

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
 Data File : BG049477.D
 Acq On : 26 Jul 2021 13:50
 Operator : CG/JU
 Sample : M3082-08
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0006

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-BG072421.M
 Title : SVOA CALIBRATION

Signal : TIC: BG049477.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.085	3	8	16	rBV2	77364	188662	44.26%	2.607%
2	3.162	16	21	35	rVB	148123	287033	67.33%	3.967%
3	3.867	137	141	148	rBV2	16195	32968	7.73%	0.456%
4	3.966	154	158	165	rVB5	6547	11111	2.61%	0.154%
5	4.096	176	180	185	rVB5	3039	4959	1.16%	0.069%
6	5.165	358	362	367	rVB4	3174	5295	1.24%	0.073%
7	5.218	367	371	374	rBV3	4764	6405	1.50%	0.089%
8	5.670	444	448	454	rBV4	3636	6304	1.48%	0.087%
9	5.981	493	501	502	rBV4	4096	5201	1.22%	0.072%
10	6.040	506	511	516	rBV4	7795	13337	3.13%	0.184%
11	6.299	550	555	557	rBV4	3572	5575	1.31%	0.077%
12	6.422	572	576	580	rVB2	3203	5304	1.24%	0.073%
13	6.522	588	593	595	rBV4	3380	4958	1.16%	0.069%
14	6.586	599	604	611	rVB	4362	7874	1.85%	0.109%
15	6.675	615	619	622	rBV2	3547	5744	1.35%	0.079%
16	6.851	644	649	655	rVB3	11015	20352	4.77%	0.281%
17	7.021	675	678	684	rBV5	3936	5206	1.22%	0.072%
18	7.203	700	709	716	rBV2	59266	117781	27.63%	1.628%
19	7.256	716	718	722	rVB3	5332	7282	1.71%	0.101%
20	7.368	731	737	743	rBV	36474	61688	14.47%	0.853%
21	7.573	767	772	782	rVV	52480	97145	22.79%	1.343%
22	7.661	782	787	797	rVB4	11707	28251	6.63%	0.390%
23	8.049	843	853	859	rBV3	76844	151967	35.65%	2.100%
24	8.102	860	862	867	rVB5	3935	6221	1.46%	0.086%
25	8.331	898	901	907	rVV3	4188	6938	1.63%	0.096%
26	8.572	936	942	946	rBV6	3746	9366	2.20%	0.129%
27	8.701	958	964	967	rBV6	6554	10627	2.49%	0.147%
28	8.754	967	973	980	rVB	63093	108619	25.48%	1.501%
29	8.889	991	996	1000	rVB5	7160	13628	3.20%	0.188%
30	9.165	1032	1043	1045	rBV4	4619	12708	2.98%	0.176%
31	9.212	1045	1051	1057	rVV	51739	100661	23.61%	1.391%
32	9.277	1058	1062	1067	rVB2	21768	35045	8.22%	0.484%
33	9.371	1071	1078	1080	rBV2	4725	7215	1.69%	0.100%
34	9.506	1098	1101	1104	rBV3	4034	5201	1.22%	0.072%
35	9.700	1128	1134	1139	rVV3	5374	9402	2.21%	0.130%
36	9.782	1144	1148	1151	rVB4	4542	6842	1.60%	0.095%

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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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
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37	9.941	1167	1175	1187	rBV4	49972	102689	24.09%	1.419%
38	10.041	1187	1192	1194	rBV4	3957	5172	1.21%	0.071%
39	10.205	1216	1220	1224	rVB4	3752	6268	1.47%	0.087%
40	10.281	1228	1233	1239	rBV5	7765	13947	3.27%	0.193%
41	10.387	1244	1251	1255	rBV4	3357	6669	1.56%	0.092%
42	10.487	1262	1268	1276	rVB	67969	125292	29.39%	1.732%
43	10.628	1286	1292	1300	rBV	60939	107009	25.10%	1.479%
44	10.828	1321	1326	1328	rBV3	34230	49103	11.52%	0.679%
45	10.869	1328	1333	1339	rVV	102026	192371	45.13%	2.659%
46	10.922	1339	1342	1348	rVB3	11761	18144	4.26%	0.251%
47	11.004	1348	1356	1365	rVB4	42858	102580	24.06%	1.418%
48	11.257	1395	1399	1402	rVV4	2882	5399	1.27%	0.075%
49	11.333	1408	1412	1419	rVV6	4170	7352	1.72%	0.102%
50	11.574	1447	1453	1458	rVV4	7103	14327	3.36%	0.198%
51	11.703	1468	1475	1480	rVB7	5447	12079	2.83%	0.167%
52	11.779	1480	1488	1495	rVB6	8509	17799	4.18%	0.246%
53	11.885	1501	1506	1510	rBV2	27045	44445	10.43%	0.614%
54	11.973	1519	1521	1526	rVB4	5181	5342	1.25%	0.074%
55	12.273	1567	1572	1578	rBV	48778	77622	18.21%	1.073%
56	12.414	1592	1596	1600	rVB4	6981	11561	2.71%	0.160%
57	12.496	1605	1610	1611	rVV2	6402	9948	2.33%	0.137%
58	12.520	1611	1614	1621	rVB3	18419	31886	7.48%	0.441%
59	12.743	1647	1652	1659	rVB2	18275	33257	7.80%	0.460%
60	12.907	1674	1680	1683	rBV6	4603	6804	1.60%	0.094%
61	12.966	1683	1690	1691	rBV6	7069	11777	2.76%	0.163%
62	13.042	1700	1703	1705	rBV4	4597	5113	1.20%	0.071%
63	13.078	1706	1709	1715	rVB6	9160	16090	3.77%	0.222%
64	13.224	1729	1734	1740	rVB3	18925	31221	7.32%	0.431%
65	13.295	1741	1746	1750	rVB5	4846	9152	2.15%	0.126%
66	13.407	1759	1765	1767	rBV6	2990	5326	1.25%	0.074%
67	13.512	1775	1783	1789	rBV2	82025	126862	29.76%	1.753%
68	13.577	1790	1794	1802	rVV2	38499	61110	14.33%	0.845%
69	13.724	1816	1819	1820	rBV3	5361	6103	1.43%	0.084%
70	13.853	1837	1841	1844	rBV3	16933	29303	6.87%	0.405%
71	13.965	1856	1860	1864	rBV6	12833	21109	4.95%	0.292%
72	14.018	1864	1869	1873	rVV3	38632	72774	17.07%	1.006%
73	14.094	1874	1882	1890	rVV	135415	268737	63.04%	3.714%
74	14.164	1890	1894	1898	rVV3	30194	46648	10.94%	0.645%
75	14.211	1898	1902	1905	rVV4	14858	22337	5.24%	0.309%
76	14.247	1905	1908	1912	rVV2	18830	30404	7.13%	0.420%

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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	14.282	1912	1914	1919	rVB2	11878	14416	3.38%	0.199%
78	14.388	1927	1932	1941	rVB	135480	211283	49.56%	2.920%
79	14.523	1951	1955	1960	rBV8	10827	18855	4.42%	0.261%
80	14.581	1960	1965	1974	rVB	101034	136895	32.11%	1.892%
81	14.693	1974	1984	1990	rBV	163038	299158	70.17%	4.135%
82	14.893	2013	2018	2028	rVB4	57252	127606	29.93%	1.764%
83	15.093	2047	2052	2058	rVB3	45130	78333	18.37%	1.083%
84	15.169	2062	2065	2068	rVV2	21021	23768	5.58%	0.328%
85	15.310	2084	2089	2094	rBV2	31750	58202	13.65%	0.804%
86	15.363	2094	2098	2104	rVB2	35563	53085	12.45%	0.734%
87	15.486	2115	2119	2122	rBV3	22532	32413	7.60%	0.448%
88	15.533	2122	2127	2132	rVB	129147	178710	41.92%	2.470%
89	15.686	2149	2153	2158	rVB	166725	249417	58.51%	3.447%
90	15.803	2169	2173	2178	rVB2	43434	67928	15.93%	0.939%
91	16.150	2230	2232	2236	rBV4	18412	22007	5.16%	0.304%
92	16.397	2269	2274	2282	rBV2	180240	360538	84.57%	4.983%
93	16.796	2339	2342	2348	rVB2	50287	64719	15.18%	0.894%
94	17.190	2405	2409	2413	rVB	190693	240482	56.41%	3.324%
95	17.448	2448	2453	2457	rBV	181498	287448	67.43%	3.973%
96	17.542	2465	2469	2478	rVB	177710	285578	66.99%	3.947%
97	17.930	2531	2535	2541	rVB	234501	305380	71.63%	4.221%
98	18.329	2599	2603	2606	rBV3	171605	243352	57.08%	3.363%
99	18.623	2649	2653	2658	rBV	209732	265691	62.32%	3.672%
100	19.827	2854	2858	2863	rBV2	278313	426303	100.00%	5.892%

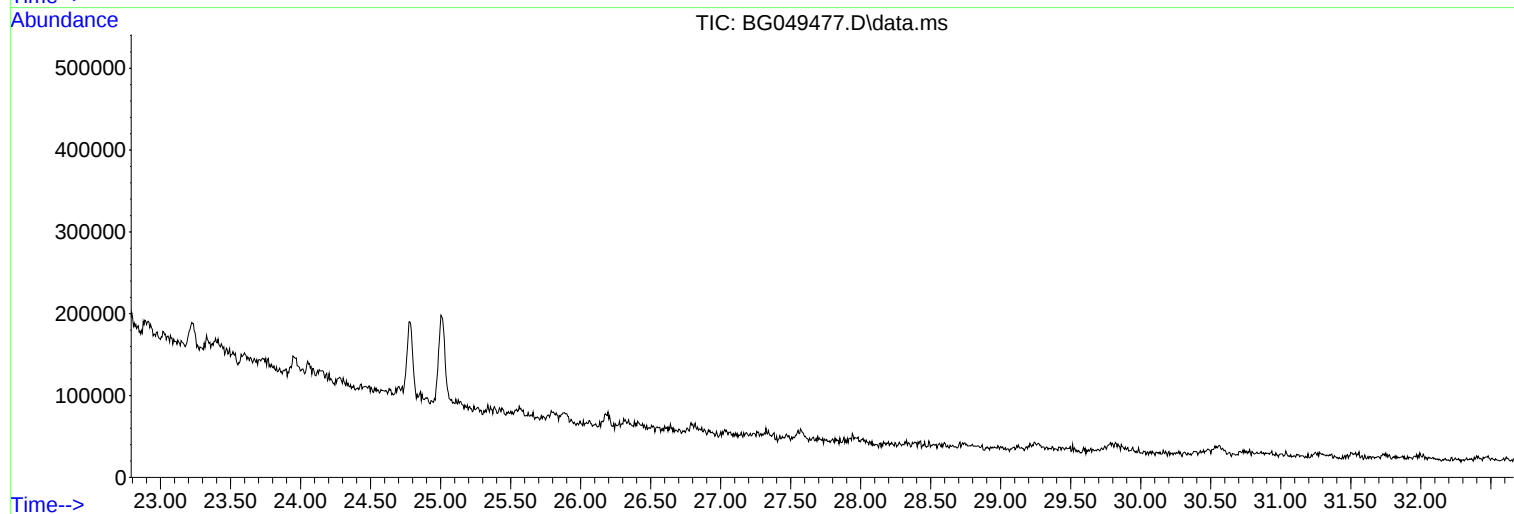
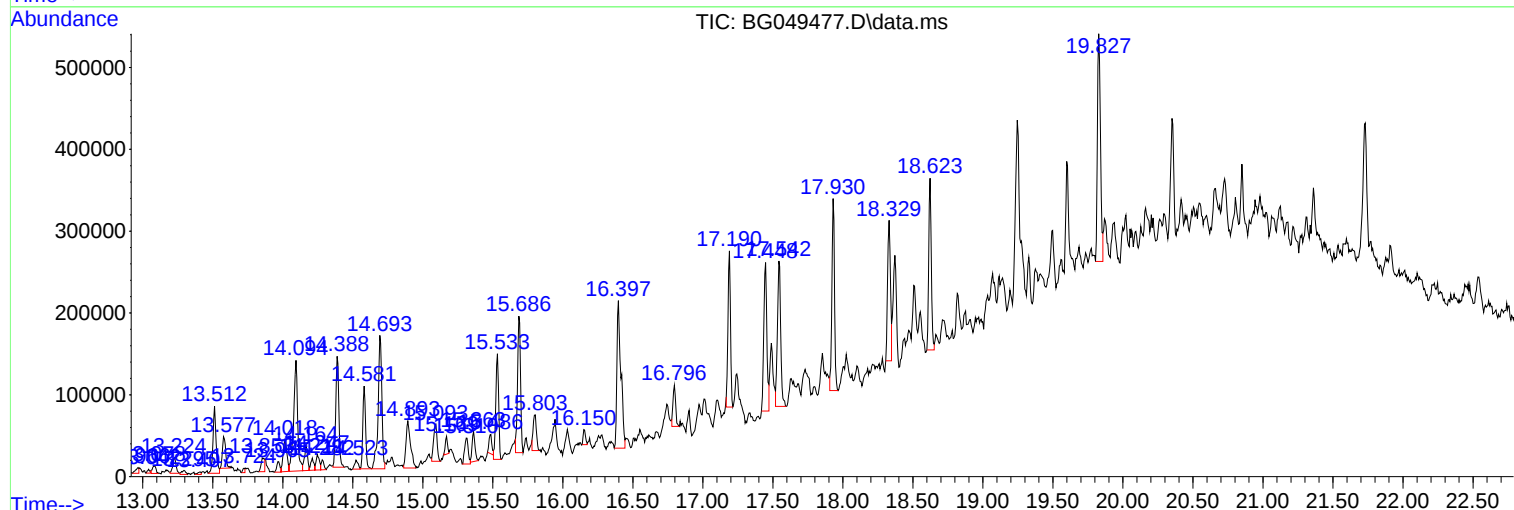
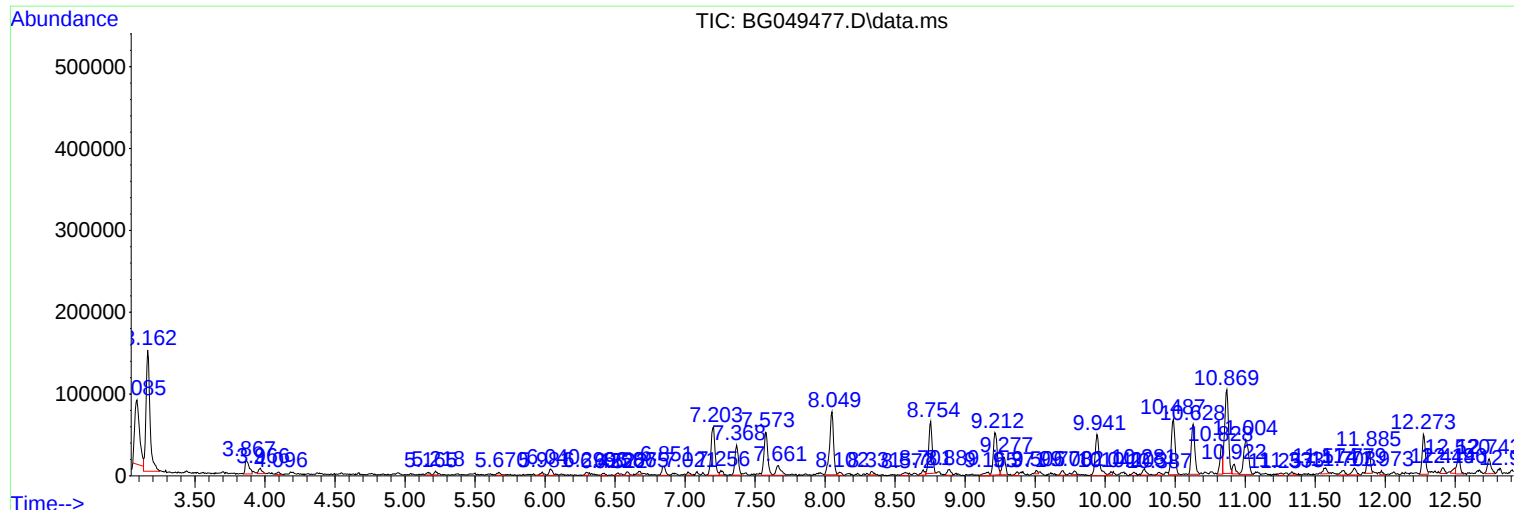
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION

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



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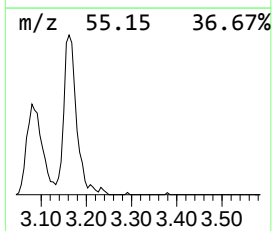
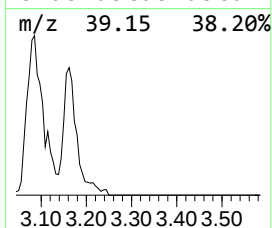
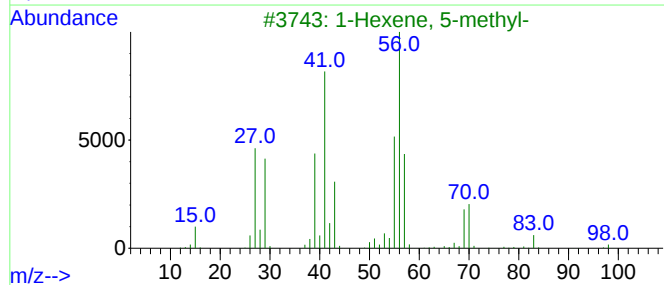
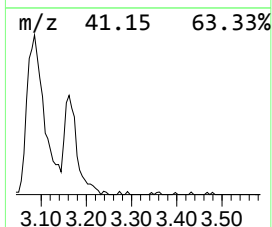
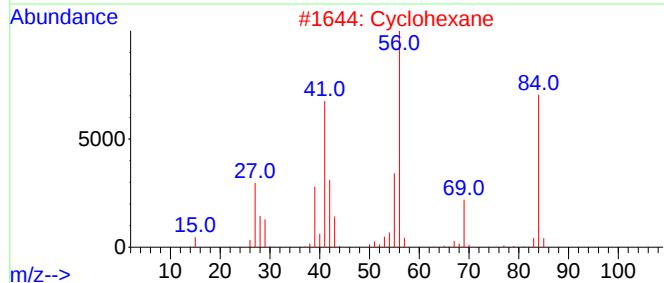
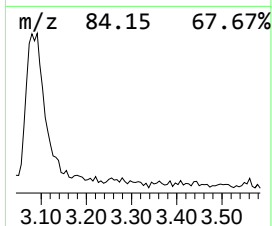
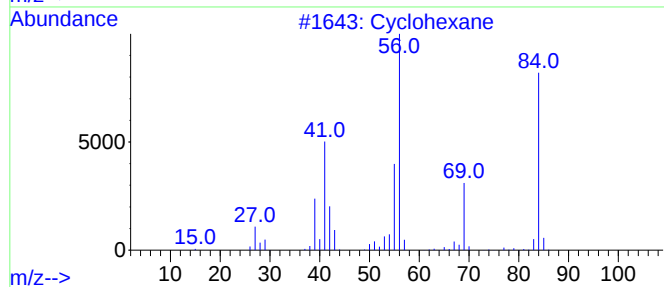
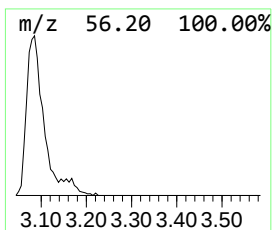
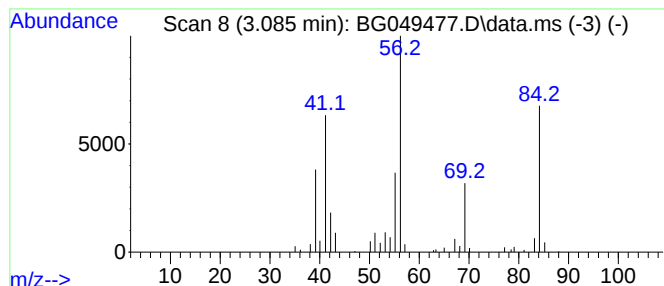
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic3.085 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.085	24.83 ng/ul	188662	1,4-Dichlorobenzene-d4	8.049

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane	84	C6H12	000110-82-7	81
2			Cyclohexane	84	C6H12	000110-82-7	62
3			1-Hexene, 5-methyl-	98	C7H14	003524-73-0	59
4			Cyclohexane	84	C6H12	000110-82-7	58
5			Cyclohexane	84	C6H12	000110-82-7	52



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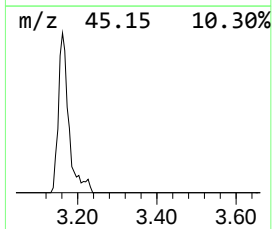
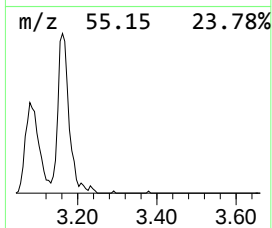
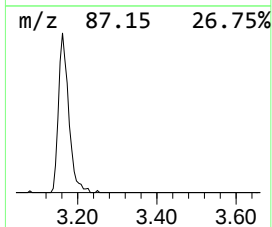
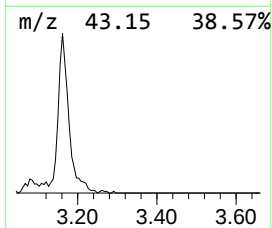
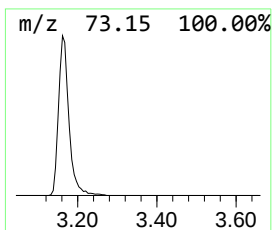
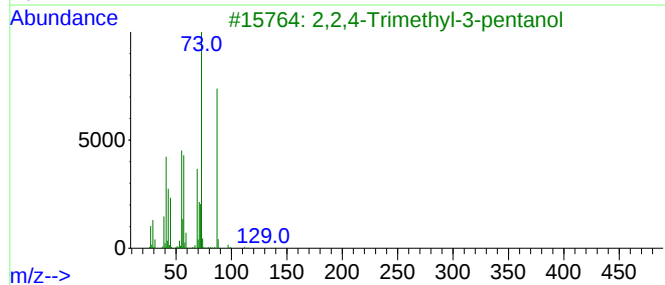
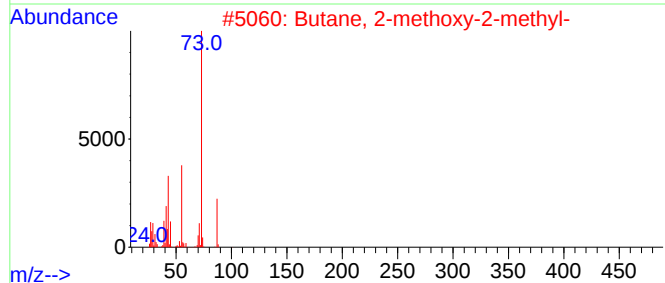
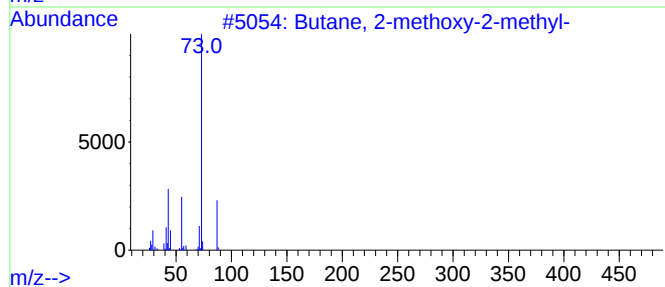
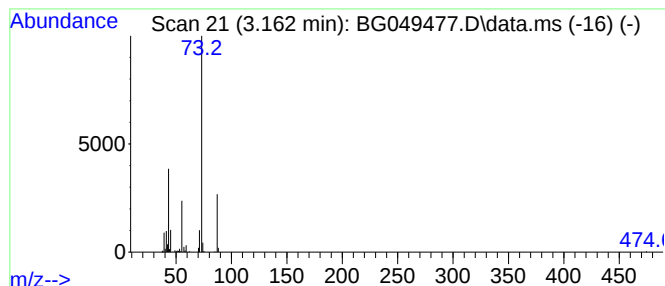
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.162	37.78 ng/ul	287033	1,4-Dichlorobenzene-d4	8.049

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	64
4			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	64
5			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	33



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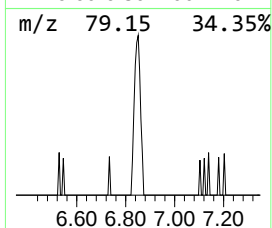
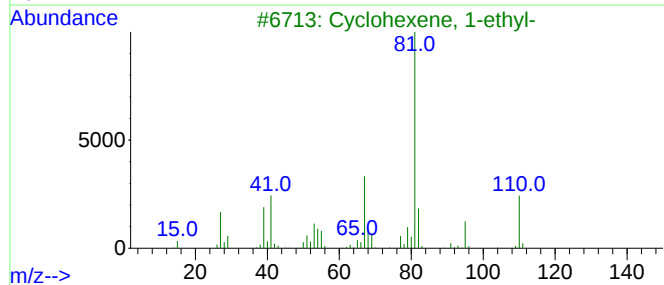
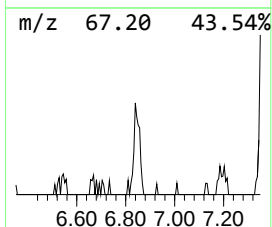
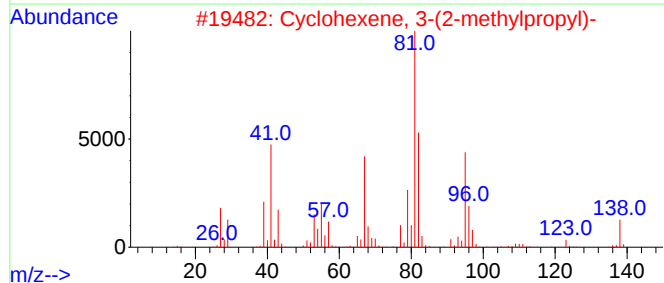
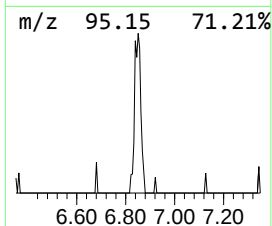
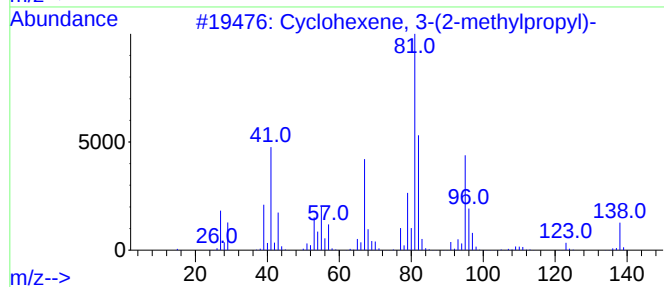
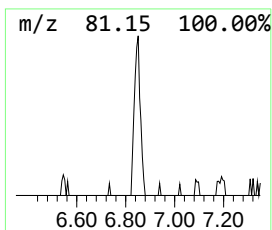
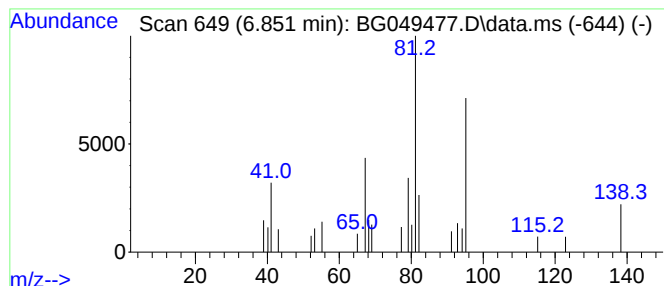
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Cyclohexene, 3-(2-methylpro... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.851	2.68 ng/ul	20352	1,4-Dichlorobenzene-d4	8.049

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene, 3-(2-methylpropyl)-	138	C10H18	004104-56-7	64
2			Cyclohexene, 3-(2-methylpropyl)-	138	C10H18	004104-56-7	64
3			Cyclohexene, 1-ethyl-	110	C8H14	001453-24-3	50
4			2-Chloro-1-isopropyl-4-methylcyc...	174	C10H19Cl	080483-50-7	50
5			2-(Hydroxyimino)cyclohexan-1-amine	128	C6H12N2O	034546-14-0	50



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Operator : CG/JU
Sample : M3082-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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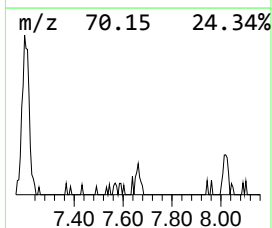
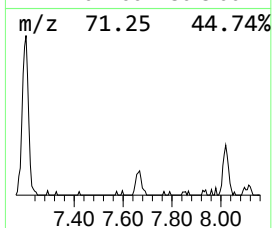
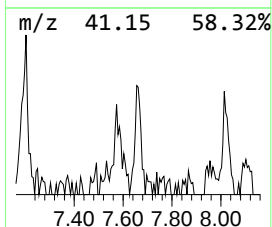
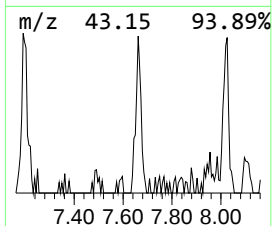
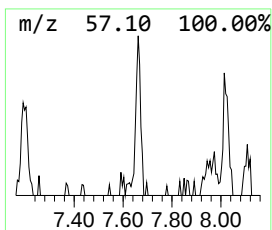
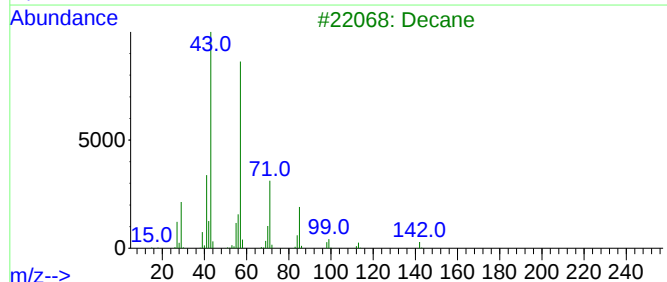
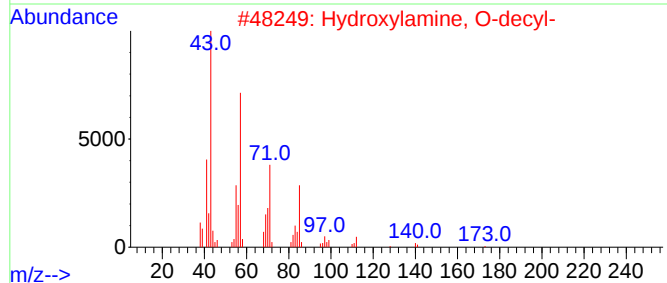
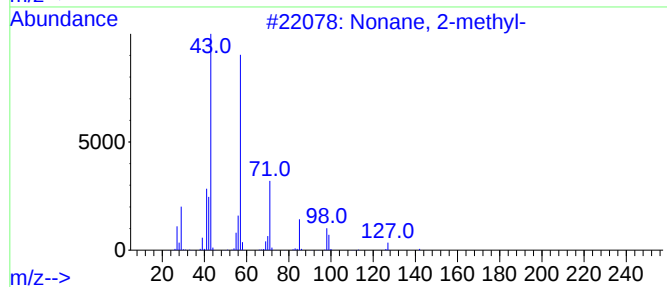
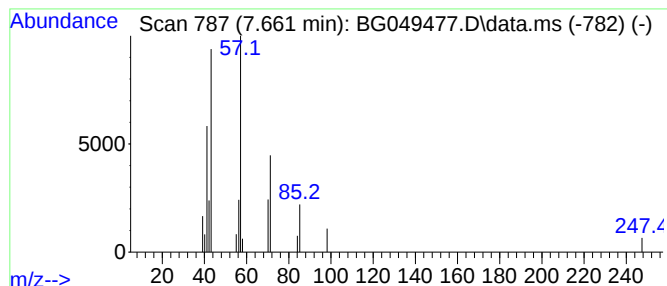
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.661	3.72 ng/ul	28251	1,4-Dichlorobenzene-d4	8.049

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonane, 2-methyl-	142	C10H22	000871-83-0	72
2		Hydroxylamine, O-decyl-	173	C10H23NO	029812-79-1	72
3		Decane	142	C10H22	000124-18-5	64
4		Decane, 2,9-dimethyl-	170	C12H26	001002-17-1	64
5		Decane	142	C10H22	000124-18-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
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Misc :
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Instrument :
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ClientSampleId :
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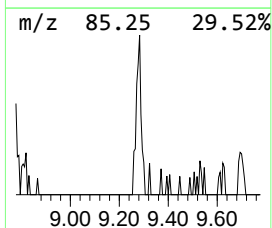
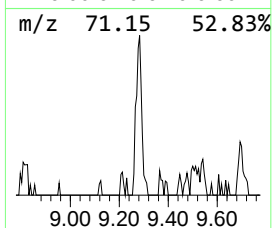
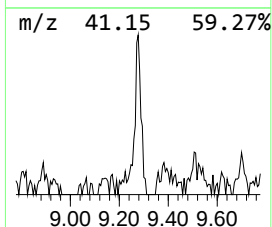
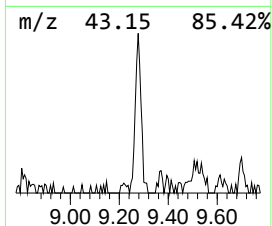
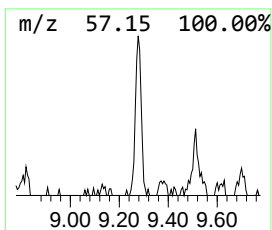
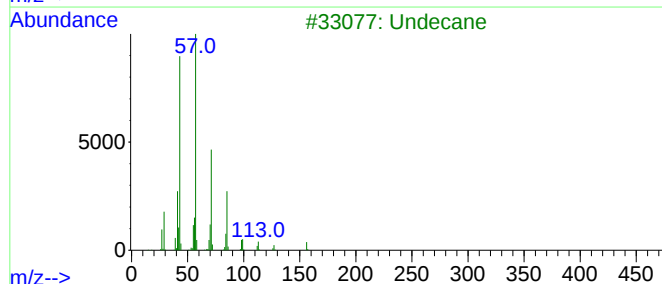
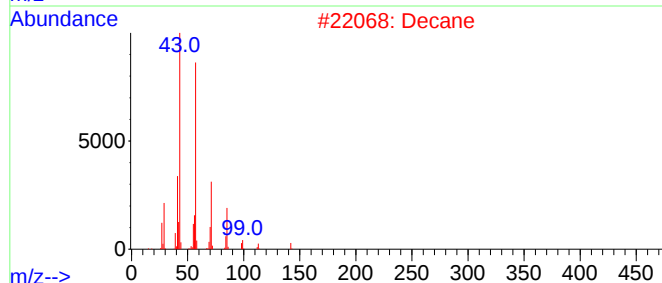
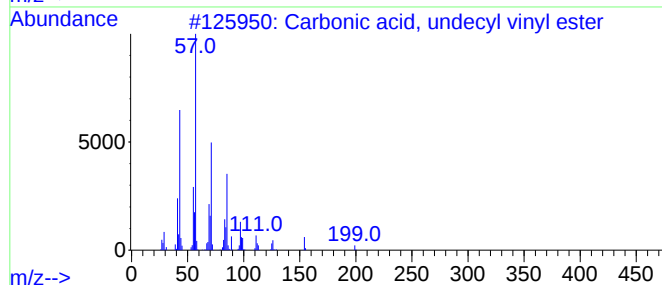
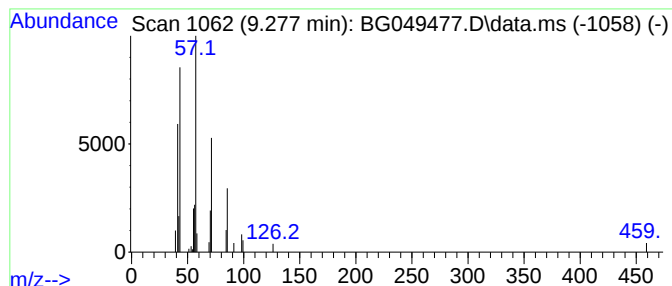
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Carbonic acid, undecyl vinyl... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.277	4.61 ng/ul	35045	1,4-Dichlorobenzene-d4	8.049

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, undecyl vinyl ester	242	C14H26O3	1000382-54-9	83
2			Decane	142	C10H22	000124-18-5	72
3			Undecane	156	C11H24	001120-21-4	72
4			Tetradecane	198	C14H30	000629-59-4	72
5			Tridecane	184	C13H28	000629-50-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049477.D
Acq On : 26 Jul 2021 13:50
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Sample : M3082-08
Misc :
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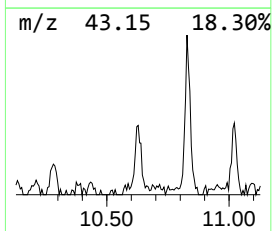
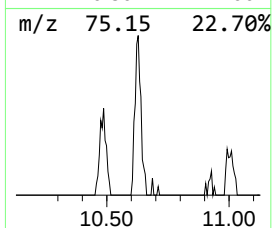
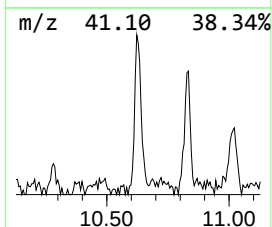
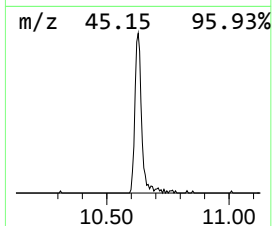
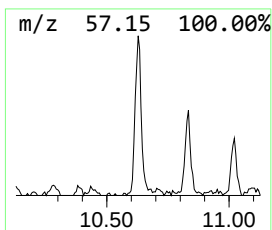
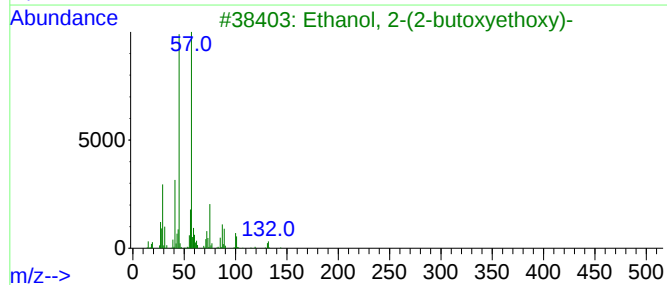
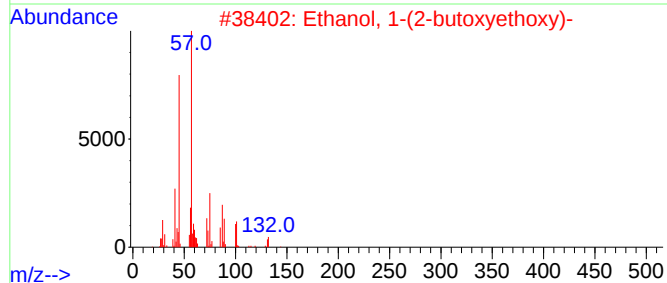
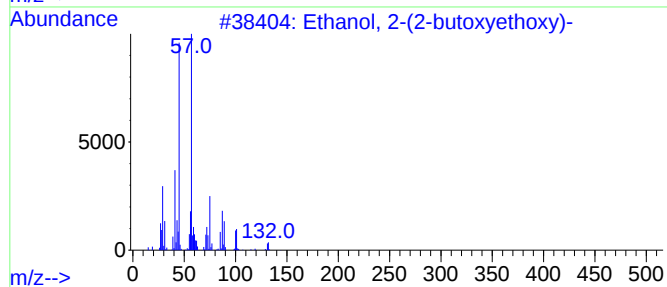
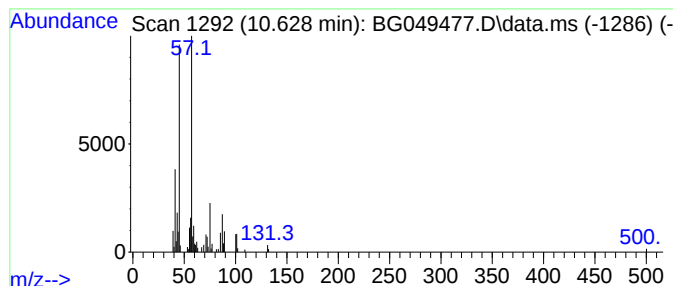
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.628	11.13 ng/ul	107009	Naphthalene-d8	10.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
5			Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049477.D
Acq On : 26 Jul 2021 13:50
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Sample : M3082-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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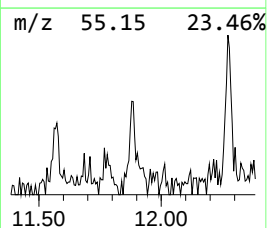
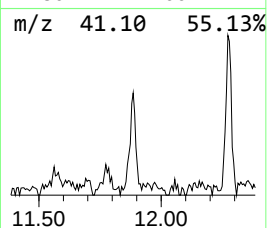
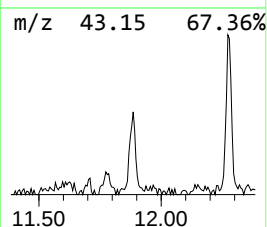
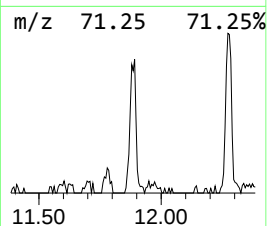
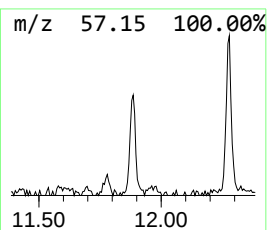
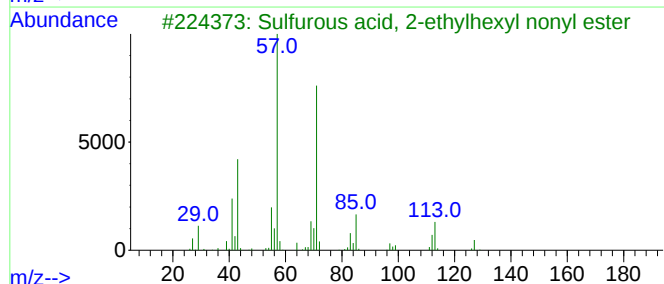
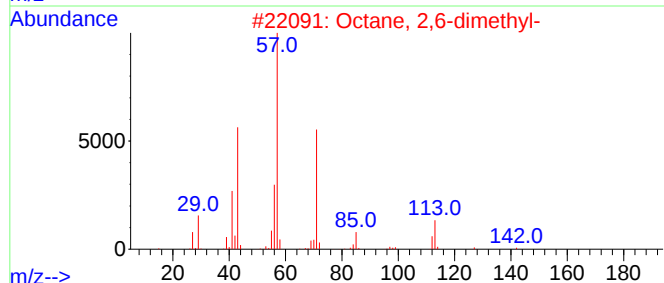
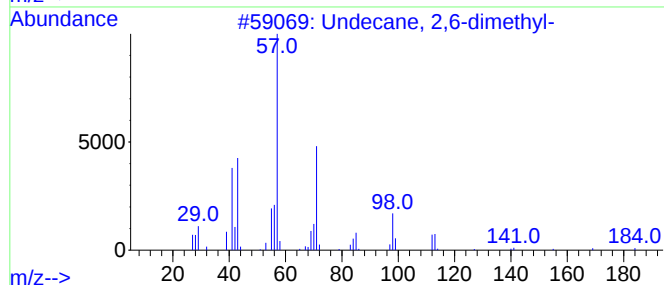
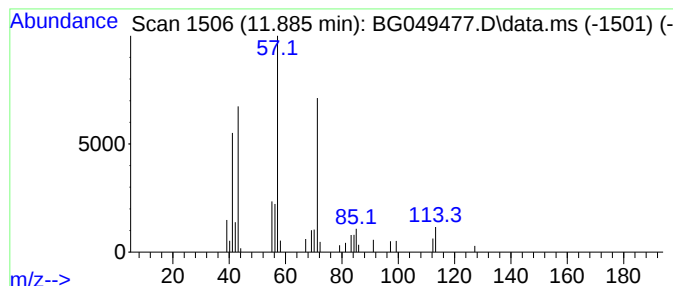
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.885	4.62 ng/ul	44445	Naphthalene-d8	10.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	83
2			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	72
3			Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	64
4			Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	64
5			Sulfurous acid, 2-ethylhexyl hep...	432	C25H52O3S	1000309-20-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049477.D
Acq On : 26 Jul 2021 13:50
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Misc :
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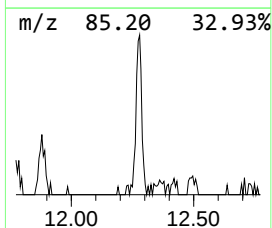
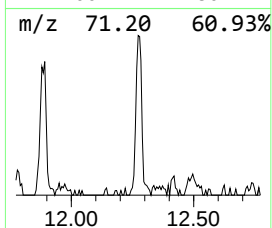
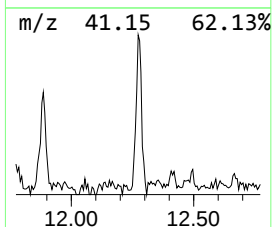
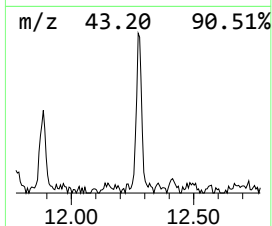
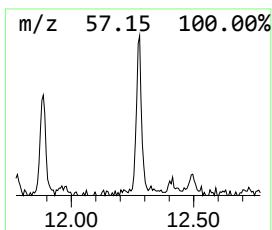
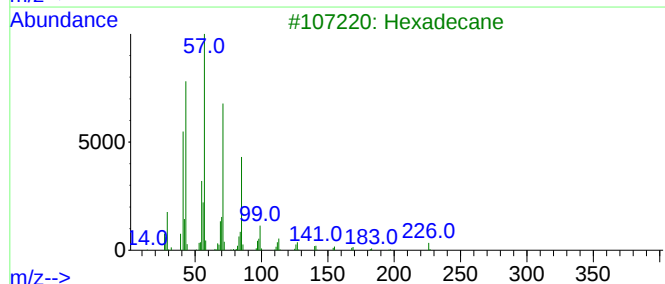
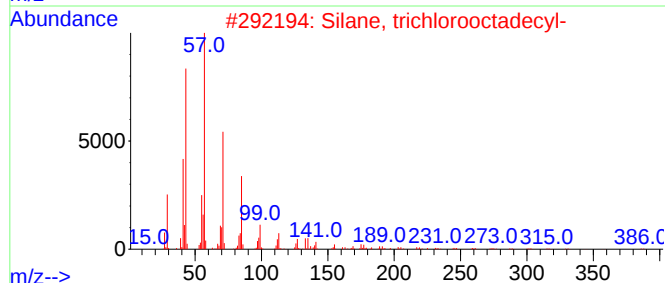
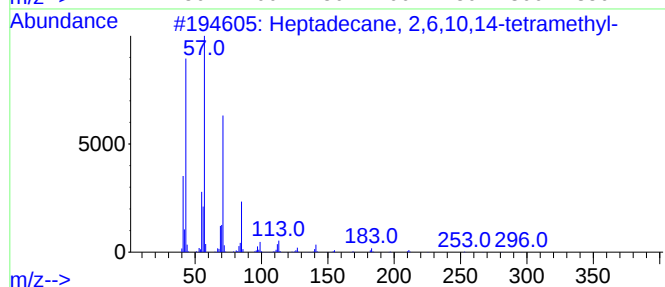
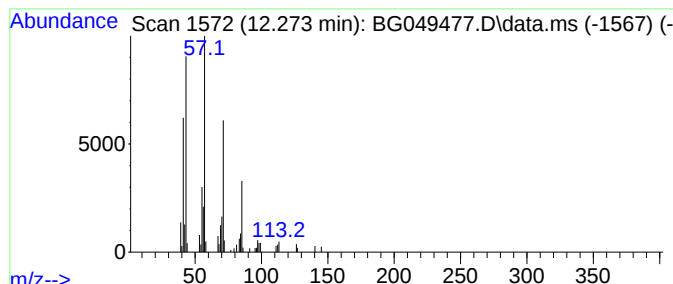
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.273	8.07 ng/ul	77622	Naphthalene-d8	10.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	86
2			Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	86
3			Hexadecane	226	C16H34	000544-76-3	86
4			Tetradecane	198	C14H30	000629-59-4	80
5			Heneicosane	296	C21H44	000629-94-7	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049477.D
Acq On : 26 Jul 2021 13:50
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Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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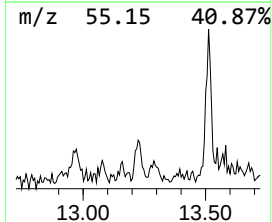
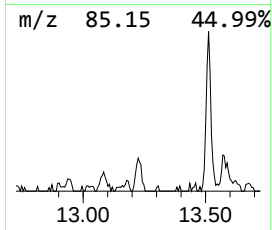
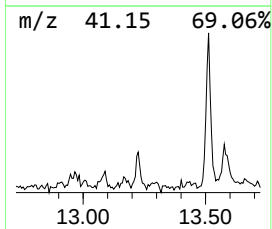
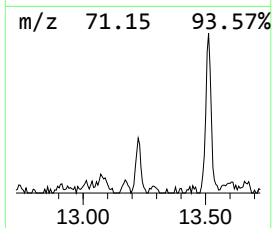
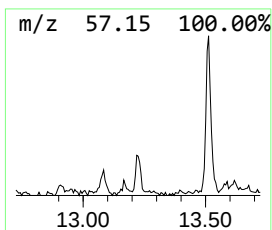
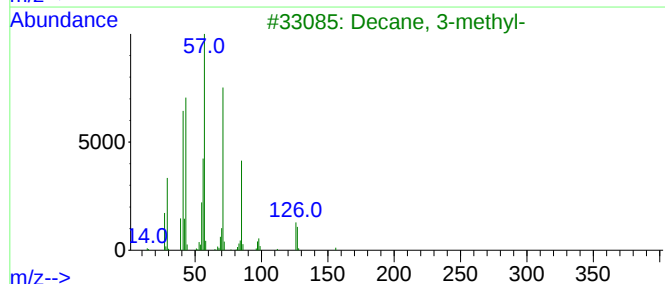
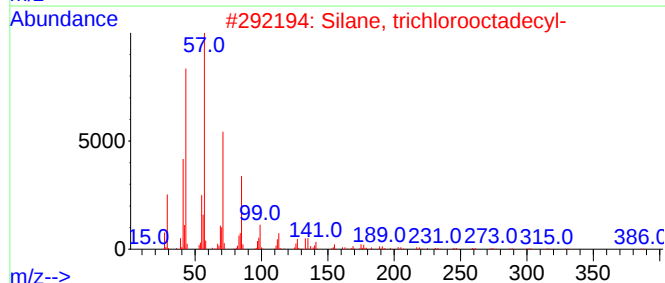
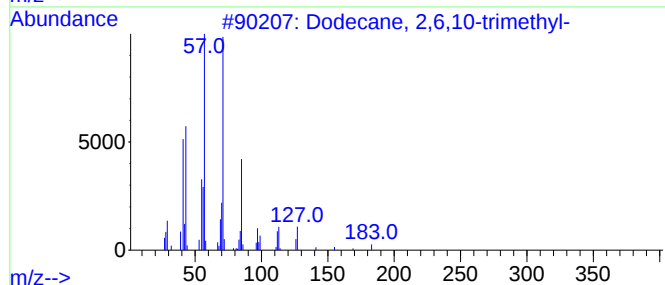
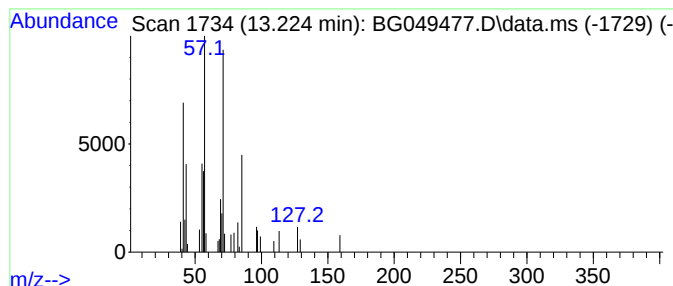
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.225	2.09 ng/ul	31221	Acenaphthene-d10	14.693

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	78
2			Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	64
3			Decane, 3-methyl-	156	C11H24	013151-34-3	64
4			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	64
5			Tetradecane	198	C14H30	000629-59-4	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049477.D
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Sample : M3082-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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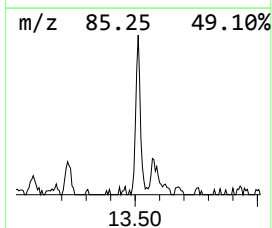
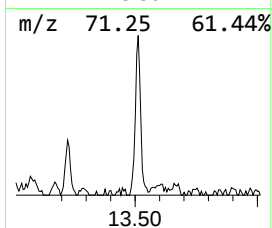
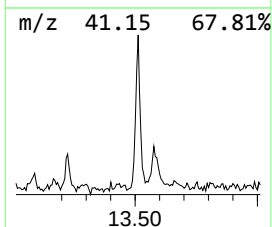
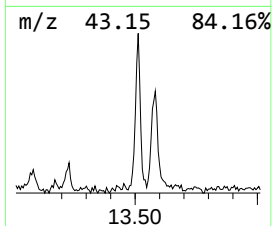
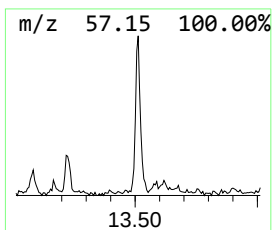
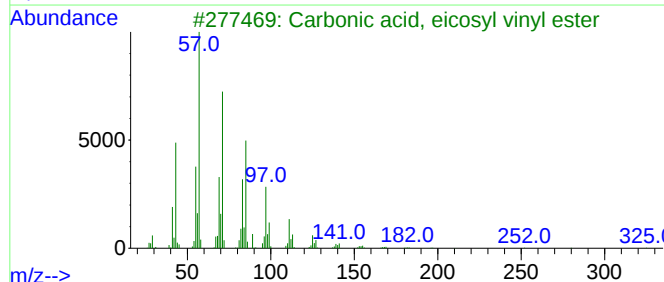
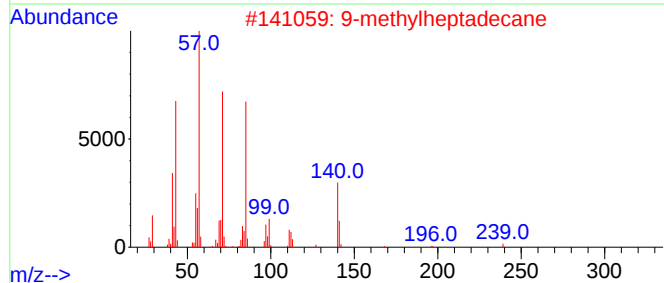
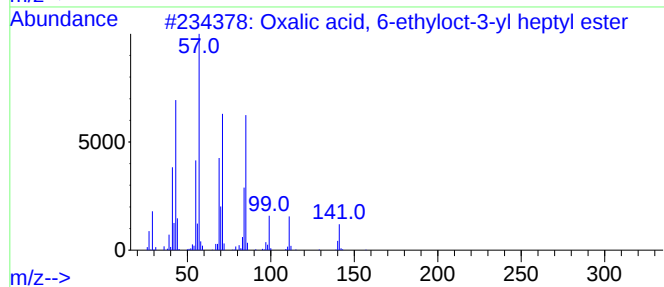
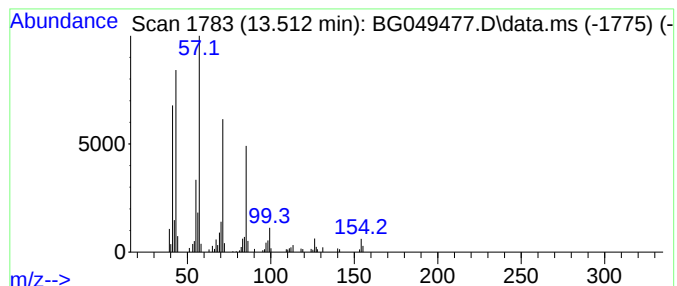
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Oxalic acid, 6-ethyloct-3-y... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.512	8.48 ng/ul	126862	Acenaphthene-d10	14.693

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, 6-ethyloct-3-yl hep...	328	C19H36O4	1000309-34-5	80
2			9-methylheptadecane	254	C18H38	026741-18-4	78
3			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	72
4			Hexadecane	226	C16H34	000544-76-3	72
5			Tetradecane	198	C14H30	000629-59-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049477.D
Acq On : 26 Jul 2021 13:50
Operator : CG/JU
Sample : M3082-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0006

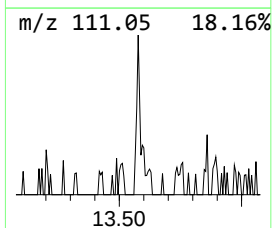
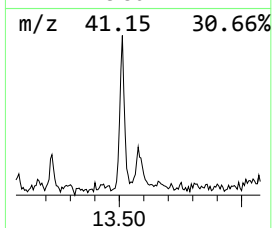
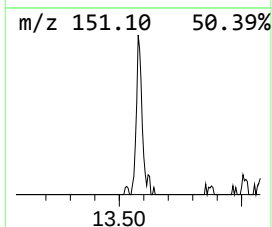
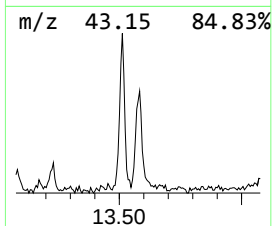
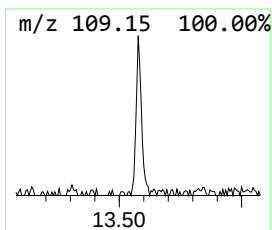
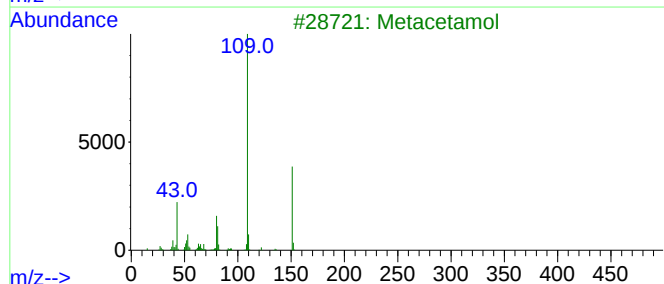
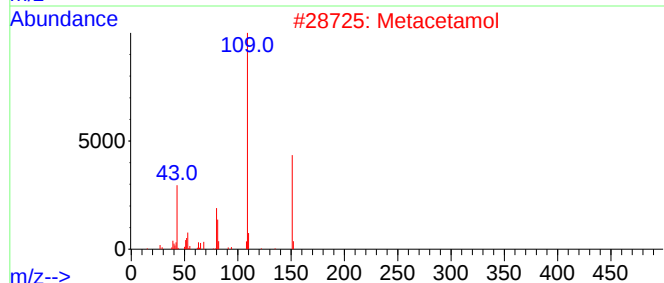
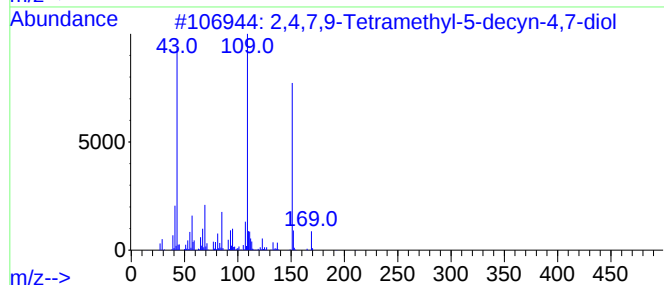
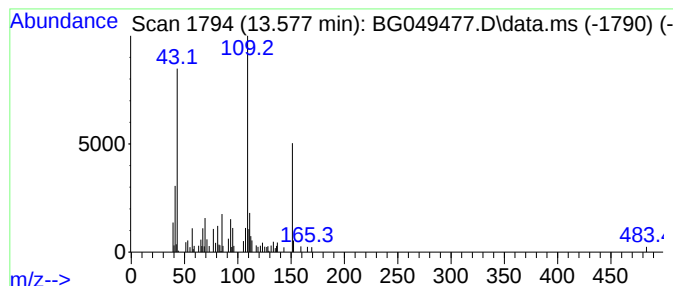
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.577	4.09 ng/ul	61110	Acenaphthene-d10		14.693

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	90	
2	Metacetamol	151	C8H9NO2	000621-42-1	58	
3	Metacetamol	151	C8H9NO2	000621-42-1	58	
4	Metacetamol	151	C8H9NO2	000621-42-1	58	
5	Acetaminophen	151	C8H9NO2	000103-90-2	58	



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049477.D
Acq On : 26 Jul 2021 13:50
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Sample : M3082-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0006

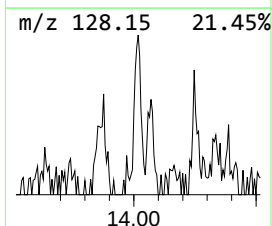
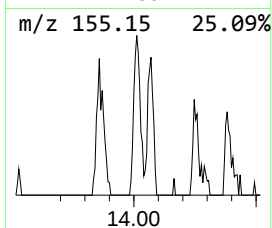
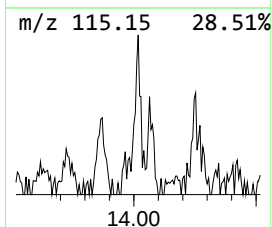
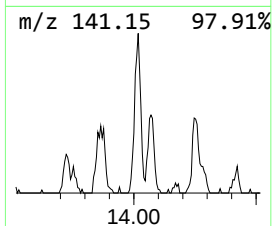
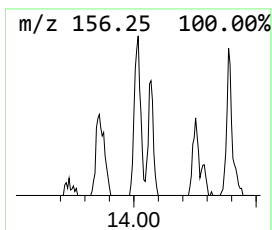
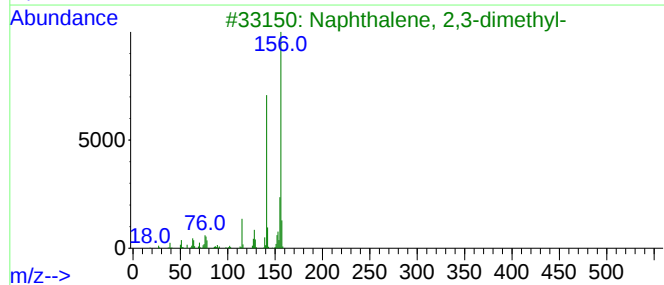
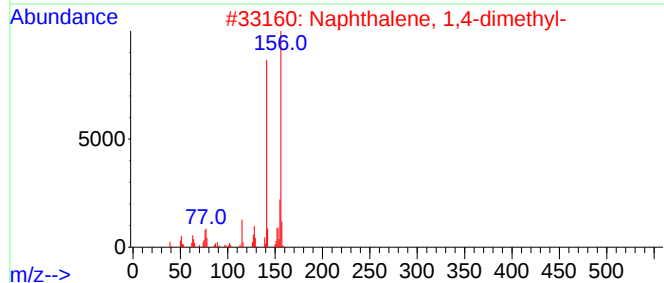
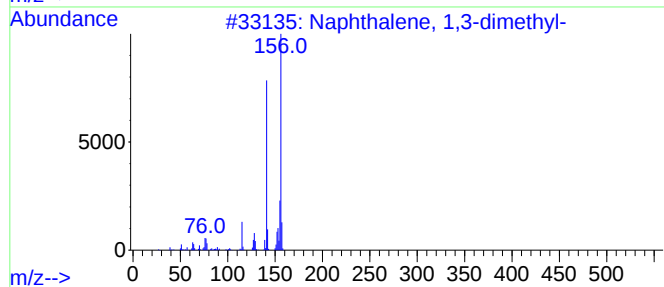
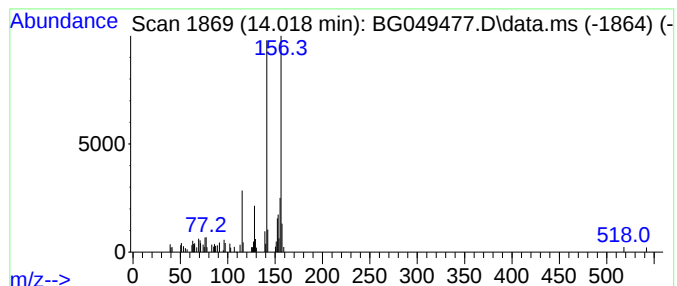
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Naphthalene, 1,3-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.018	4.87 ng/ul	72774	Acenaphthene-d10	14.693

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	94
2			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	93
3			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	93
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	93
5			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049477.D
Acq On : 26 Jul 2021 13:50
Operator : CG/JU
Sample : M3082-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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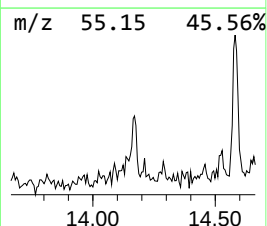
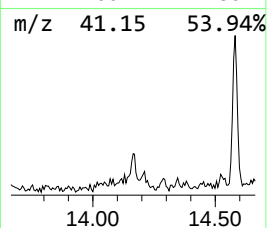
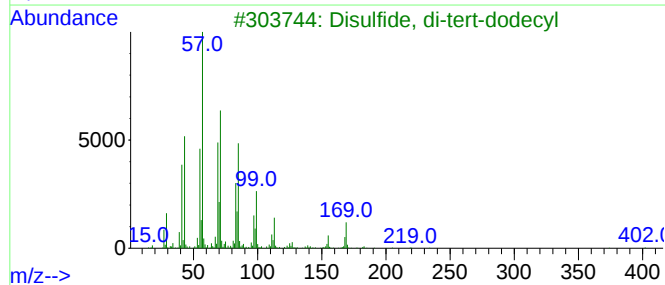
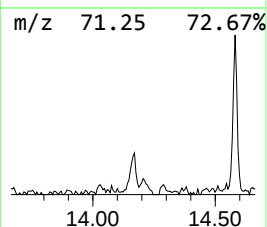
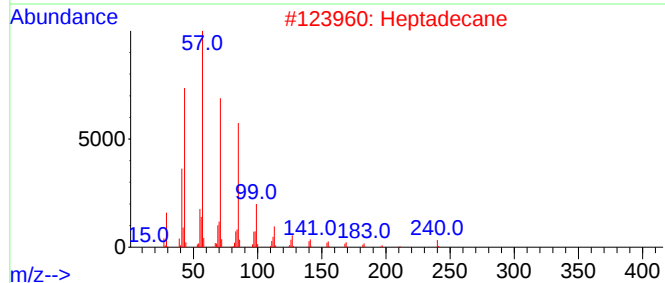
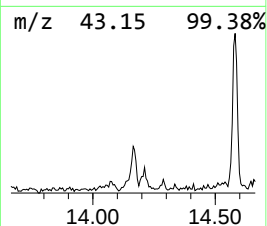
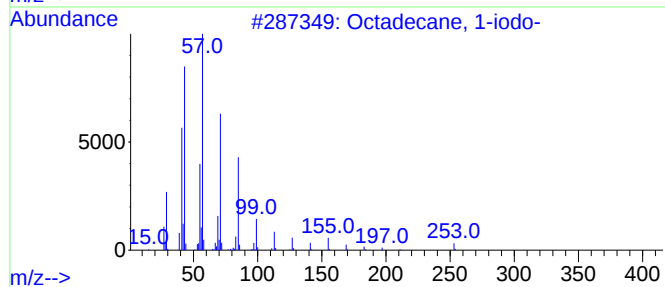
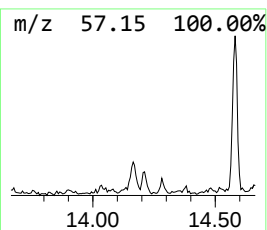
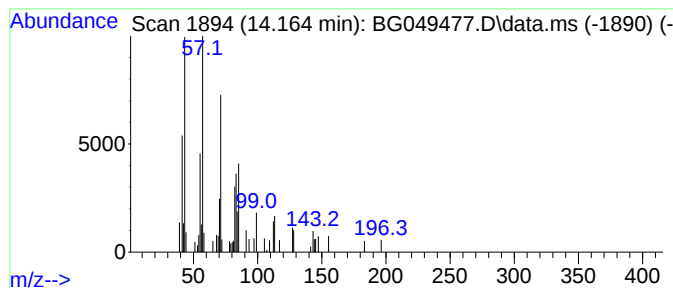
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Octadecane, 1-iodo- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.164	3.12 ng/ul	46648	Acenaphthene-d10	14.693

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	58
2			Heptadecane	240	C17H36	000629-78-7	53
3			Disulfide, di-tert-dodecyl	402	C24H50S2	027458-90-8	50
4			Undecane, 3-ethyl-	184	C13H28	017312-58-2	49
5			Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049477.D
Acq On : 26 Jul 2021 13:50
Operator : CG/JU
Sample : M3082-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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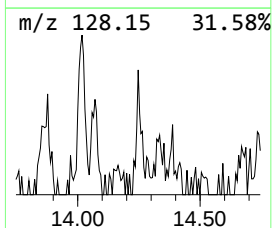
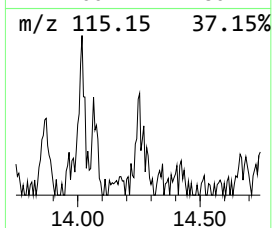
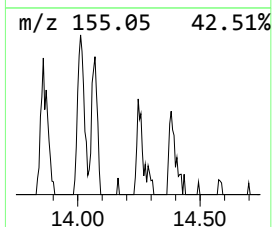
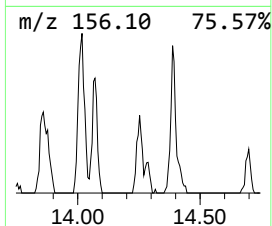
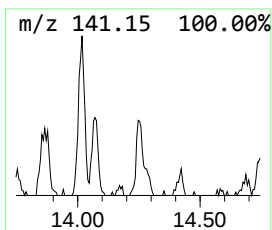
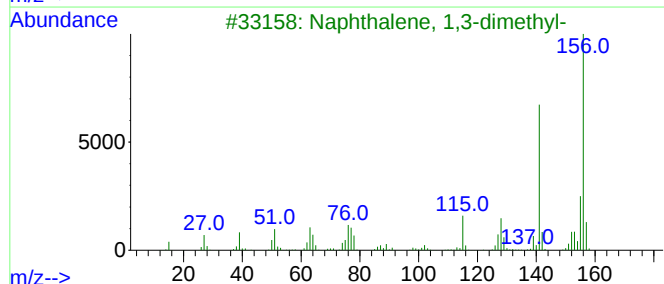
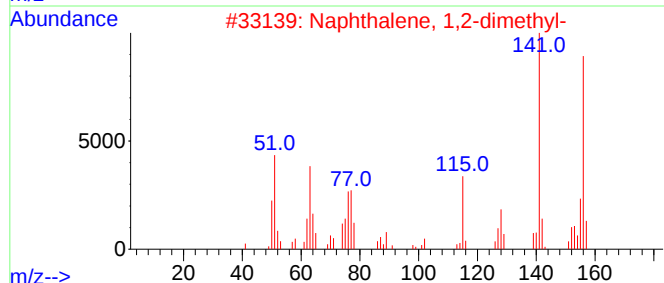
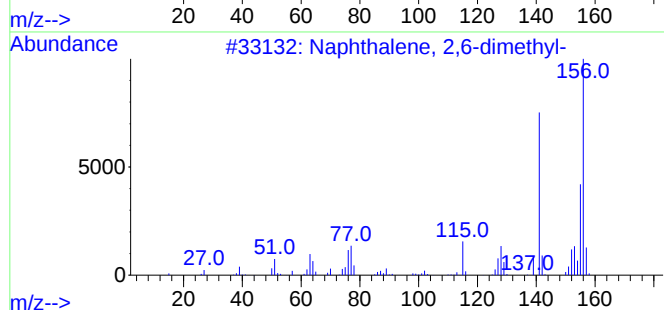
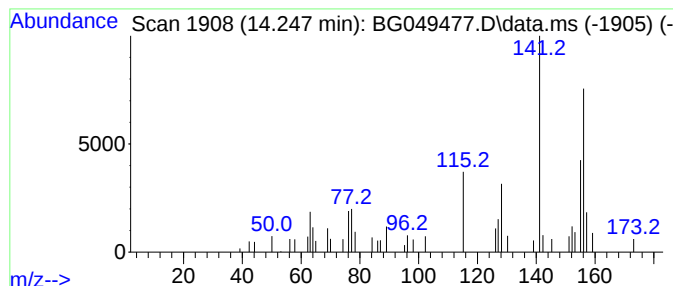
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Naphthalene, 2,6-dimethyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.247	2.03 ng/ul	30404	Acenaphthene-d10	14.693

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	90
2			Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	87
3			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	87
4			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	80
5			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049477.D
Acq On : 26 Jul 2021 13:50
Operator : CG/JU
Sample : M3082-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0006

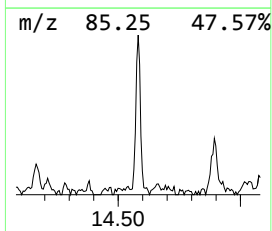
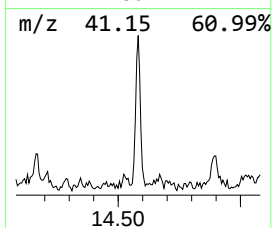
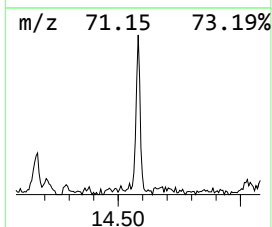
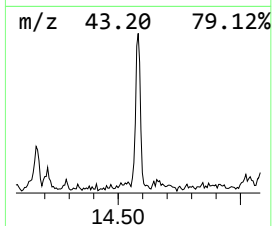
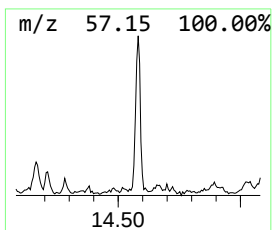
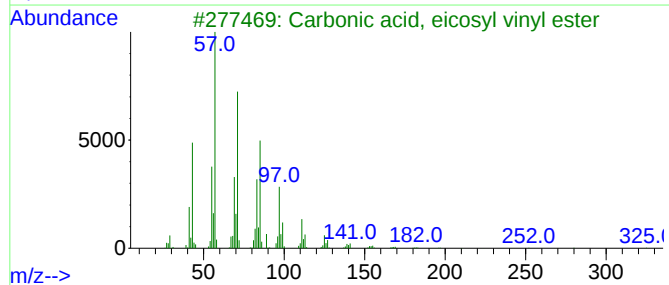
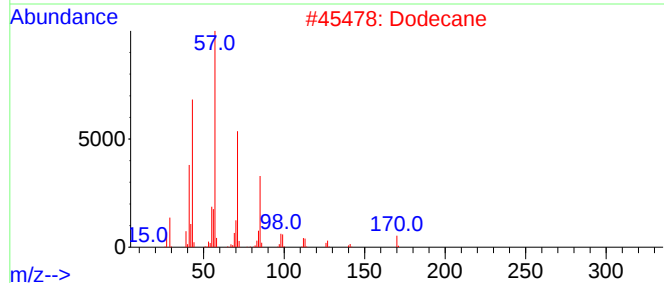
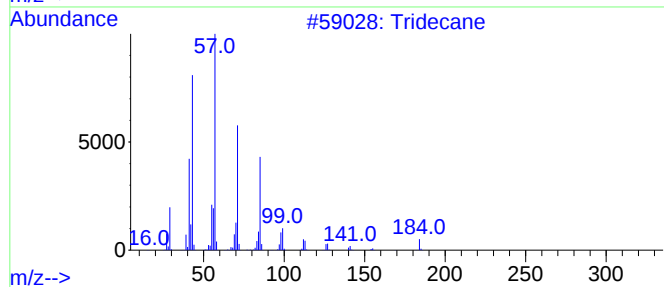
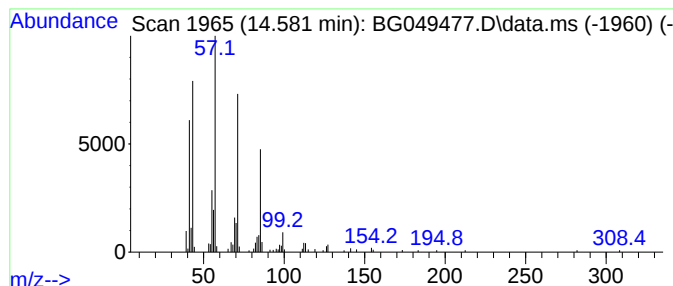
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.582	9.15 ng/ul	136895	Acenaphthene-d10	14.693

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane	184	C13H28	000629-50-5	91
2		Dodecane	170	C12H26	000112-40-3	90
3		Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	78
4		Pentadecane	212	C15H32	000629-62-9	78
5		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
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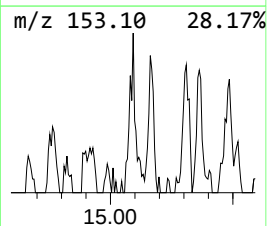
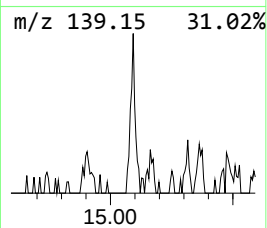
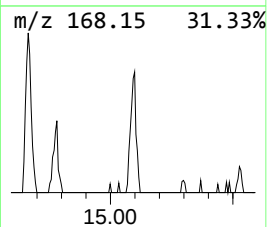
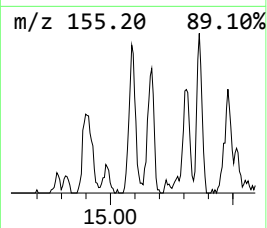
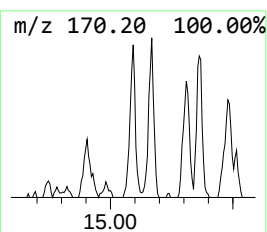
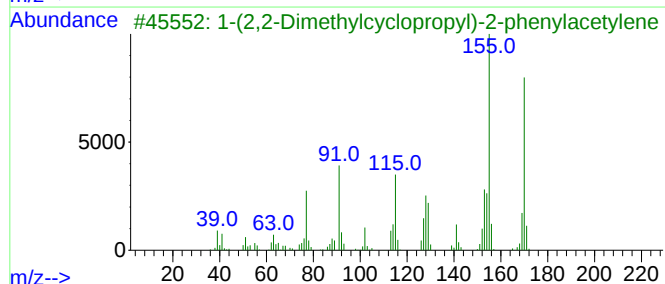
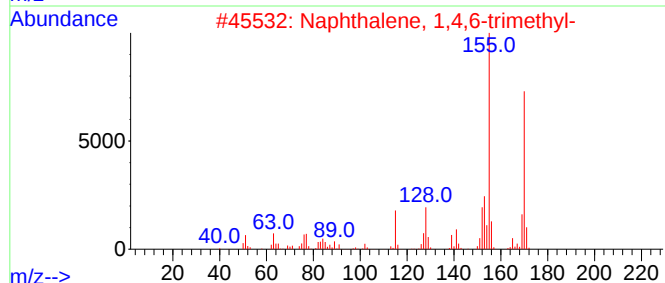
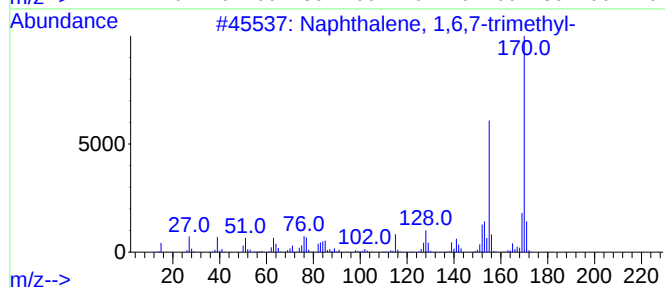
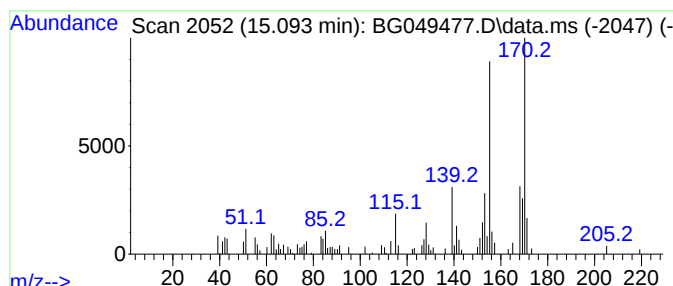
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ClientSampled :
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Naphthalene, 1,6,7-trimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.093	5.24 ng/ul	78333	Acenaphthene-d10	14.693	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	95
2	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	94
3	1-(2,2-Dimethylcyclopropyl)-2-ph...	170	C13H14	1000294-91-1	89
4	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	89
5	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049477.D
Acq On : 26 Jul 2021 13:50
Operator : CG/JU
Sample : M3082-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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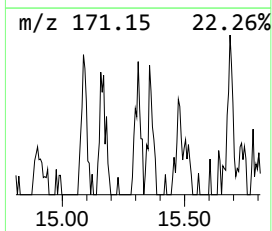
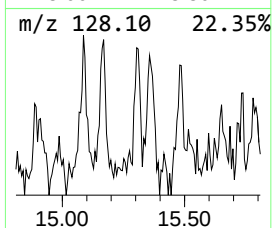
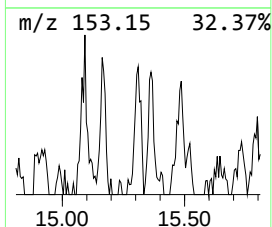
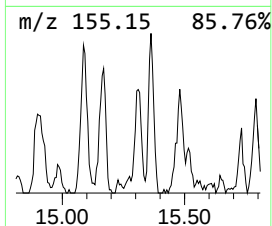
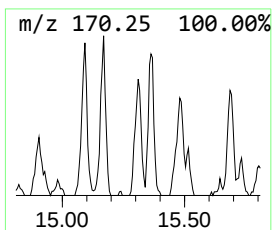
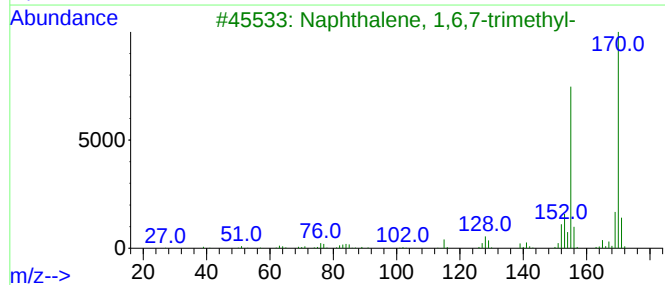
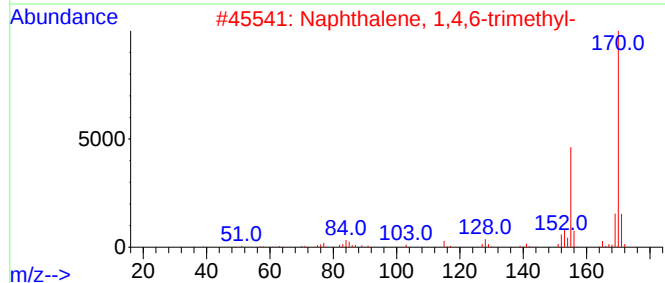
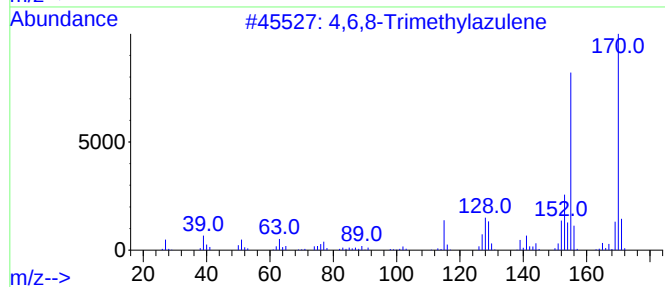
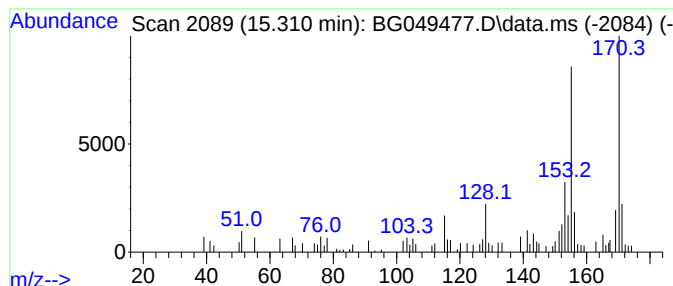
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 4,6,8-Trimethylazulene Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.310	3.89 ng/ul	58202	Acenaphthene-d10	14.693

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		4,6,8-Trimethylazulene	170	C13H14	000941-81-1	90
2		Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	87
3		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	83
4		Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	81
5		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049477.D
Acq On : 26 Jul 2021 13:50
Operator : CG/JU
Sample : M3082-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

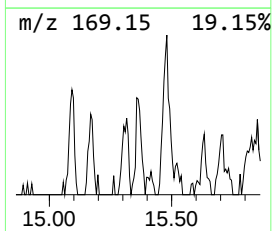
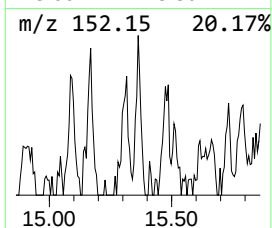
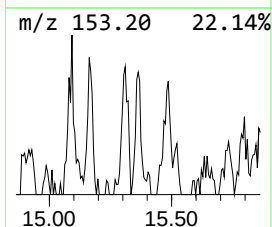
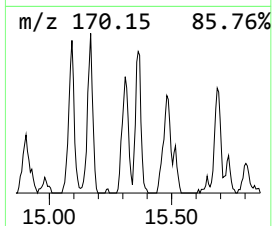
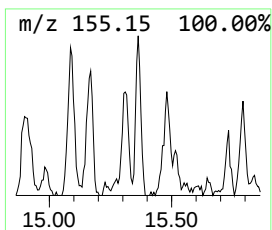
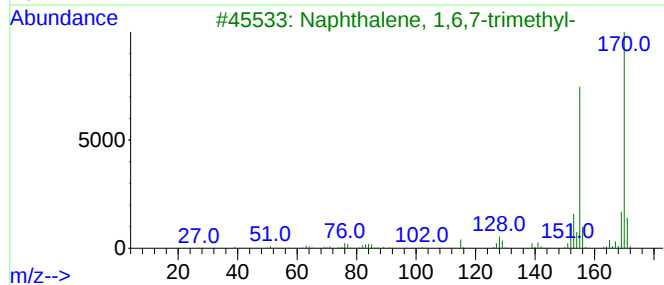
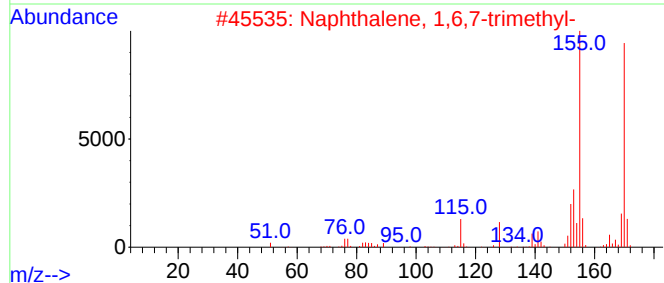
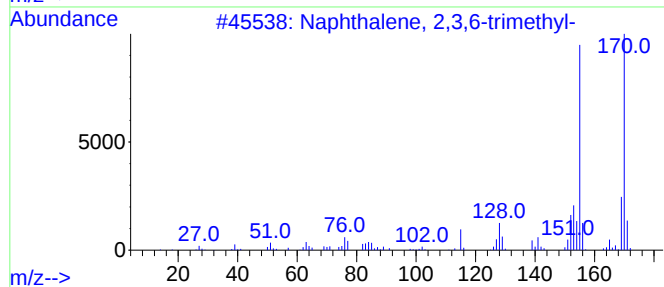
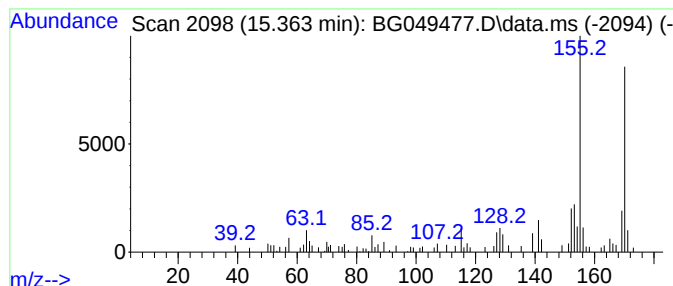
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 Naphthalene, 2,3,6-trimethyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.363	3.55 ng/ul	53085	Acenaphthene-d10	14.693	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
2	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
3	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
4	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	96
5	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049477.D
Acq On : 26 Jul 2021 13:50
Operator : CG/JU
Sample : M3082-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

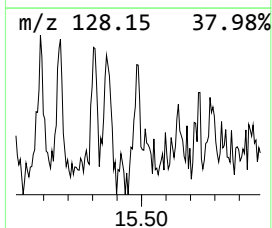
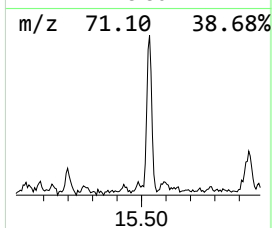
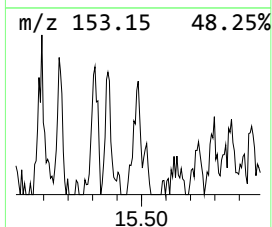
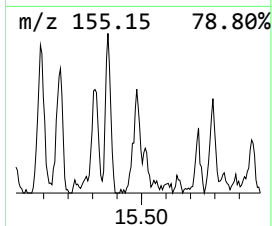
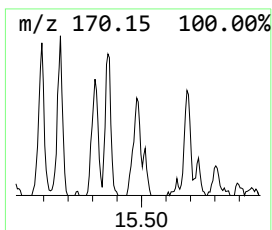
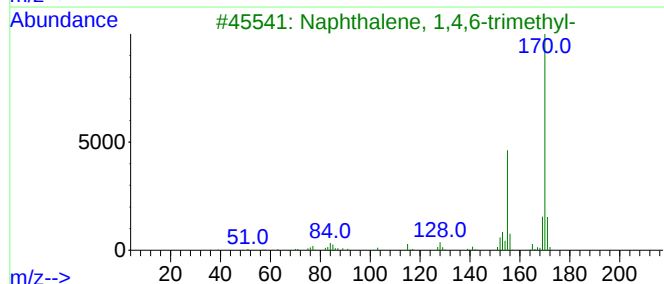
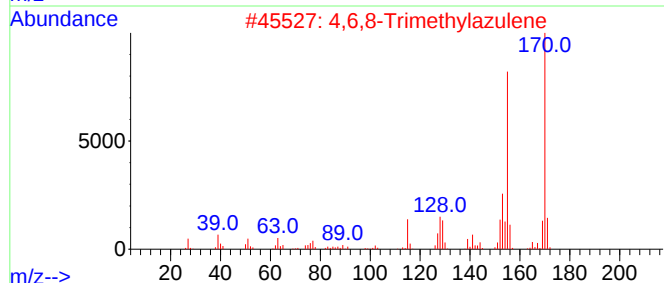
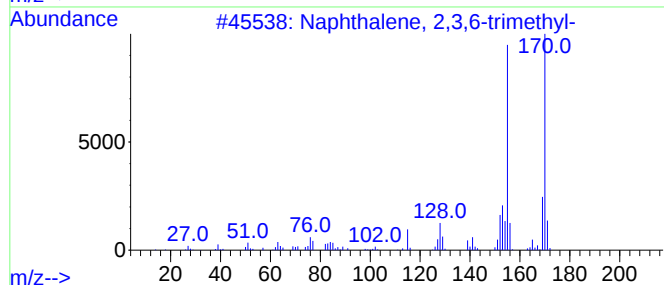
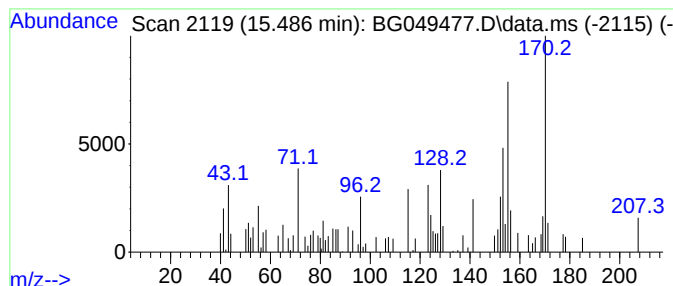
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 unknown-01 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.486	2.17 ng/ul	32413	Acenaphthene-d10	14.693	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	55
2	4,6,8-Trimethylazulene	170	C13H14	000941-81-1	53
3	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	49
4	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	49
5	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049477.D
Acq On : 26 Jul 2021 13:50
Operator : CG/JU
Sample : M3082-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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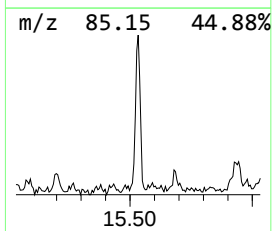
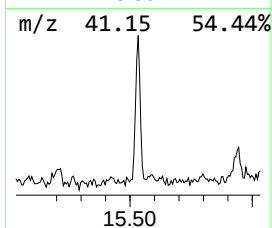
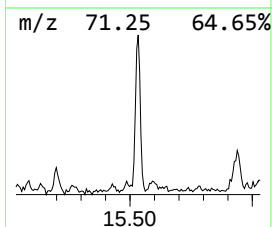
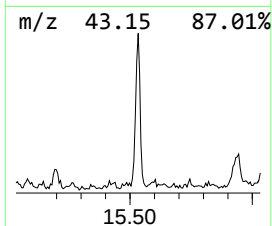
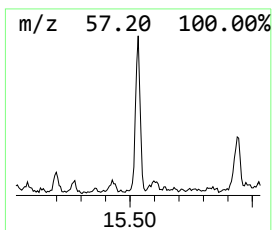
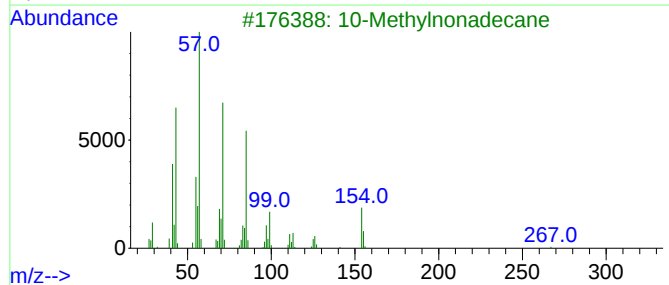
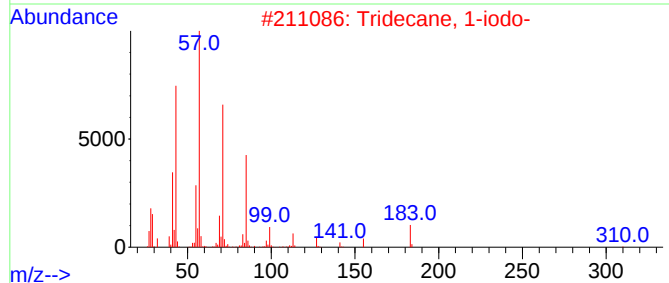
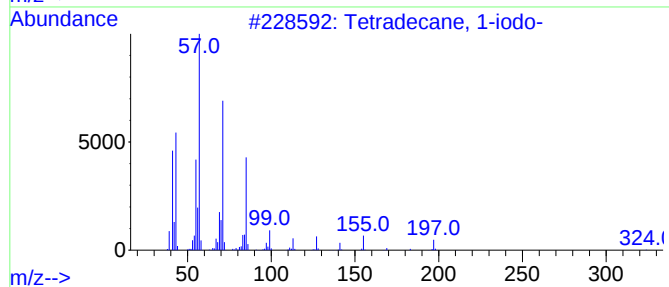
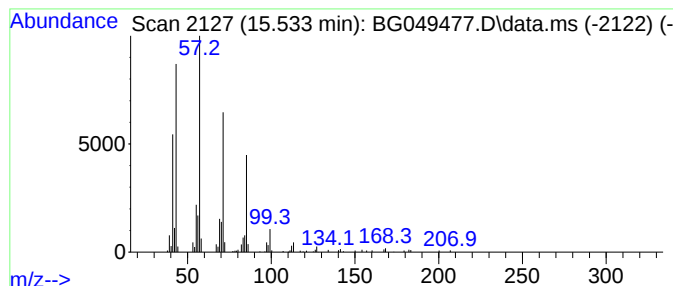
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 Tetradecane, 1-iodo- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.533	11.95 ng/ul	178710	Acenaphthene-d10	14.693

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	90
2			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	90
3			10-Methylnonadecane	282	C20H42	056862-62-5	90
4			Heneicosane	296	C21H44	000629-94-7	90
5			Nonadecane	268	C19H40	000629-92-5	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049477.D
Acq On : 26 Jul 2021 13:50
Operator : CG/JU
Sample : M3082-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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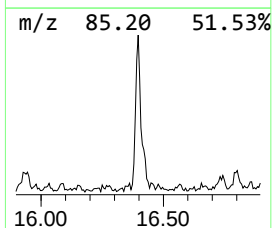
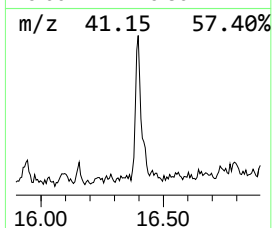
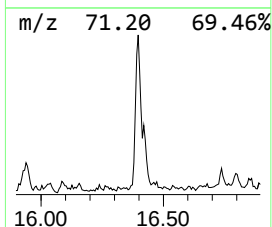
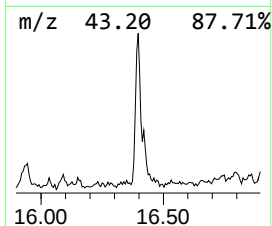
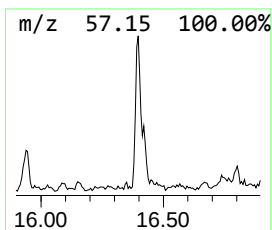
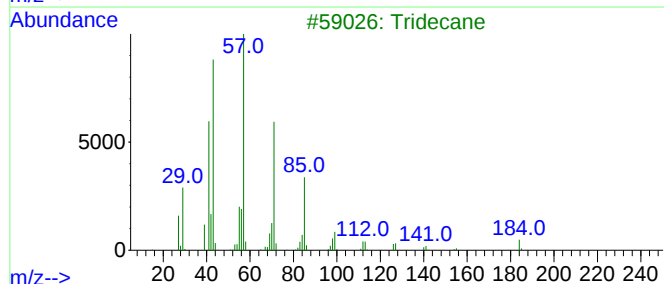
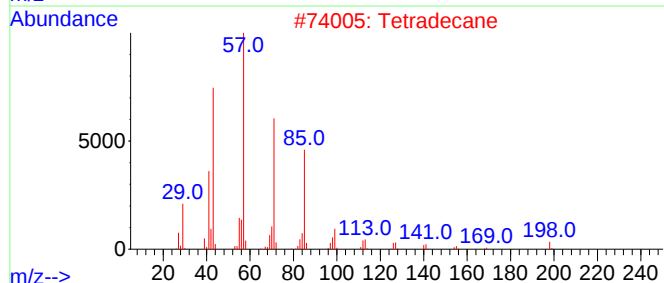
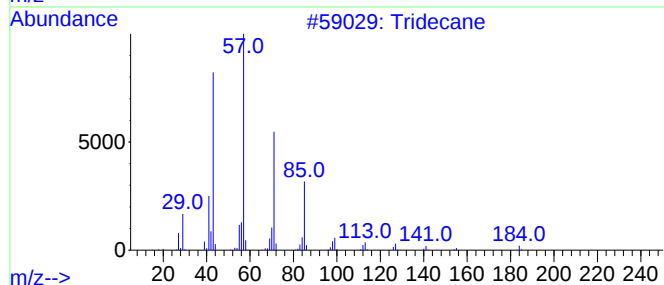
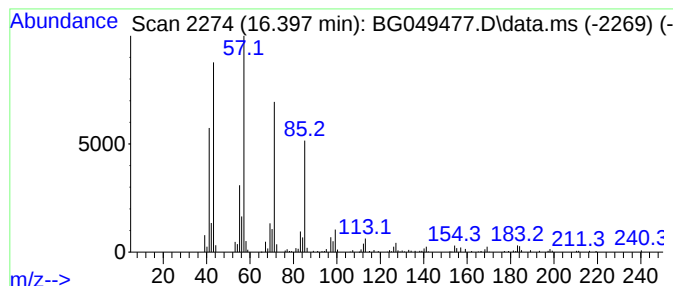
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.397	25.09 ng/ul	360538	Phenanthrene-d10	17.448

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane	184	C13H28	000629-50-5	94
2		Tetradecane	198	C14H30	000629-59-4	91
3		Tridecane	184	C13H28	000629-50-5	90
4		Nonadecane	268	C19H40	000629-92-5	90
5		Hexadecane	226	C16H34	000544-76-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049477.D
Acq On : 26 Jul 2021 13:50
Operator : CG/JU
Sample : M3082-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0006

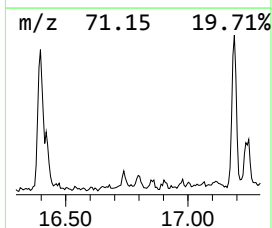
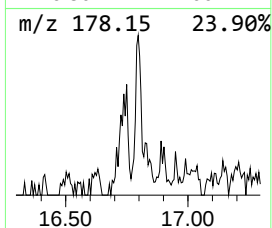
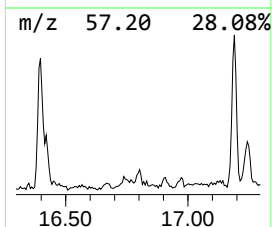
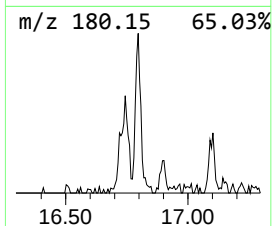
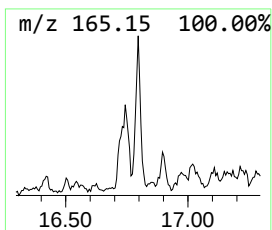
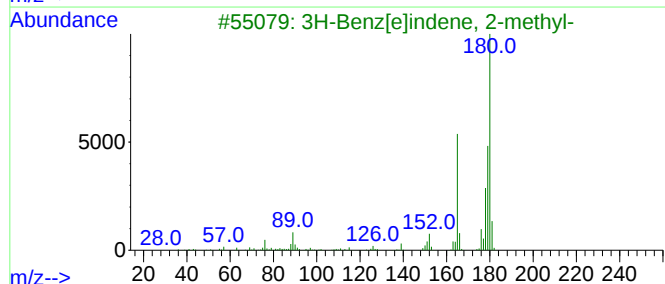
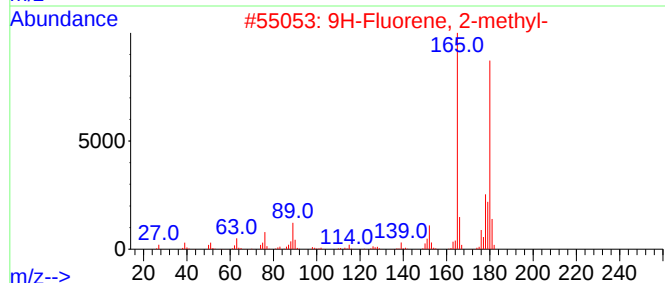
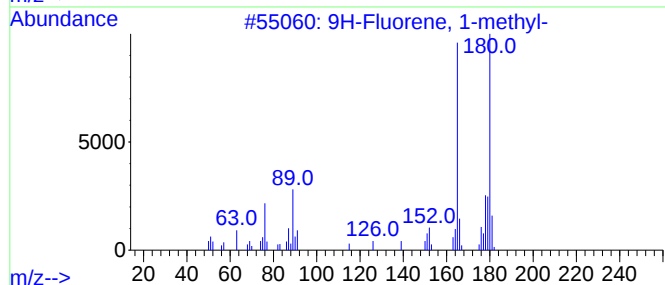
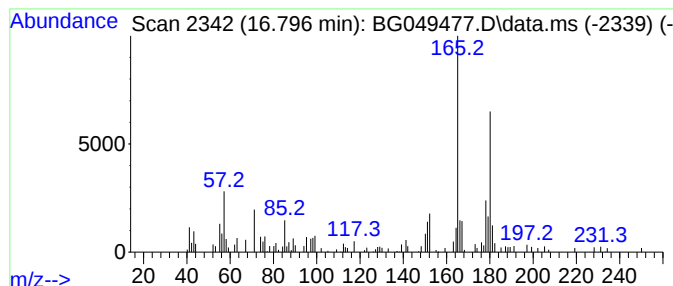
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 9H-Fluorene, 1-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.796	4.50 ng/ul	64719	Phenanthrene-d10	17.448

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	87
2			9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	81
3			3H-Benz[e]indene, 2-methyl-	180	C14H12	150096-60-9	74
4			9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	68
5			Benzenethiol, 4-(1,1-dimethyleth...	180	C11H16S	015570-10-2	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049477.D
Acq On : 26 Jul 2021 13:50
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Sample : M3082-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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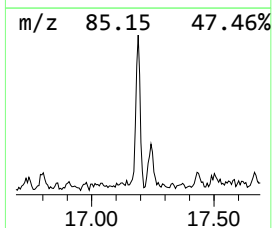
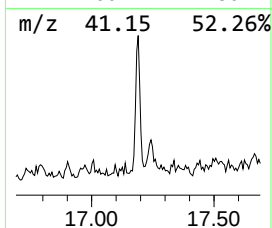
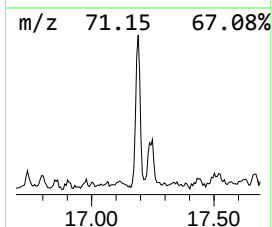
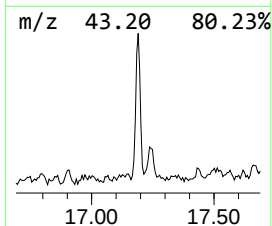
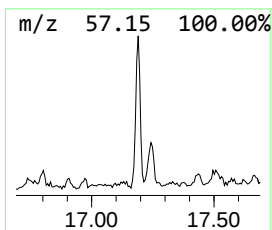
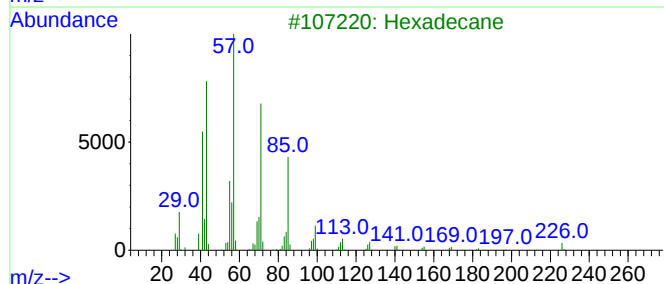
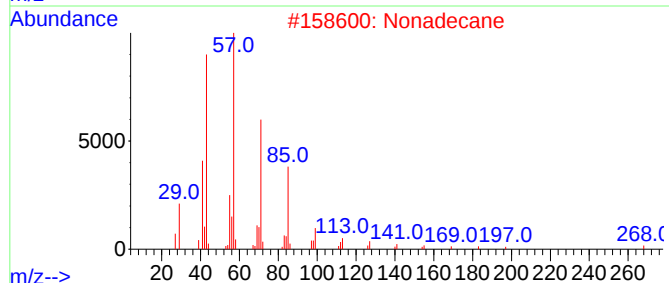
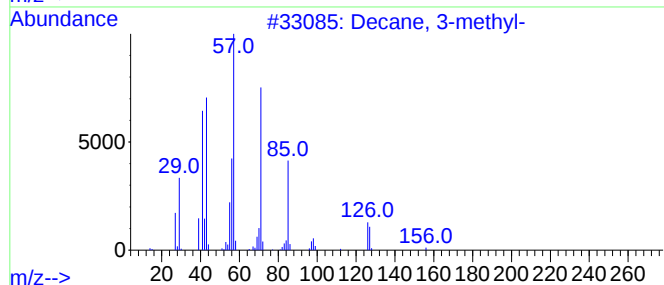
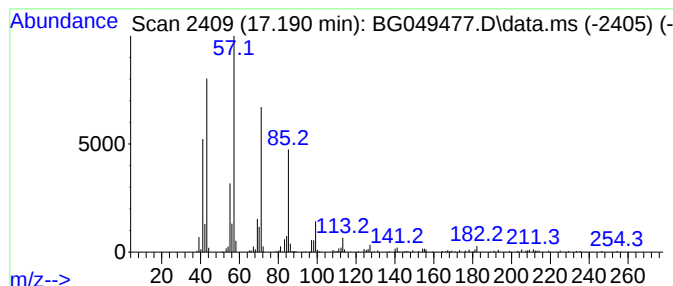
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.190	16.73 ng/ul	240482	Phenanthrene-d10	17.448

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 3-methyl-	156	C11H24	013151-34-3	87
2			Nonadecane	268	C19H40	000629-92-5	87
3			Hexadecane	226	C16H34	000544-76-3	86
4			Tetradecane	198	C14H30	000629-59-4	83
5			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049477.D
Acq On : 26 Jul 2021 13:50
Operator : CG/JU
Sample : M3082-08
Misc :
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Instrument :
BNA_G
ClientSampleId :
C0006

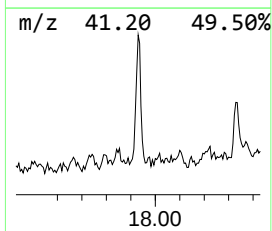
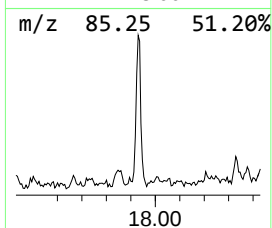
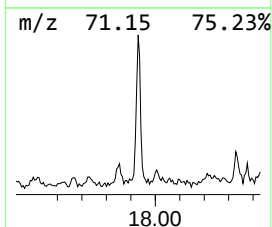
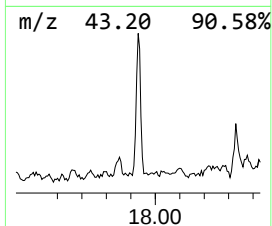
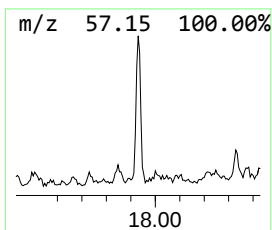
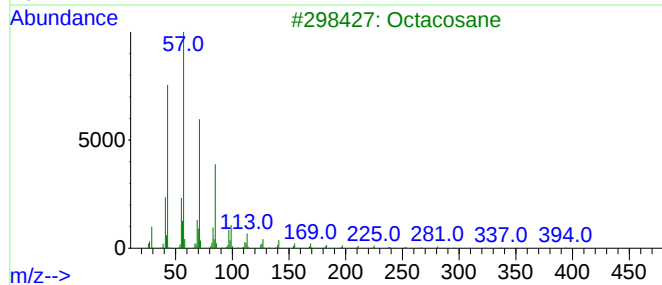
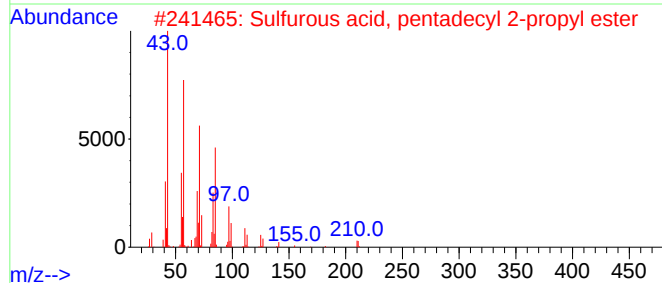
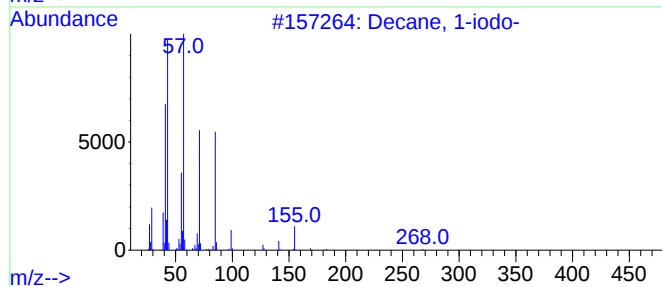
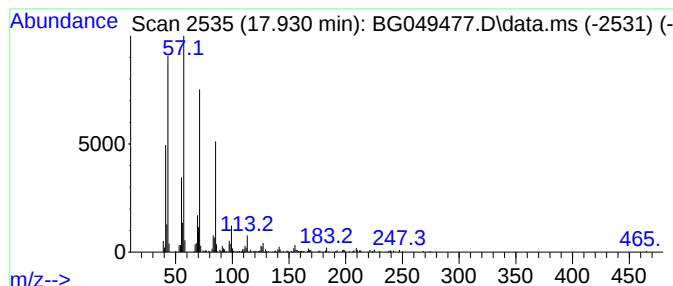
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 Decane, 1-iodo- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.930	21.25 ng/ul	305380	Phenanthrene-d10	17.448

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 1-iodo-	268	C10H21I	002050-77-3	93
2			Sulfurous acid, pentadecyl 2-pro...	334	C18H38O3S	1000309-12-6	90
3			Octacosane	394	C28H58	000630-02-4	90
4			Undecane	156	C11H24	001120-21-4	90
5			Docosane, 1-iodo-	436	C22H45I	1000406-31-9	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049477.D
Acq On : 26 Jul 2021 13:50
Operator : CG/JU
Sample : M3082-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0006

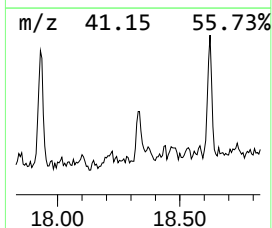
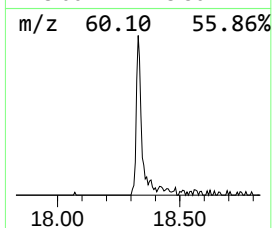
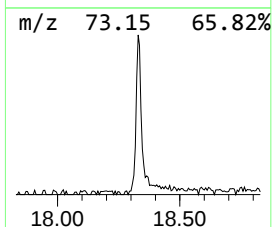
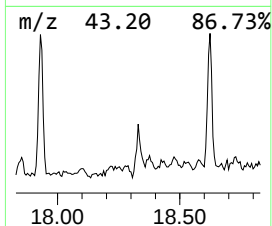
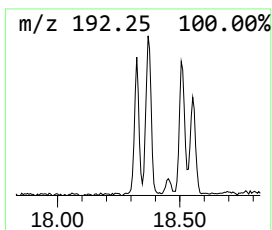
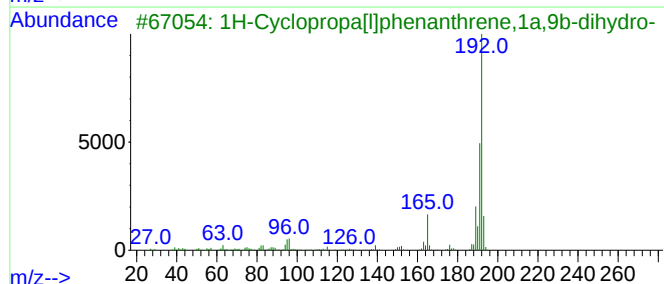
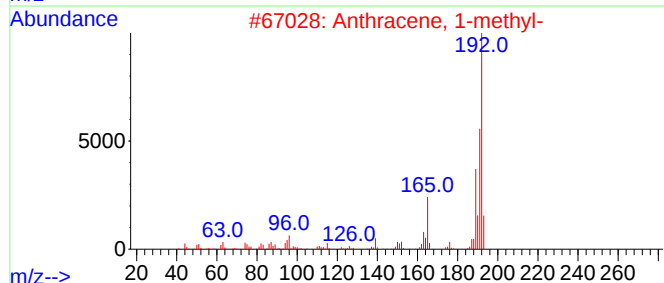
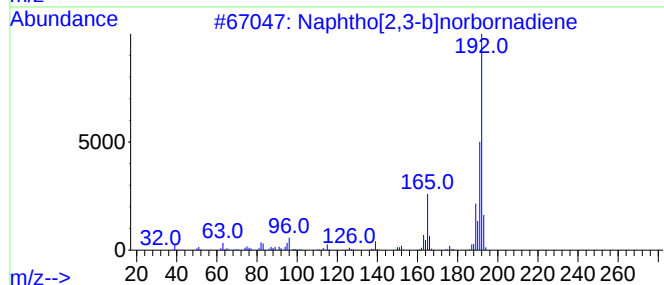
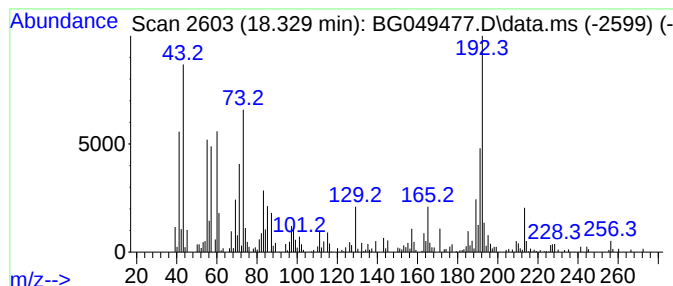
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 Naphtho[2,3-b]norbornadiene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.329	16.93 ng/ul	243352	Phenanthrene-d10	17.448

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	53
2			Anthracene, 1-methyl-	192	C15H12	000610-48-0	46
3			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	46
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	46
5			Anthracene, 9-methyl-	192	C15H12	000779-02-2	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049477.D
Acq On : 26 Jul 2021 13:50
Operator : CG/JU
Sample : M3082-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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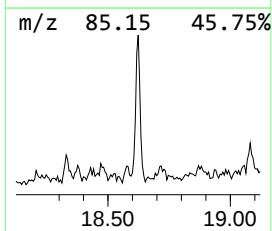
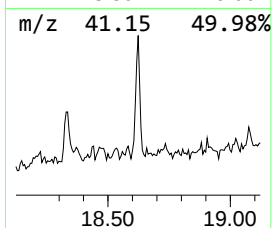
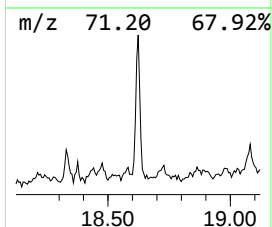
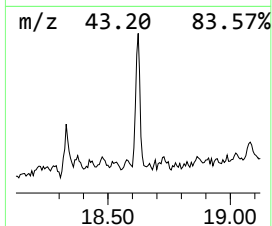
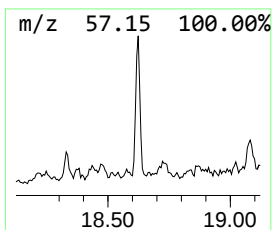
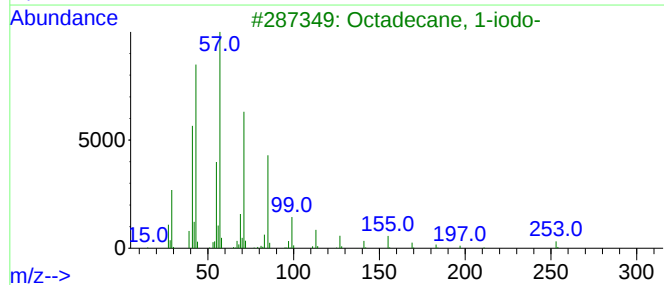
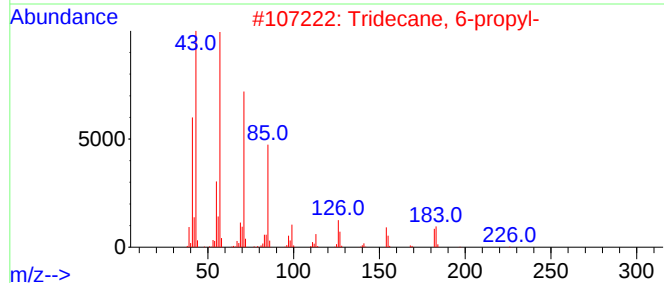
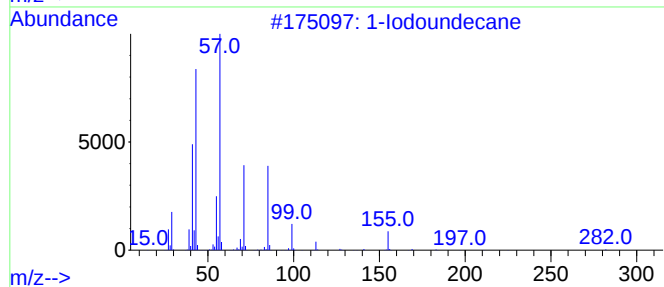
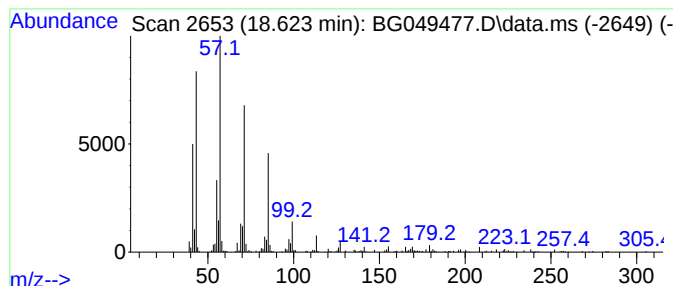
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 1-Iodoundecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.623	18.49 ng/ul	265691	Phenanthrene-d10	17.448

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Iodoundecane	282	C11H23I	004282-44-4	91
2		Tridecane, 6-propyl-	226	C16H34	055045-10-8	91
3		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	90
4		Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	90
5		Nonadecane	268	C19H40	000629-92-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
 Data File : BG049477.D
 Acq On : 26 Jul 2021 13:50
 Operator : CG/JU
 Sample : M3082-08
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0006

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.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
(DEL) Alkane: C...	3.085	24.8	ng/ul	188662	1	8.049	151967	20.0
Butane, 2-metho...	3.162	37.8	ng/ul	287033	1	8.049	151967	20.0
Cyclohexene, 3-...	6.851	2.7	ng/ul	20352	1	8.049	151967	20.0
(DEL) Alkane: S...	7.661	3.7	ng/ul	28251	1	8.049	151967	20.0
Carbonic acid, ...	9.277	4.6	ng/ul	35045	1	8.049	151967	20.0
Ethanol, 2-(2-b...	10.628	11.1	ng/ul	107009	2	10.869	192371	20.0
(DEL) Alkane: S...	11.885	4.6	ng/ul	44445	2	10.869	192371	20.0
(DEL) Alkane: S...	12.273	8.1	ng/ul	77622	2	10.869	192371	20.0
(DEL) Alkane: S...	13.225	2.1	ng/ul	31221	3	14.693	299158	20.0
Oxalic acid, 6-...	13.512	8.5	ng/ul	126862	3	14.693	299158	20.0
2,4,7,9-Tetrame...	13.577	4.1	ng/ul	61110	3	14.693	299158	20.0
Naphthalene, 1,...	14.018	4.9	ng/ul	72774	3	14.693	299158	20.0
Octadecane, 1-i...	14.164	3.1	ng/ul	46648	3	14.693	299158	20.0
Naphthalene, 2,...	14.247	2.0	ng/ul	30404	3	14.693	299158	20.0
(DEL) Alkane: S...	14.582	9.2	ng/ul	136895	3	14.693	299158	20.0
Naphthalene, 1,...	15.093	5.2	ng/ul	78333	3	14.693	299158	20.0
4,6,8-Trimethyl...	15.310	3.9	ng/ul	58202	3	14.693	299158	20.0
Naphthalene, 2,...	15.363	3.5	ng/ul	53085	3	14.693	299158	20.0
unknown-01	15.486	2.2	ng/ul	32413	3	14.693	299158	20.0
Tetradecane, 1-...	15.533	11.9	ng/ul	178710	3	14.693	299158	20.0
(DEL) Alkane: S...	16.397	25.1	ng/ul	360538	4	17.448	287448	20.0
9H-Fluorene, 1-...	16.796	4.5	ng/ul	64719	4	17.448	287448	20.0
(DEL) Alkane: S...	17.190	16.7	ng/ul	240482	4	17.448	287448	20.0
Decane, 1-iodo-	17.930	21.3	ng/ul	305380	4	17.448	287448	20.0
Naphtho[2,3-b]n...	18.329	16.9	ng/ul	243352	4	17.448	287448	20.0
1-Iodoundecane	18.623	18.5	ng/ul	265691	4	17.448	287448	20.0