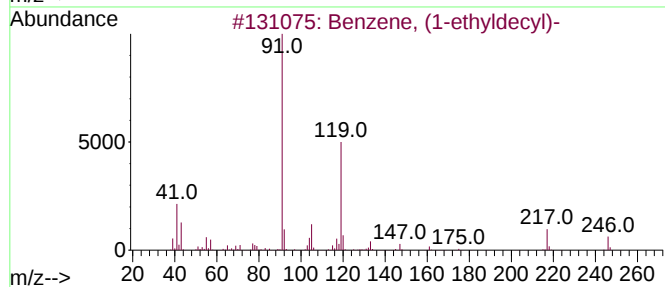
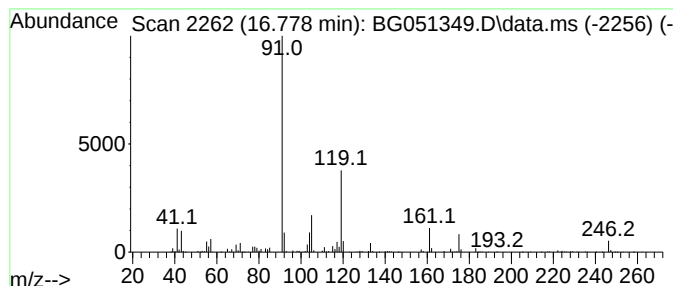
Instrument :
BNA_G
ClientSampleId :
BGKQ1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 39 Benzene, (1-ethyldecyl)- Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.779	19.78 ng/ul	927186	Phenanthrene-d10		17.583	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, (1-ethyldecyl)-		246	C18H30	002400-00-2	93
2	Benzene, (1-ethyldecyl)-		246	C18H30	002400-00-2	87
3	Benzene, (1-pentylheptyl)-		246	C18H30	002719-62-2	76
4	Benzene, (1-pentylheptyl)-		246	C18H30	002719-62-2	70
5	Benzene, (1-pentylheptyl)-		246	C18H30	002719-62-2	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

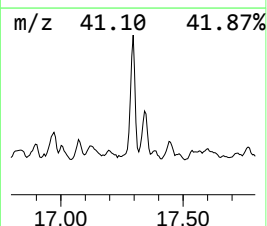
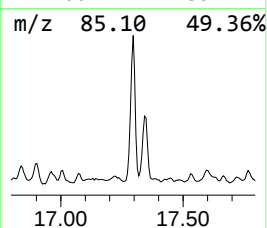
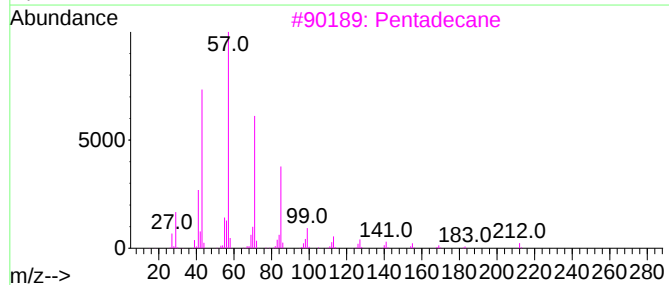
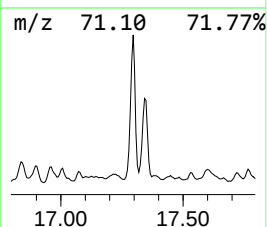
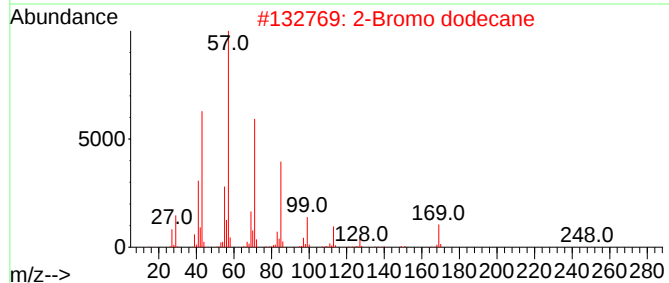
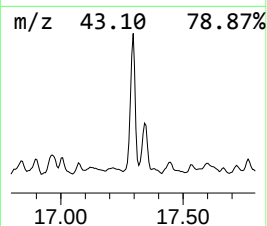
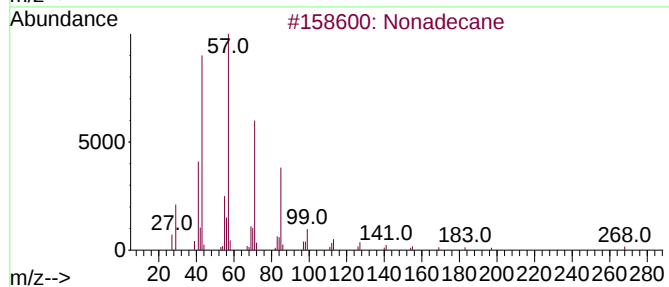
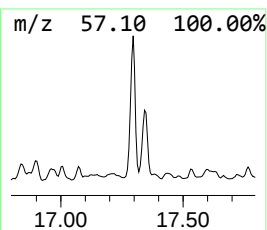
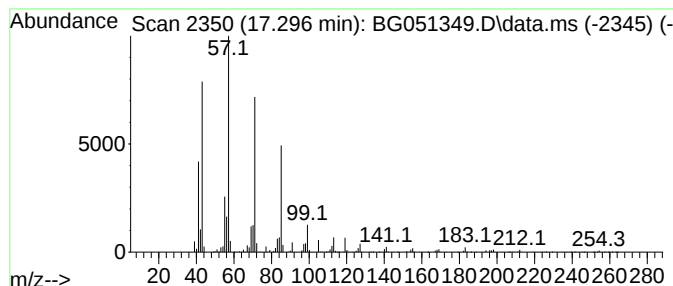
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD			R.T.
17.296	66.85 ng/ul	3133450	Phenanthrene-d10			17.583
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	90
2	2-Bromo dodecane		248	C12H25Br	013187-99-0	89
3	Pentadecane		212	C15H32	000629-62-9	86
4	Dodecane, 2-methyl-6-propyl-		226	C16H34	055045-08-4	86
5	Octadecane		254	C18H38	000593-45-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
BGKQ1

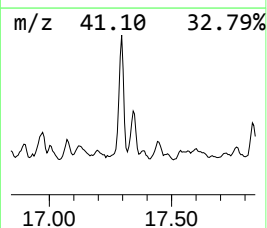
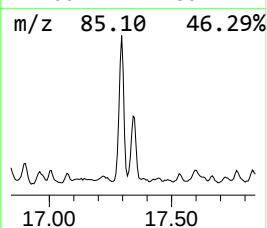
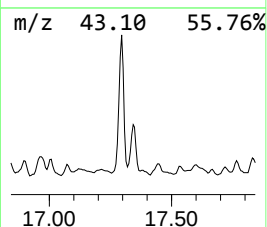
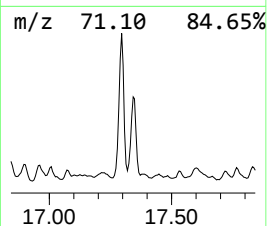
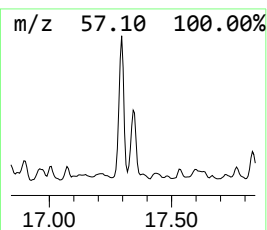
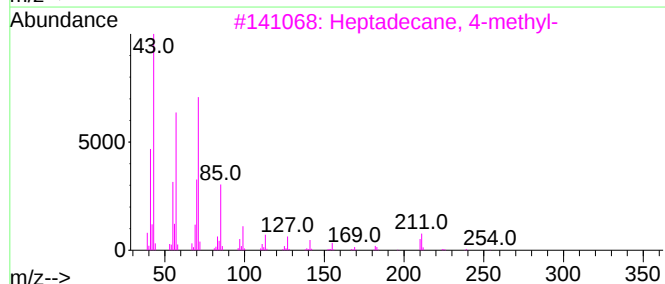
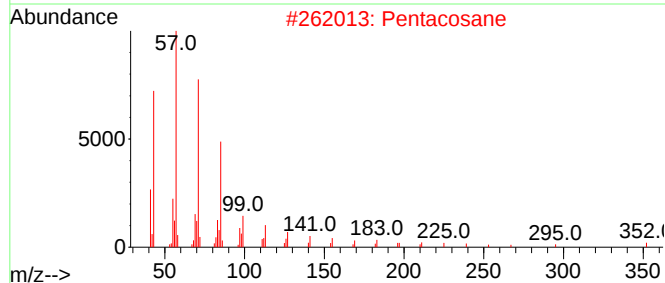
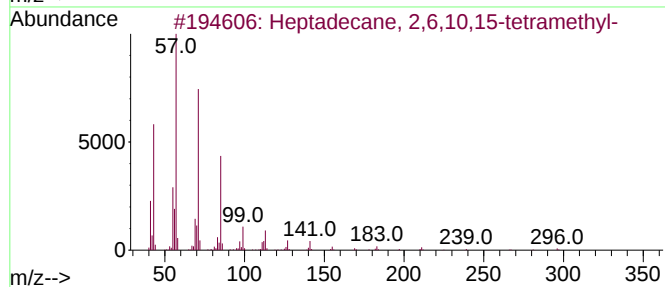
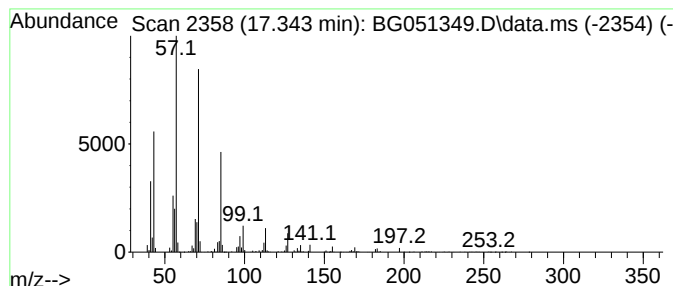
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Quant Title : SVOA CALIBRATION

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

Peak Number 41 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.343	31.25 ng/ul	1464900	Phenanthrene-d10	17.583

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
2		Pentacosane	352	C25H52	000629-99-2	90
3		Heptadecane, 4-methyl-	254	C18H38	026429-11-8	90
4		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	86
5		Hexadecane	226	C16H34	000544-76-3	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

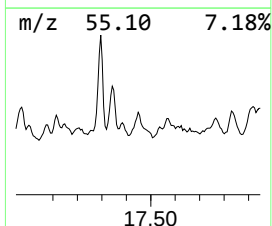
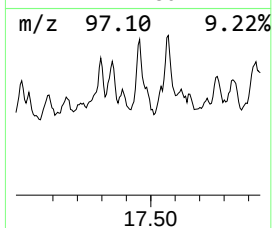
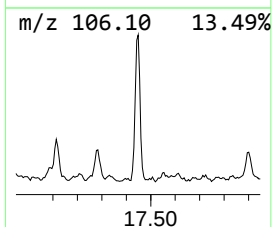
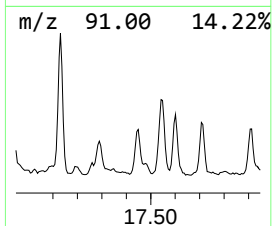
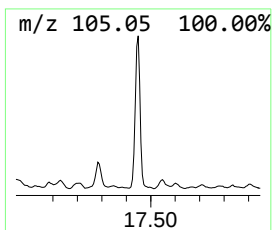
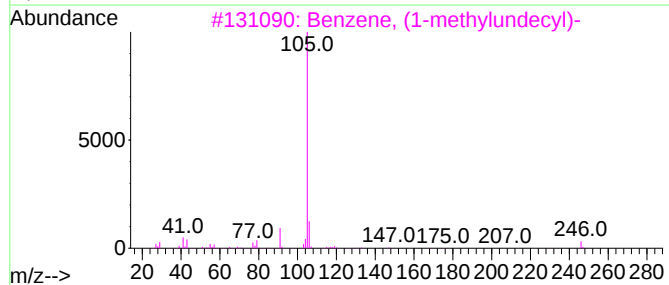
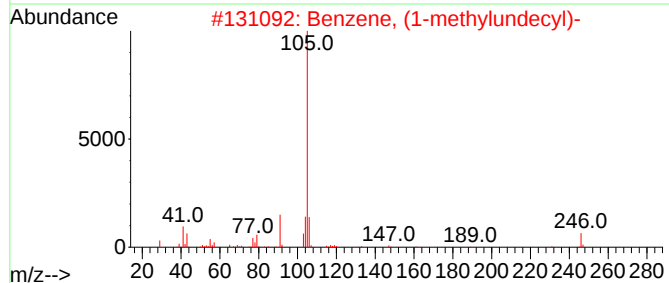
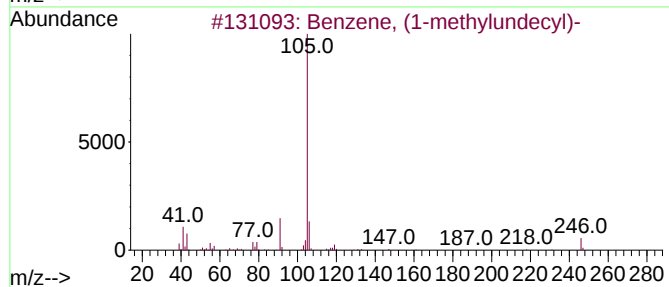
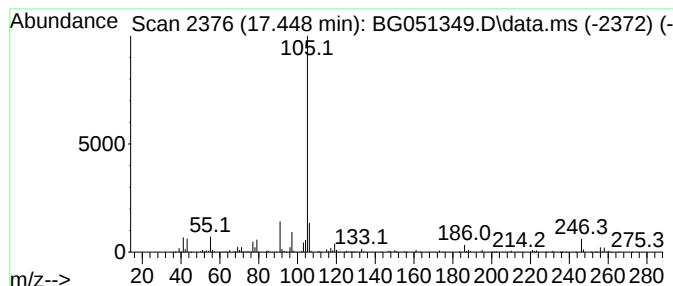
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Quant Title : SVOA CALIBRATION

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

Peak Number 42 Benzene, (1-methylundecyl)- Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.448	20.00 ng/ul	937490	Phenanthrene-d10	17.583

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methylundecyl)-	246	C18H30	002719-61-1	70
2			Benzene, (1-methylundecyl)-	246	C18H30	002719-61-1	70
3			Benzene, (1-methylundecyl)-	246	C18H30	002719-61-1	58
4			Benzene, (1,3,3-trimethylnonyl)-	246	C18H30	054986-44-6	53
5			Benzene, (1-methylundecyl)-	246	C18H30	002719-61-1	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

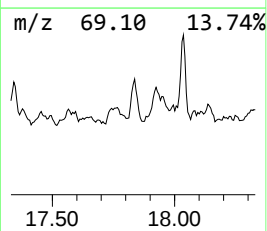
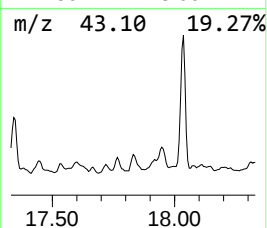
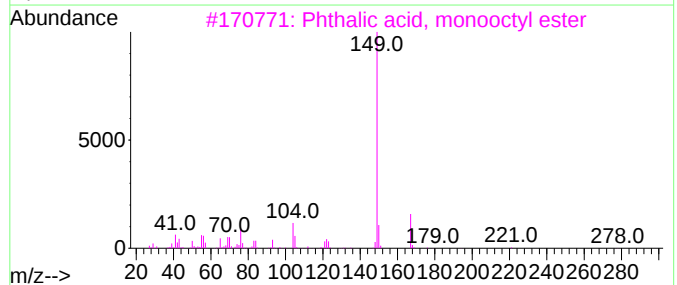
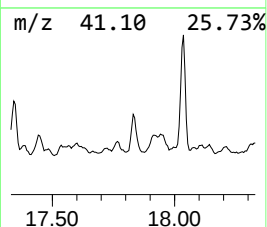
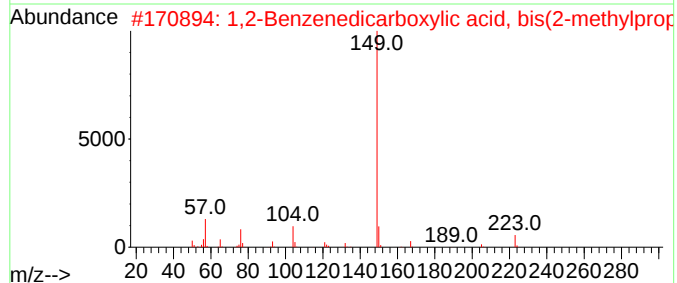
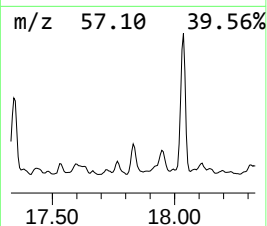
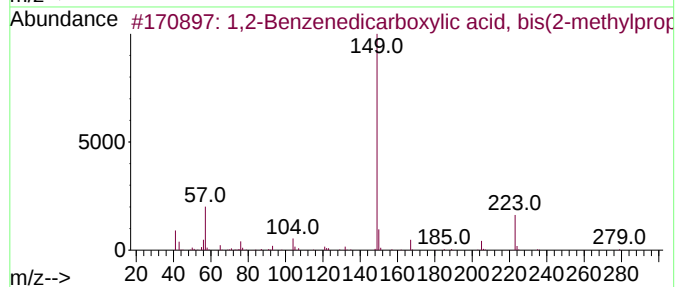
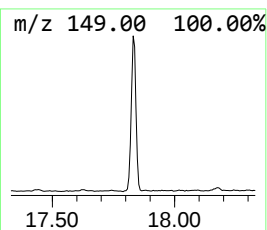
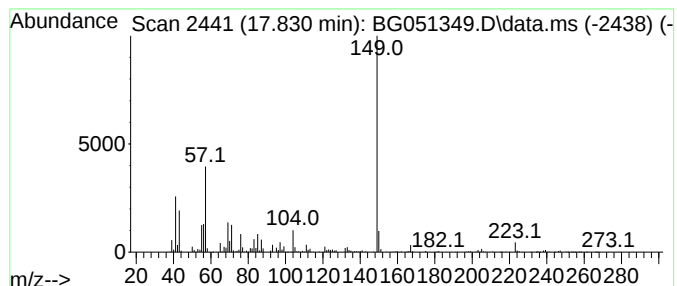
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Quant Title : SVOA CALIBRATION

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

Peak Number 43 1,2-Benzenedicarboxylic aci... Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.830	25.09 ng/ul	1176310	Phenanthrene-d10	17.583

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Benzenedicarboxylic acid, bi...	278	C16H22O4	000084-69-5	76
2			1,2-Benzenedicarboxylic acid, bi...	278	C16H22O4	000084-69-5	68
3			Phthalic acid, monoethyl ester	278	C16H22O4	005393-19-1	64
4			Phthalic acid, isopropyl propyl ...	250	C14H18O4	1010314-99-7	64
5			1,2-Benzenedicarboxylic acid, bi...	278	C16H22O4	000084-69-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

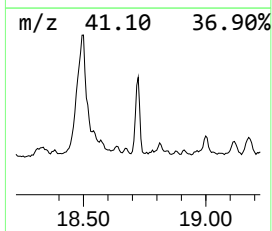
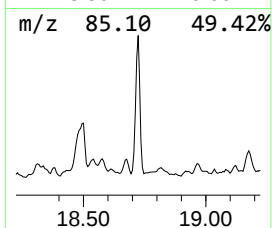
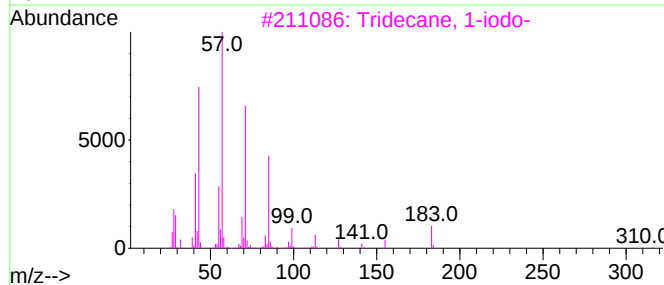
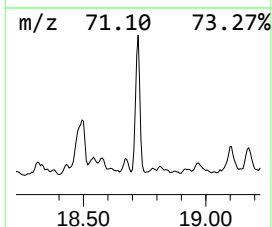
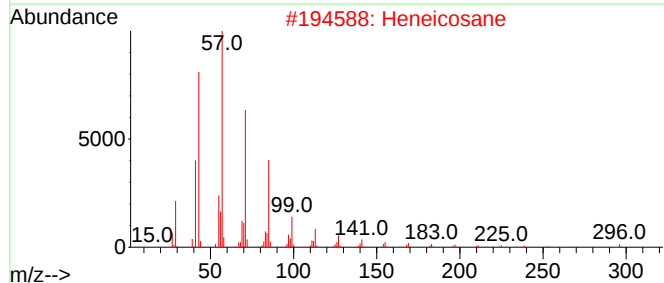
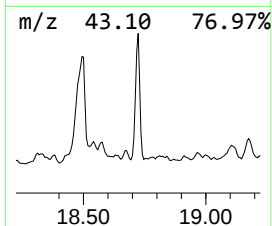
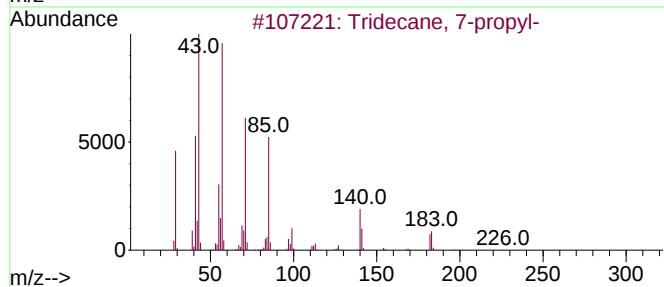
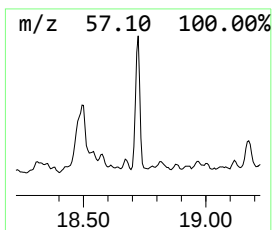
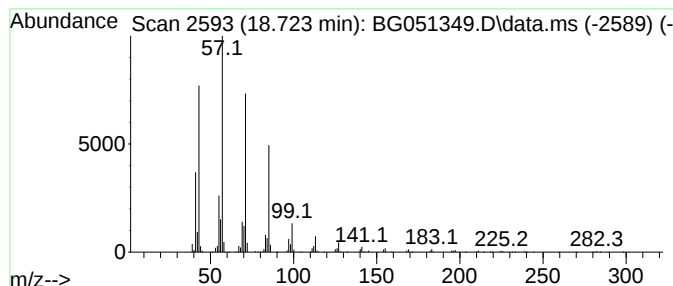
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Quant Title : SVOA CALIBRATION

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

Peak Number 44 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.723	41.39 ng/ul	1940210	Phenanthrene-d10	17.583

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 7-propyl-	226	C16H34	055045-09-5	93
2			Heneicosane	296	C21H44	000629-94-7	91
3			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	90
4			Octacosane	394	C28H58	000630-02-4	90
5			Sulfurous acid, 2-propyl tetradec...	320	C17H36O3S	1010309-12-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

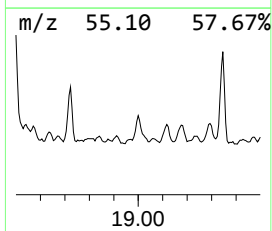
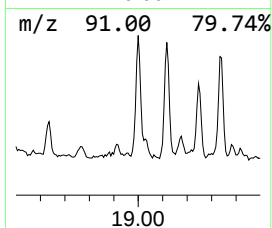
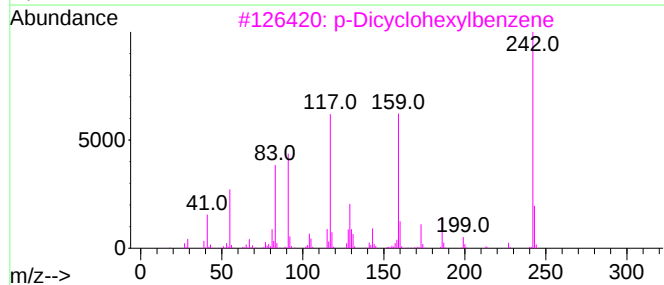
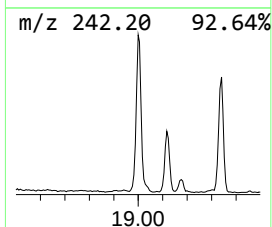
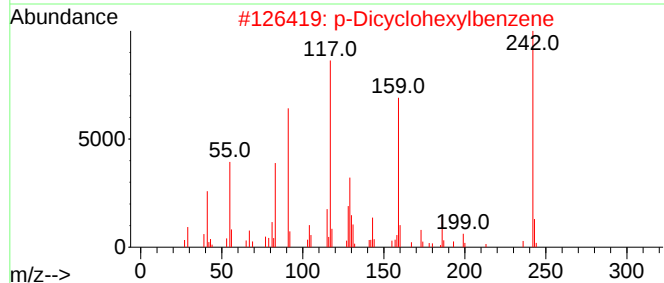
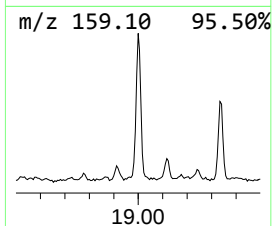
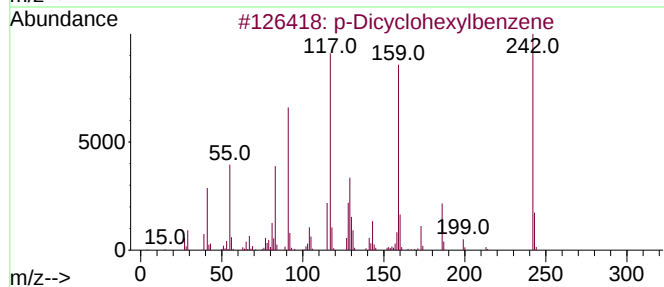
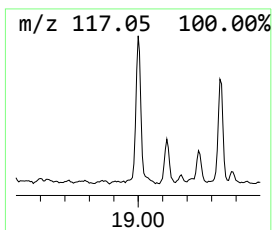
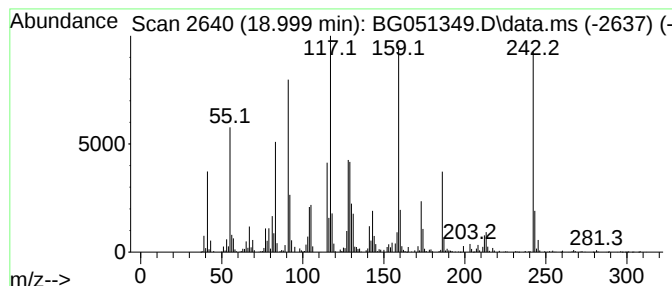
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 45 p-Dicyclohexylbenzene Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.999	27.48 ng/ul	1288330	Phenanthrene-d10		17.583	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	p-Dicyclohexylbenzene		242	C18H26	001087-02-1	93
2	p-Dicyclohexylbenzene		242	C18H26	001087-02-1	91
3	p-Dicyclohexylbenzene		242	C18H26	001087-02-1	91
4	Bicyclohexyl, 4-phenyl-		242	C18H26	020273-27-2	76
5	2-(4-(2-Methylallyl)phenyl)propa...		204	C13H16O2	131734-08-2	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

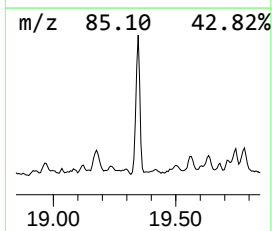
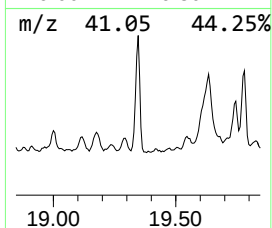
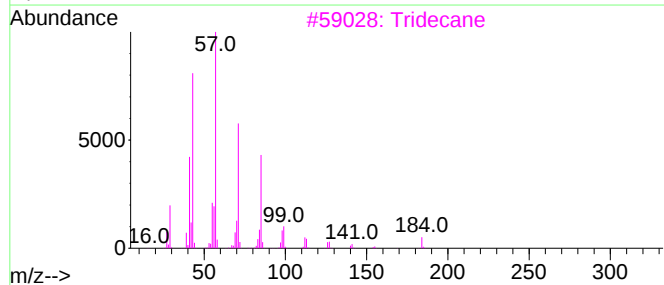
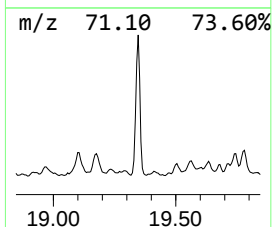
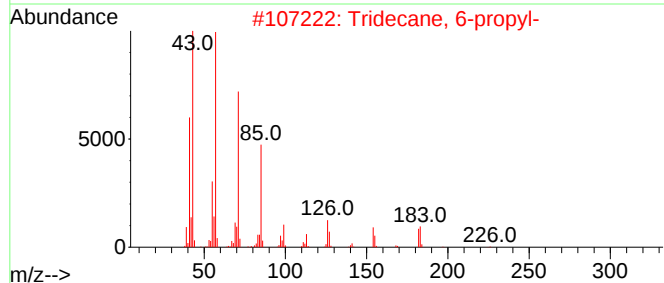
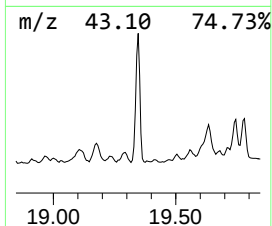
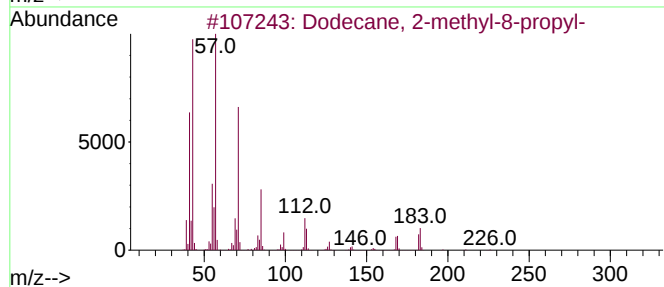
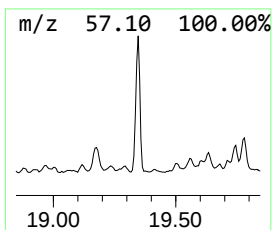
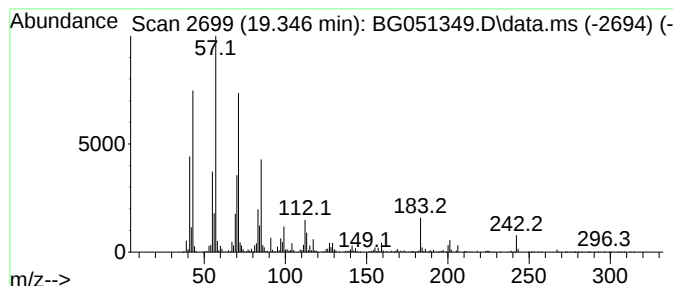
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Quant Title : SVOA CALIBRATION

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

Peak Number 46 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.346	77.35 ng/ul	3625530	Phenanthrene-d10	17.583

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Dodecane, 2-methyl-8-propyl-	226	C16H34	055045-07-3	86
2		Tridecane, 6-propyl-	226	C16H34	055045-10-8	76
3		Tridecane	184	C13H28	000629-50-5	76
4		Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	68
5		Hexadecane	226	C16H34	000544-76-3	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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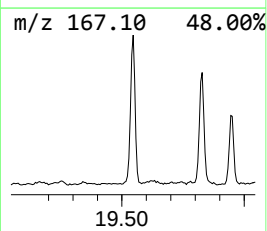
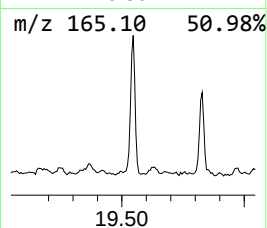
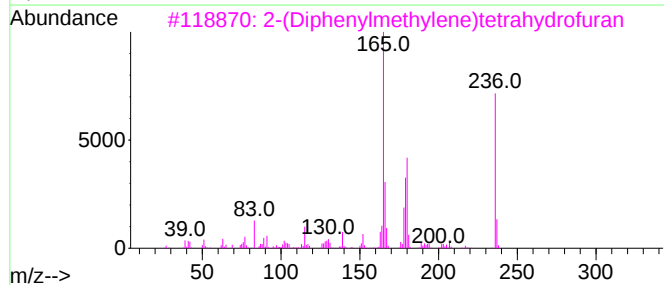
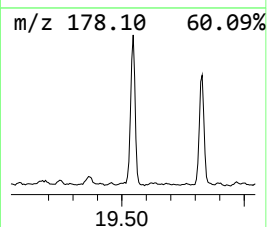
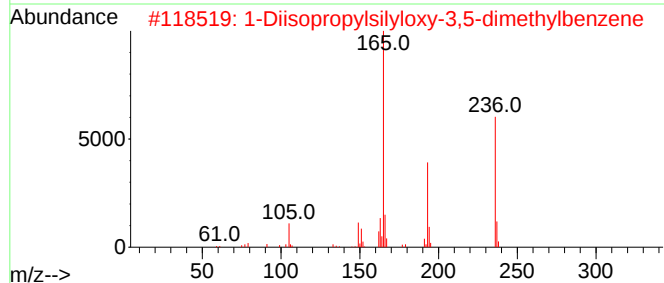
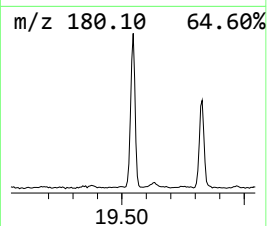
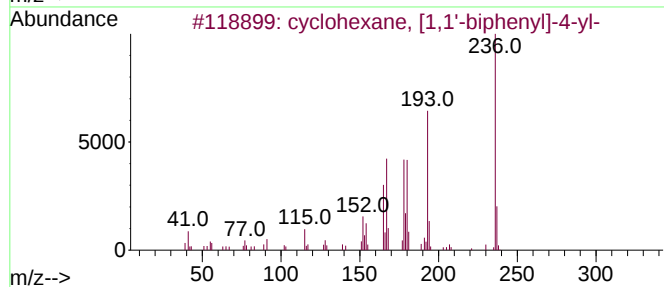
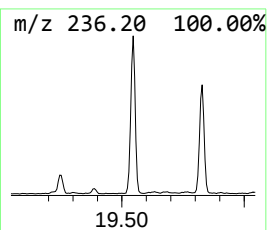
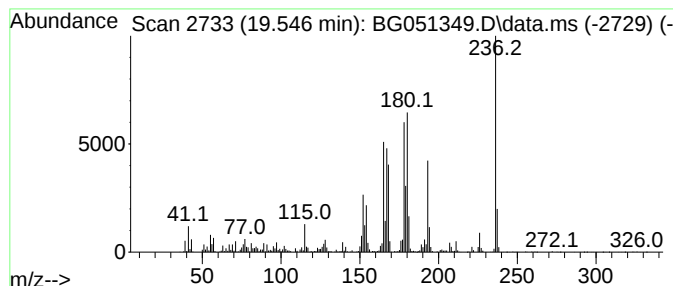
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Quant Title : SVOA CALIBRATION

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

Peak Number 47 cyclohexane, [1,1'-biphenyl]... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.546	34.20 ng/ul	1603170	Phenanthrene-d10	17.583

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cyclohexane, [1,1'-biphenyl]-4-yl-	236	C18H20	1000401-12-4	94
2			1-Diisopropylsilyloxy-3,5-dimeth...	236	C14H24OSi	1000307-97-3	45
3			2-(Diphenylmethylene)tetrahydrof...	236	C17H16O	039077-35-5	38
4			2-Methyl-2-phenyl-1H-indene-1,3(...	236	C16H12O2	1010332-18-2	27
5			1-[(2-Hydroxyethyl)amino]anthra...	267	C16H13NO3	004465-58-1	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

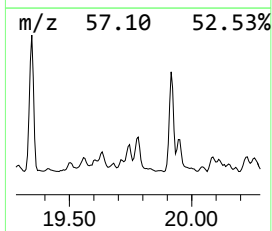
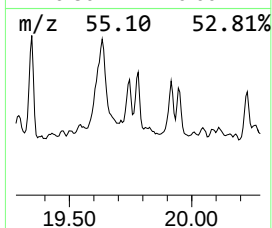
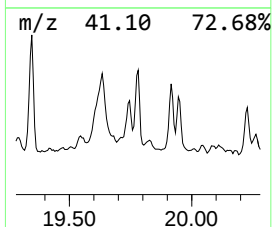
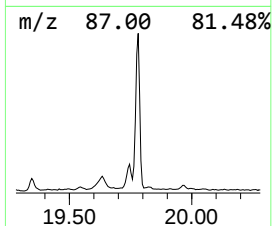
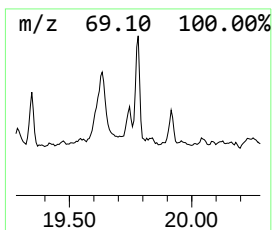
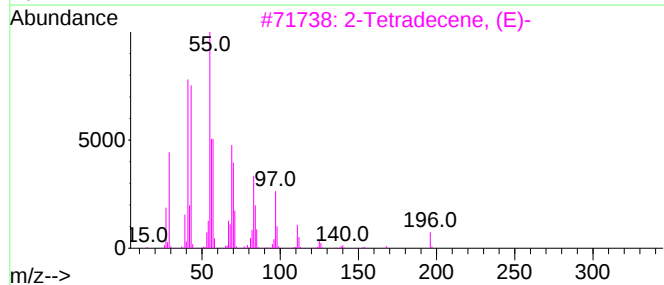
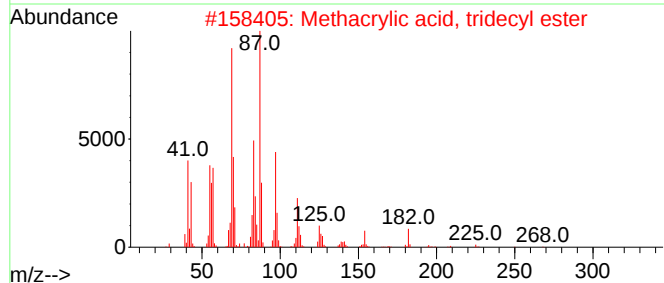
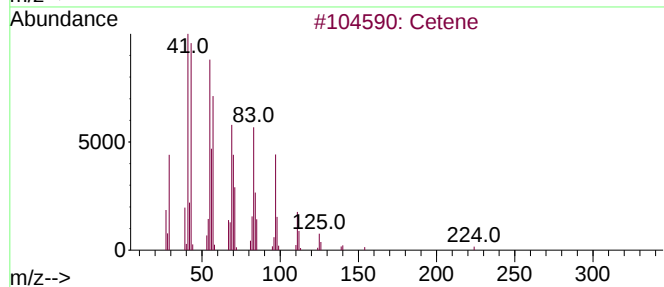
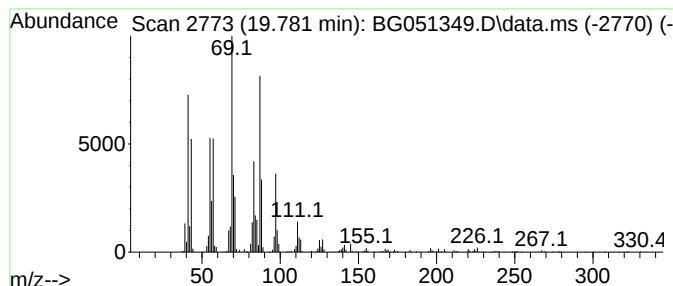
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 48 (DEL) Alkane: Straight-Chai... Concentration Rank 59

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.781	12.70 ng/ul	1339030	Chrysene-d12		21.902	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Cetene		224	C16H32	000629-73-2	92
2	Methacrylic acid, tridecyl ester		268	C17H32O2	1000340-28-9	83
3	2-Tetradecene, (E)-		196	C14H28	035953-54-9	74
4	Pentadecafluorooctanoic acid, te...		610	C22H29F15O2	1000406-04-4	42
5	Dichloroacetic acid, tetradecyl ...		324	C16H30Cl2O2	083005-02-1	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

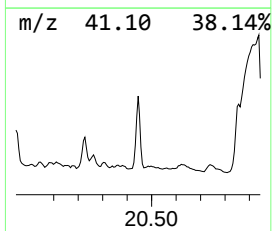
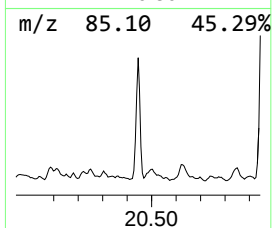
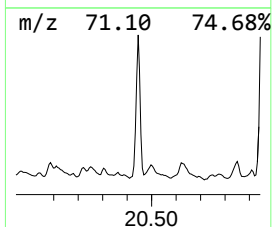
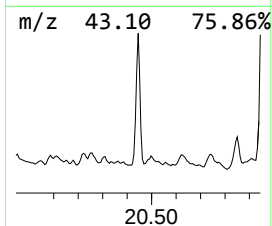
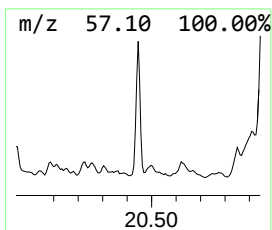
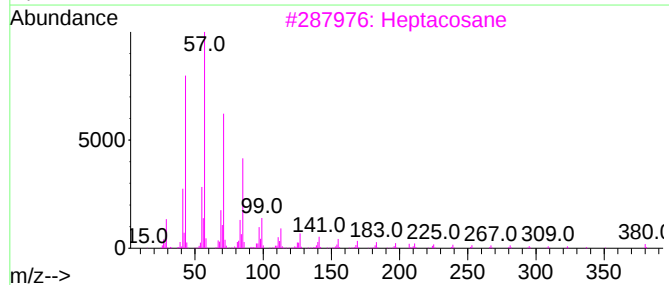
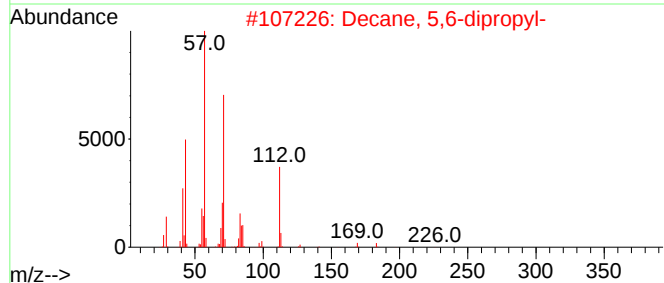
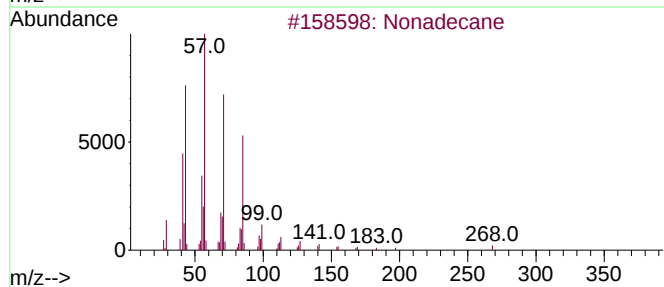
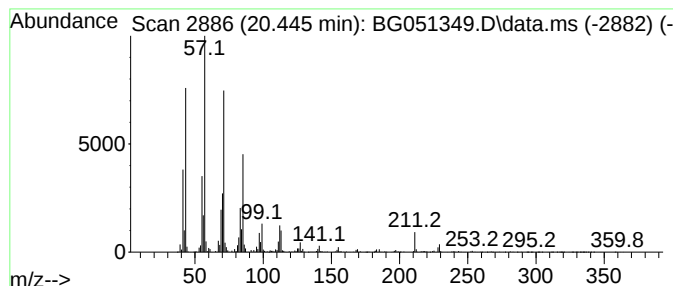
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Quant Title : SVOA CALIBRATION

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

Peak Number 52 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.445	28.36 ng/ul	2990360	Chrysene-d12	21.902

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	81
2			Decane, 5,6-dipropyl-	226	C16H34	119209-20-0	76
3			Heptacosane	380	C27H56	000593-49-7	72
4			Octacosane	394	C28H58	000630-02-4	72
5			Heptacosane	380	C27H56	000593-49-7	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

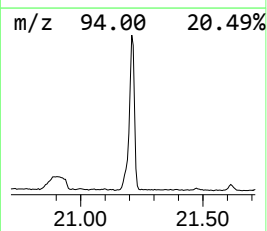
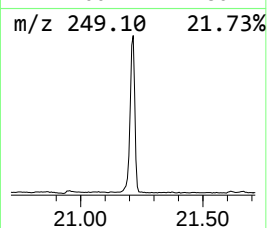
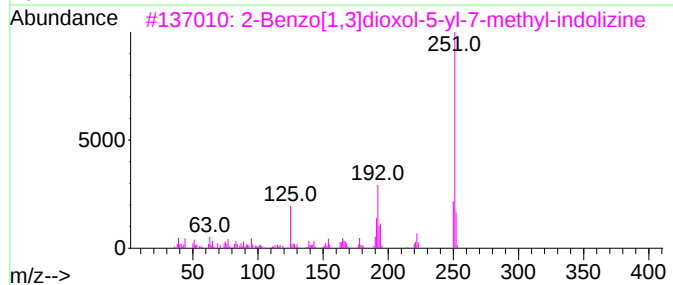
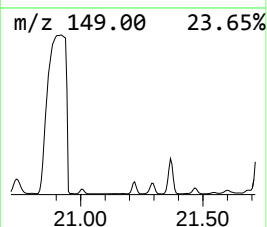
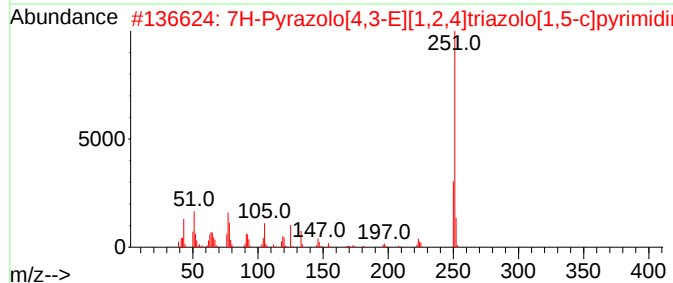
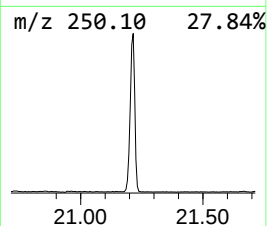
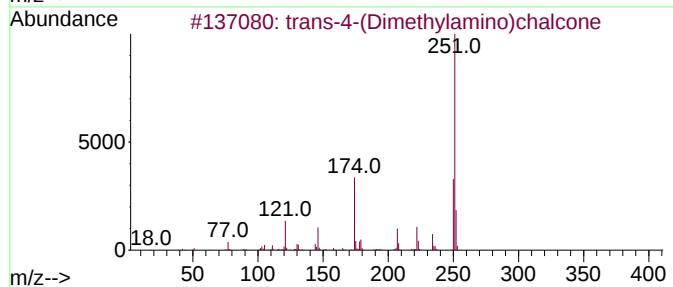
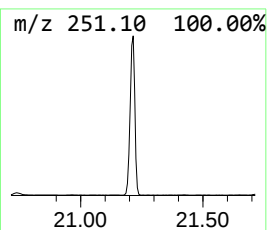
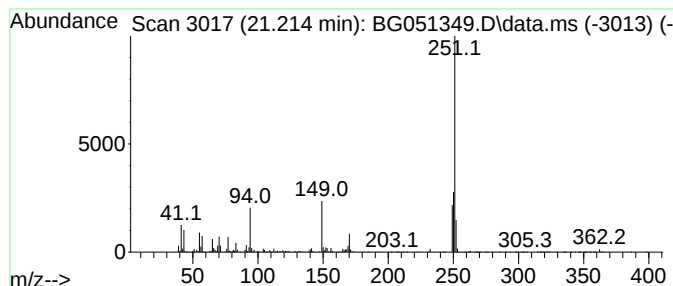
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Quant Title : SVOA CALIBRATION

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

Peak Number 53 unknown-03 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.215	31.31 ng/ul	3301070	Chrysene-d12	21.902

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-4-(Dimethylamino)chalcone	251	C17H17NO	022965-98-6	47
2			7H-Pyrazolo[4,3-E][1,2,4]triazol...	251	C12H9N7	1000310-01-6	47
3			2-Benzo[1,3]dioxol-5-yl-7-methyl...	251	C16H13NO2	1000274-75-9	47
4			N,N-Dimethyl-4-(6-methyl-1H-benz...	251	C16H17N3	069570-95-2	47
5			Phosphoric acid, isodecyl diphen...	390	C22H31O4P	1346599-15-2	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

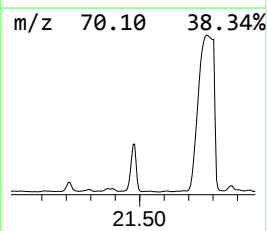
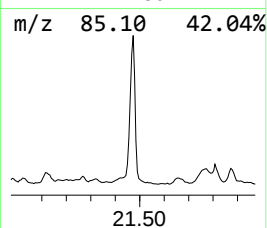
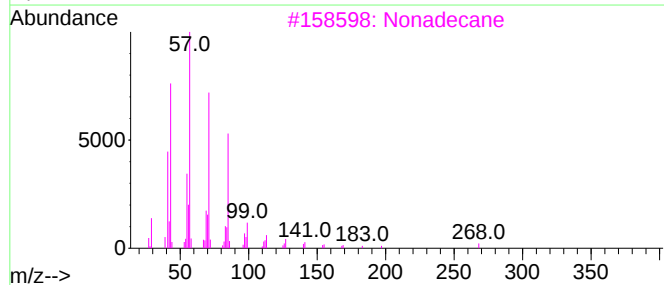
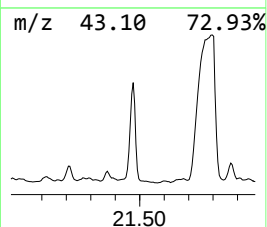
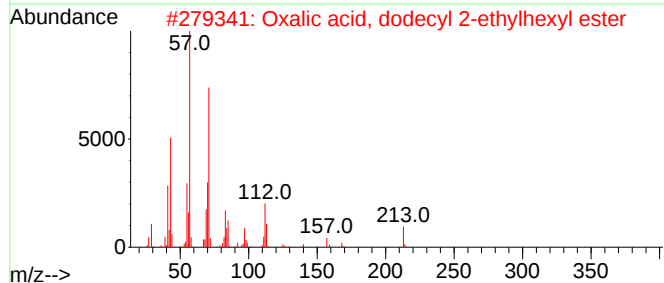
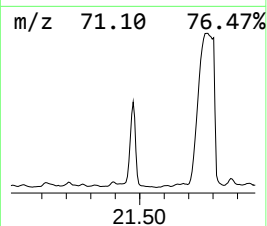
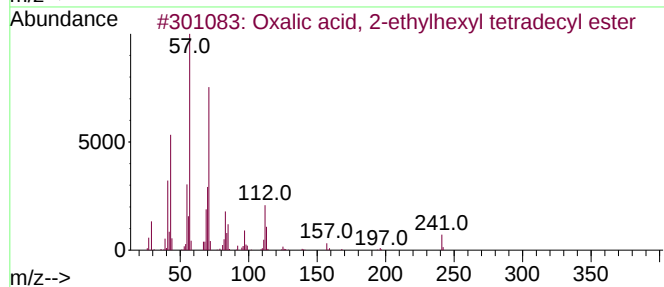
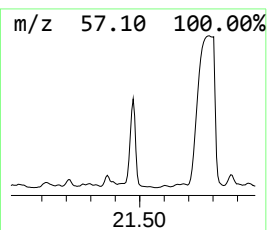
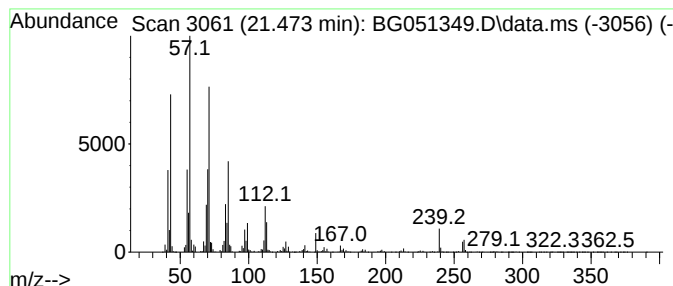
Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 55 Oxalic acid, 2-ethylhexyl t... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.473	58.23 ng/ul	6138430	Chrysene-d12	21.902
Hit#	of 5	Tentative ID	MW MolForm	CAS# Qual
1		Oxalic acid, 2-ethylhexyl tetrad...	398 C24H46O4	1000309-39-7 74
2		Oxalic acid, dodecyl 2-ethylhexy...	370 C22H42O4	1000309-39-5 74
3		Nonadecane	268 C19H40	000629-92-5 58
4		Heneicosane	296 C21H44	000629-94-7 58
5		Nonane, 4-methyl-5-propyl-	184 C13H28	062185-55-1 53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

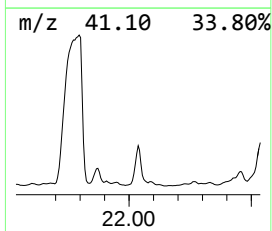
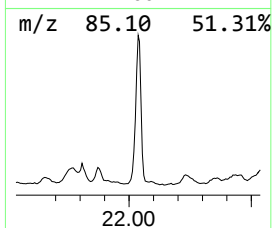
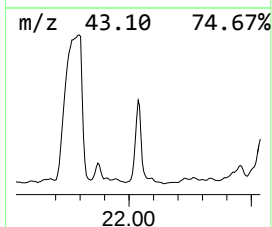
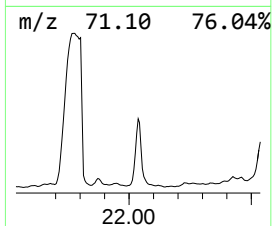
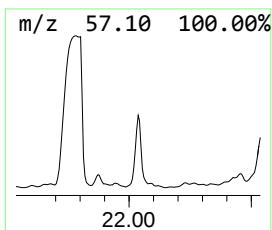
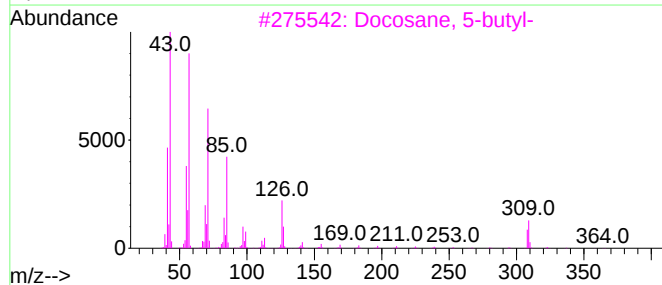
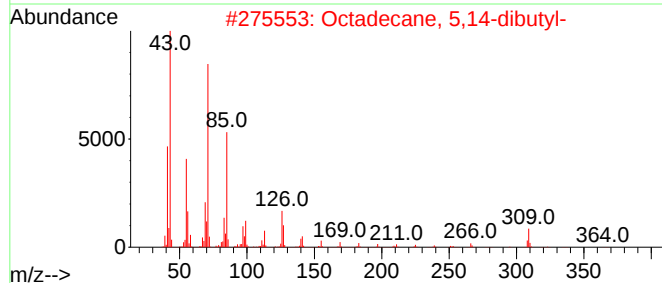
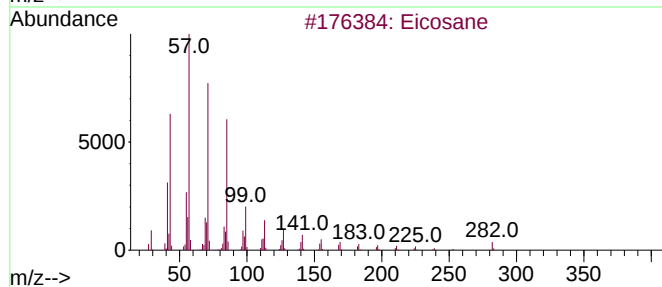
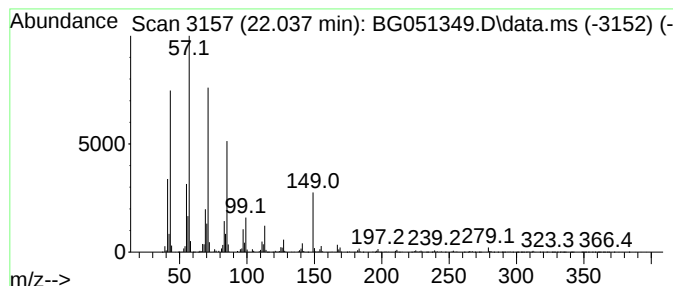
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Quant Title : SVOA CALIBRATION

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

Peak Number 56 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.037	47.25 ng/ul	4981180	Chrysene-d12	21.902

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	96
2		Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	93
3		Docosane, 5-butyl-	366	C26H54	055282-16-1	90
4		Heneicosane	296	C21H44	000629-94-7	87
5		Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

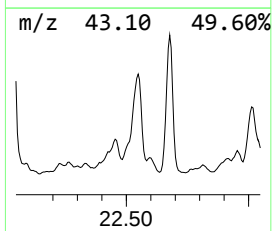
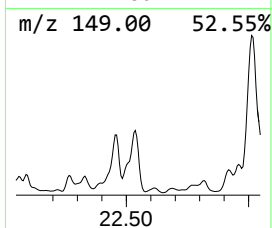
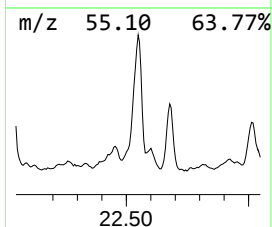
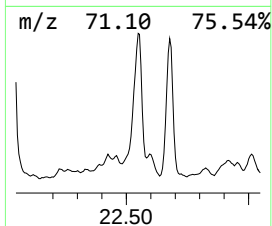
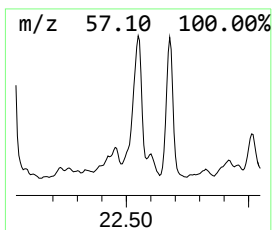
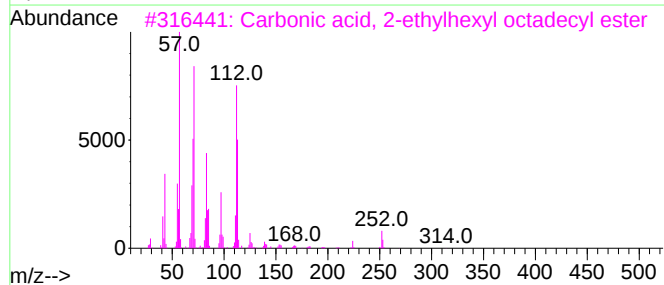
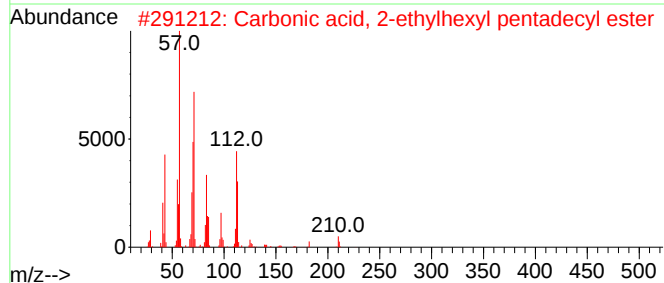
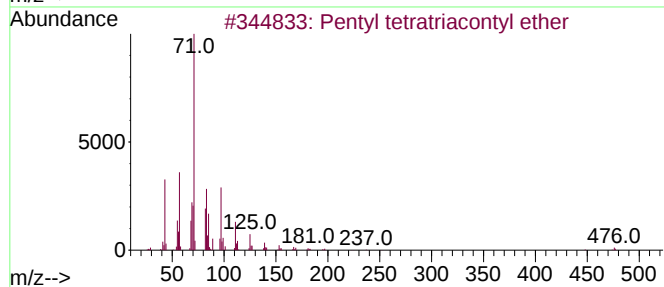
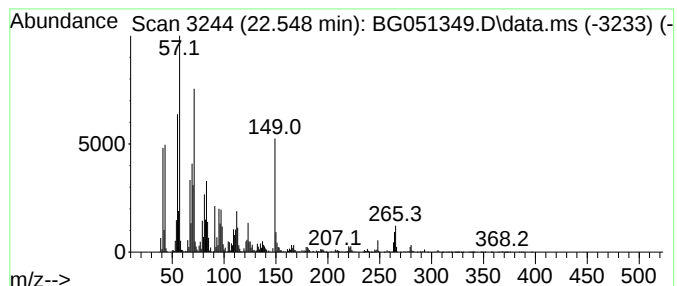
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Quant Title : SVOA CALIBRATION

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

Peak Number 59 unknown-04 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.548	105.35 ng/ul	11106200	Chrysene-d12	21.902

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentyl tetratriacontyl ether	565	C39H80O	1010406-42-7	38
2			Carbonic acid, 2-ethylhexyl pent...	384	C24H48O3	1000383-14-2	38
3			Carbonic acid, 2-ethylhexyl octa...	426	C27H54O3	1000383-14-5	38
4			Dodecyl pentyl ether	256	C17H36O	1010406-41-6	35
5			1-Decene, 4-methyl-	154	C11H22	013151-29-6	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

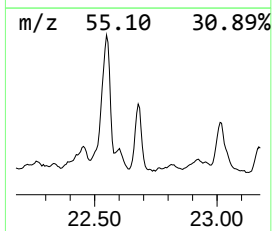
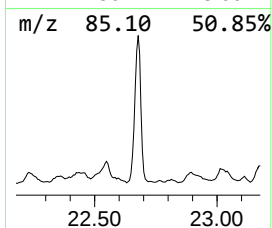
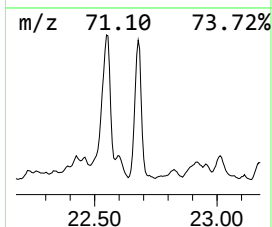
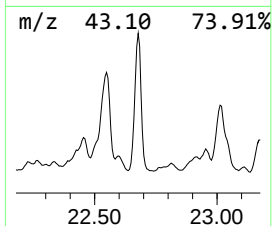
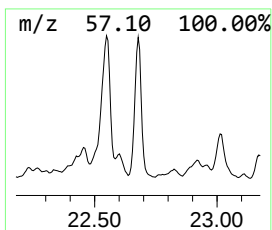
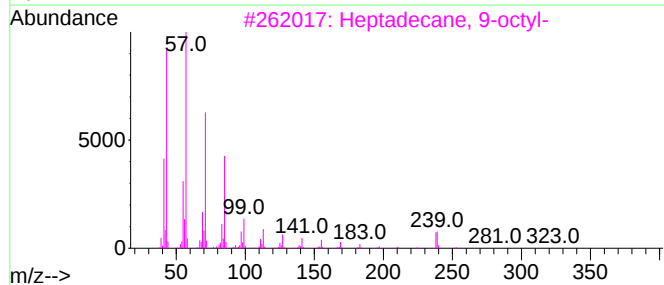
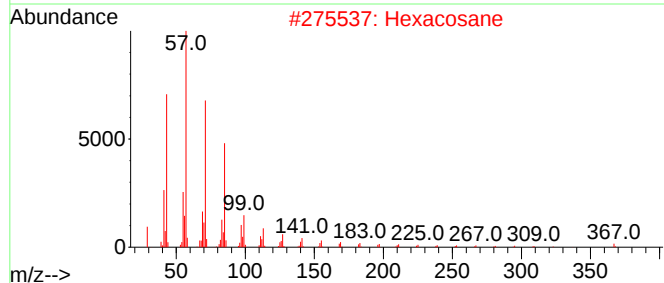
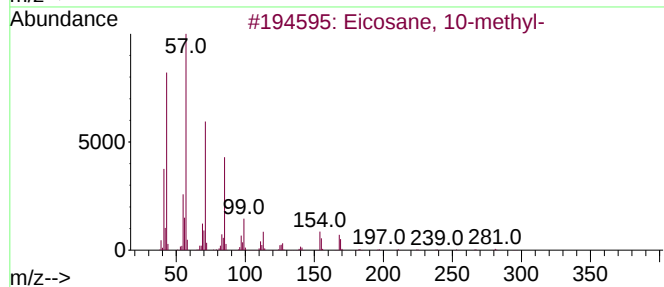
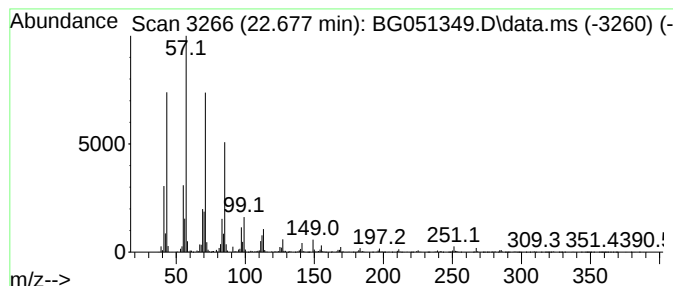
Instrument :
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ClientSampleId :
BGKQ1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 60 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
22.677	48.49 ng/ul	5112540	Chrysene-d12		21.902	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Eicosane, 10-methyl-		296	C21H44	054833-23-7	93
2	Hexacosane		366	C26H54	000630-01-3	91
3	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91
4	Pentadecane, 8-hexyl-		296	C21H44	013475-75-7	91
5	Heneicosane		296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

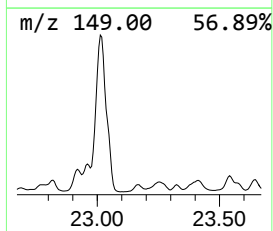
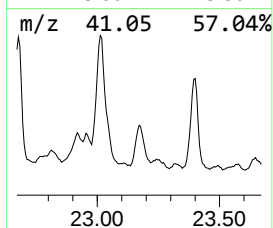
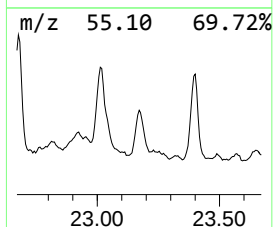
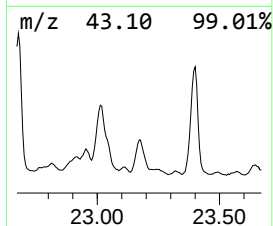
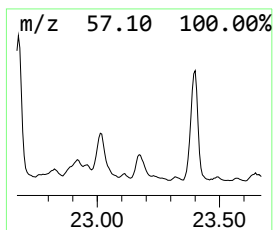
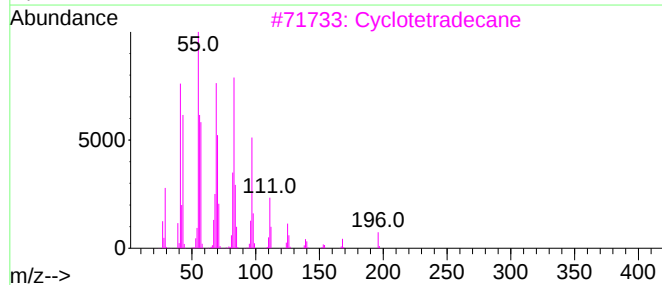
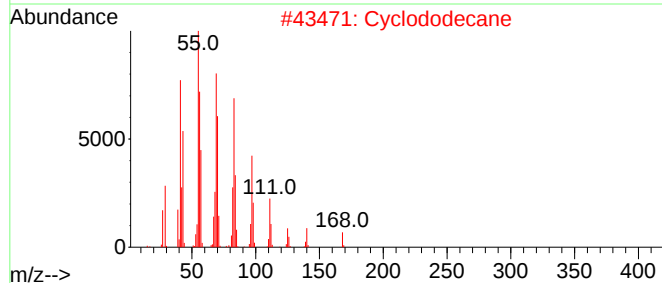
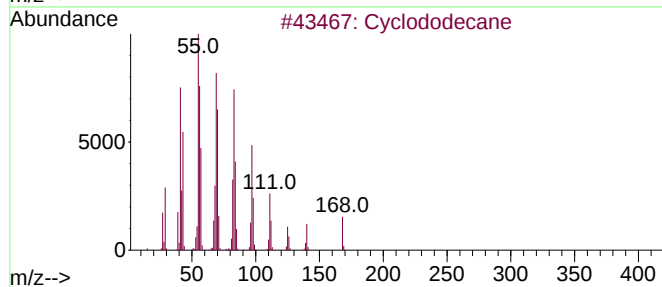
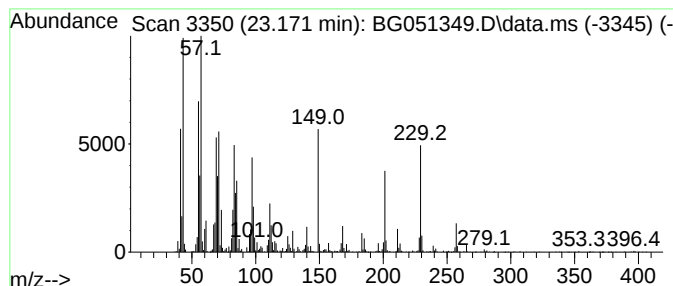
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Quant Title : SVOA CALIBRATION

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

Peak Number 62 (DEL) Alkane: Cyclic23.171 Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.171	18.61 ng/ul	1962200	Chrysene-d12	21.902

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclododecane	168	C12H24	000294-62-2	89
2			Cyclododecane	168	C12H24	000294-62-2	86
3			Cyclotetradecane	196	C14H28	000295-17-0	60
4			1-Tricosene	322	C23H46	018835-32-0	52
5			Dichloroacetic acid, tridecyl ester	310	C15H28Cl2O2	1000280-48-3	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

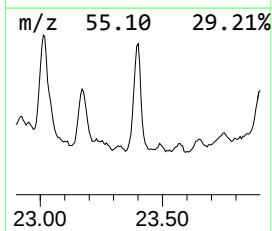
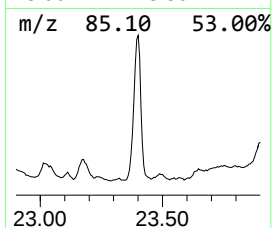
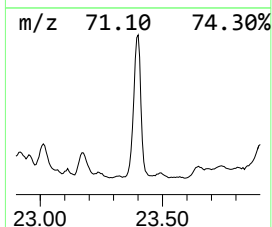
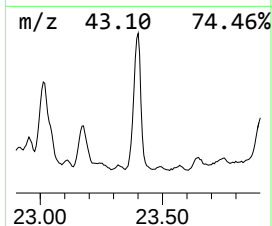
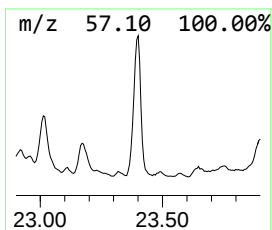
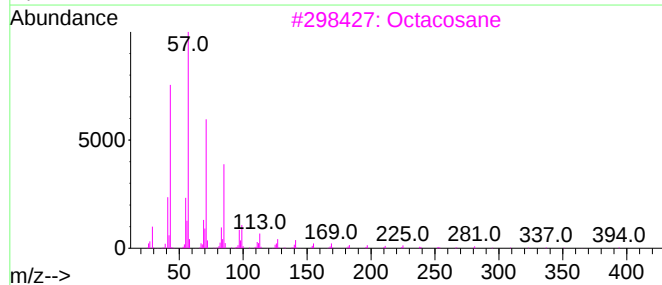
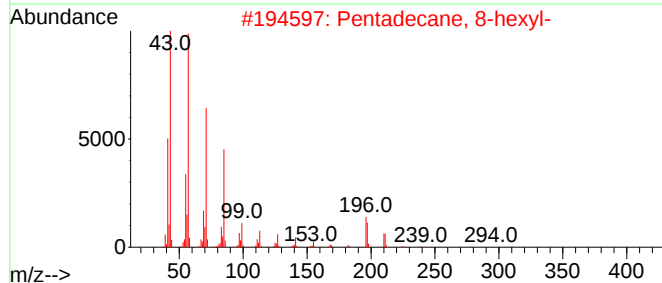
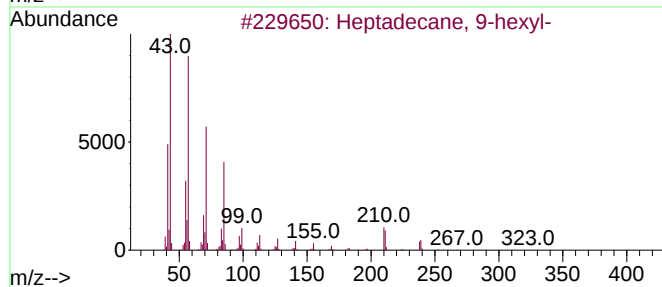
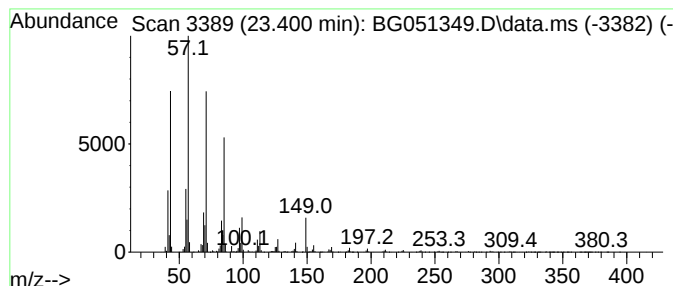
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Quant Title : SVOA CALIBRATION

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

Peak Number 63 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.400	40.23 ng/ul	4241370	Chrysene-d12	21.902

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
2		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
3		Octacosane	394	C28H58	000630-02-4	93
4		Hexacosane	366	C26H54	000630-01-3	91
5		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

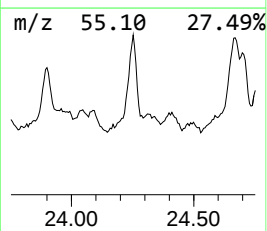
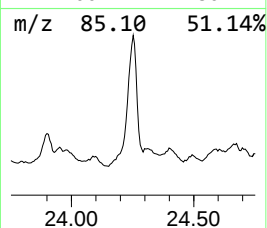
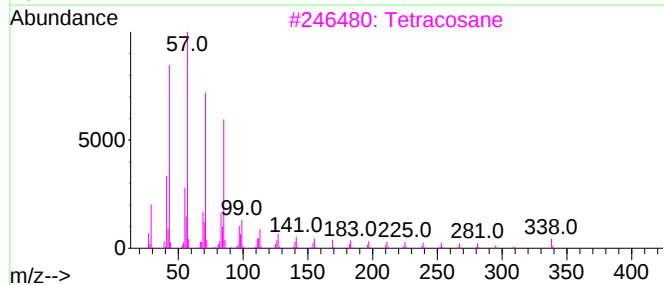
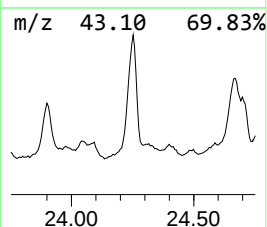
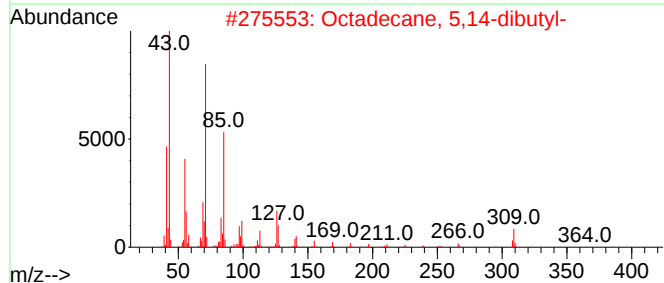
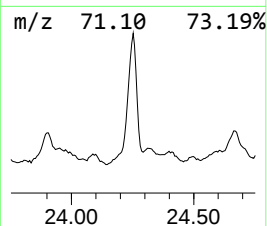
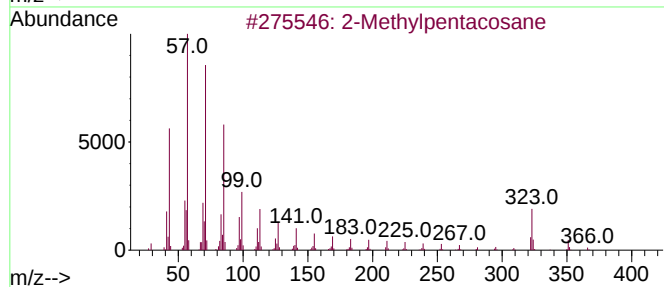
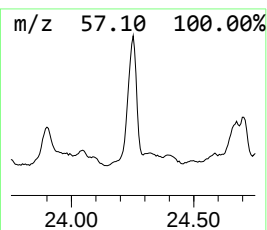
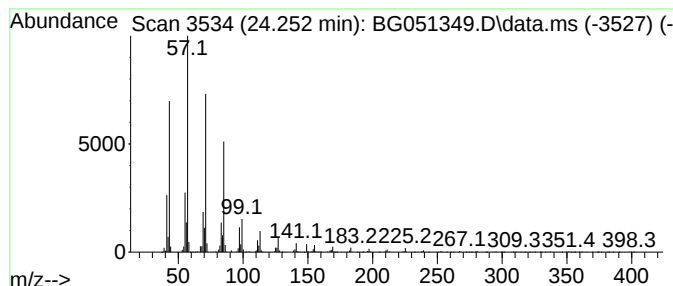
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 66 (DEL) Alkane: Straight-Chai... Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD		R.T.	
24.252	20.32 ng/ul	2755960	Perylene-d12		25.321	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Methylpentacosane		366	C26H54	000629-87-8	95
2	Octadecane, 5,14-dibutyl-		366	C26H54	055282-13-8	94
3	Tetracosane		338	C24H50	000646-31-1	94
4	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	94
5	Hexacosane		366	C26H54	000630-01-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

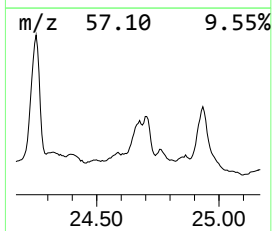
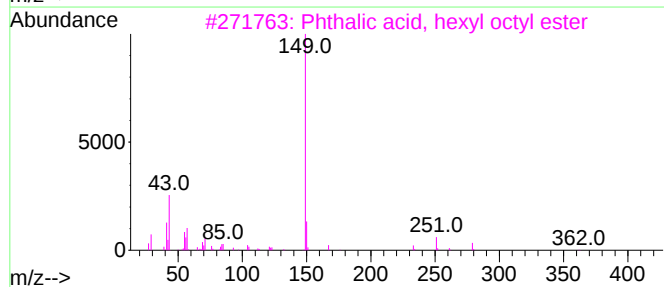
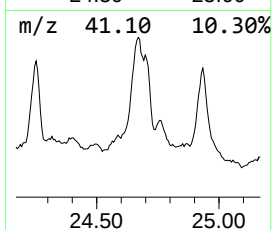
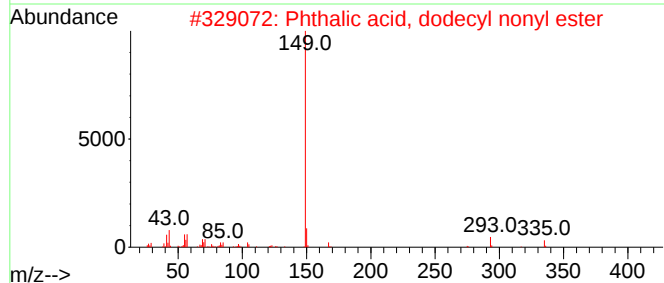
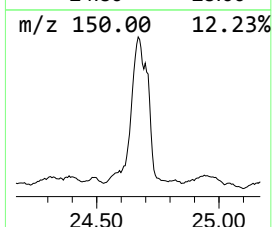
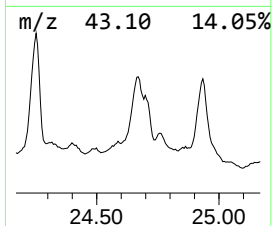
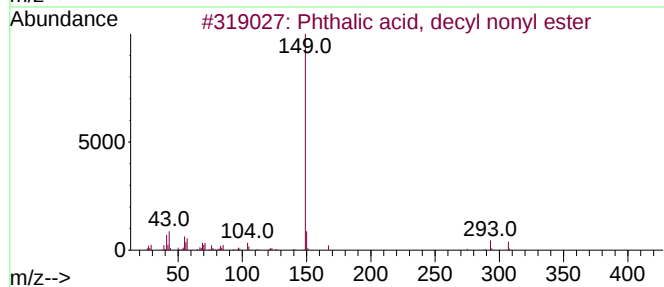
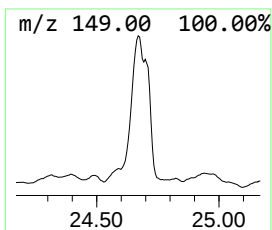
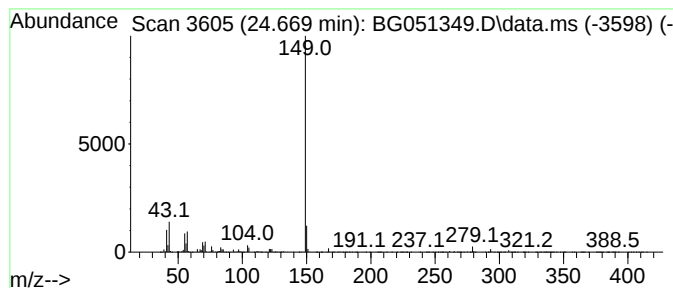
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Quant Title : SVOA CALIBRATION

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

Peak Number 67 Phthalic acid, decyl nonyl ... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.669	26.48 ng/ul	3591010	Perylene-d12	25.321

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, decyl nonyl ester	432	C27H44O4	1000308-89-8	87
2			Phthalic acid, dodecyl nonyl ester	460	C29H48O4	1000308-92-6	86
3			Phthalic acid, hexyl octyl ester	362	C22H34O4	1000424-04-5	86
4			Di-n-octyl phthalate	390	C24H38O4	000117-84-0	86
5			1,2-Benzenedicarboxylic acid, de...	418	C26H42O4	000119-07-3	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

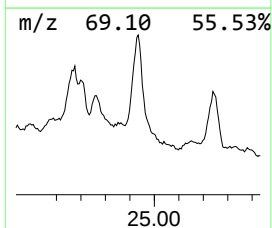
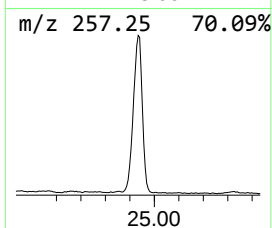
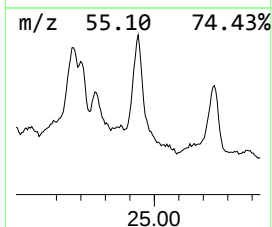
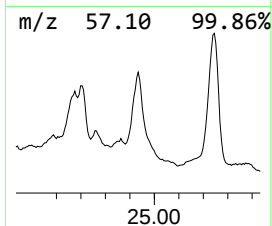
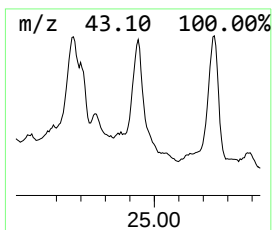
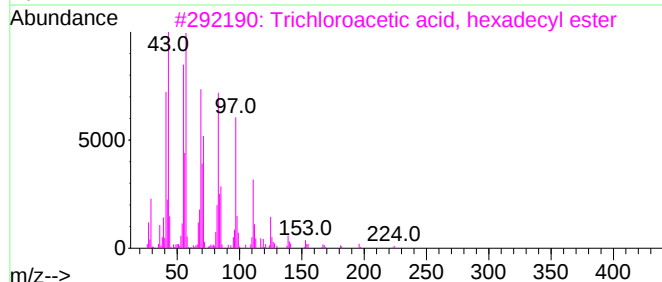
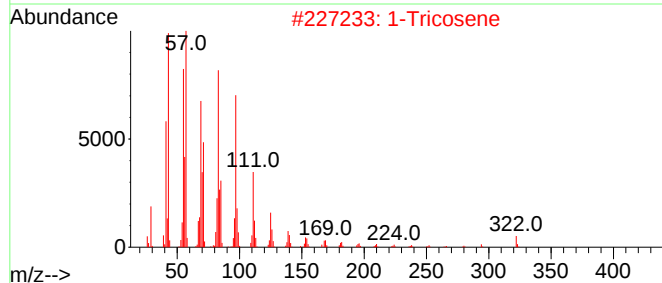
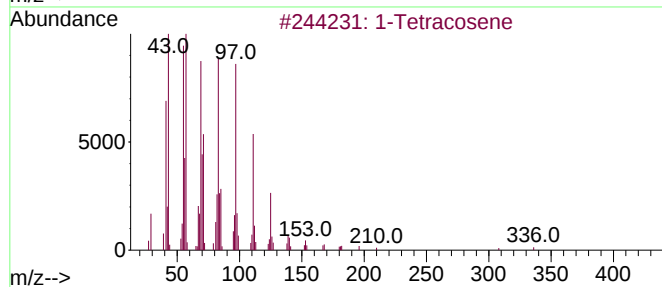
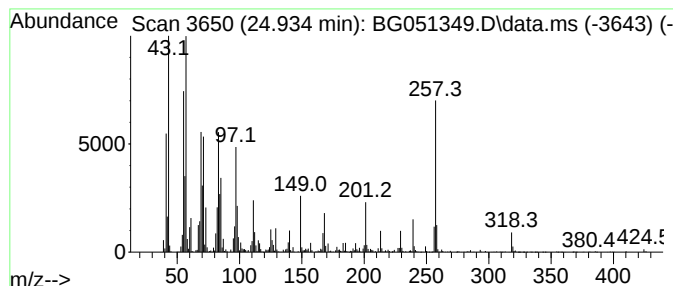
Instrument :
BNA_G
ClientSampleId :
BGKQ1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 68 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.934	36.82 ng/ul	4993310	Perylene-d12	25.321	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetracosene	336	C24H48	010192-32-2	78
2	1-Tricosene	322	C23H46	018835-32-0	64
3	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	64
4	Trifluoroacetic acid,n-tridecyl ...	296	C15H27F3O2	053800-02-5	50
5	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

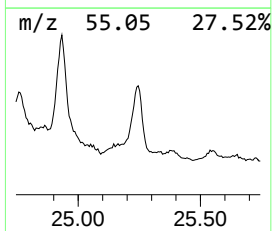
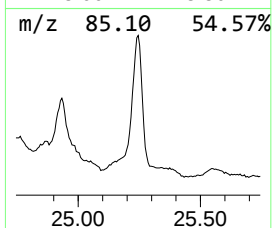
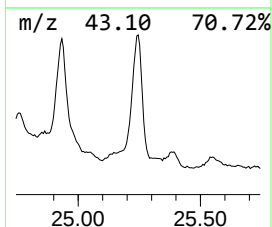
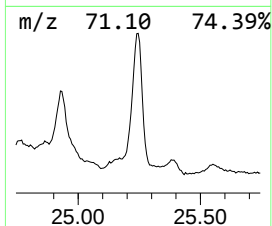
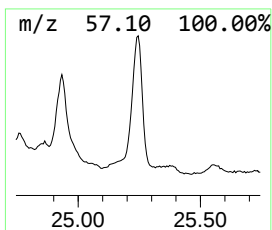
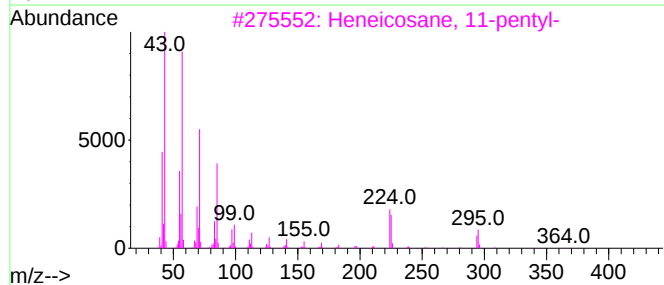
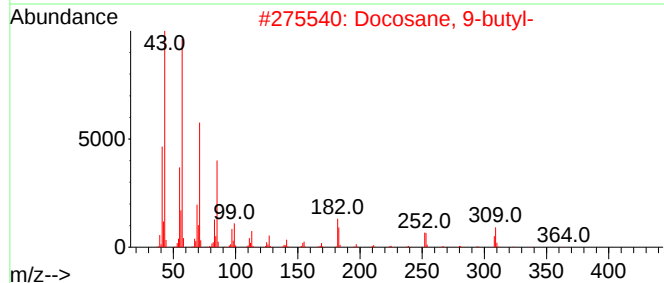
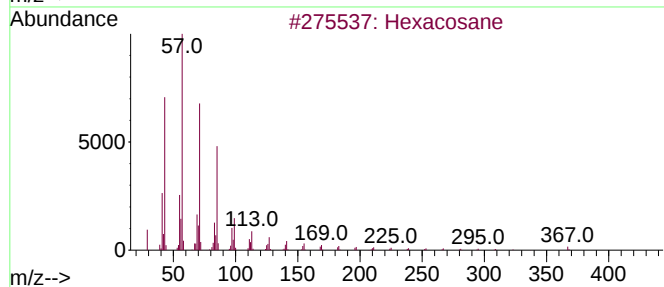
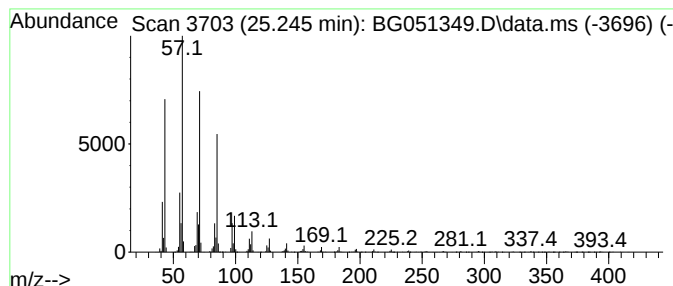
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Quant Title : SVOA CALIBRATION

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

Peak Number 69 (DEL) Alkane: Straight-Chai... Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.245	20.00 ng/ul	2711960	Perylene-d12	25.321

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	94
2			Docosane, 9-butyl-	366	C26H54	055282-14-9	94
3			Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	94
4			Docosane, 11-butyl-	366	C26H54	013475-76-8	93
5			Docosane, 5-butyl-	366	C26H54	055282-16-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

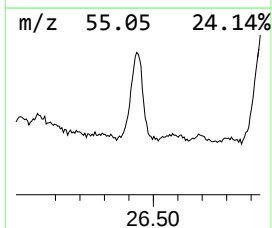
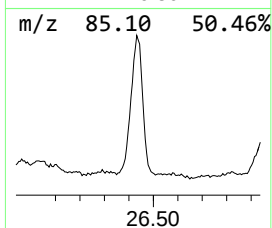
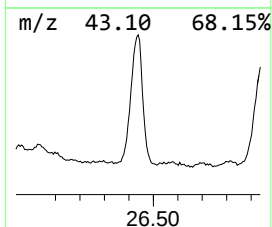
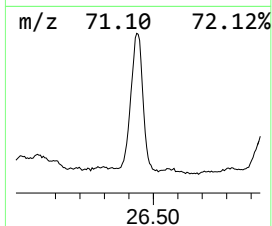
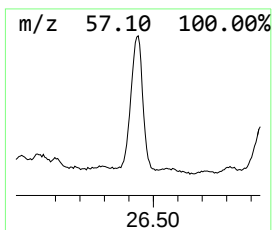
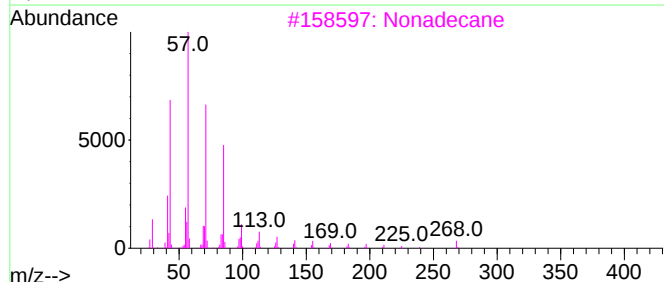
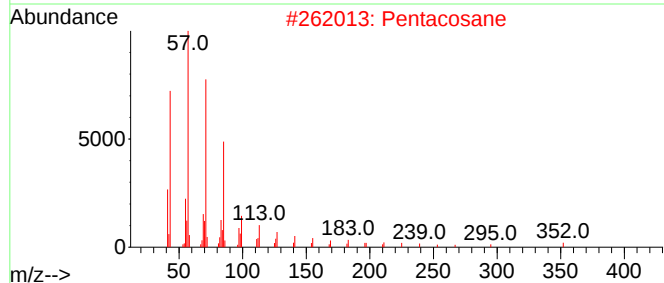
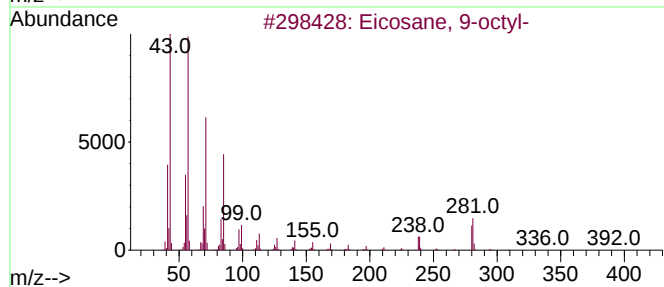
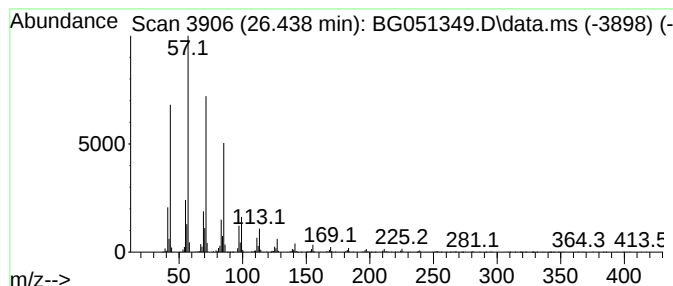
Instrument :
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ClientSampleId :
BGKQ1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

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

Peak Number 70 (DEL) Alkane: Straight-Chai... Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD		R.T.	
26.438	17.57 ng/ul	2382290	Perylene-d12		25.321	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Eicosane, 9-octyl-		394	C28H58	013475-77-9	94
2	Pentacosane		352	C25H52	000629-99-2	94
3	Nonadecane		268	C19H40	000629-92-5	93
4	Pentadecane, 8-heptyl-		310	C22H46	071005-15-7	93
5	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

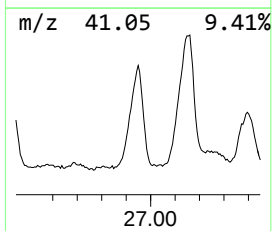
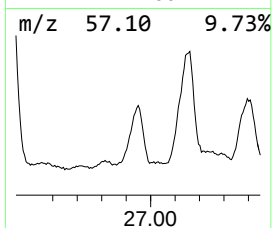
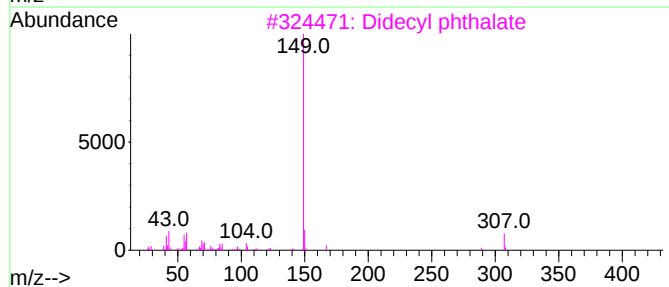
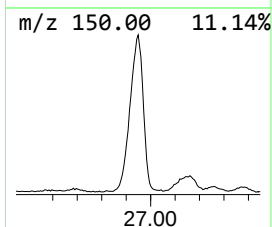
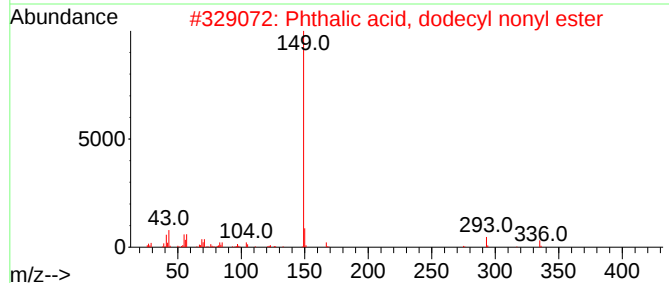
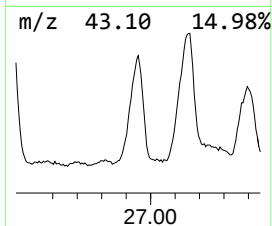
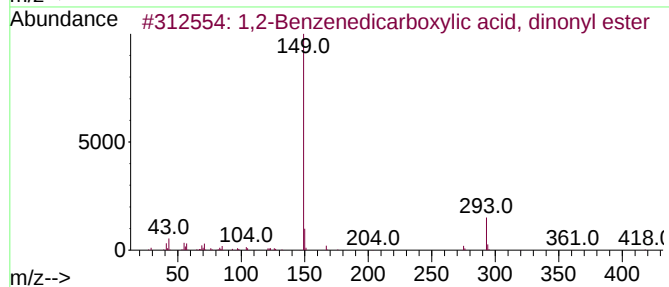
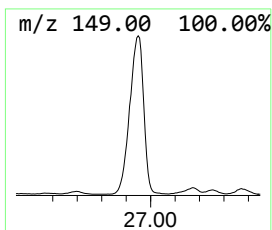
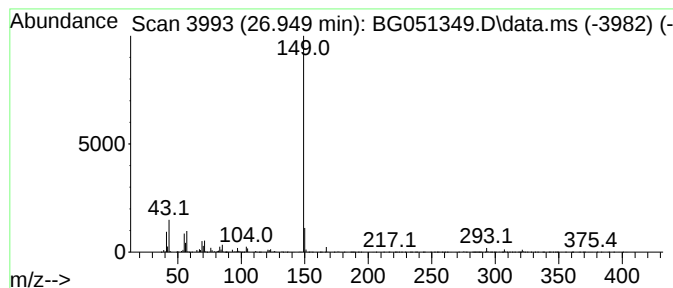
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Quant Title : SVOA CALIBRATION

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

Peak Number 71 1,2-Benzenedicarboxylic aci... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.949	30.30 ng/ul	4108190	Perylene-d12	25.321

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Benzenedicarboxylic acid, di...	418	C26H42O4	000084-76-4	86
2			Phthalic acid, dodecyl nonyl ester	460	C29H48O4	1000308-92-6	86
3			Didecyl phthalate	446	C28H46O4	000084-77-5	86
4			1,2-Benzenedicarboxylic acid, di...	474	C30H50O4	003648-20-2	83
5			1,2-Benzenedicarboxylic acid, di...	418	C26H42O4	000084-76-4	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

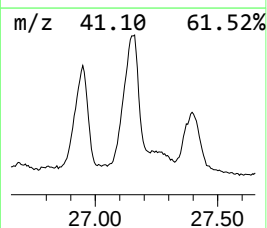
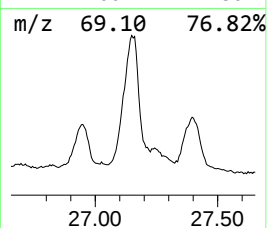
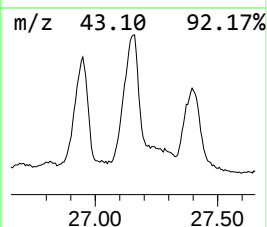
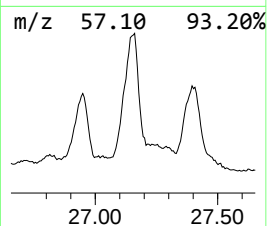
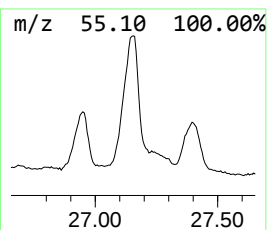
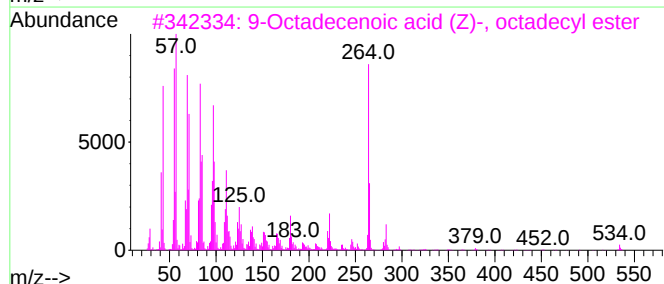
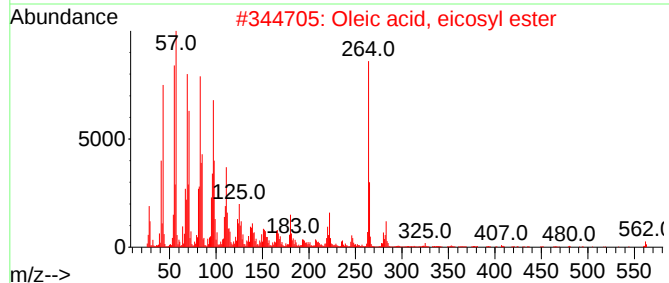
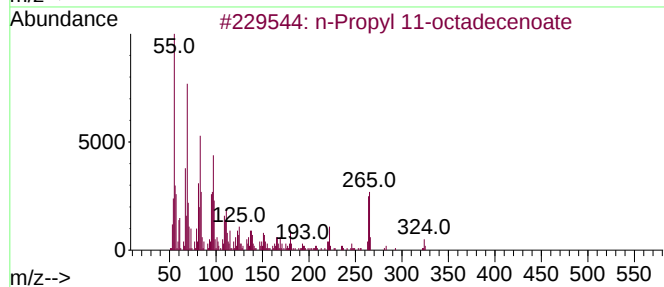
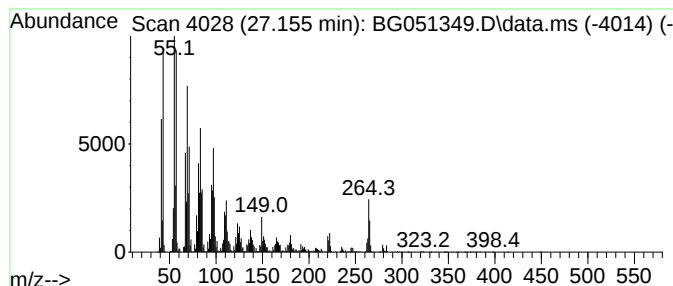
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Quant Title : SVOA CALIBRATION

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

Peak Number 72 n-Propyl 11-octadecenoate Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.155	47.19 ng/ul	6399000	Perylene-d12	25.321

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Propyl 11-octadecenoate	324	C21H40O2	1000336-71-7	91
2			Oleic acid, eicosyl ester	563	C38H74O2	022393-88-0	91
3			9-Octadecenoic acid (Z)-, octade...	535	C36H70O2	017673-49-3	91
4			3-Eicosene, (E)-	280	C20H40	074685-33-9	91
5			Octyl cis-vaccenate	394	C26H50O2	180252-12-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051349.D
Acq On : 6 Dec 2021 14:04
Operator : CG/JU
Sample : M4942-03 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1

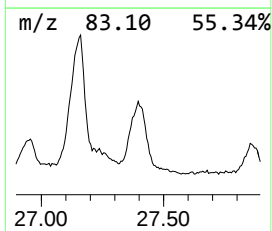
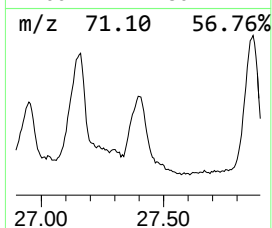
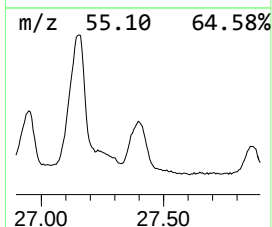
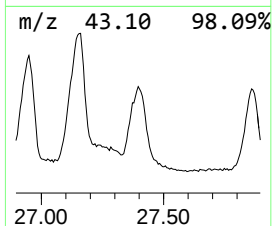
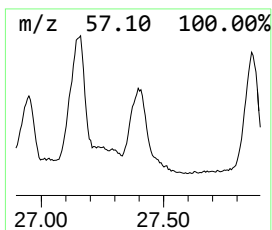
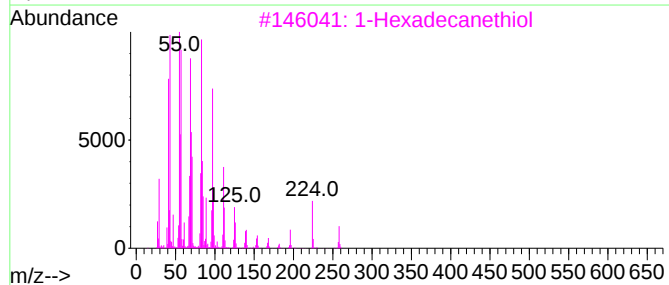
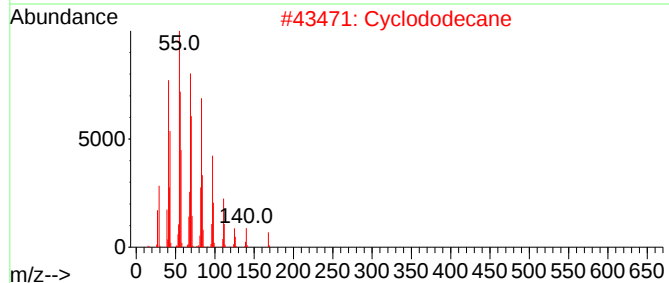
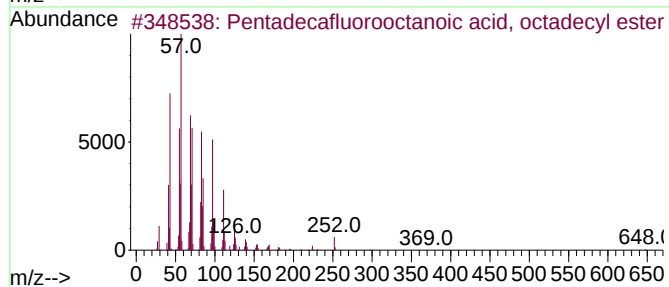
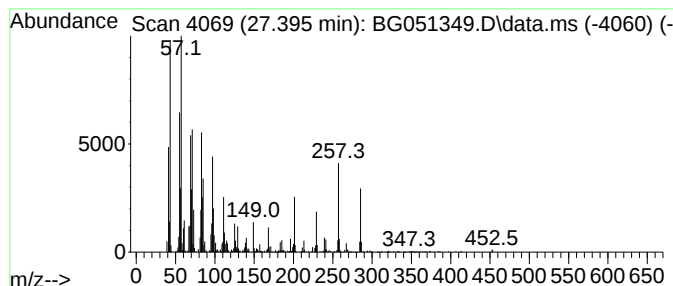
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Quant Title : SVOA CALIBRATION

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

Peak Number 73 Pentadecafluorooctanoic aci... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.395	23.42 ng/ul	3176320	Perylene-d12	25.321

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	89
2			Cyclododecane	168	C12H24	000294-62-2	86
3			1-Hexadecanethiol	258	C16H34S	002917-26-2	86
4			2- Chloropropionic acid, hexadec...	332	C19H37ClO2	086711-81-1	55
5			Carbonic acid, tetradecyl 2,2,2-...	388	C17H31Cl3O3	1000314-56-0	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
 Data File : BG051349.D
 Acq On : 6 Dec 2021 14:04
 Operator : CG/JU
 Sample : M4942-03 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGKQ1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.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-met...	6.649	10.5	ng/ul	866112	1	8.194	1656080	20.0
Benzene, propyl-	7.143	16.6	ng/ul	1370420	1	8.194	1656080	20.0
Benzene, 1-ethy...	7.260	23.1	ng/ul	1911320	1	8.194	1656080	20.0
Benzene, 1-ethy...	7.566	11.9	ng/ul	986617	1	8.194	1656080	20.0
(DEL) Alkane: S...	7.789	56.1	ng/ul	4642290	1	8.194	1656080	20.0
Mesitylene	7.836	45.2	ng/ul	3745170	1	8.194	1656080	20.0
(DEL) Alkane: S...	8.077	10.8	ng/ul	897065	1	8.194	1656080	20.0
(DEL) Alkane: S...	8.142	19.2	ng/ul	1592730	1	8.194	1656080	20.0
(DEL) Alkane: S...	8.694	16.6	ng/ul	1378760	1	8.194	1656080	20.0
4-Methyl-3-hept...	8.753	23.9	ng/ul	1976090	1	8.194	1656080	20.0
Benzene, 1,4-di...	8.829	39.0	ng/ul	3227350	1	8.194	1656080	20.0
Benzene, 1-ethy...	9.293	22.0	ng/ul	1821170	1	8.194	1656080	20.0
unknown-01	9.540	20.4	ng/ul	1692320	1	8.194	1656080	20.0
(DEL) Alkane: S...	10.956	20.0	ng/ul	3977720	2	11.026	3977720	20.0
(DEL) Alkane: S...	11.996	7.1	ng/ul	1418760	2	11.026	3977720	20.0
(DEL) Alkane: S...	12.395	20.0	ng/ul	3971580	2	11.026	3977720	20.0
Benzenamine, 2,...	13.277	34.9	ng/ul	1600400	3	14.828	917054	20.0
(DEL) Alkane: S...	13.330	20.9	ng/ul	959958	3	14.828	917054	20.0
Naphthalene, 2-...	13.858	18.0	ng/ul	823373	3	14.828	917054	20.0
Naphthalene, 2,...	14.005	44.6	ng/ul	2044740	3	14.828	917054	20.0
Naphthalene, 2,...	14.152	46.0	ng/ul	2106870	3	14.828	917054	20.0
(DEL) Alkane: S...	14.270	26.5	ng/ul	1214340	3	14.828	917054	20.0
Naphthalene, 1,...	14.387	19.9	ng/ul	914367	3	14.828	917054	20.0
(DEL) Alkane: S...	14.693	93.5	ng/ul	4288660	3	14.828	917054	20.0
Naphthalene, 2,...	15.298	35.0	ng/ul	1602490	3	14.828	917054	20.0
Naphthalene, 1,...	15.439	20.3	ng/ul	931695	3	14.828	917054	20.0
unknown-02	15.744	21.0	ng/ul	964230	3	14.828	917054	20.0
(DEL) Alkane: S...	16.502	90.6	ng/ul	4245510	4	17.583	937490	20.0
Benzene, (1-met...	16.632	35.2	ng/ul	1648640	4	17.583	937490	20.0
Benzene, (1-eth...	16.779	19.8	ng/ul	927186	4	17.583	937490	20.0
(DEL) Alkane: S...	17.296	66.8	ng/ul	3133450	4	17.583	937490	20.0
(DEL) Alkane: S...	17.343	31.3	ng/ul	1464900	4	17.583	937490	20.0
Benzene, (1-met...	17.448	20.0	ng/ul	937490	4	17.583	937490	20.0
1,2-Benzenedica...	17.830	25.1	ng/ul	1176310	4	17.583	937490	20.0
(DEL) Alkane: S...	18.723	41.4	ng/ul	1940210	4	17.583	937490	20.0
p-Dicyclohexylb...	18.999	27.5	ng/ul	1288330	4	17.583	937490	20.0
(DEL) Alkane: S...	19.346	77.3	ng/ul	3625530	4	17.583	937490	20.0
cyclohexane, [1...	19.546	34.2	ng/ul	1603170	4	17.583	937490	20.0
(DEL) Alkane: S...	19.781	12.7	ng/ul	1339030	5	21.902	2108510	20.0
(DEL) Alkane: S...	20.445	28.4	ng/ul	2990360	5	21.902	2108510	20.0
unknown-03	21.215	31.3	ng/ul	3301070	5	21.902	2108510	20.0
Oxalic acid, 2-...	21.473	58.2	ng/ul	6138430	5	21.902	2108510	20.0
(DEL) Alkane: S...	22.037	47.3	ng/ul	4981180	5	21.902	2108510	20.0
unknown-04	22.548	105.3	ng/ul	11106200	5	21.902	2108510	20.0
(DEL) Alkane: S...	22.677	48.5	ng/ul	5112540	5	21.902	2108510	20.0
(DEL) Alkane: C...	23.171	18.6	ng/ul	1962200	5	21.902	2108510	20.0
(DEL) Alkane: S...	23.400	40.2	ng/ul	4241370	5	21.902	2108510	20.0
(DEL) Alkane: S...	24.252	20.3	ng/ul	2755960	6	25.321	2711960	20.0
Phthalic acid, ...	24.669	26.5	ng/ul	3591010	6	25.321	2711960	20.0
(DEL) Alkane: S...	24.934	36.8	ng/ul	4993310	6	25.321	2711960	20.0
(DEL) Alkane: S...	25.245	20.0	ng/ul	2711960	6	25.321	2711960	20.0
(DEL) Alkane: S...	26.438	17.6	ng/ul	2382290	6	25.321	2711960	20.0

1,2-Benzenedica...	26.949	30.3	ng/ul	4108190	6	25.321	2711960	20.0
n-Propyl 11-oct...	27.155	47.2	ng/ul	6399000	6	25.321	2711960	20.0
Pentadecafluoro...	27.395	23.4	ng/ul	3176320	6	25.321	2711960	20.0

Instrument :

BNA_G

ClientSampleId :

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