

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
 Data File : BG050107.D
 Acq On : 2 Sep 2021 21:53
 Operator : CG/JU
 Sample : M3526-14
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW7C4

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-BG081921.M
 Title : SVOA CALIBRATION

Signal : TIC: BG050107.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.754	117	122	130	rBV3	12161	25524	0.50%	0.089%
2	3.848	134	138	147	rVV2	22827	38438	0.75%	0.134%
3	4.071	168	176	185	rBV	99038	178390	3.47%	0.621%
4	6.192	528	537	544	rBV8	10405	27055	0.53%	0.094%
5	6.404	569	573	581	rBV7	8039	20274	0.39%	0.071%
6	6.592	604	605	614	rVV6	12250	22752	0.44%	0.079%
7	6.803	637	641	650	rVV5	9027	21878	0.43%	0.076%
8	6.921	656	661	667	rVV6	11262	21145	0.41%	0.074%
9	7.085	683	689	692	rVV6	17026	42583	0.83%	0.148%
10	7.120	692	695	707	rVB2	46290	96529	1.88%	0.336%
11	7.273	713	721	726	rBV4	38409	72186	1.41%	0.251%
12	7.432	742	748	752	rBV4	9900	19425	0.38%	0.068%
13	7.484	752	757	766	rVB3	50697	101308	1.97%	0.353%
14	7.843	810	818	819	rBV6	11901	19772	0.39%	0.069%
15	7.919	826	831	832	rBV3	16029	20256	0.39%	0.071%
16	7.949	832	836	843	rVB	101736	167072	3.25%	0.582%
17	8.007	843	846	851	rVB6	14329	21596	0.42%	0.075%
18	8.131	862	867	871	rVB4	25346	38735	0.75%	0.135%
19	8.225	877	883	887	rVB7	14456	25299	0.49%	0.088%
20	8.295	891	895	903	rVB6	12845	28154	0.55%	0.098%
21	8.389	903	911	917	rVB7	14408	33288	0.65%	0.116%
22	8.495	927	929	935	rVB3	12844	20713	0.40%	0.072%
23	8.642	947	954	955	rBV5	11414	21949	0.43%	0.076%
24	8.677	955	960	967	rVB3	39296	75626	1.47%	0.263%
25	8.789	974	979	983	rBV4	27235	45834	0.89%	0.160%
26	9.059	1014	1025	1030	rBV4	38375	94634	1.84%	0.330%
27	9.123	1031	1036	1048	rVV3	52842	138162	2.69%	0.481%
28	9.212	1048	1051	1056	rVB6	16303	29885	0.58%	0.104%
29	9.270	1056	1061	1063	rBV3	29030	45771	0.89%	0.159%
30	9.311	1063	1068	1077	rVB5	39357	80811	1.57%	0.281%
31	9.423	1077	1087	1096	rVB6	29654	86372	1.68%	0.301%
32	9.529	1098	1105	1110	rBV10	15783	41599	0.81%	0.145%
33	9.587	1110	1115	1120	rBV6	38420	82653	1.61%	0.288%
34	9.646	1123	1125	1127	rVV3	30907	39012	0.76%	0.136%
35	9.681	1127	1131	1135	rVV4	50352	88230	1.72%	0.307%
36	9.717	1135	1137	1143	rVB6	14056	23030	0.45%	0.080%

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Smoothing : OFF

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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-BG081921.M
 Title : SVOA CALIBRATION

37	9.852	1148	1160	1162	rBV7	68659	173160	3.37%	0.603%
38	9.940	1169	1175	1181	rBV5	59336	110302	2.15%	0.384%
39	10.075	1194	1198	1200	rBV4	19250	29044	0.57%	0.101%
40	10.175	1208	1215	1217	rBV6	19736	42490	0.83%	0.148%
41	10.257	1227	1229	1235	rVB4	27736	32575	0.63%	0.113%
42	10.339	1236	1243	1249	rBV8	34310	99304	1.93%	0.346%
43	10.410	1249	1255	1260	rVB3	56359	103218	2.01%	0.359%
44	10.469	1261	1265	1267	rBV4	14183	22449	0.44%	0.078%
45	10.539	1272	1277	1281	rBV7	29057	56610	1.10%	0.197%
46	10.657	1288	1297	1304	rBV8	41627	151938	2.96%	0.529%
47	10.727	1304	1309	1311	rVV4	45694	73512	1.43%	0.256%
48	10.774	1312	1317	1323	rVV	151049	281868	5.49%	0.981%
49	10.833	1323	1327	1331	rVB3	29988	48190	0.94%	0.168%
50	10.886	1332	1336	1337	rBV4	21084	31604	0.62%	0.110%
51	10.921	1337	1342	1347	rVB	161884	266320	5.19%	0.927%
52	10.980	1349	1352	1356	rBV7	33916	62869	1.22%	0.219%
53	11.138	1372	1379	1384	rBV5	45124	119044	2.32%	0.415%
54	11.250	1394	1398	1405	rBV9	26276	54776	1.07%	0.191%
55	11.315	1405	1409	1418	rVB6	40966	83975	1.64%	0.292%
56	11.391	1418	1422	1424	rBV5	19981	32338	0.63%	0.113%
57	11.420	1424	1427	1430	rBV2	47734	66967	1.30%	0.233%
58	11.467	1430	1435	1441	rVB7	80911	148735	2.90%	0.518%
59	11.532	1443	1446	1451	rVB3	43118	69578	1.36%	0.242%
60	11.602	1453	1458	1466	rVB8	70373	150782	2.94%	0.525%
61	11.696	1467	1474	1480	rBV7	62975	153209	2.98%	0.533%
62	11.743	1480	1482	1484	rBV3	20605	24219	0.47%	0.084%
63	11.785	1484	1489	1493	rBV	331199	493554	9.61%	1.719%
64	11.826	1493	1496	1500	rVB2	88758	119378	2.32%	0.416%
65	12.090	1538	1541	1546	rVB4	53636	71706	1.40%	0.250%
66	12.172	1552	1555	1558	rVB2	39283	37355	0.73%	0.130%
67	12.319	1577	1580	1585	rBV4	58673	102780	2.00%	0.358%
68	12.401	1587	1594	1601	rVB6	140132	395758	7.71%	1.378%
69	12.501	1607	1611	1615	rBV5	49234	91607	1.78%	0.319%
70	12.619	1628	1631	1635	rBV3	52802	99003	1.93%	0.345%
71	12.660	1635	1638	1650	rVB7	107795	220309	4.29%	0.767%
72	12.766	1652	1656	1660	rBV4	88795	132328	2.58%	0.461%
73	12.818	1660	1665	1668	rVV3	110421	170092	3.31%	0.592%
74	12.877	1672	1675	1684	rVB	159742	294890	5.74%	1.027%
75	13.136	1715	1719	1725	rBV	526339	748474	14.58%	2.606%
76	13.412	1762	1766	1769	rBV4	63983	120971	2.36%	0.421%

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 Stop Thrs : 0 Peak Location: TOP

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Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
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77	13.453	1769	1773	1777	rVB3	190719	301505	5.87%	1.050%
78	13.559	1788	1791	1794	rBV3	92002	91263	1.78%	0.318%
79	13.635	1801	1804	1805	rBV2	71349	75240	1.47%	0.262%
80	13.664	1806	1809	1815	rVB4	158737	264364	5.15%	0.921%
81	13.817	1831	1835	1840	rBV2	200841	319794	6.23%	1.114%
82	14.087	1876	1881	1885	rVB	1221014	1679054	32.70%	5.847%
83	14.140	1886	1890	1895	rVB4	186262	353986	6.89%	1.233%
84	14.258	1906	1910	1913	rBV5	151131	233175	4.54%	0.812%
85	14.305	1914	1918	1923	rVB2	166099	298019	5.80%	1.038%
86	14.440	1937	1941	1947	rVB4	229329	399410	7.78%	1.391%
87	14.487	1947	1949	1953	rBV4	110436	155898	3.04%	0.543%
88	14.575	1960	1964	1966	rBV2	246012	330537	6.44%	1.151%
89	14.704	1982	1986	1989	rBV2	238128	325381	6.34%	1.133%
90	14.839	2005	2009	2015	rVB2	283996	424505	8.27%	1.478%
91	15.010	2034	2038	2042	rVB4	268134	378000	7.36%	1.316%
92	15.127	2054	2058	2063	rVB3	332835	497124	9.68%	1.731%
93	15.867	2178	2184	2189	rVB	1734968	2729225	53.15%	9.503%
94	16.349	2260	2266	2278	rVB	2707419	5134667	100.00%	17.879%
95	17.171	2401	2406	2416	rVB	1856501	3338969	65.03%	11.626%
96	17.365	2436	2439	2440	rBV	342022	396289	7.72%	1.380%
97	17.776	2505	2509	2514	rBV	818125	1222633	23.81%	4.257%
98	19.004	2715	2718	2723	rVB2	431185	602941	11.74%	2.099%
99	19.439	2789	2792	2796	rVB	447839	557390	10.86%	1.941%
100	19.803	2851	2854	2862	rVB2	643309	1066214	20.77%	3.713%

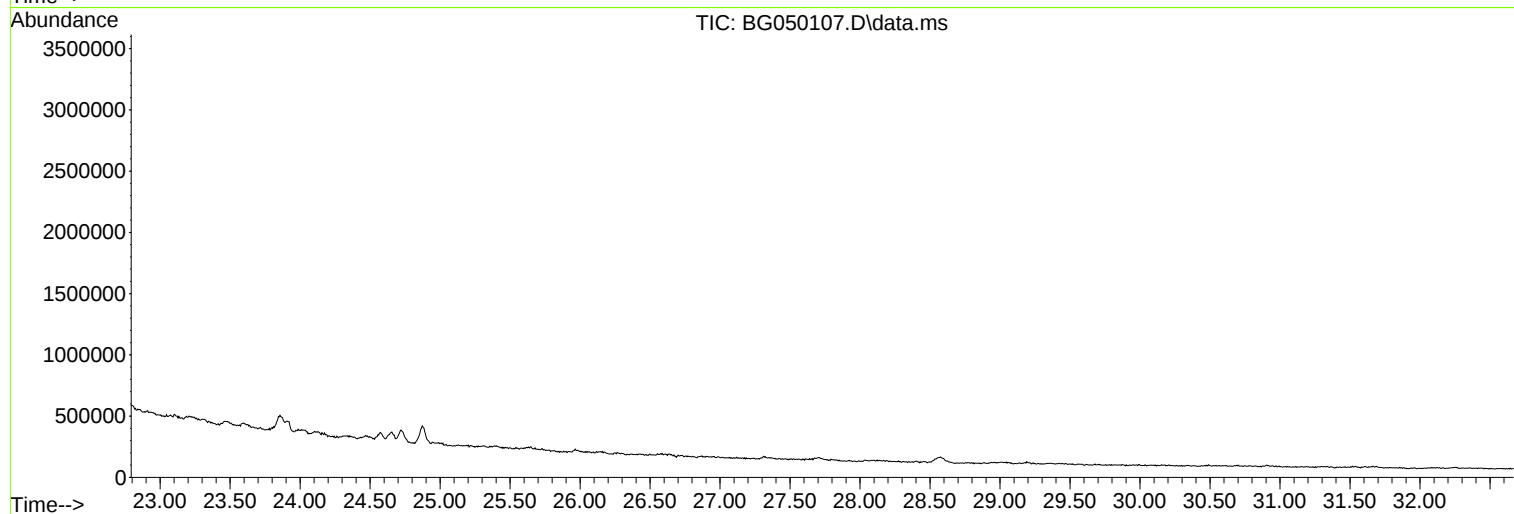
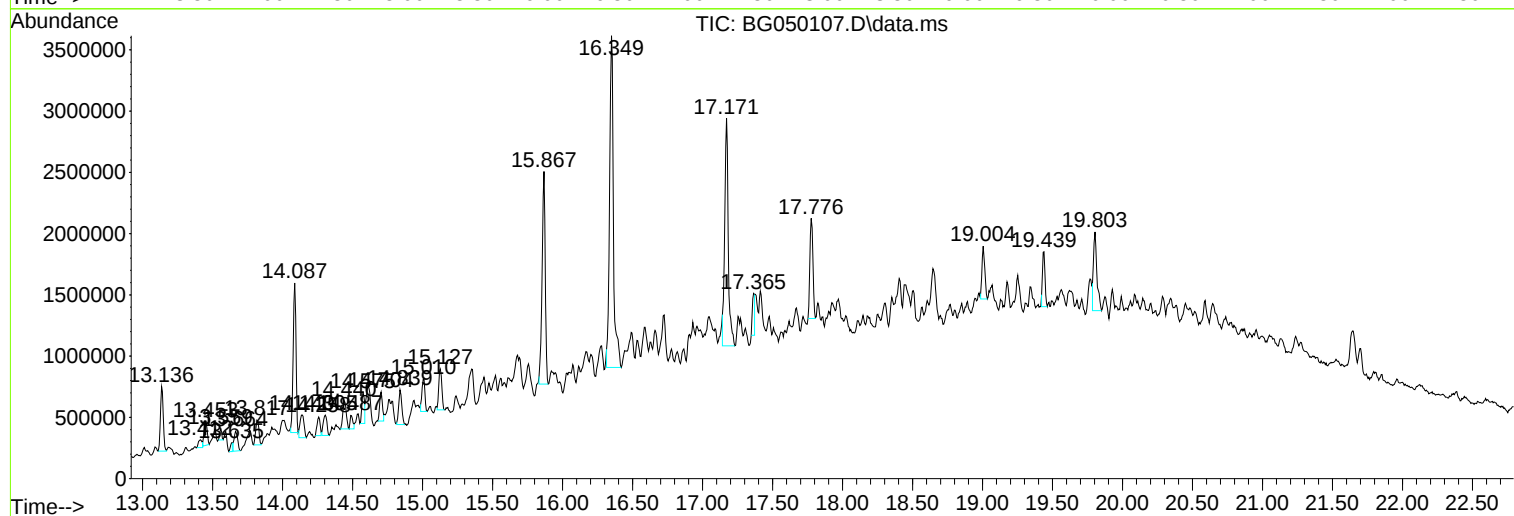
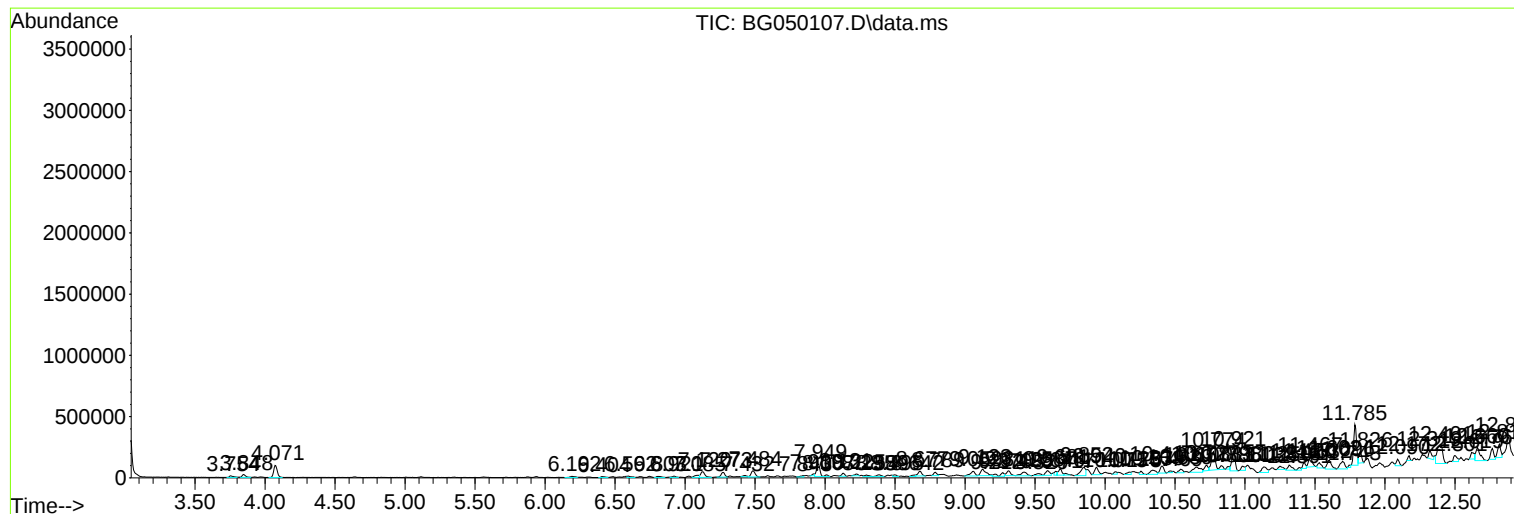
Sum of corrected areas: 28718804

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

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



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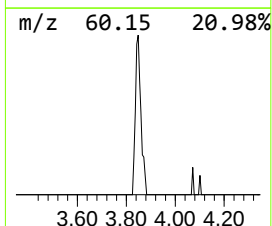
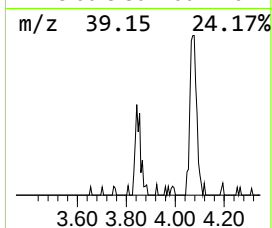
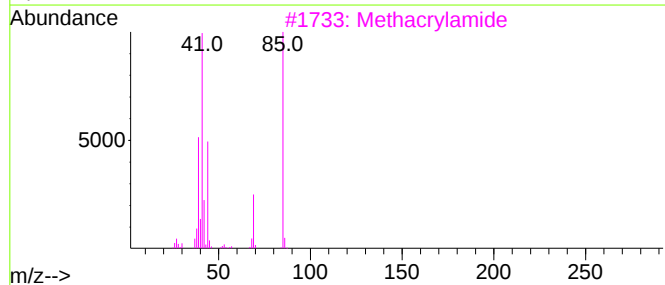
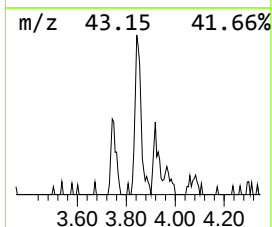
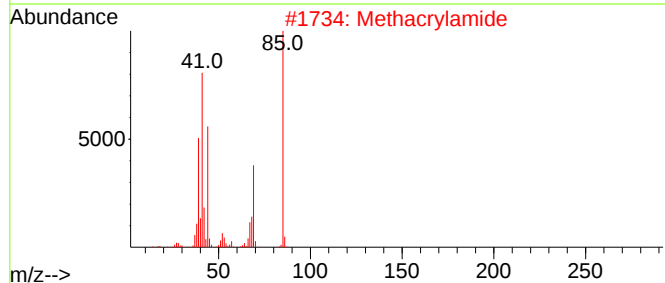
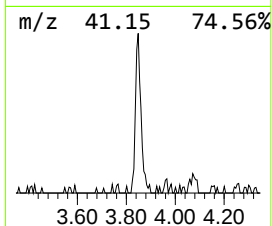
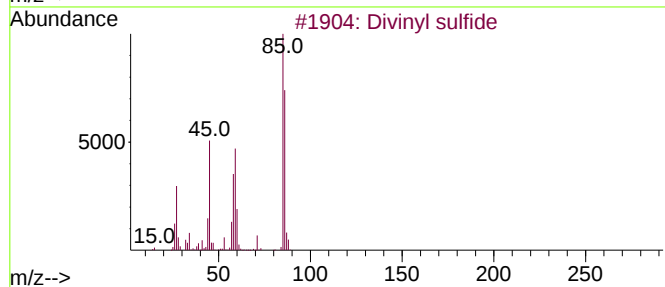
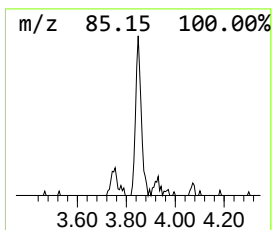
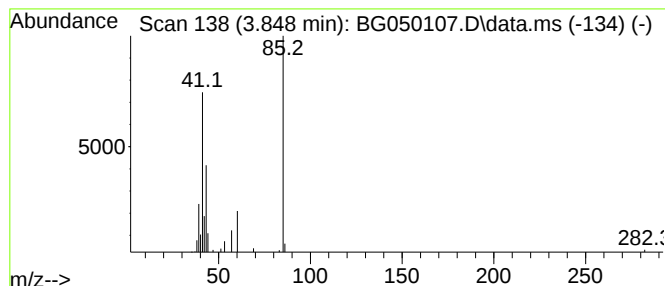
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Divinyl sulfide Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.848	4.60 ng/ul	38438	1,4-Dichlorobenzene-d4	7.949

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Divinyl sulfide	86	C4H6S	000627-51-0	50
2			Methacrylamide	85	C4H7NO	000079-39-0	38
3			Methacrylamide	85	C4H7NO	000079-39-0	38
4			Piperazine, 1-nitroso-	115	C4H9N3O	005632-47-3	38
5			2H-Pyran, 2-(4-chlorobutoxy)tetr...	192	C9H17ClO2	041302-05-0	36



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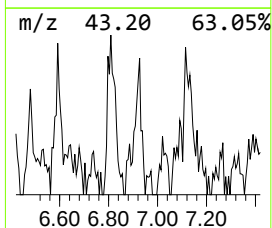
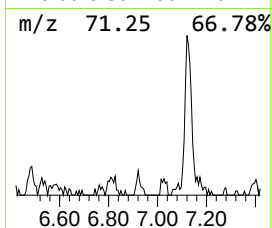
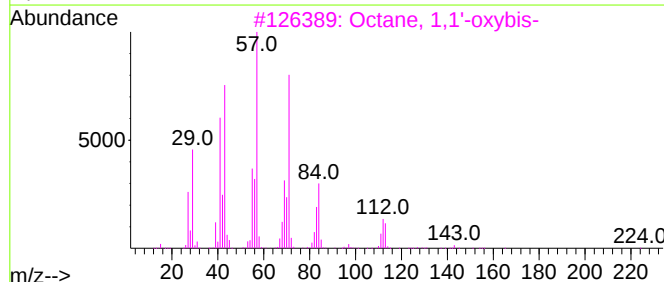
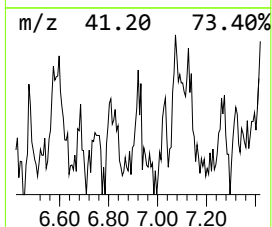
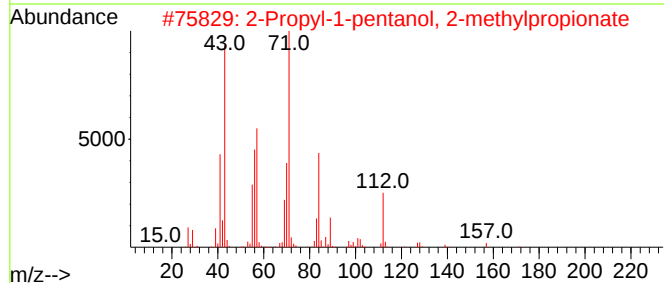
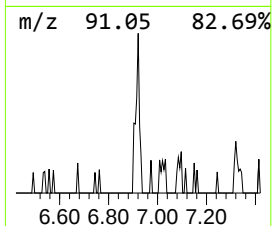
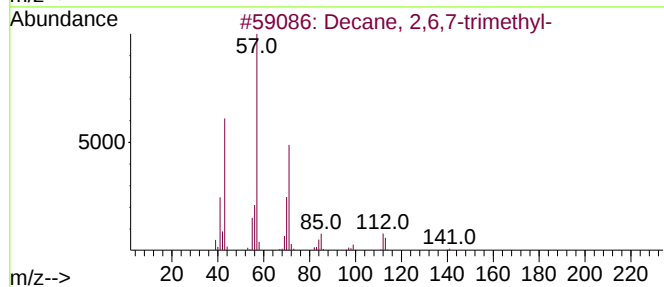
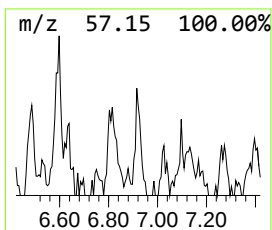
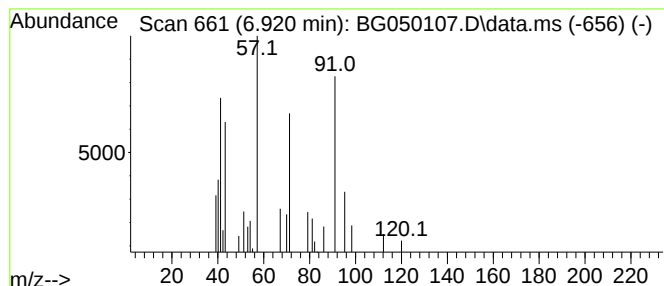
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 61

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.920	2.53 ng/ul	21145	1,4-Dichlorobenzene-d4	7.949

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 2,6,7-trimethyl-	184	C13H28	062108-25-2	16
2			2-Propyl-1-pentanol, 2-methylpro...	200	C12H24O2	1000447-36-4	16
3			Octane, 1,1'-oxybis-	242	C16H34O	000629-82-3	16
4			Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	14
5			Oxalic acid, 2-ethylhexyl hexyl ...	286	C16H30O4	1000309-38-9	12



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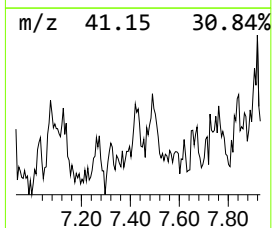
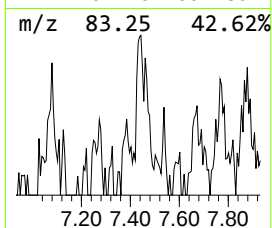
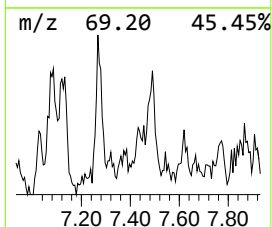
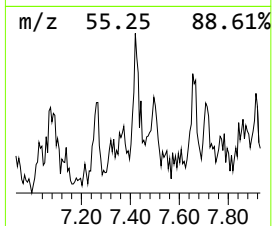
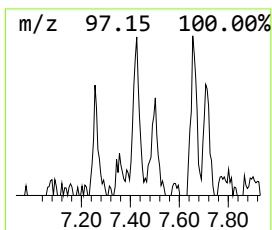
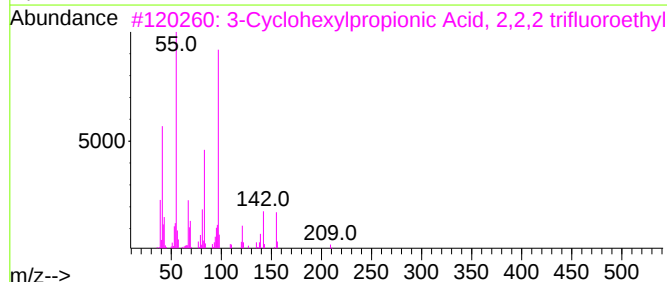
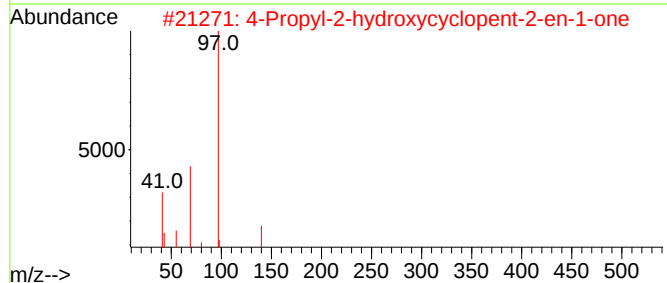
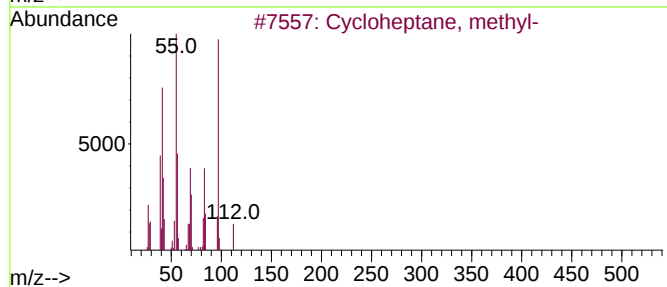
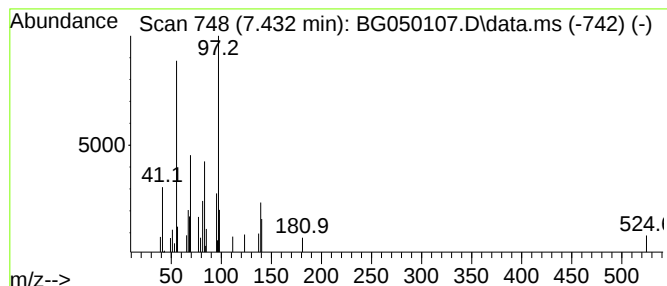
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Cyclic7.432 Concentration Rank 65

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.432	2.33 ng/ul	19425	1,4-Dichlorobenzene-d4	7.949

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cycloheptane, methyl-	112	C8H16	004126-78-7	37
2			4-Propyl-2-hydroxycyclopent-2-en...	140	C8H12O2	082147-26-0	35
3			3-Cyclohexylpropionic Acid, 2,2,...	238	C11H17F3O2	1000376-54-1	28
4			Cyclohexane, 1-(cyclohexylmethyl...	194	C14H26	054823-96-0	28
5			6-Nitroundec-5-ene	199	C11H21NO2	1000192-40-3	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampled :
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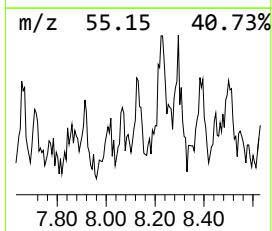
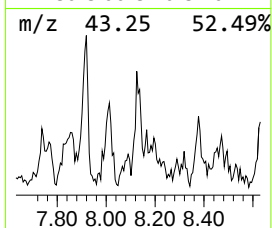
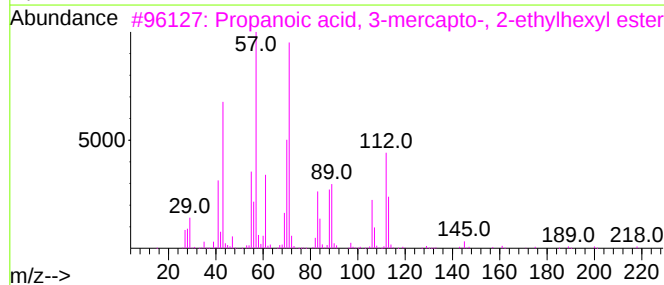
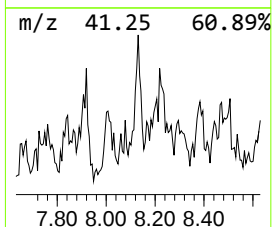
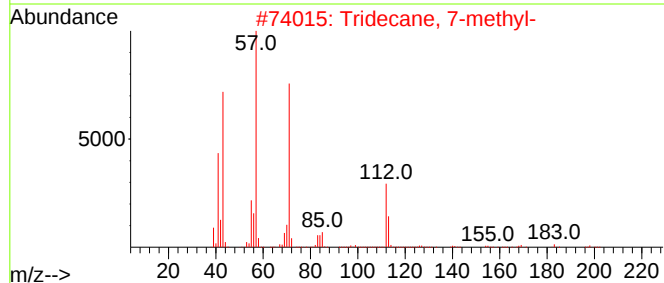
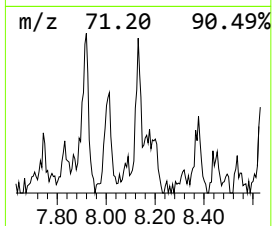
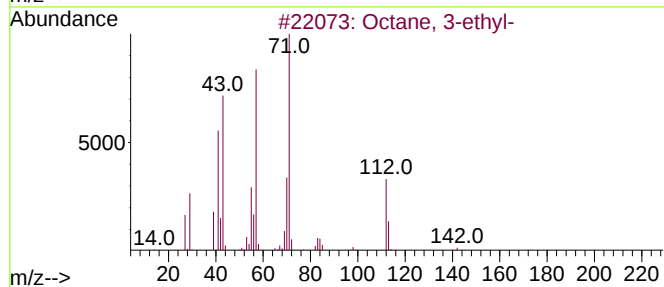
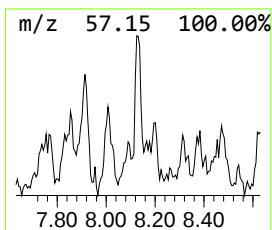
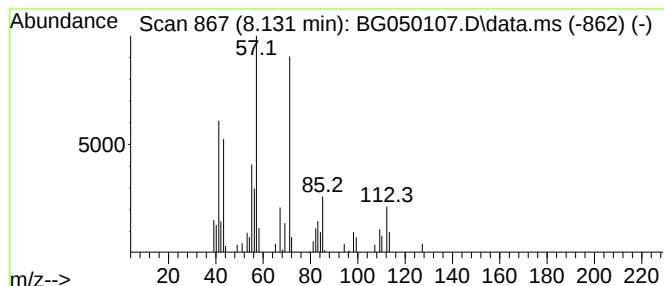
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.131	4.64 ng/ul	38735	1,4-Dichlorobenzene-d4	7.949

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 3-ethyl-			142	C10H22	005881-17-4	53
2	Tridecane, 7-methyl-			198	C14H30	026730-14-3	53
3	Propanoic acid, 3-mercapto-, 2-e...			218	C11H22O2S	050448-95-8	53
4	Dodecane, 2,6,10-trimethyl-			212	C15H32	003891-98-3	50
5	Decane, 2-methyl-			156	C11H24	006975-98-0	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
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Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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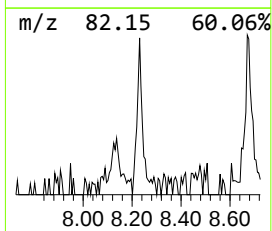
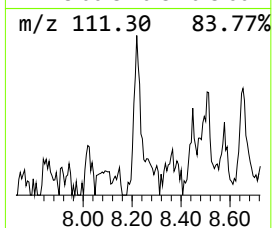
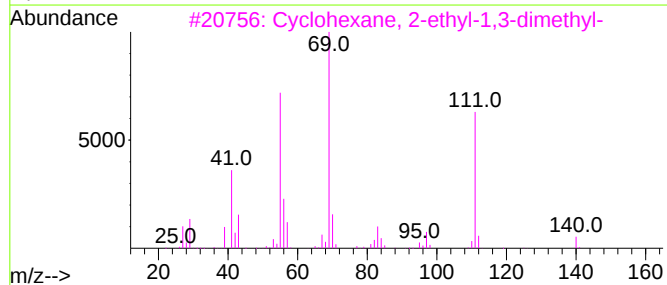
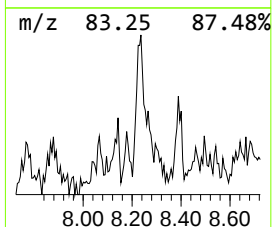
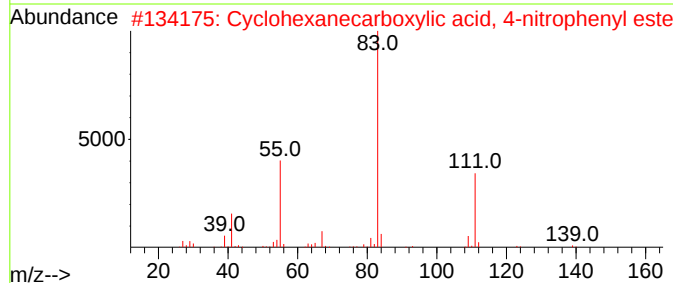
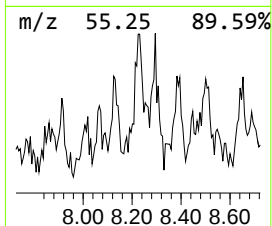
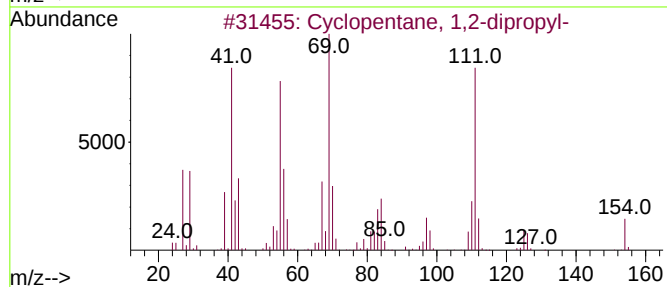
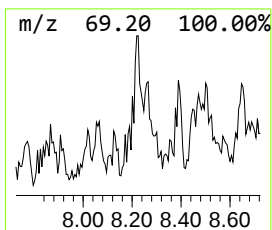
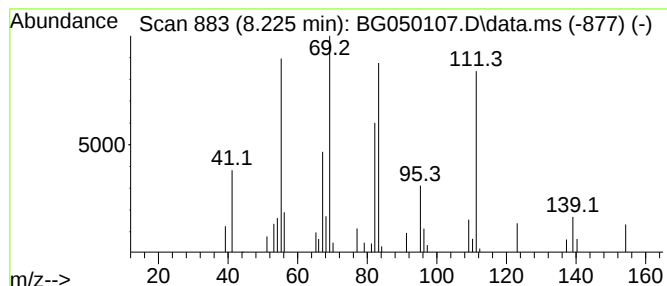
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Cyclic8.225 Concentration Rank 53

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.225	3.03 ng/ul	25299	1,4-Dichlorobenzene-d4	7.949

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, 1,2-dipropyl-	154	C11H22	091242-57-8	38
2		Cyclohexanecarboxylic acid, 4-ni...	249	C13H15NO4	1000307-70-8	38
3		Cyclohexane, 2-ethyl-1,3-dimethyl-	140	C10H20	007045-67-2	35
4		Cyclohexane, 1-ethyl-2,4-dimethyl-	140	C10H20	061142-69-6	35
5		Cyclooctane, ethyl-	140	C10H20	013152-02-8	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
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Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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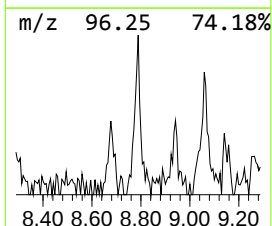
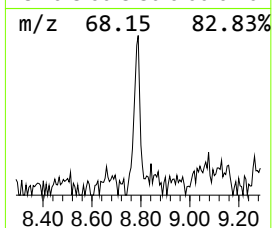
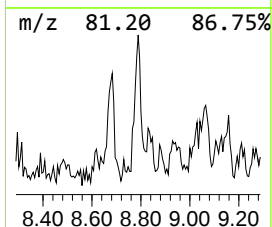
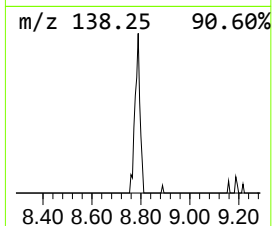
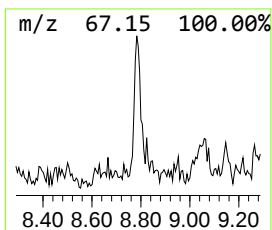
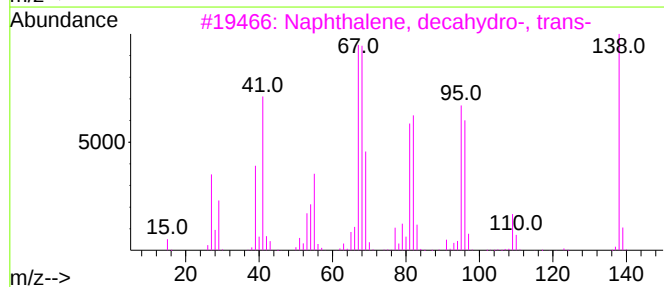
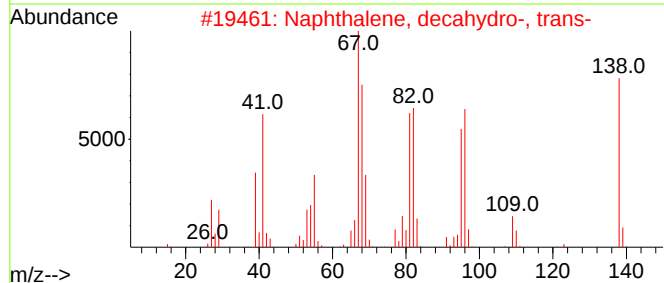
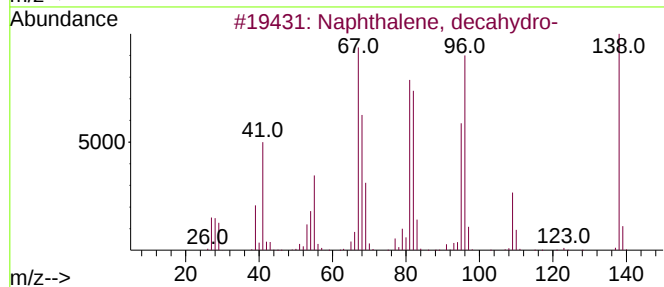
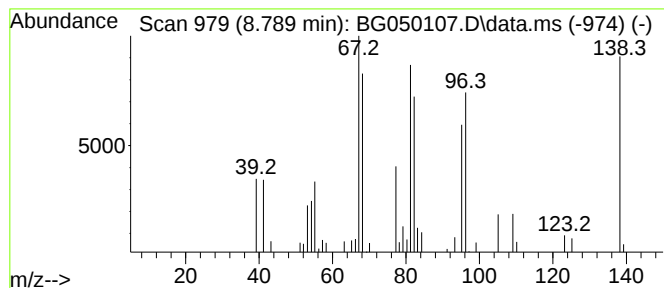
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Naphthalene, decahydro- Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.789	5.49 ng/ul	45834	1,4-Dichlorobenzene-d4	7.949

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, decahydro-	138	C10H18	000091-17-8	93
2			Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	89
3			Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	89
4			Naphthalene, decahydro-	138	C10H18	000091-17-8	87
5			Naphthalene, decahydro-	138	C10H18	000091-17-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
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Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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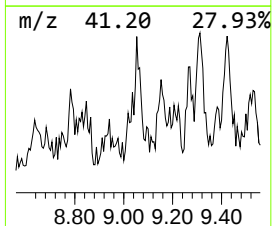
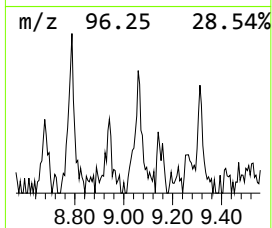
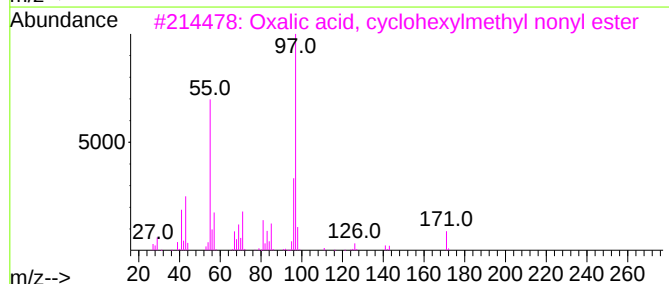
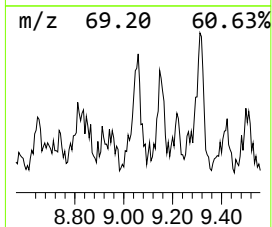
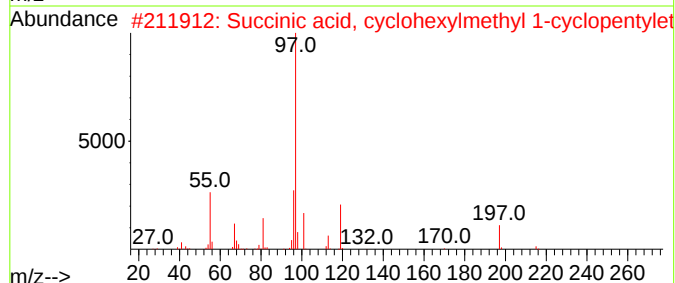
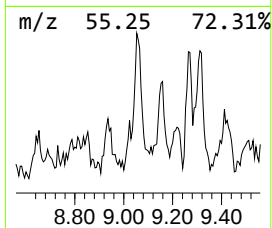
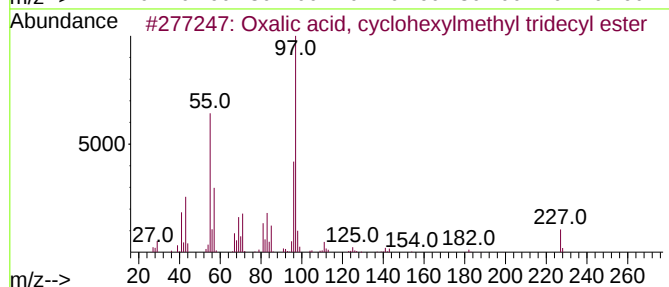
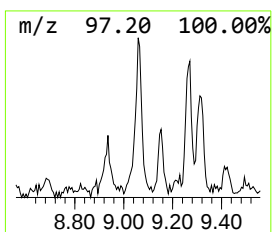
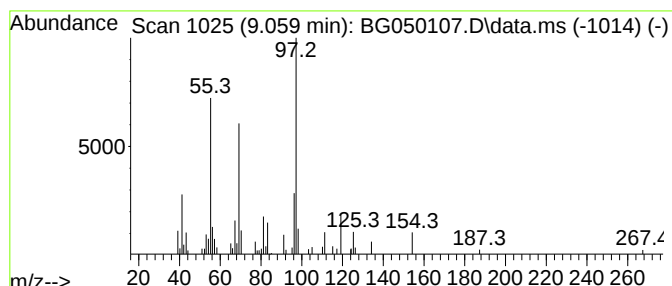
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Oxalic acid, cyclohexylmeth... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.059	11.33 ng/ul	94634	1,4-Dichlorobenzene-d4	7.949

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, cyclohexylmethyl tr...	368	C22H40O4	1000309-69-1	53
2			Succinic acid, cyclohexylmethyl ...	310	C18H30O4	1000391-41-3	53
3			Oxalic acid, cyclohexylmethyl no...	312	C18H32O4	1000309-68-6	53
4			Diborane(6), tetrapropyl-	196	C12H30B2	022784-01-6	53
5			1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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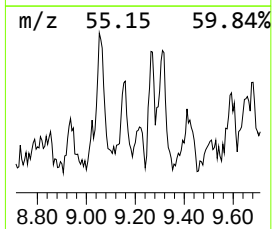
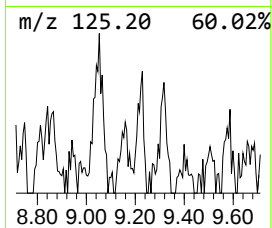
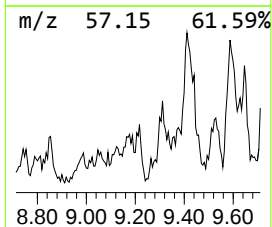
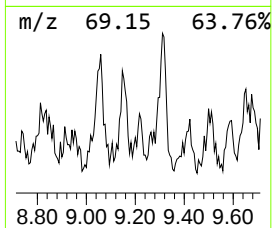
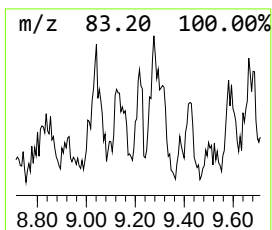
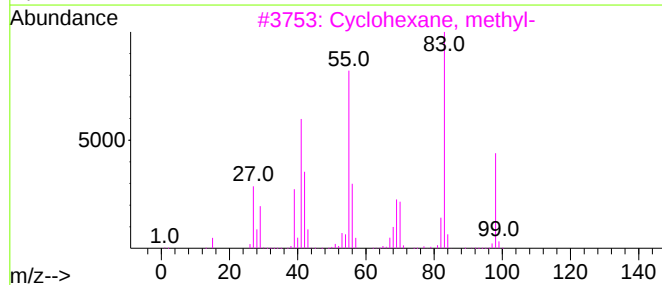
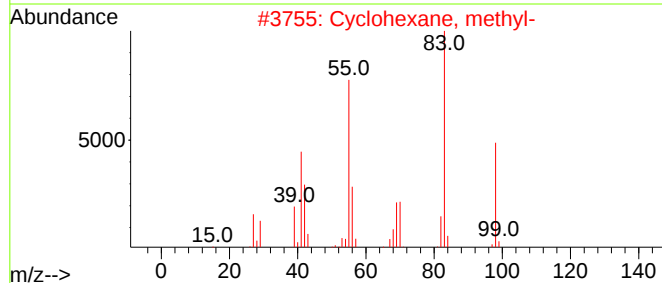
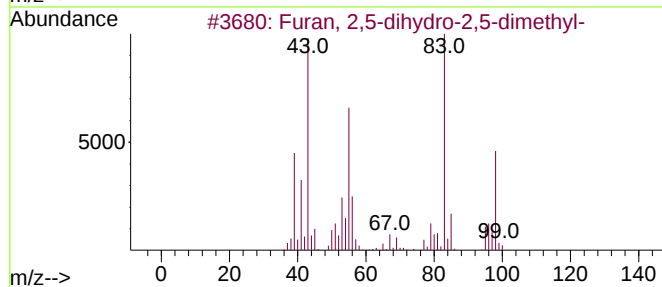
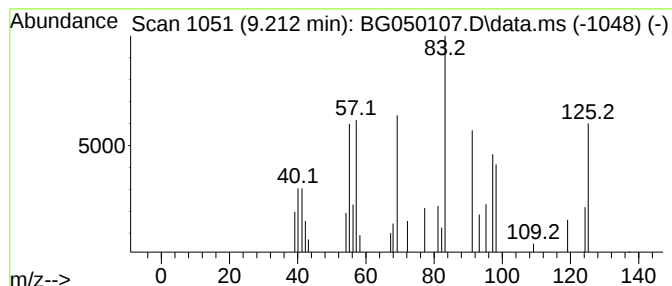
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 unknown-01 Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.212	3.58 ng/ul	29885	1,4-Dichlorobenzene-d4	7.949

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Furan, 2,5-dihydro-2,5-dimethyl-	98	C6H10O	059242-27-2	38
2			Cyclohexane, methyl-	98	C7H14	000108-87-2	38
3			Cyclohexane, methyl-	98	C7H14	000108-87-2	35
4			Cyclohexane, methyl-	98	C7H14	000108-87-2	35
5			1-Butyl-1H-1,2,4-triazole	125	C6H11N3	006086-22-2	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
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Misc :
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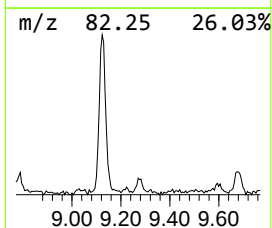
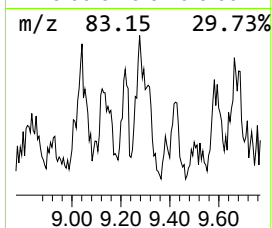
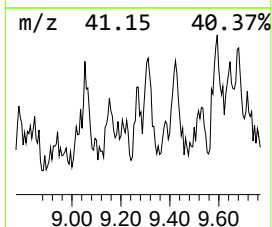
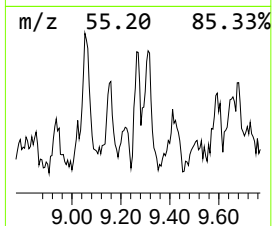
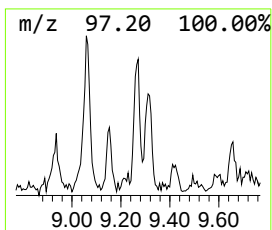
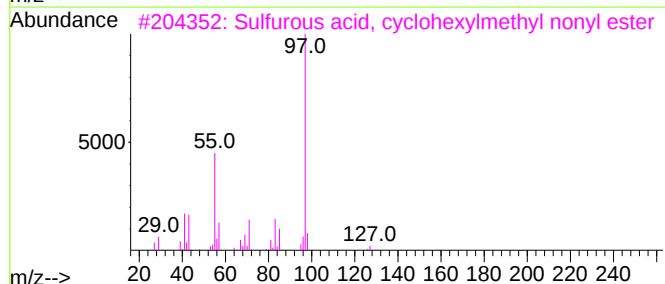
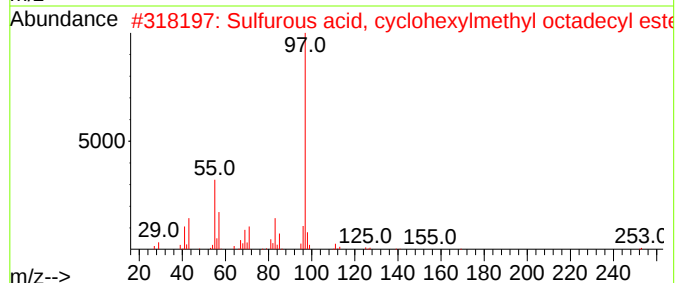
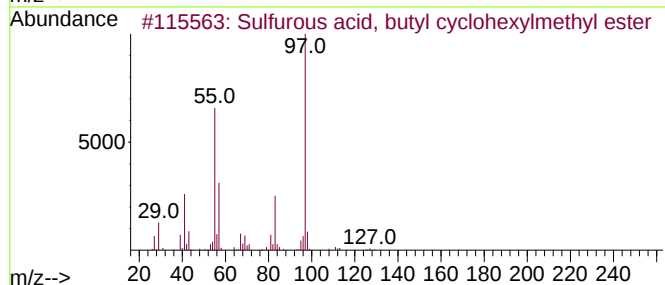
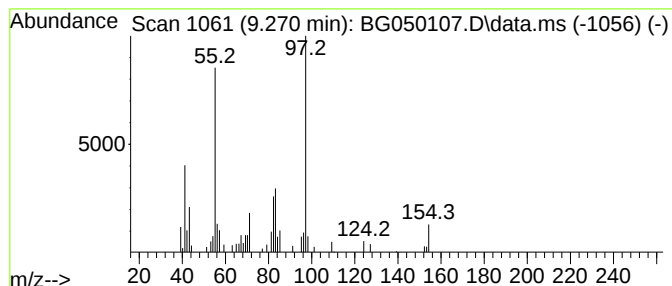
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 Sulfurous acid, butyl cyclo... Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.270	5.48 ng/ul	45771	1,4-Dichlorobenzene-d4	7.949

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, butyl cyclohexyl...	234	C11H22O3S	1000309-21-4	53
2			Sulfurous acid, cyclohexylmethyl...	430	C25H50O3S	1000309-22-6	53
3			Sulfurous acid, cyclohexylmethyl...	304	C16H32O3S	1000309-21-8	50
4			Sulfurous acid, cyclohexylmethyl...	402	C23H46O3S	1000309-22-4	50
5			Cyclohexane, 1,1-dimethyl-	112	C8H16	000590-66-9	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
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Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

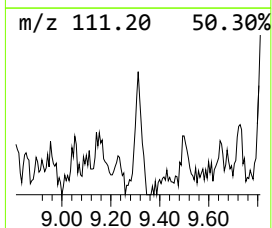
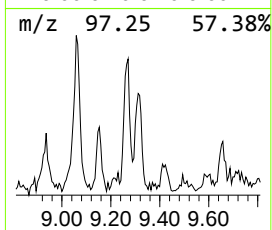
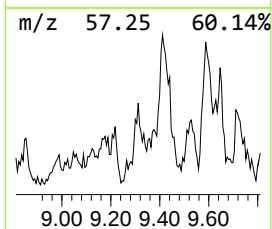
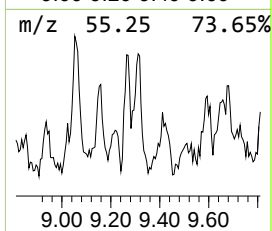
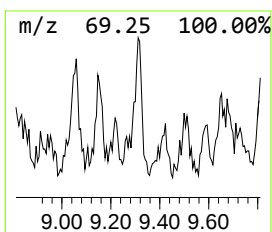
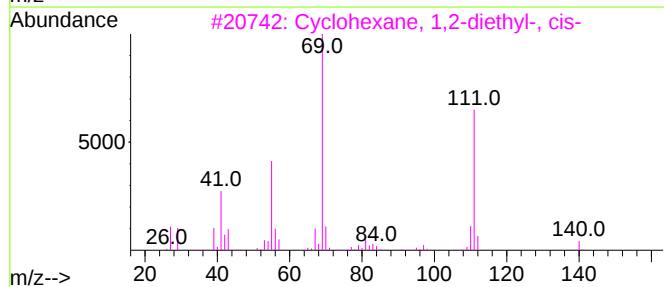
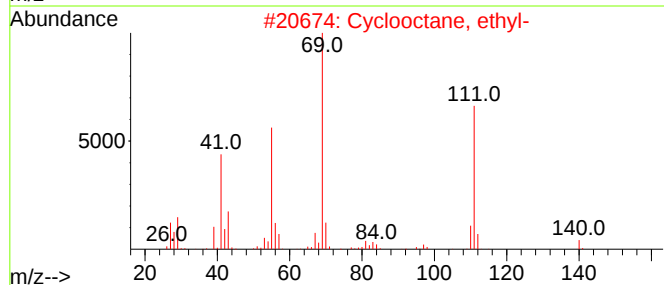
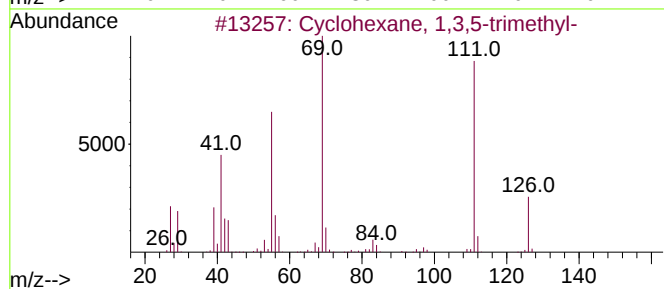
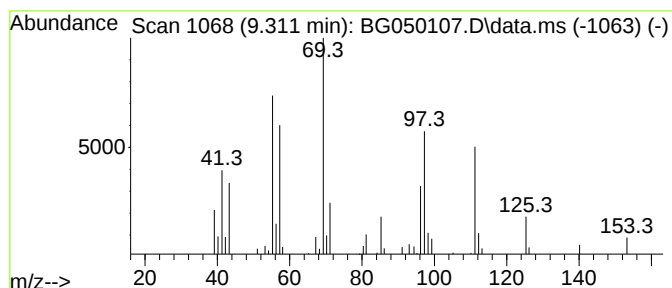
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

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

Peak Number 20 (DEL) Alkane: Cyclic9.311 Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.311	9.67 ng/ul	80811	1,4-Dichlorobenzene-d4	7.949	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	43
2	Cyclooctane, ethyl-	140	C10H20	013152-02-8	43
3	Cyclohexane, 1,2-diethyl-, cis-	140	C10H20	000824-43-1	43
4	Cyclooctane, ethyl-	140	C10H20	013152-02-8	38
5	2,4-Dimethyl-1,5-diazabicyclo[3....	112	C6H12N2	1000283-20-9	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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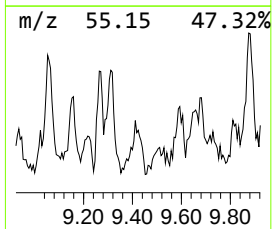
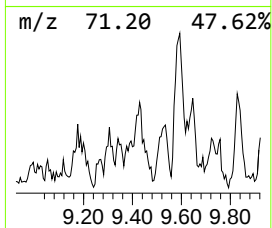
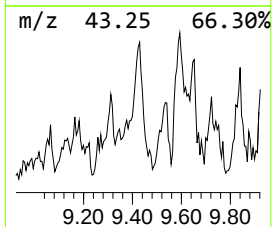
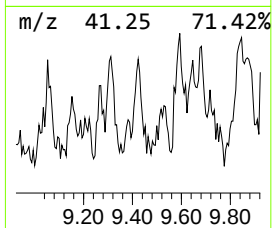
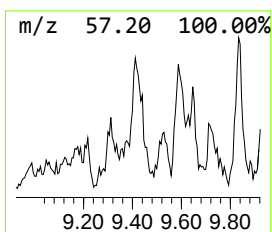
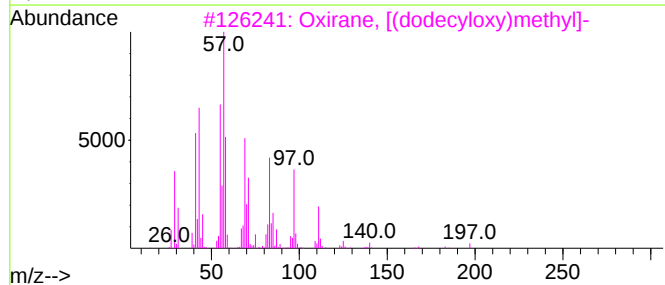
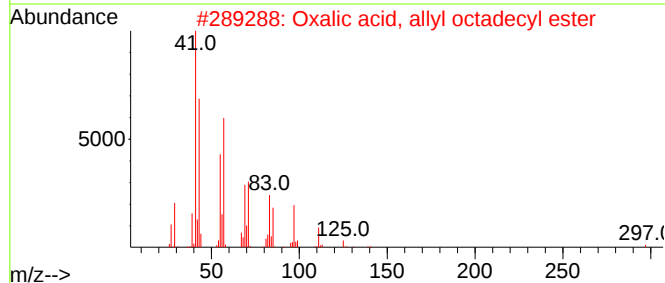
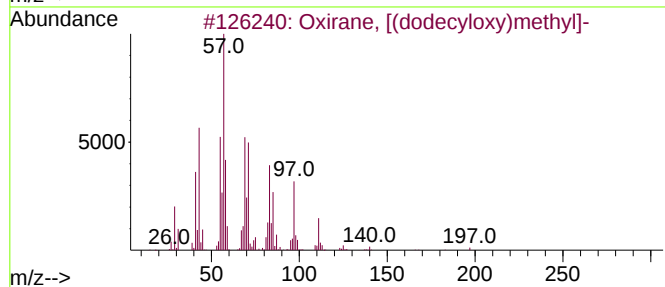
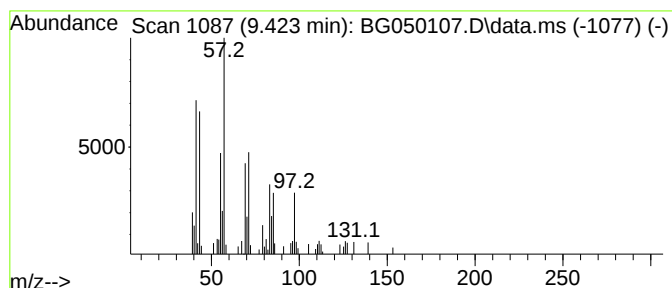
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 Oxirane, [(dodecyloxy)methyl]- Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.423	6.13 ng/ul	86372	Naphthalene-d8	10.774

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Oxirane, [(dodecyloxy)methyl]-	242	C15H30O2	002461-18-9	87
2		Oxalic acid, allyl octadecyl ester	382	C23H42O4	1000309-24-5	80
3		Oxirane, [(dodecyloxy)methyl]-	242	C15H30O2	002461-18-9	53
4		Butyl decyl ether	214	C14H30O	1000406-40-2	53
5		Decane, 2,4-dimethyl-	170	C12H26	002801-84-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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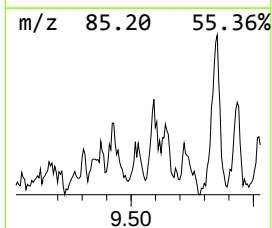
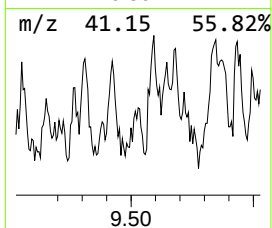
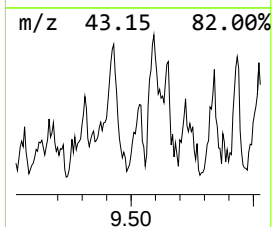
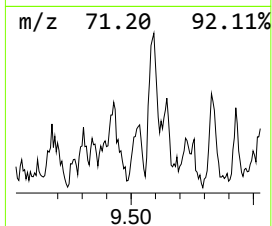
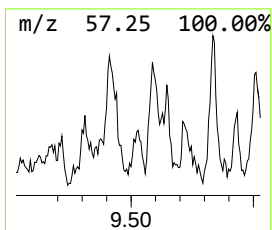
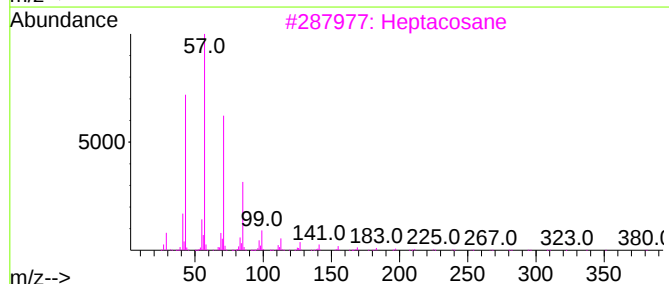
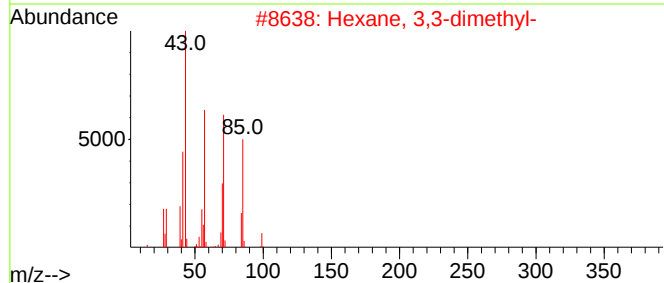
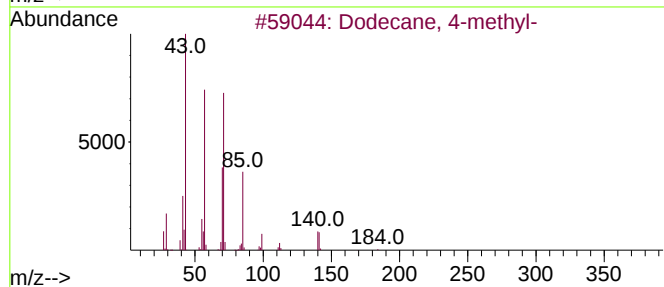
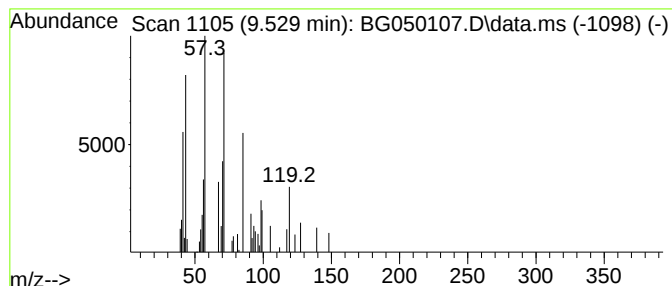
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 55

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.529	2.95 ng/ul	41599	Naphthalene-d8	10.774

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Dodecane, 4-methyl-	184	C13H28	006117-97-1	50
2		Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	47
3		Heptacosane	380	C27H56	000593-49-7	47
4		Nonane, 4,5-dimethyl-	156	C11H24	017302-23-7	47
5		Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW7C4

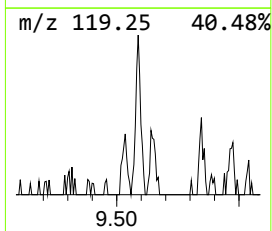
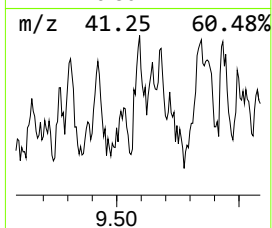
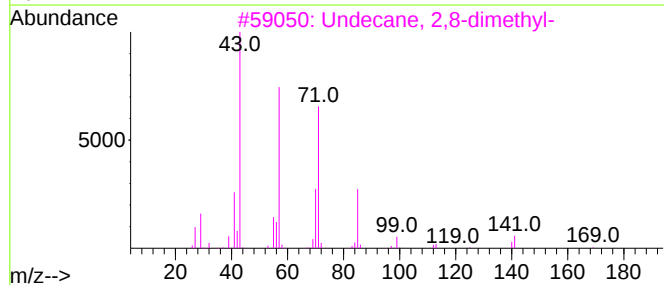
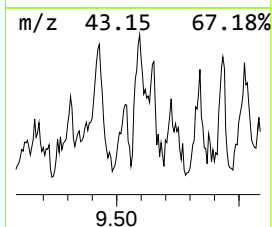
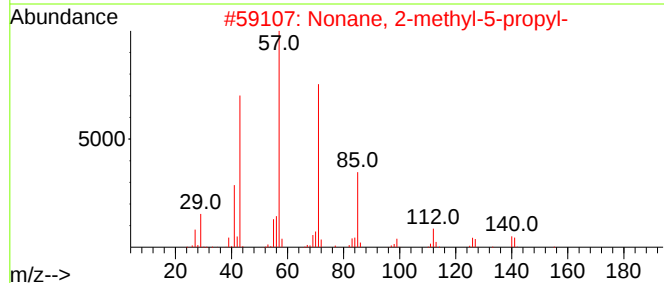
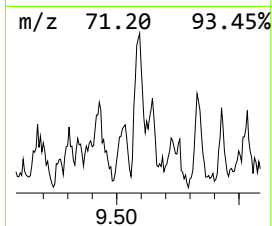
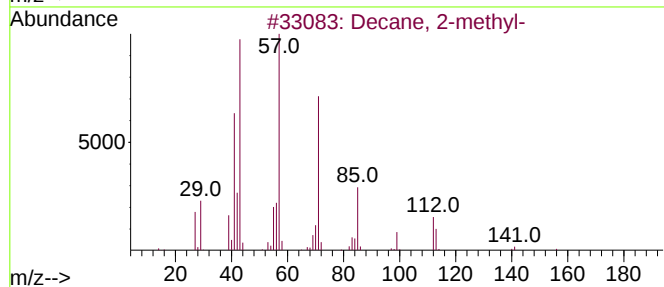
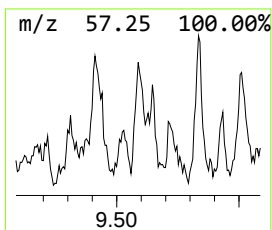
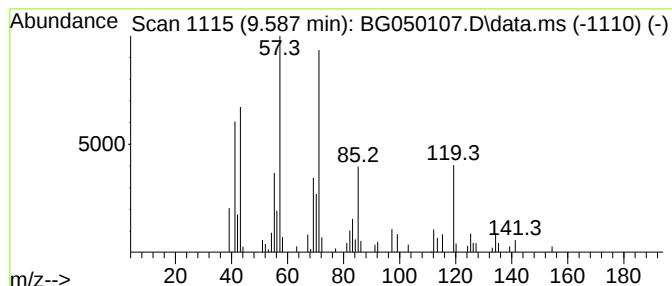
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.587	5.86 ng/ul	82653	Naphthalene-d8	10.774

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane, 2-methyl-	156	C11H24	006975-98-0	58
2		Nonane, 2-methyl-5-propyl-	184	C13H28	031081-17-1	58
3		Undecane, 2,8-dimethyl-	184	C13H28	017301-25-6	50
4		Decane, 2,3,8-trimethyl-	184	C13H28	062238-14-6	50
5		Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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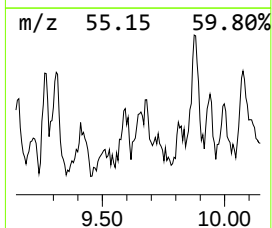
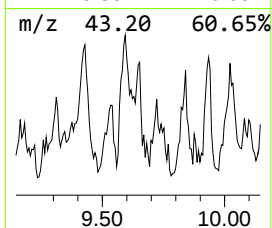
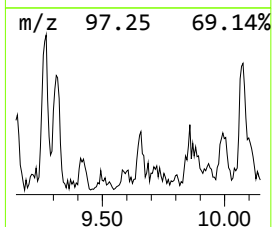
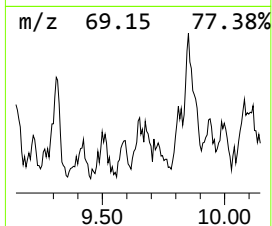
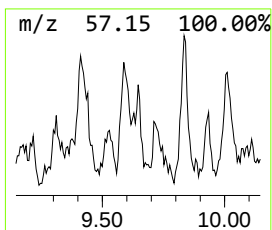
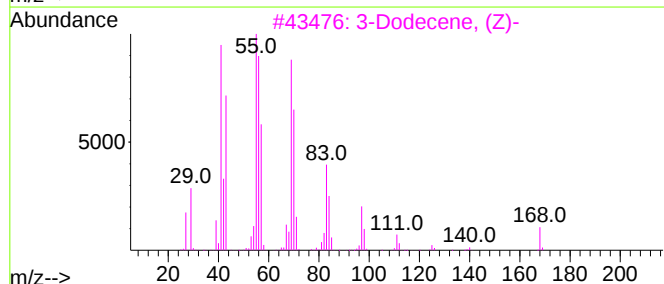
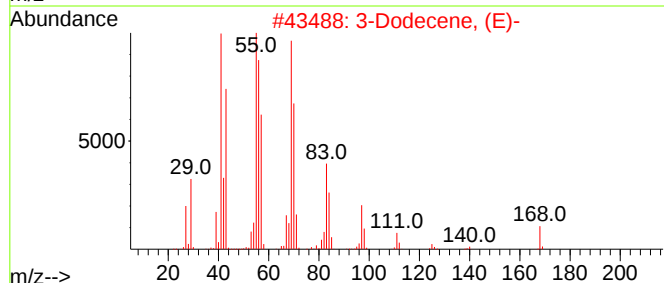
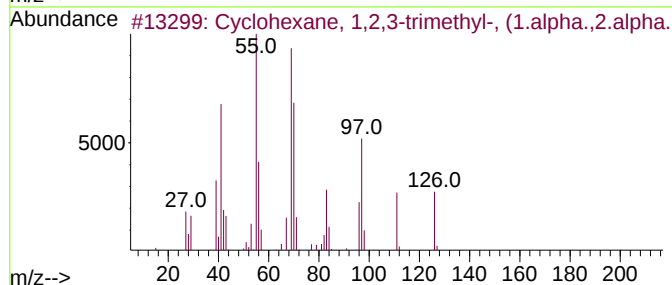
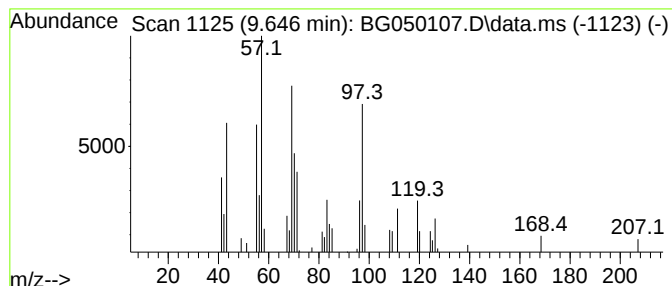
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 (DEL) Alkane: Cyclic9.646 Concentration Rank 56

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.646	2.77 ng/ul	39012	Naphthalene-d8	10.774

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2,3-trimethyl-, (...)	126	C9H18	007667-55-2	76
2			3-Dodecene, (E)-	168	C12H24	007206-14-6	43
3			3-Dodecene, (Z)-	168	C12H24	007239-23-8	43
4			2,2,4-Trimethyl-3-hexene	126	C9H18	059643-72-0	43
5			5-Nonanol, pentafluoropropionate	290	C12H19F5O2	1000463-47-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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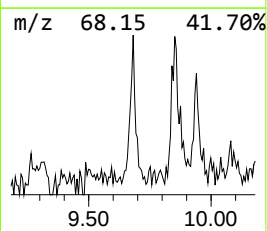
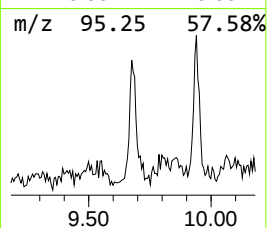
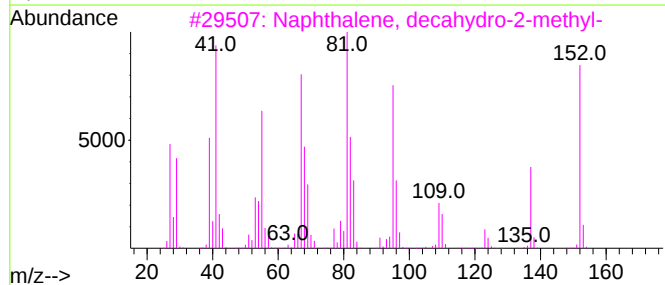
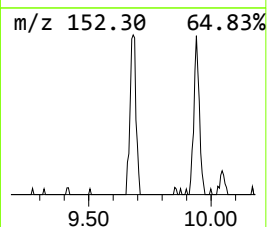
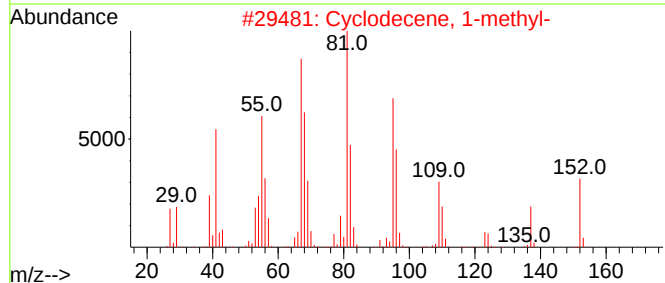
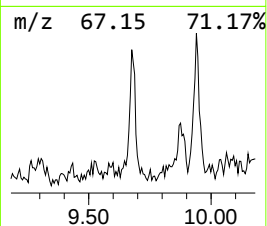
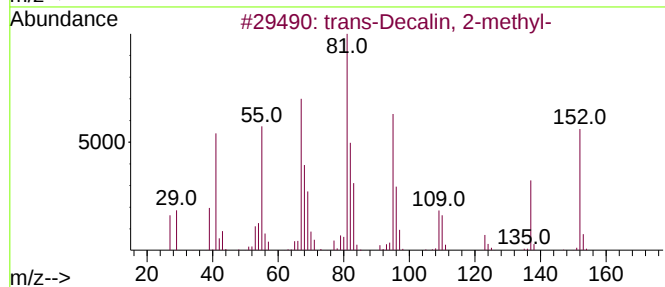
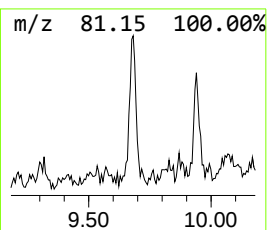
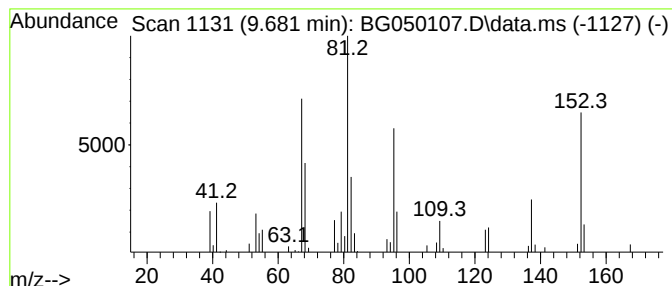
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 trans-Decalin, 2-methyl- Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.681	6.26 ng/ul	88230	Naphthalene-d8	10.774

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	91
2			Cyclodecene, 1-methyl-	152	C11H20	066633-38-3	86
3			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	86
4			Furan, 2-hexyl-	152	C10H16O	003777-70-6	68
5			Cyclohexanone, 2-methyl-5-(1-met...	152	C10H16O	007764-50-3	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
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Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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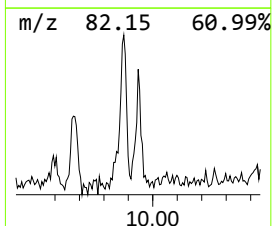
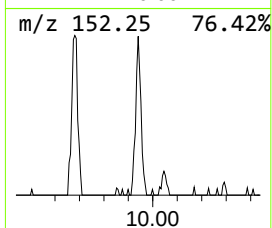
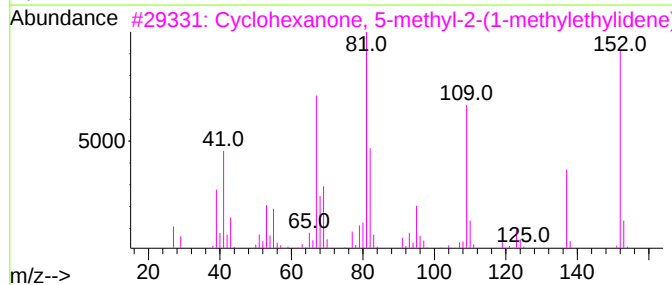
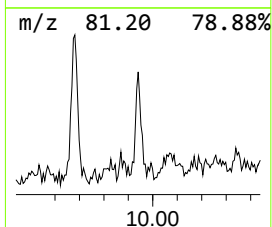
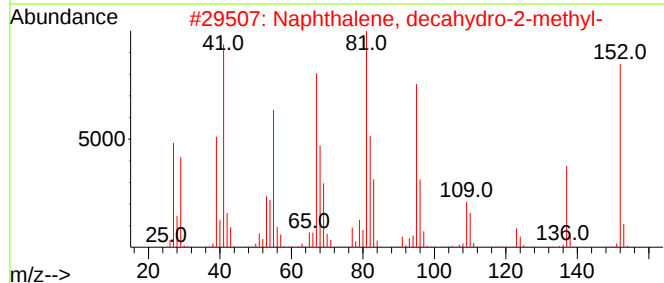
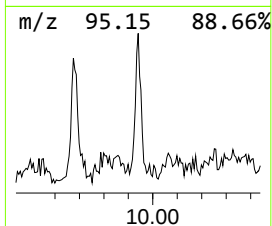
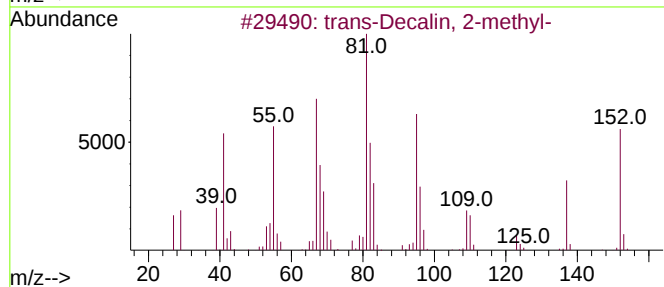
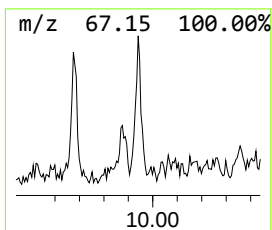
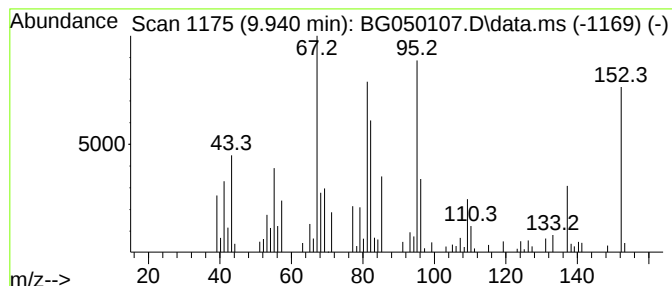
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 Naphthalene, decahydro-2-me... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.940	7.83 ng/ul	110302	Naphthalene-d8	10.774

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	62
2			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	58
3			Cyclohexanone, 5-methyl-2-(1-met...	152	C10H16O	015932-80-6	55
4			4,5-Nonadiene, 2-methyl-	138	C10H18	055956-32-6	52
5			Cyclohexanone, 5-methyl-2-(1-met...	152	C10H16O	015932-80-6	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW7C4

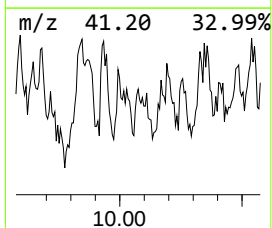
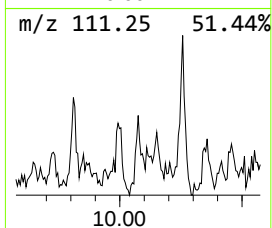
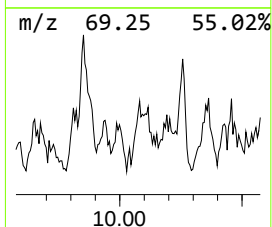
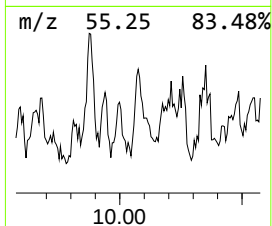
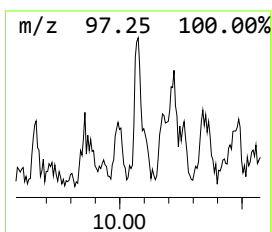
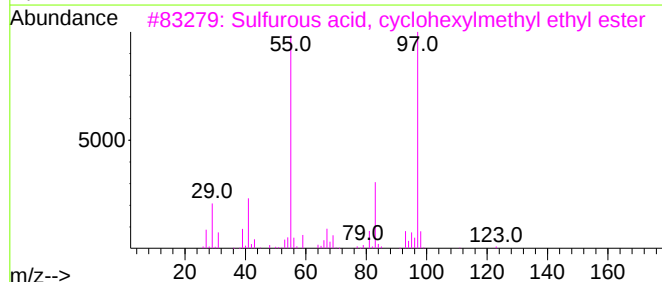
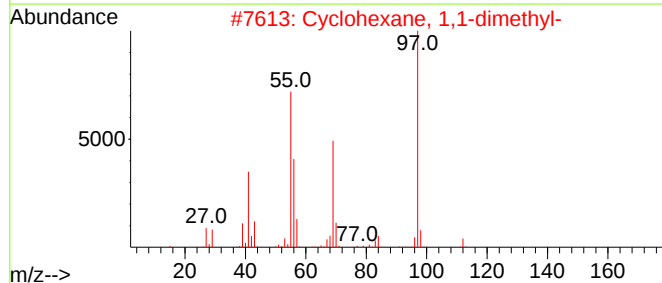
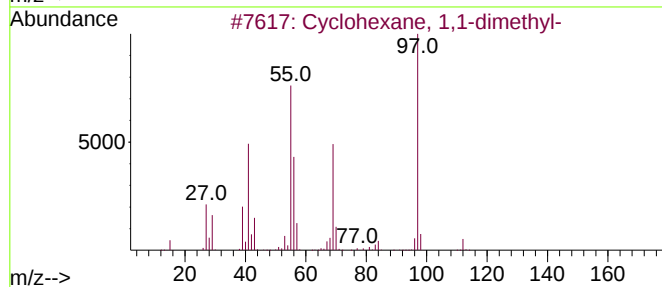
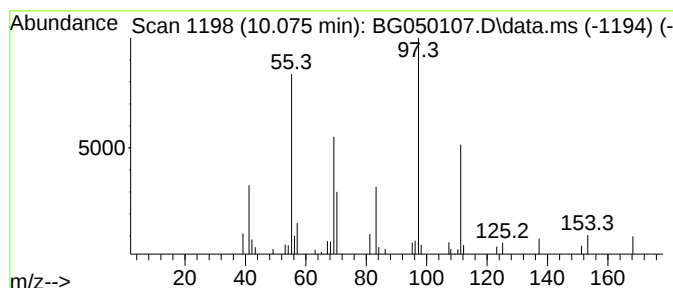
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 27 (DEL) Alkane: Cyclic10.075 Concentration Rank 68

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.075	2.06 ng/ul	29044	Naphthalene-d8	10.774	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,1-dimethyl-	112	C8H16	000590-66-9	50
2	Cyclohexane, 1,1-dimethyl-	112	C8H16	000590-66-9	45
3	Sulfurous acid, cyclohexylmethyl...	206	C9H18O3S	1000309-21-2	43
4	Cyclohexane, 1,1-dimethyl-	112	C8H16	000590-66-9	41
5	Sulfurous acid, cyclohexylmethyl...	346	C19H38O3S	1000309-22-0	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW7C4

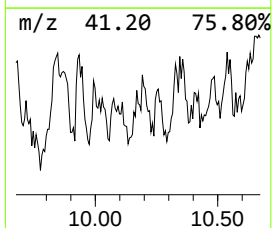
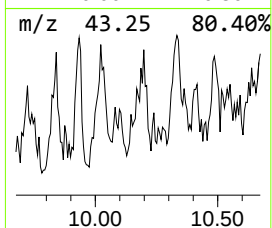
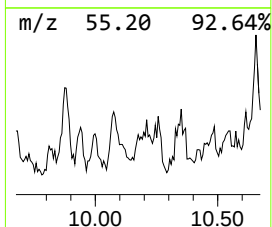
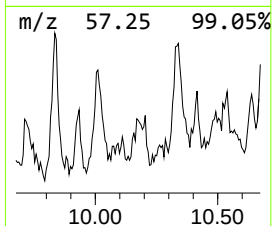
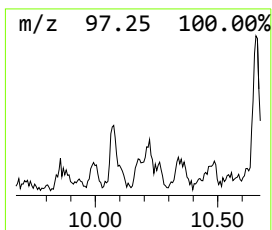
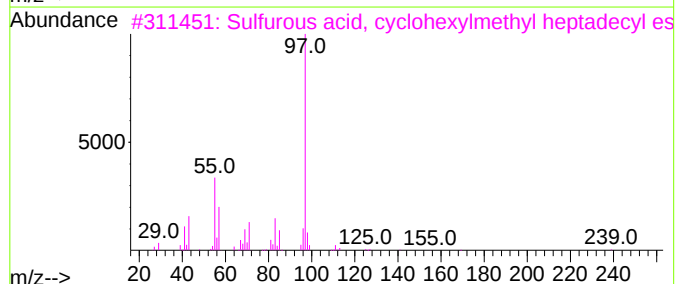
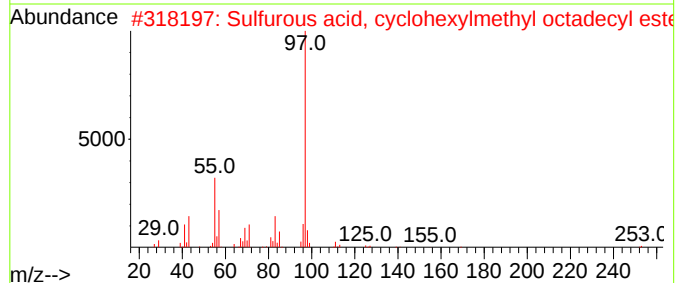
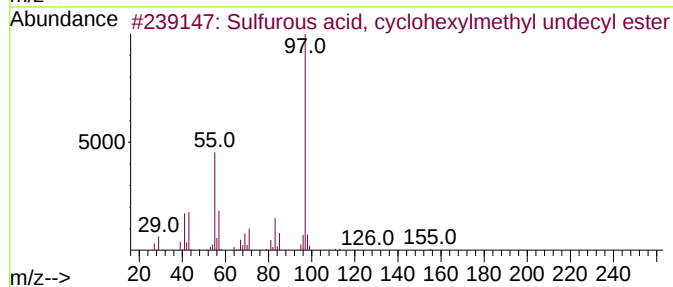
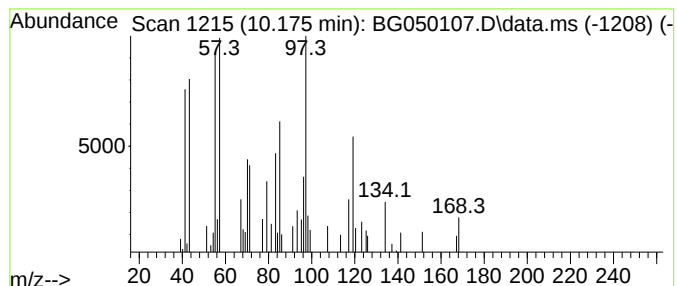
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Quant Title : SVOA CALIBRATION

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

Peak Number 28 unknown-02 Concentration Rank 54

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.175	3.01 ng/ul	42490	Naphthalene-d8	10.774

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, cyclohexylmethyl...	332	C18H36O3S	1000309-21-9	46
2			Sulfurous acid, cyclohexylmethyl...	430	C25H50O3S	1000309-22-6	46
3			Sulfurous acid, cyclohexylmethyl...	416	C24H48O3S	1000309-22-5	46
4			Oxalic acid, cyclobutyl pentadec...	354	C21H38O4	1010309-70-5	35
5			Succinic acid, 2-methylpent-3-yl...	298	C17H30O4	1000390-06-8	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

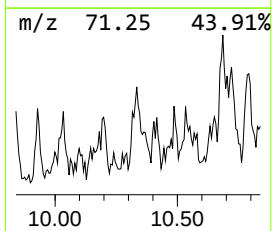
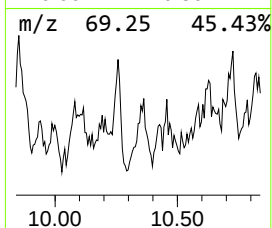
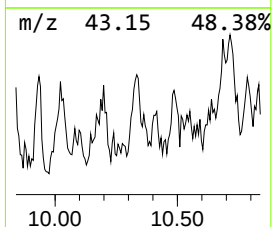
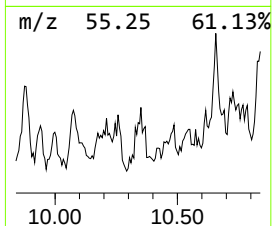
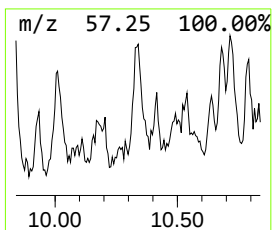
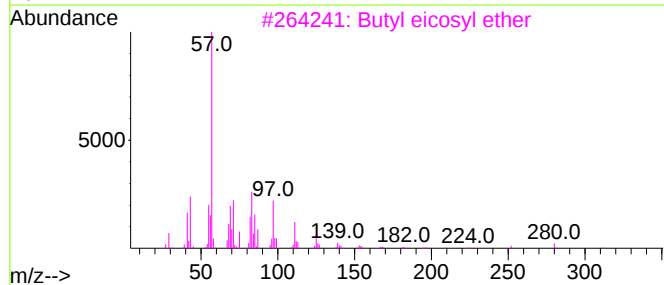
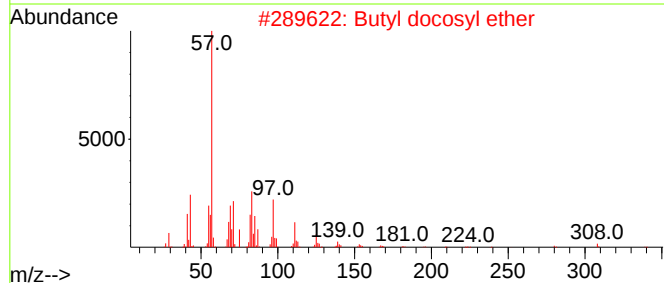
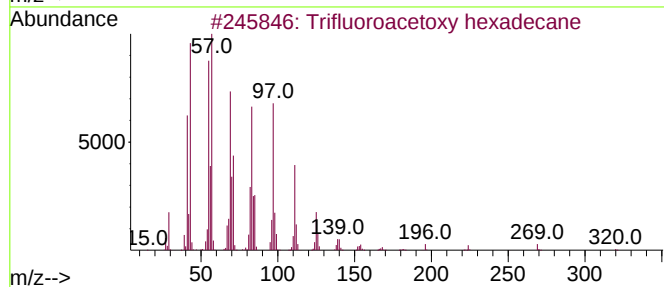
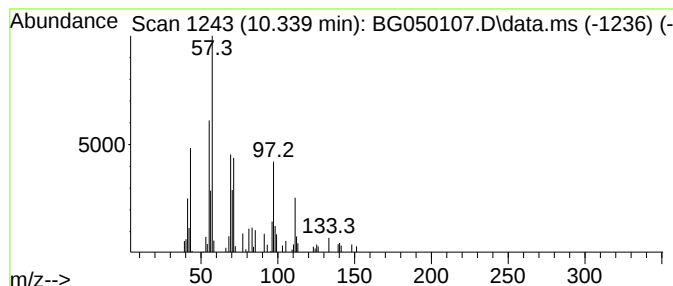
Instrument :
BNA_G
ClientSampleId :
EW7C4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

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

Peak Number 30 Trifluoroacetoxy hexadecane Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD		R.T.	
10.339	7.05 ng/ul	99304	Naphthalene-d8		10.774	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Trifluoroacetoxy hexadecane		338	C18H33F3O2	006222-03-3	64
2	Butyl docosyl ether		382	C26H54O	1000406-40-8	53
3	Butyl eicosyl ether		354	C24H50O	1000406-40-7	53
4	5-Octadecene, (E)-		252	C18H36	007206-21-5	43
5	Decyl isobutyl carbonate		258	C15H30O3	959226-07-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW7C4

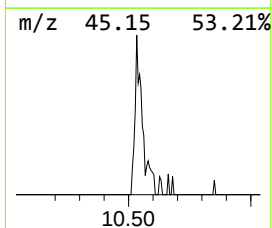
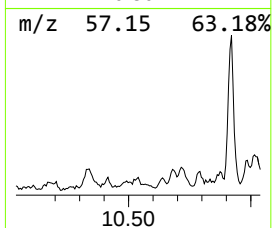
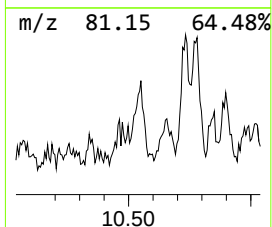
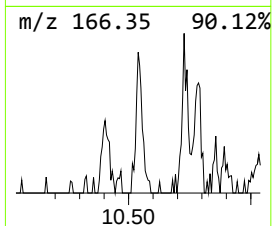
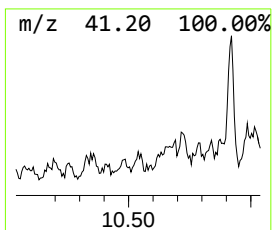
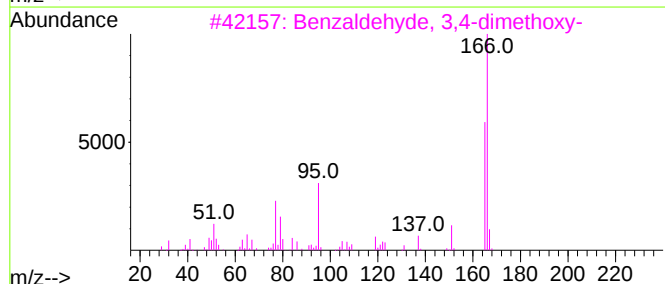
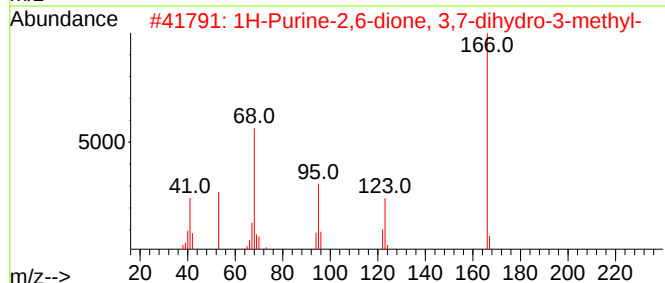
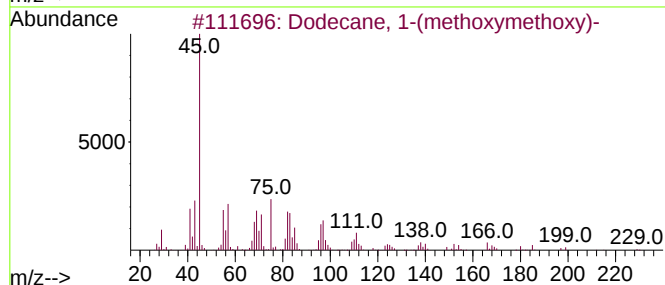
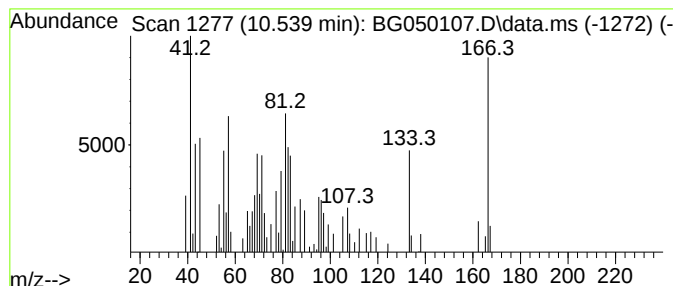
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Quant Title : SVOA CALIBRATION

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

Peak Number 31 unknown-03 Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.539	4.02 ng/ul	56610	Naphthalene-d8	10.774

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 1-(methoxymethoxy)-	230	C14H30O2	034458-41-8	16
2			1H-Purine-2,6-dione, 3,7-dihydro...	166	C6H6N4O2	001076-22-8	14
3			Benzaldehyde, 3,4-dimethoxy-	166	C9H10O3	000120-14-9	14
4			Benzaldehyde, 2,5-dimethoxy-	166	C9H10O3	000093-02-7	14
5			2,5-Cyclohexadiene-1,4-dione, 2,...	166	C8H10N2O2	075083-35-1	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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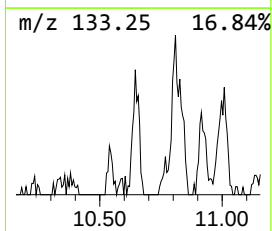
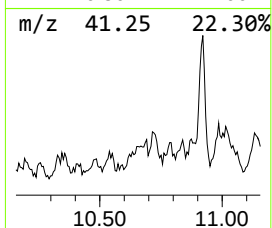
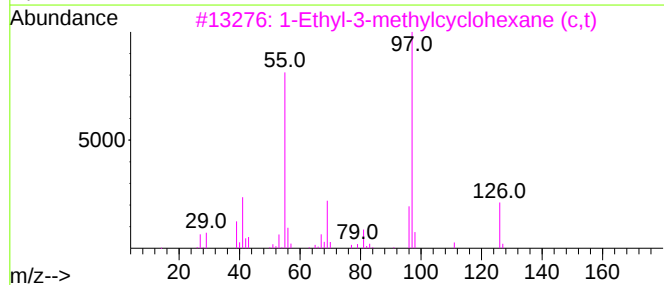
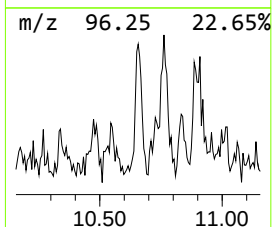
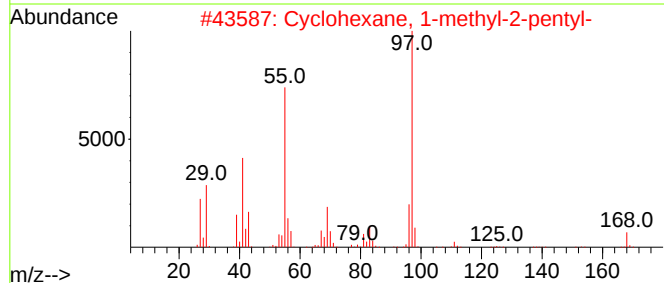
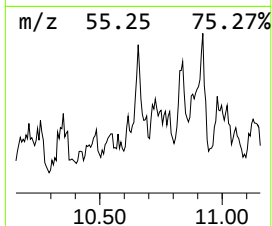
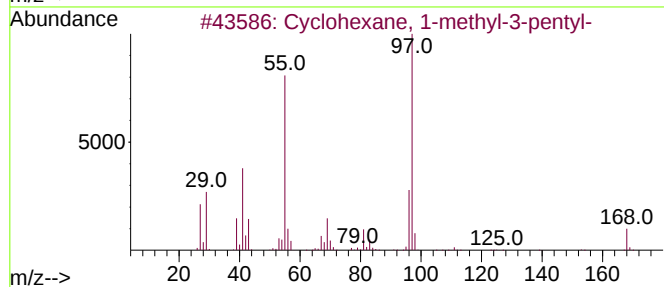
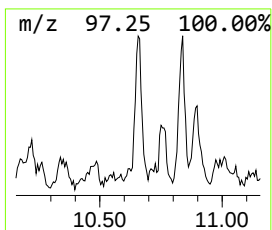
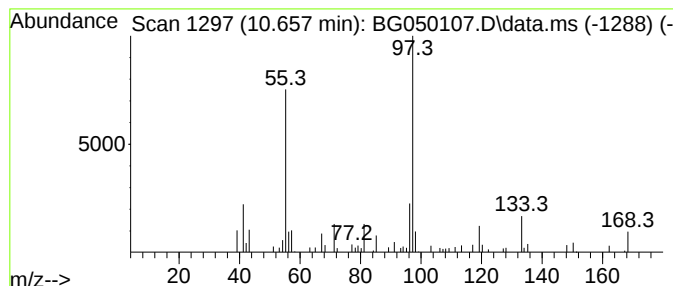
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Quant Title : SVOA CALIBRATION

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

Peak Number 32 (DEL) Alkane: Cyclic10.657 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.657	10.78 ng/ul	151938	Naphthalene-d8	10.774

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, 1-methyl-3-pentyl-	168	C12H24	054411-02-8	74
2		Cyclohexane, 1-methyl-2-pentyl-	168	C12H24	054411-01-7	72
3		1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	59
4		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	59
5		Ethanone, 1-(1-methylcyclohexyl)-	140	C9H16O	002890-62-2	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
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Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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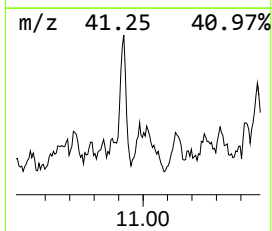
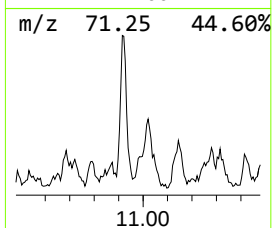
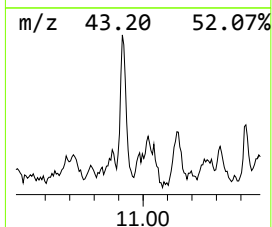
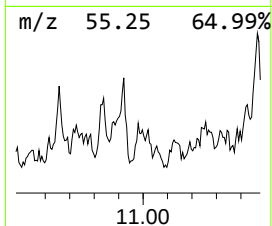
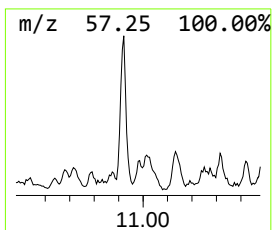
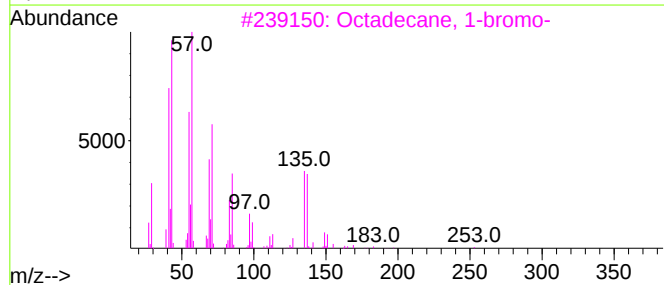
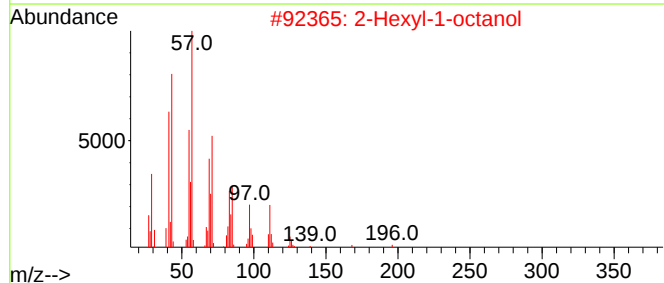
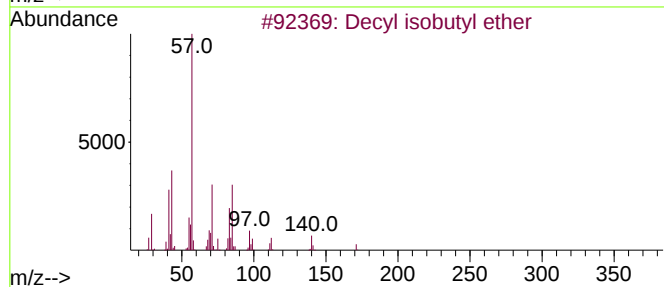
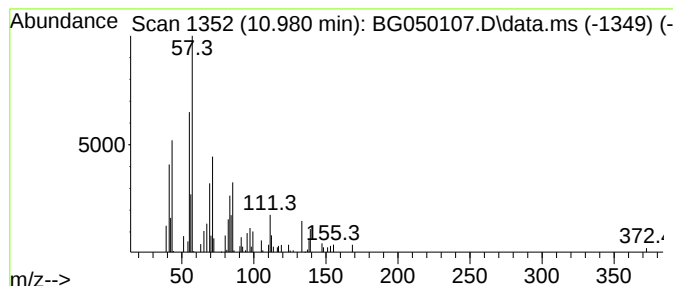
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Quant Title : SVOA CALIBRATION

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

Peak Number 33 Decyl isobutyl ether Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.980	4.46 ng/ul	62869	Naphthalene-d8	10.774

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decyl isobutyl ether	214	C14H30O	1010406-32-5	53
2			2-Hexyl-1-octanol	214	C14H30O	019780-79-1	53
3			Octadecane, 1-bromo-	332	C18H37Br	000112-89-0	50
4			Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	47
5			Hexadecane, 1-chloro-	260	C16H33Cl	004860-03-1	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
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ALS Vial : 9 Sample Multiplier: 1

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ClientSampleId :
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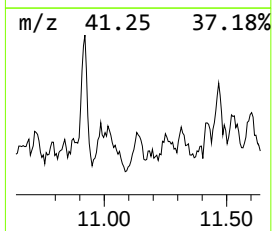
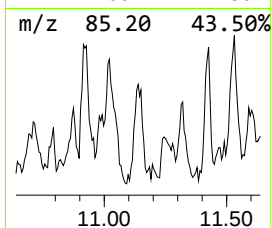
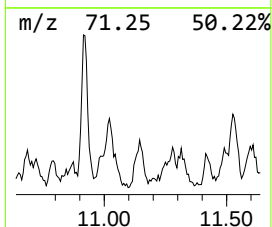
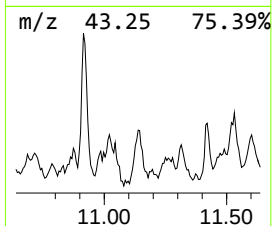
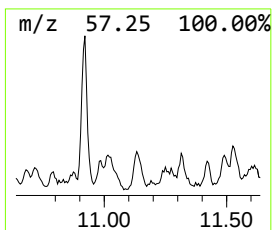
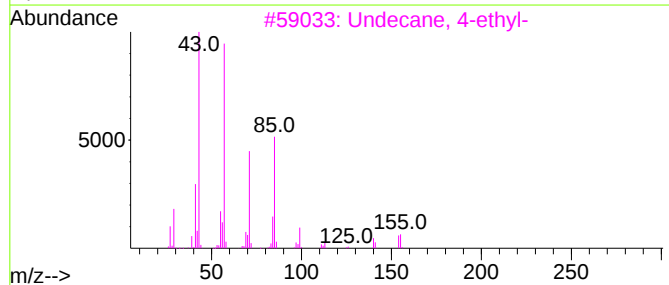
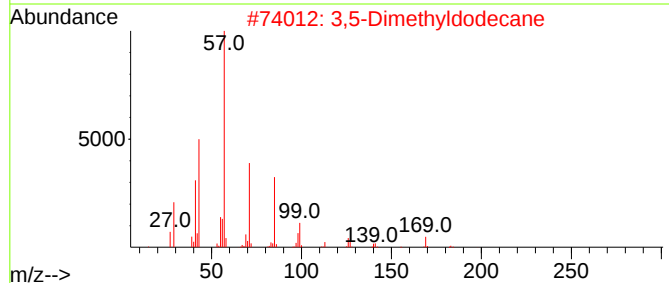
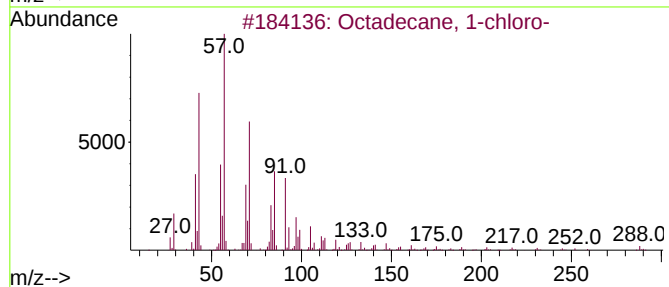
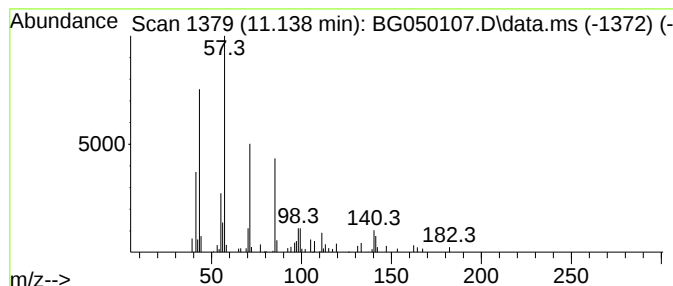
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Quant Title : SVOA CALIBRATION

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

Peak Number 34 Octadecane, 1-chloro- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.138	8.45 ng/ul	119044	Naphthalene-d8	10.774

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane, 1-chloro-		288	C18H37Cl		003386-33-2	78
2	3,5-Dimethyldodecane		198	C14H30		107770-99-0	64
3	Undecane, 4-ethyl-		184	C13H28		017312-59-3	59
4	Tetracosane		338	C24H50		000646-31-1	59
5	Oxalic acid, bis(6-ethyloct-3-yl...		370	C22H42O4		1000309-34-6	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW7C4

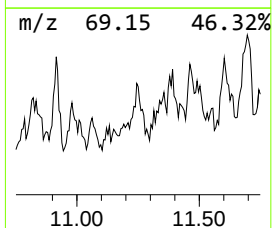
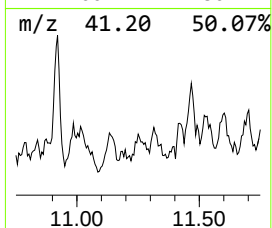
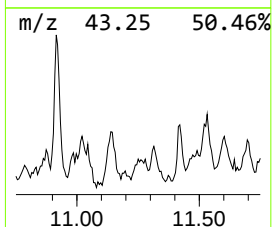
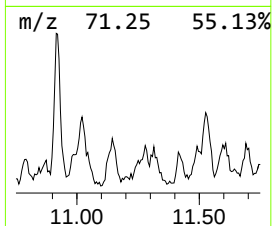
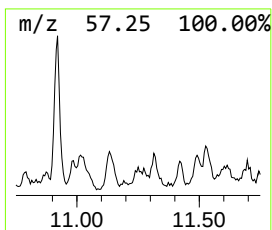
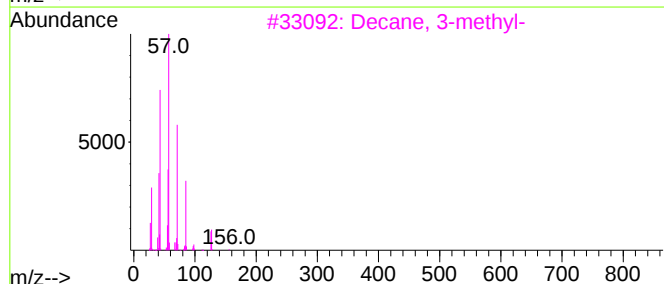
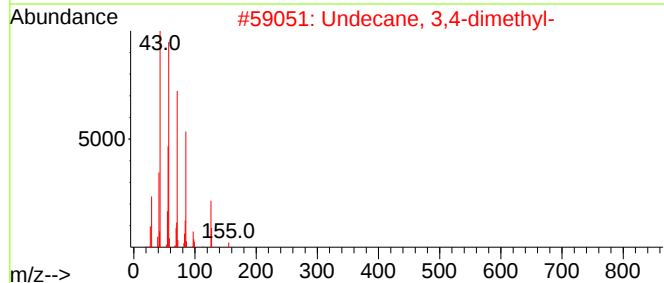
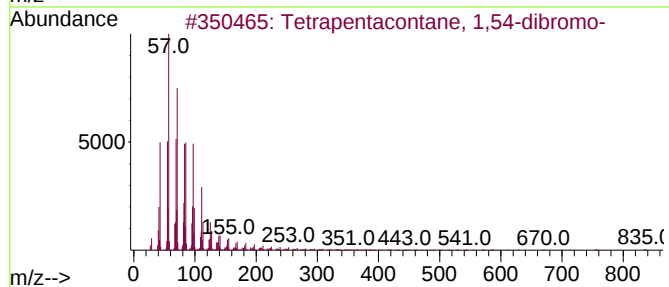
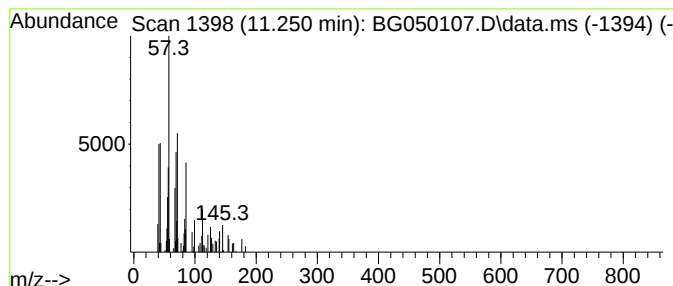
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Quant Title : SVOA CALIBRATION

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

Peak Number 35 Tetrapentacontane, 1,54-dib... Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.250	3.89 ng/ul	54776	Naphthalene-d8	10.774

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetrapentacontane, 1,54-dibromo-	915	C54H108Br2	1000156-09-4	52
2		Undecane, 3,4-dimethyl-	184	C13H28	017312-78-6	50
3		Decane, 3-methyl-	156	C11H24	013151-34-3	47
4		1-Hexanol, 5-methyl-2-(1-methyle...	158	C10H22O	002051-33-4	43
5		Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

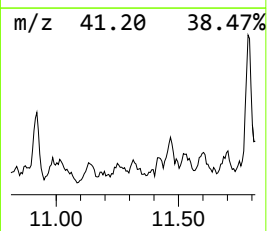
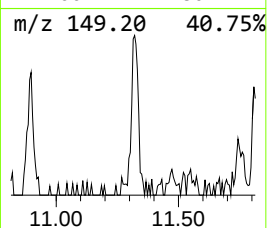
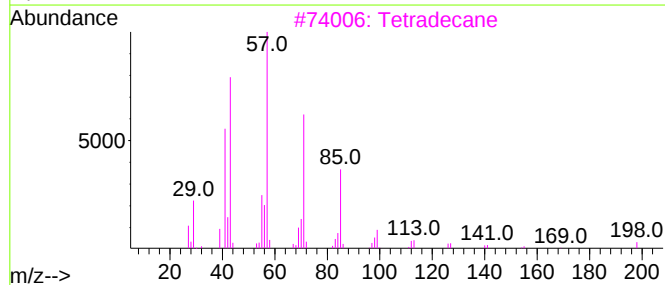
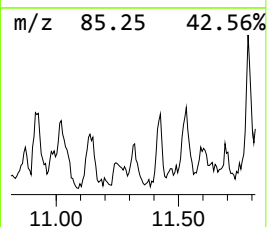
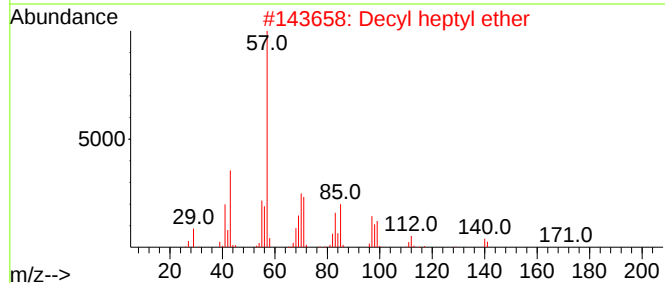
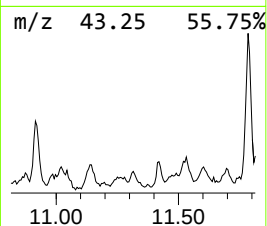
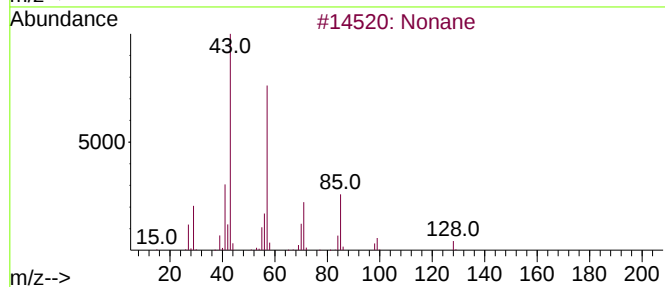
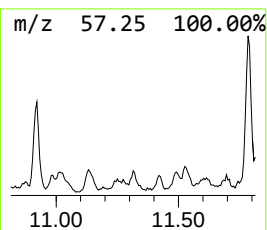
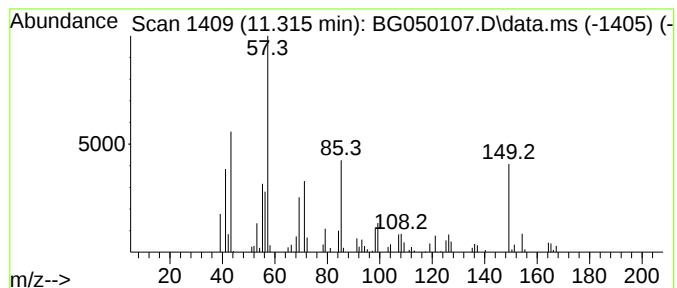
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Quant Title : SVOA CALIBRATION

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

Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.315	5.96 ng/ul	83975	Naphthalene-d8	10.774	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane	128	C9H20	000111-84-2	53
2	Decyl heptyl ether	256	C17H36O	1000406-39-1	47
3	Tetradecane	198	C14H30	000629-59-4	47
4	Nonane	128	C9H20	000111-84-2	47
5	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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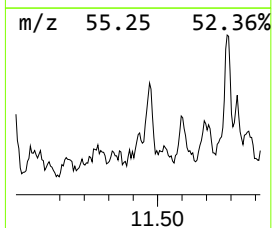
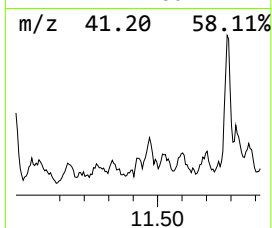
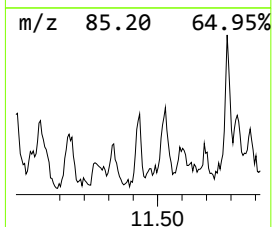
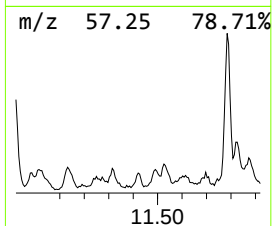
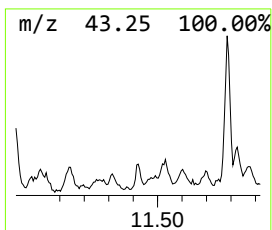
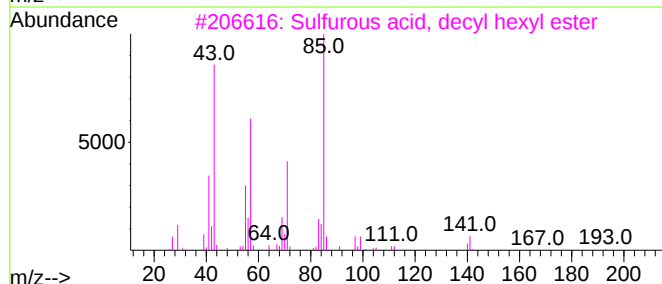
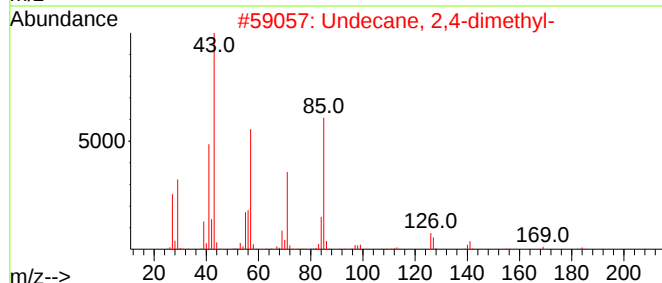
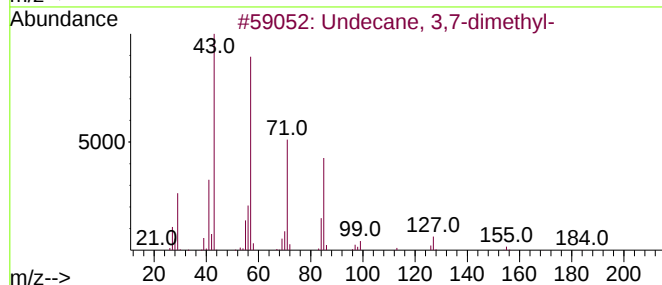
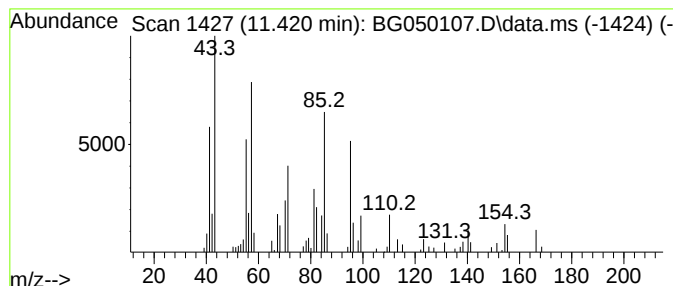
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Quant Title : SVOA CALIBRATION

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

Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.420	4.75 ng/ul	66967	Naphthalene-d8	10.774

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Undecane, 3,7-dimethyl-	184	C13H28	017301-29-0	35
2		Undecane, 2,4-dimethyl-	184	C13H28	017312-80-0	35
3		Sulfurous acid, decyl hexyl ester	306	C16H34O3S	1000309-13-2	35
4		Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	27
5		Cyclohexanol, 2,3-dimethyl-	128	C8H16O	001502-24-5	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
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Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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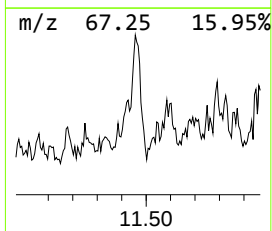
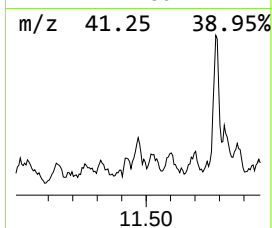
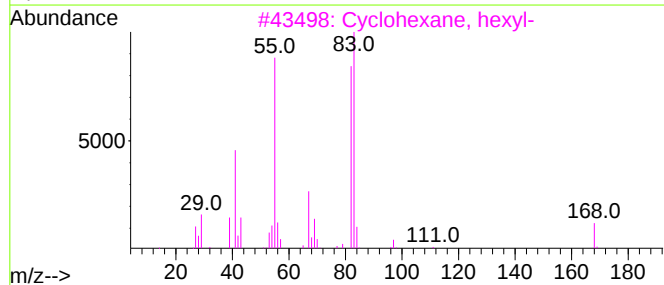
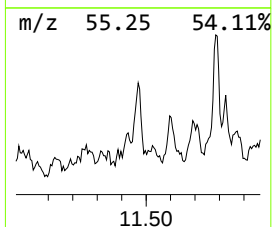
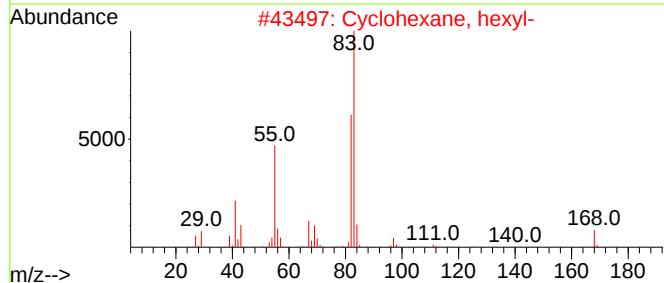
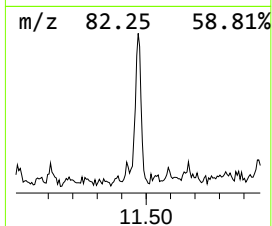
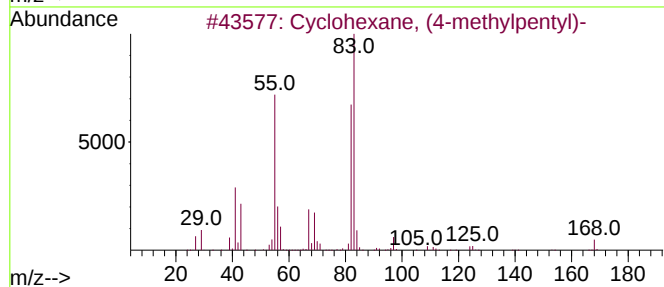
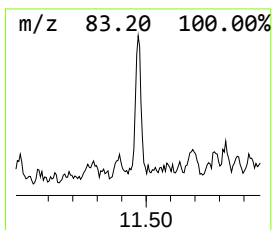
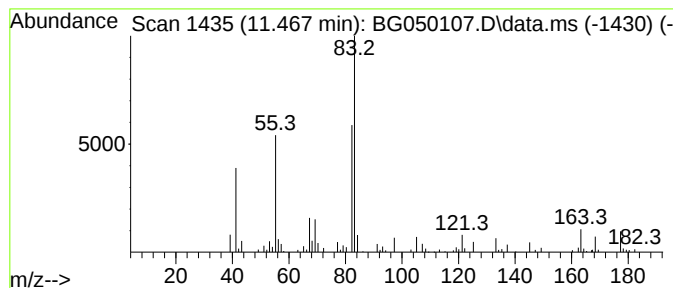
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Quant Title : SVOA CALIBRATION

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

Peak Number 39 (DEL) Alkane: Cyclic11.467 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.467	10.55 ng/ul	148735	Naphthalene-d8	10.774

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, (4-methylpentyl)-	168	C12H24	061142-20-9	64
2		Cyclohexane, hexyl-	168	C12H24	004292-75-5	62
3		Cyclohexane, hexyl-	168	C12H24	004292-75-5	58
4		Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	53
5		Cyclohexane, octyl-	196	C14H28	001795-15-9	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
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Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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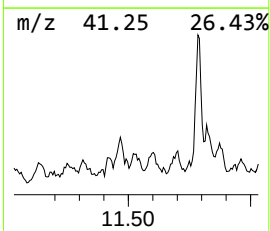
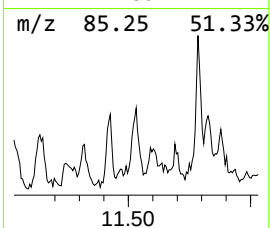
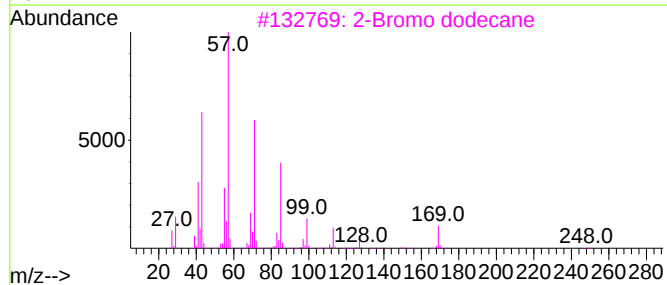
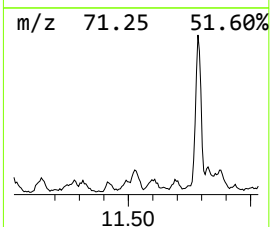
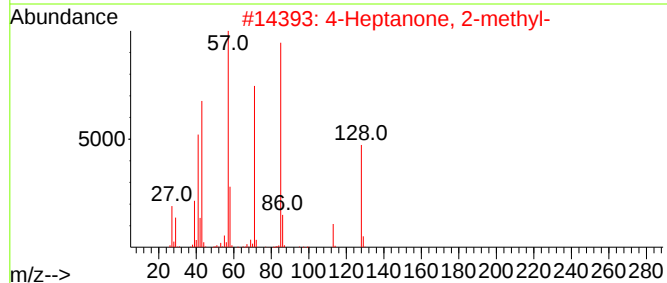
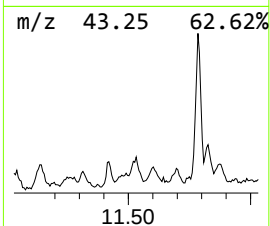
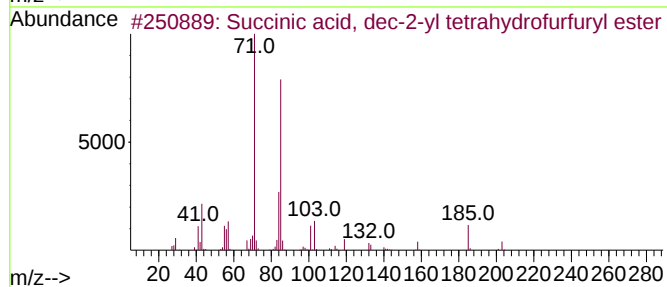
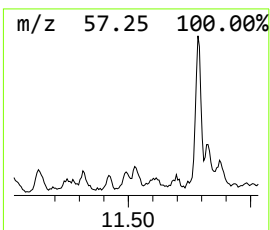
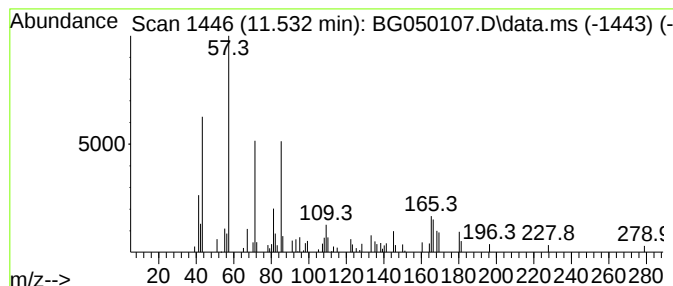
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Quant Title : SVOA CALIBRATION

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

Peak Number 40 unknown-04 Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.532	4.94 ng/ul	69578	Naphthalene-d8	10.774

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Succinic acid, dec-2-yl tetrahyd...	342	C19H34O5	1000390-72-3	38
2			4-Heptanone, 2-methyl-	128	C8H16O	000626-33-5	38
3			2-Bromo dodecane	248	C12H25Br	013187-99-0	38
4			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	38
5			4-Heptanone, 2-methyl-	128	C8H16O	000626-33-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
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Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

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ClientSampleId :
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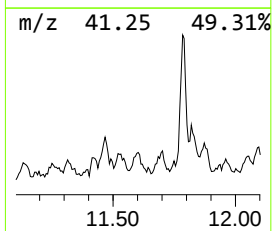
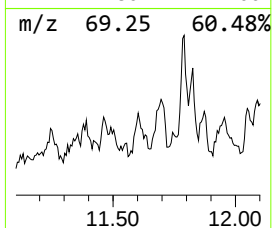
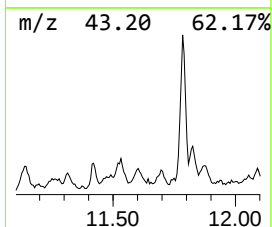
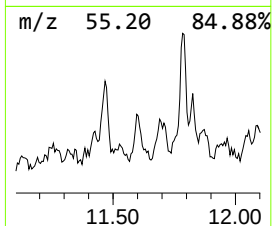
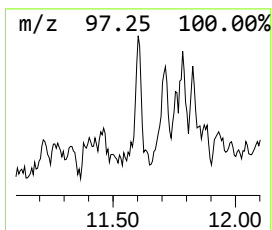
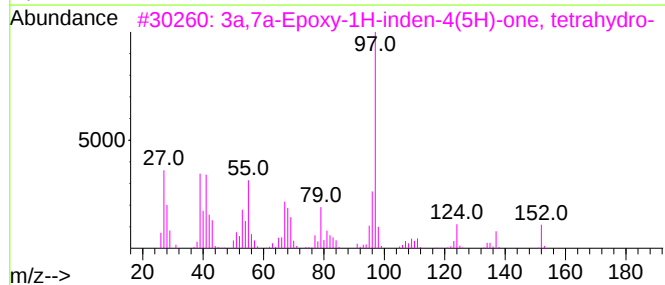
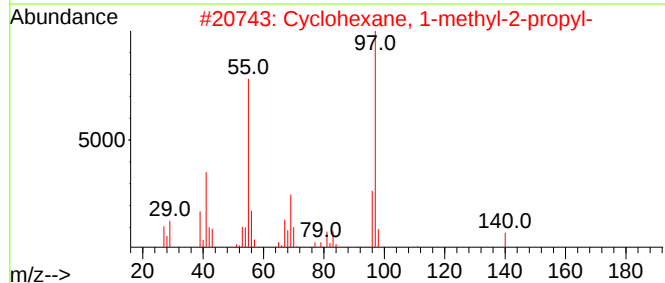
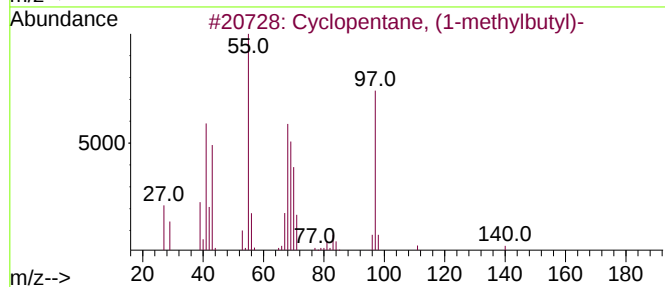
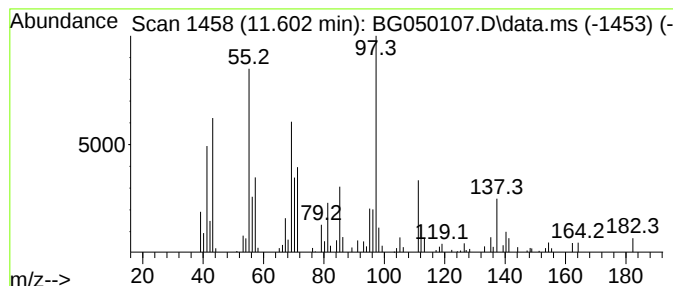
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Quant Title : SVOA CALIBRATION

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

Peak Number 41 (DEL) Alkane: Cyclic11.602 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.602	10.70 ng/ul	150782	Naphthalene-d8	10.774

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, (1-methylbutyl)-	140	C10H20	004737-43-3	45
2			Cyclohexane, 1-methyl-2-propyl-	140	C10H20	004291-79-6	38
3			3a,7a-Epoxy-1H-inden-4(5H)-one, ...	152	C9H12O2	039746-31-1	35
4			2-(3-Methyl-2,5-dihydro-1H-pyraz...	127	C6H13N3	1000387-20-9	35
5			cis-2-Oxabicyclo[4.4.0]decane	140	C9H16O	060416-19-5	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
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Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

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ClientSampleId :
EW7C4

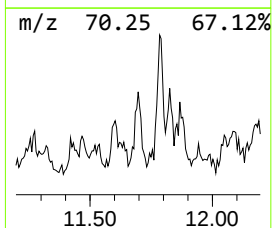
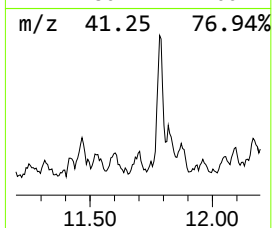
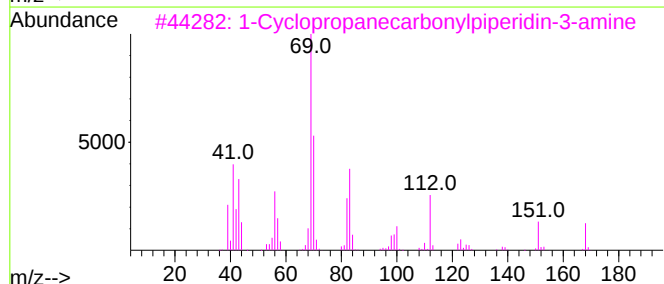
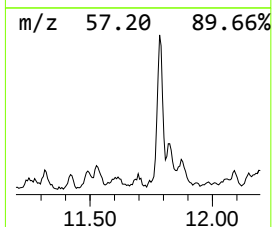
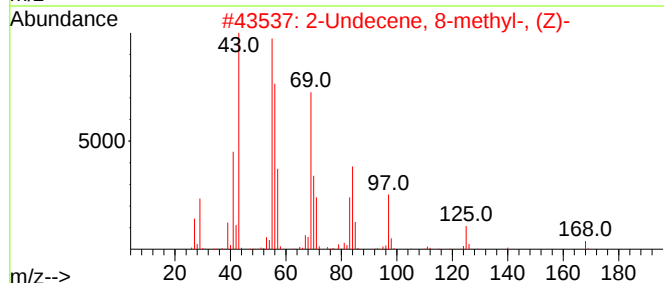
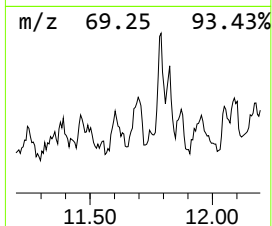
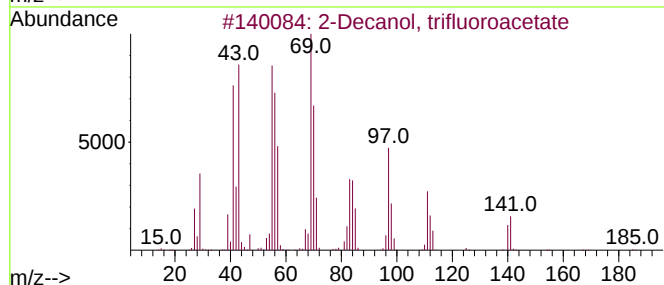
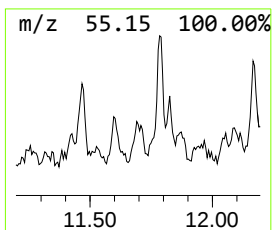
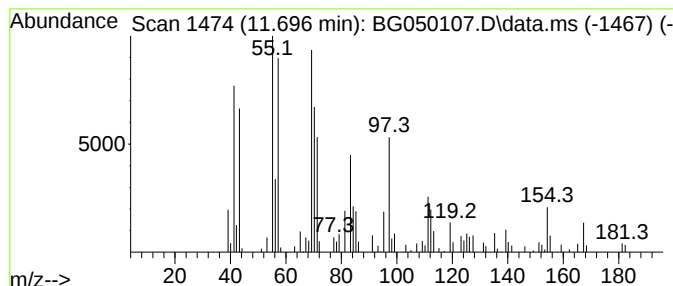
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Quant Title : SVOA CALIBRATION

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

Peak Number 42 2-Decanol, trifluoroacetate Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.696	10.87 ng/ul	153209	Naphthalene-d8	10.774

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Decanol, trifluoroacetate	254	C12H21F3O2	1010352-32-8	64
2			2-Undecene, 8-methyl-, (Z)-	168	C12H24	074630-44-7	52
3			1-Cyclopropanecarbonylpiperidin-...	168	C9H16N2O	1114596-39-2	49
4			Pentafluoropropionic acid, tride...	346	C16H27F5O2	959261-22-4	38
5			4-Octene, 2,3,7-trimethyl-, [S-(...	154	C11H22	052763-13-0	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW7C4

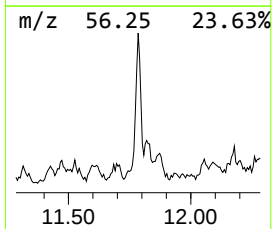
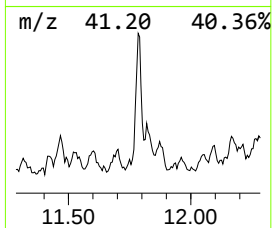
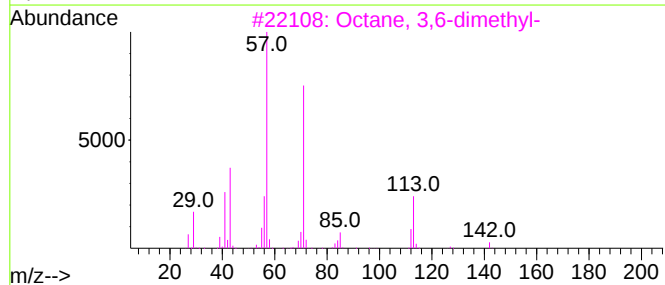
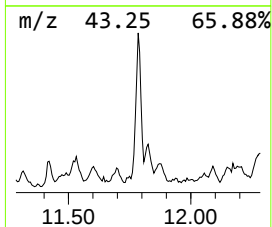
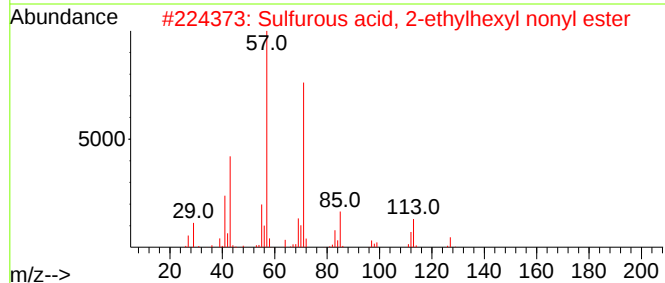
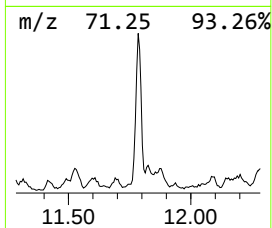
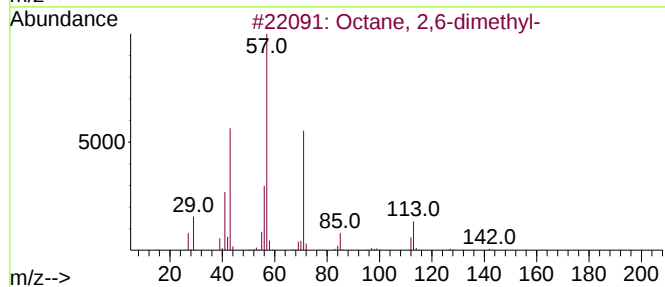
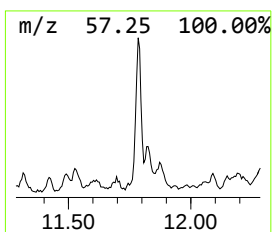
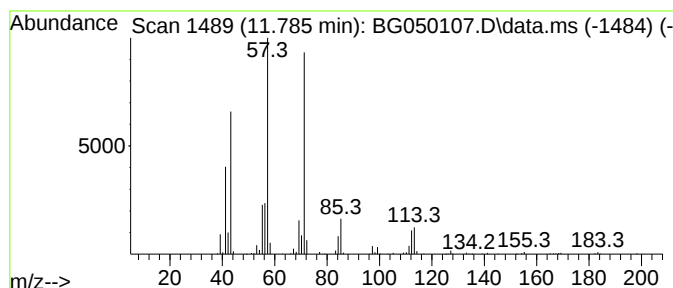
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Quant Title : SVOA CALIBRATION

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

Peak Number 43 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.785	35.02 ng/ul	493554	Naphthalene-d8	10.774

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	83
2			Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	80
3			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	72
4			Octane, 2,3,7-trimethyl-	156	C11H24	062016-34-6	72
5			Sulfurous acid, 2-ethylhexyl pen...	264	C13H28O3S	1000309-18-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

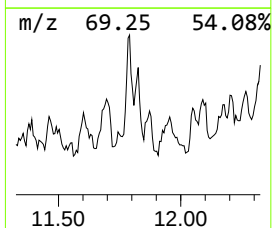
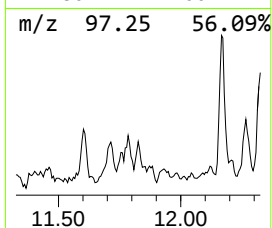
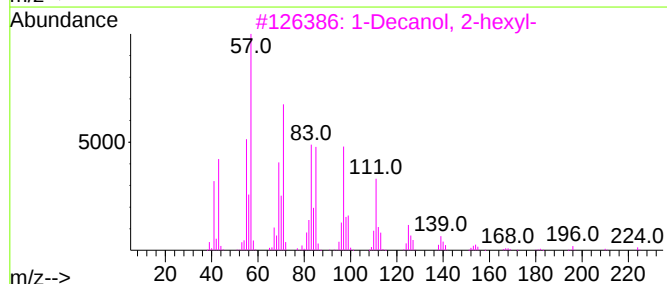
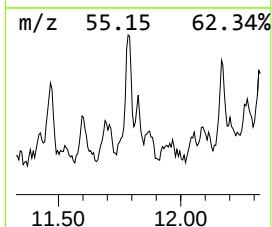
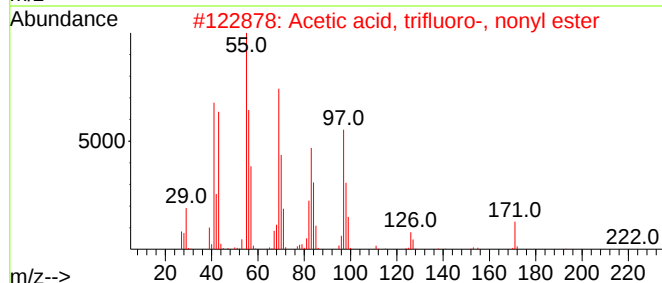
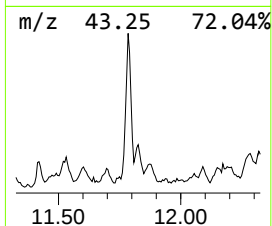
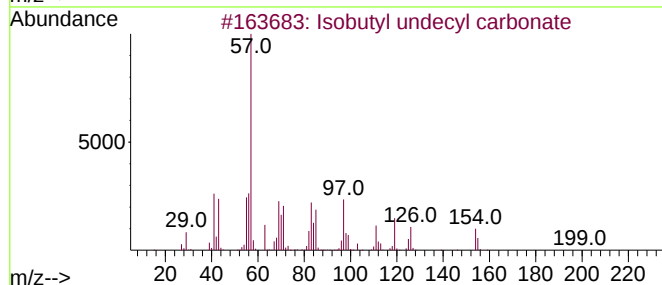
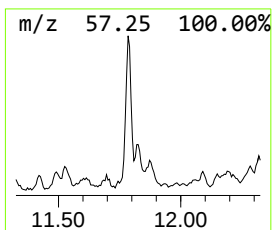
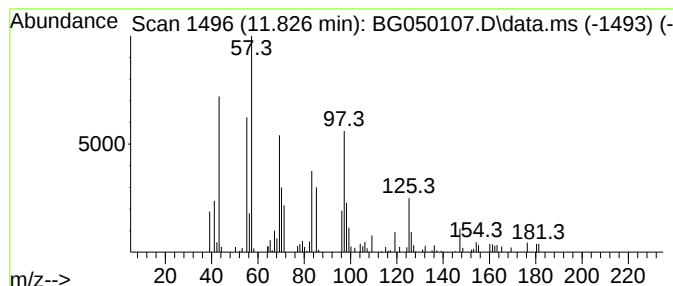
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 44 unknown-05 Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.			
11.826	8.47 ng/ul	119378	Naphthalene-d8	10.774			
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isobutyl undecyl carbonate	272	C16H32O3	1000372-78-8	43
2			Acetic acid, trifluoro-, nonyl e...	240	C11H19F3O2	030767-14-7	38
3			1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	35
4			1,7-Dimethyl-4-(1-methylethyl)cy...	210	C15H30	000645-10-3	27
5			Sulfurous acid, cyclohexylmethyl...	360	C20H40O3S	1000309-22-1	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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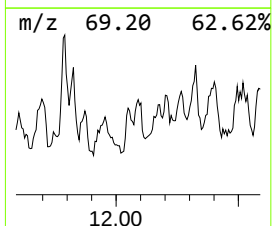
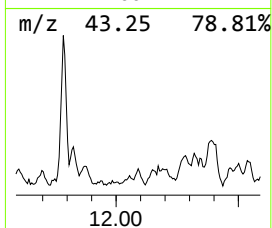
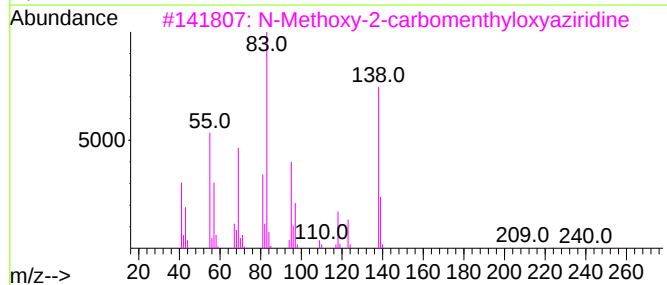
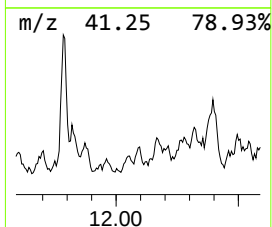
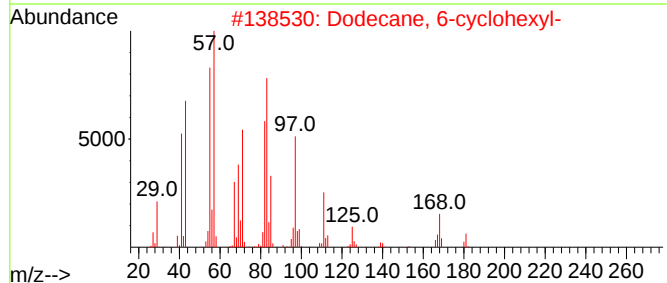
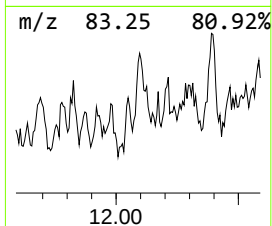
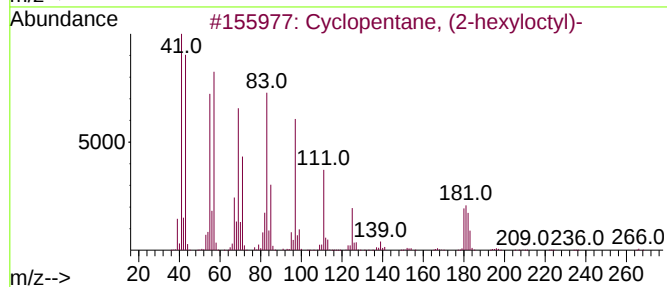
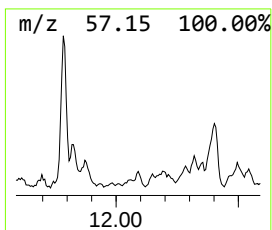
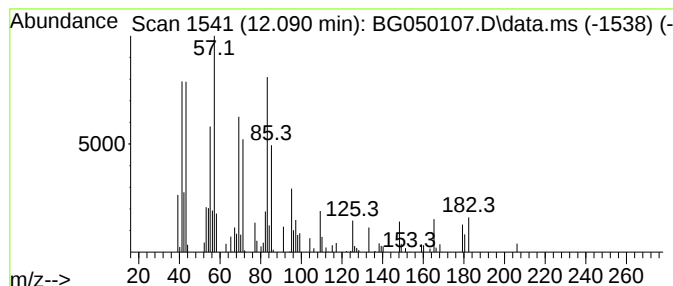
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Quant Title : SVOA CALIBRATION

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

Peak Number 45 (DEL) Alkane: Cyclic12.090 Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.090	5.09 ng/ul	71706	Naphthalene-d8	10.774

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, (2-hexyloctyl)-	266	C19H38	055044-77-4	47
2			Dodecane, 6-cyclohexyl-	252	C18H36	013151-86-5	43
3			N-Methoxy-2-carbomethyloxyaziri...	255	C14H25NO3	060999-67-9	27
4			Oxalic acid, cyclobutyl tetradec...	340	C20H36O4	1000309-70-9	27
5			Sulfurous acid, butyl heptadecyl...	376	C21H44O3S	1000309-18-4	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW7C4

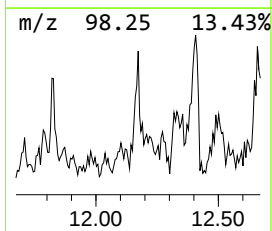
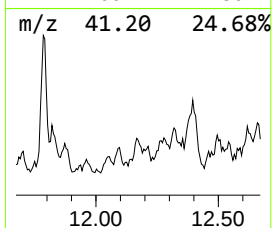
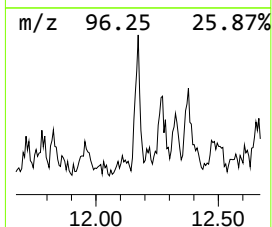
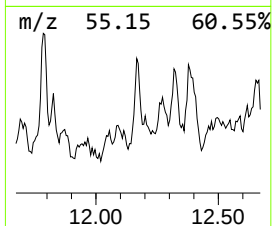
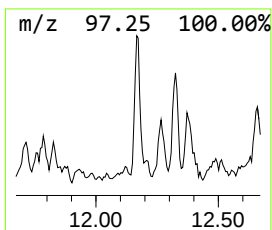
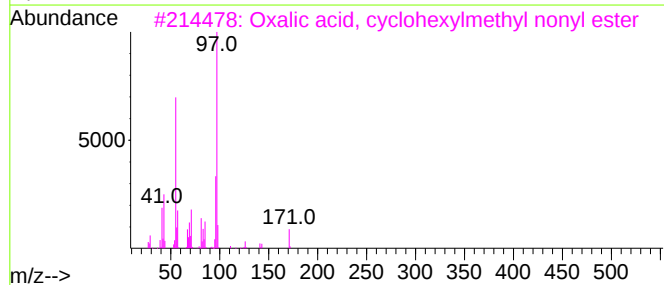
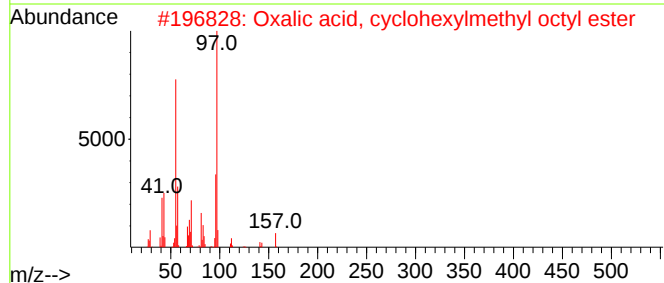
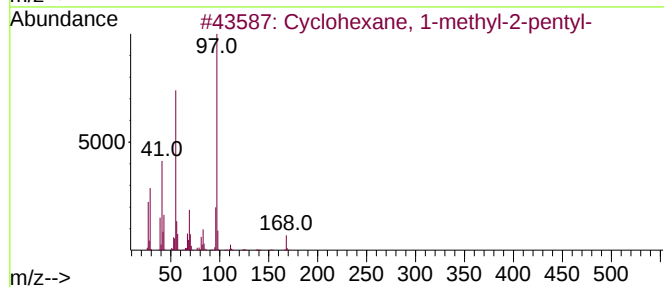
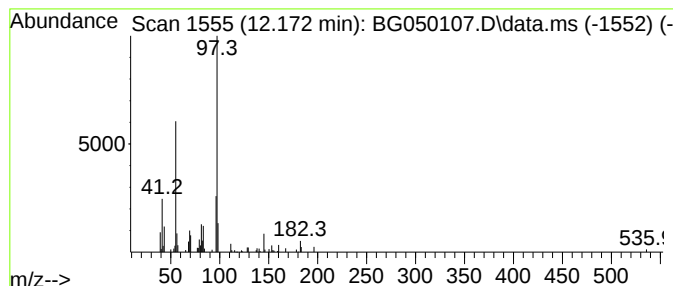
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Quant Title : SVOA CALIBRATION

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

Peak Number 46 (DEL) Alkane: Cyclic12.172 Concentration Rank 58

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.172	2.65 ng/ul	37355	Naphthalene-d8	10.774

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-methyl-2-pentyl-	168	C12H24	054411-01-7	78
2			Oxalic acid, cyclohexylmethyl oc...	298	C17H30O4	1000309-68-4	72
3			Oxalic acid, cyclohexylmethyl no...	312	C18H32O4	1000309-68-6	72
4			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	64
5			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW7C4

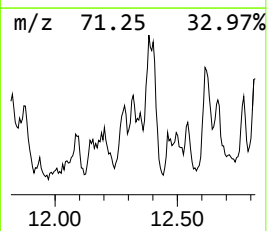
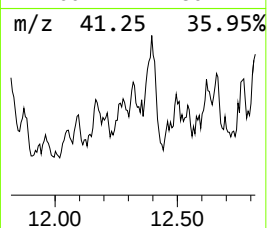
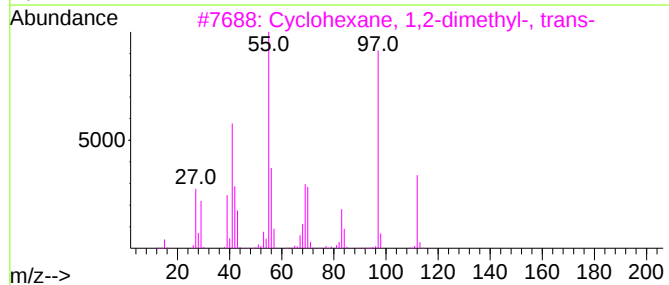
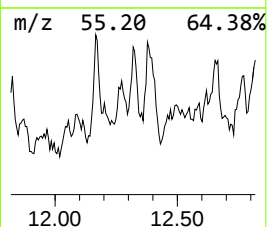
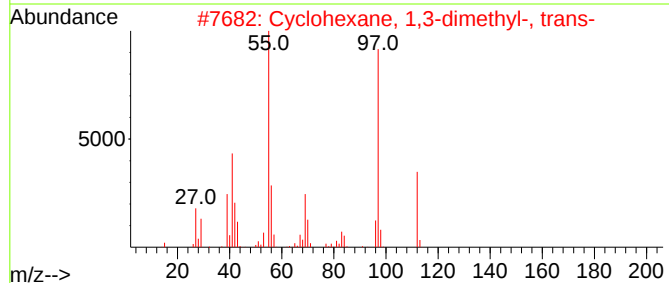
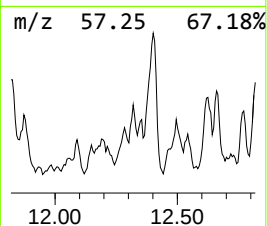
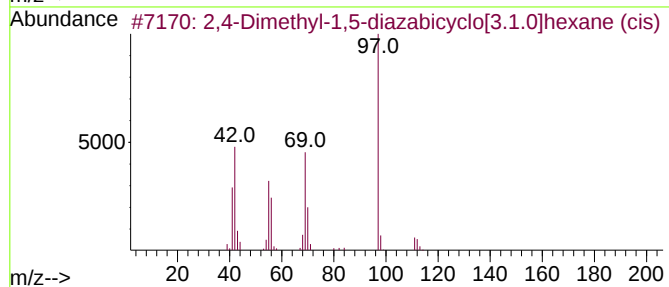
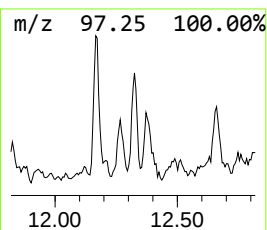
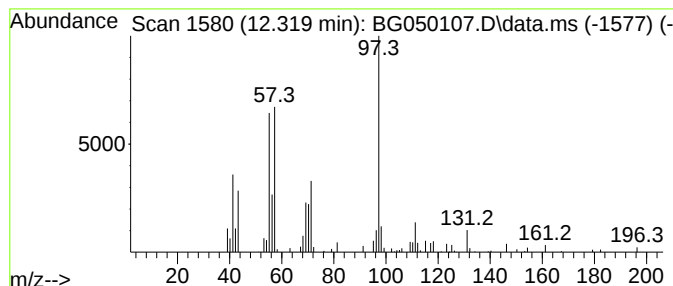
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Quant Title : SVOA CALIBRATION

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

Peak Number 47 2,4-Dimethyl-1,5-diazabicyc... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.319	7.29 ng/ul	102780	Naphthalene-d8	10.774

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,4-Dimethyl-1,5-diazabicyclo[3...]	112	C6H12N2	100463-01-2	53
2			Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	53
3			Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	53
4			Sulfurous acid, cyclohexylmethyl...	416	C24H48O3S	1000309-22-5	50
5			Cyclohexane, 1,1-dimethyl-	112	C8H16	000590-66-9	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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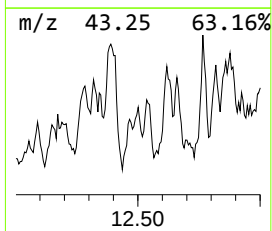
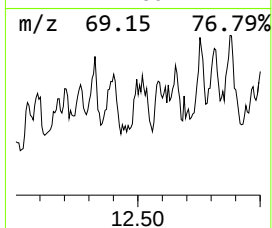
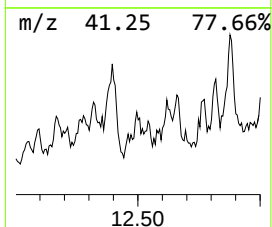
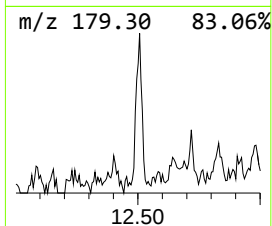
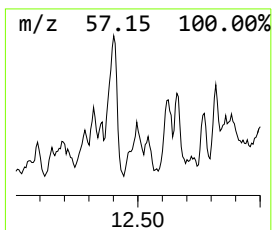
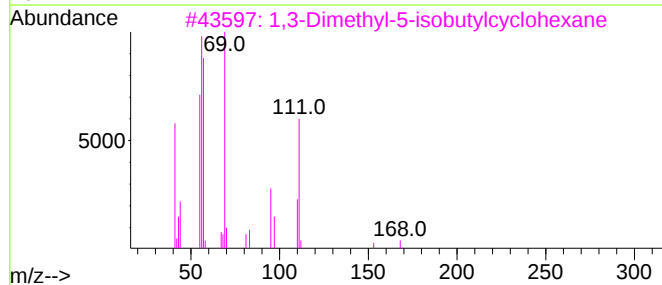
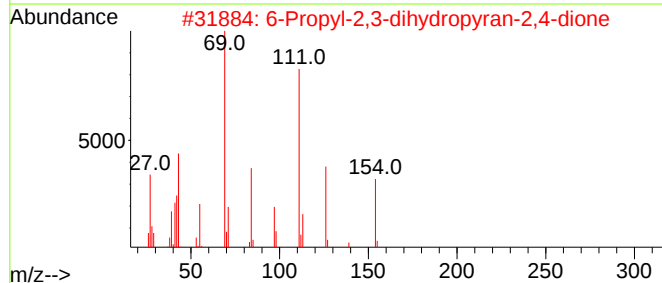
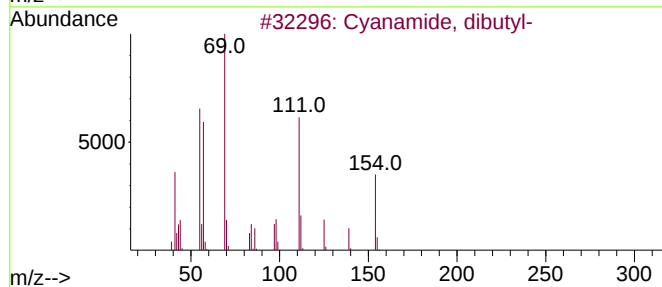
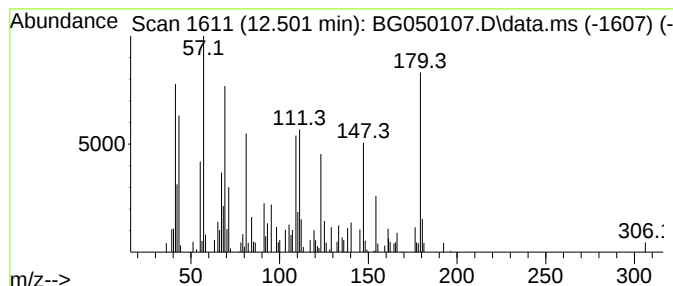
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Quant Title : SVOA CALIBRATION

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

Peak Number 48 unknown-06 Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.501	6.50 ng/ul	91607	Naphthalene-d8	10.774

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyanamide, dibutyl-	154	C9H18N2	002050-54-6	38
2			6-Propyl-2,3-dihydropyran-2,4-dione	154	C8H10O3	005698-53-3	22
3			1,3-Dimethyl-5-isobutylcyclohexane	168	C12H24	013131-76-5	22
4			Cyanamide, dibutyl-	154	C9H18N2	002050-54-6	20
5			Trans-1,4-diethylcyclohexane	140	C10H20	013990-93-7	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

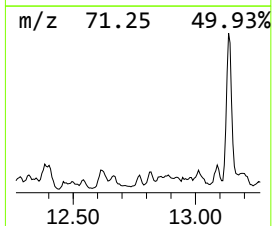
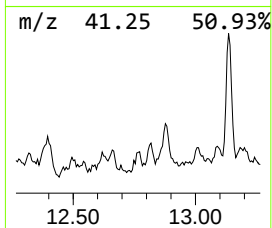
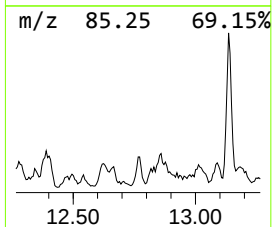
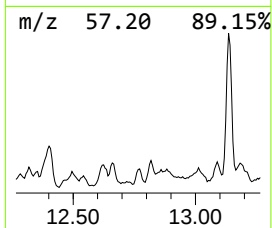
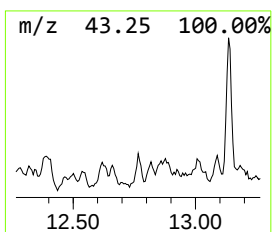
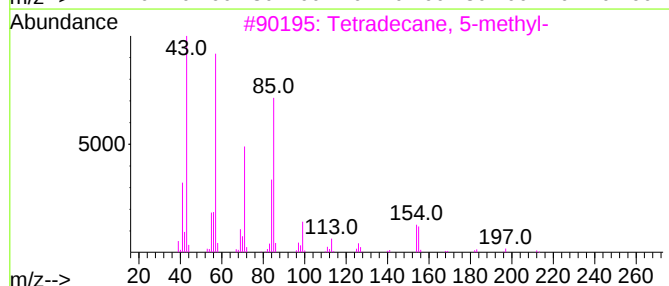
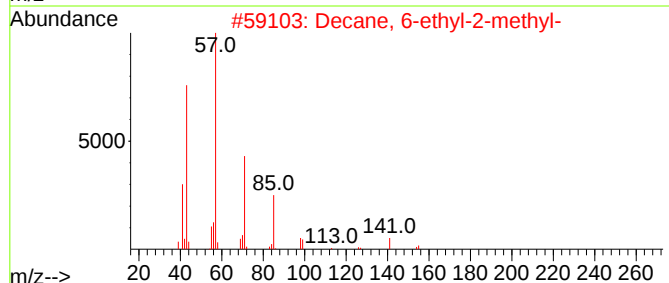
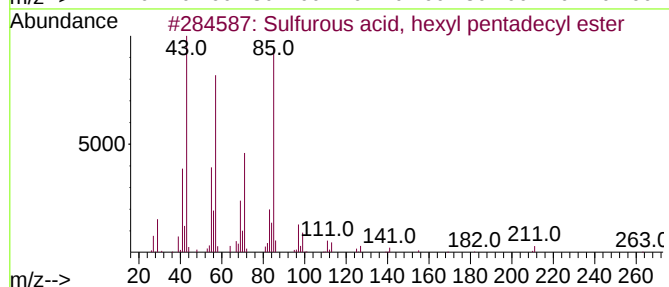
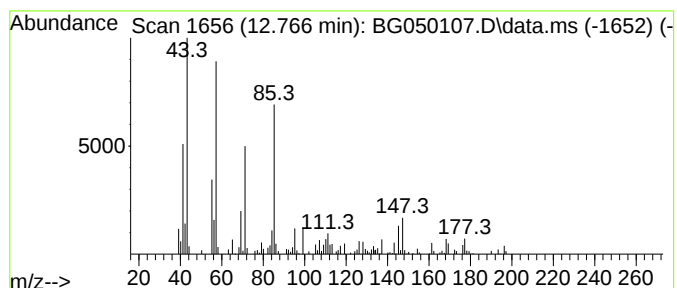
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

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

Peak Number 49 Sulfurous acid, hexyl penta... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.765	8.01 ng/ul	132328	Acenaphthene-d10	14.622	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Sulfurous acid, hexyl pentadecyl...	376	C21H44O3S	1000309-13-7	64
2	Decane, 6-ethyl-2-methyl-	184	C13H28	062108-21-8	53
3	Tetradecane, 5-methyl-	212	C15H32	025117-32-2	50
4	Decane, 2,3,5,8-tetramethyl-	198	C14H30	192823-15-7	47
5	4-Heptanone, 2-methyl-	128	C8H16O	000626-33-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

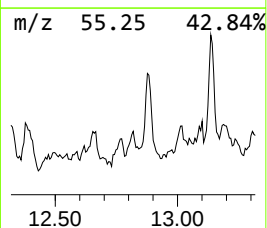
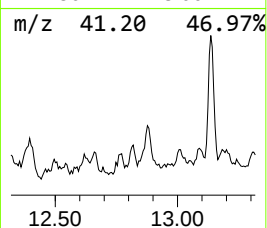
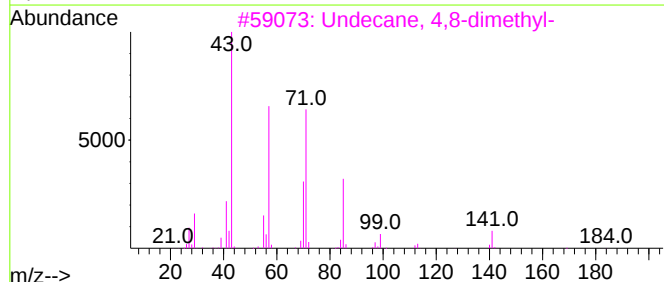
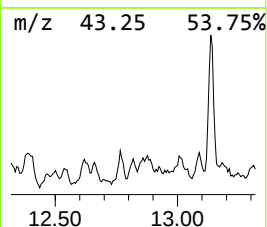
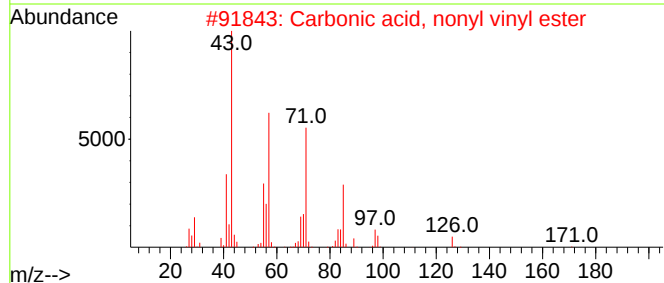
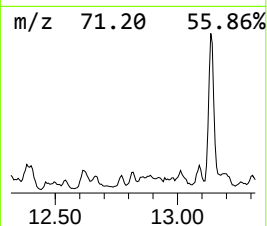
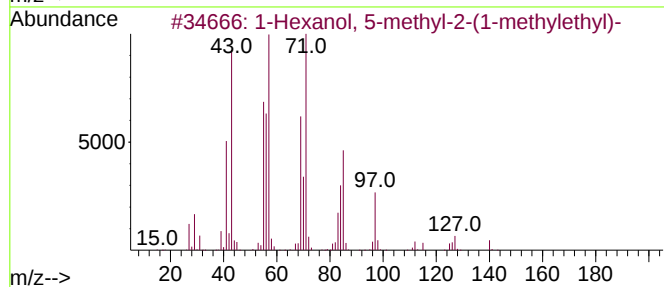
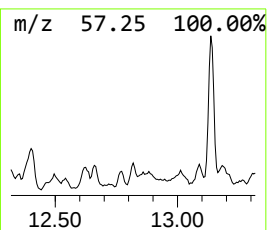
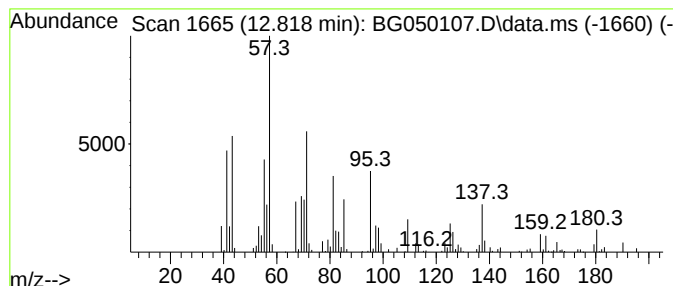
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ClientSampleId :
EW7C4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

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

Peak Number 50 unknown-07 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.818	10.29 ng/ul	170092	Acenaphthene-d10	14.622
Hit#	of 5	Tentative ID	MW MolForm	CAS# Qual
1		1-Hexanol, 5-methyl-2-(1-methyle...	158 C10H22O	002051-33-4 35
2		Carbonic acid, nonyl vinyl ester	214 C12H22O3	1000383-25-6 27
3		Undecane, 4,8-dimethyl-	184 C13H28	017301-33-6 27
4		Nonane, 3,7-dimethyl-	156 C11H24	017302-32-8 27
5		2-Bromopropionic acid, 2-pentyl ...	222 C8H15BrO2	130232-40-5 27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

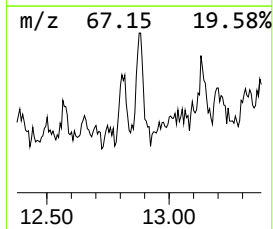
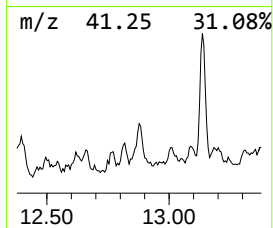
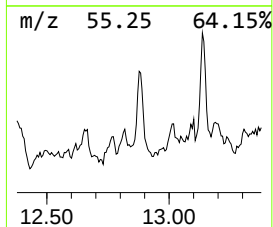
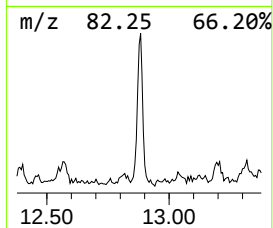
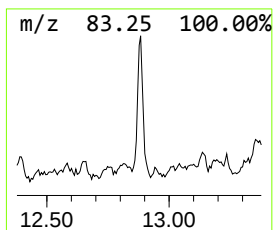
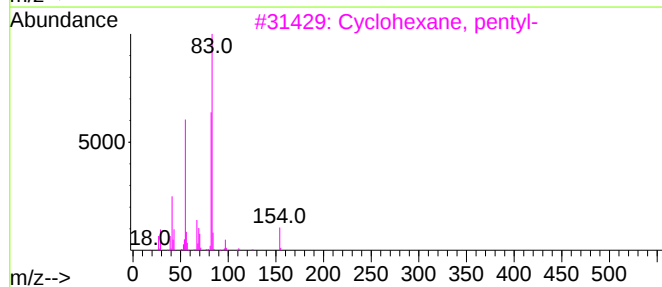
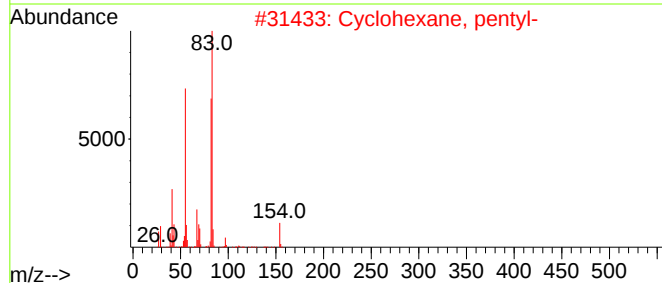
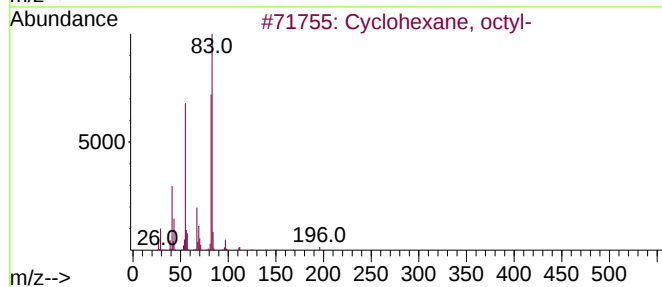
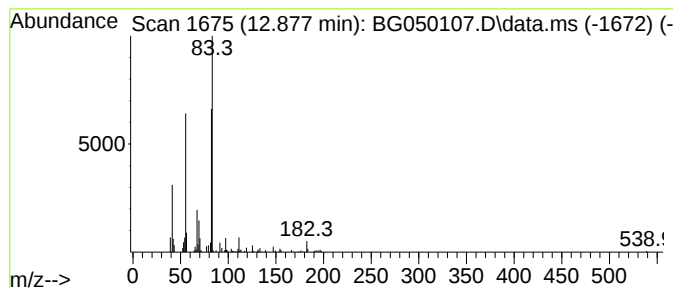
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

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

Peak Number 51 (DEL) Alkane: Cyclic12.877 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.877	17.84 ng/ul	294890	Acenaphthene-d10		14.622	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Cyclohexane, octyl-		196	C14H28	001795-15-9	90
2	Cyclohexane, pentyl-		154	C11H22	004292-92-6	87
3	Cyclohexane, pentyl-		154	C11H22	004292-92-6	87
4	Cyclohexane, pentyl-		154	C11H22	004292-92-6	86
5	Heptylcyclohexane		182	C13H26	005617-41-4	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

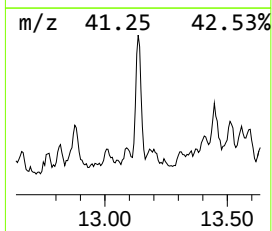
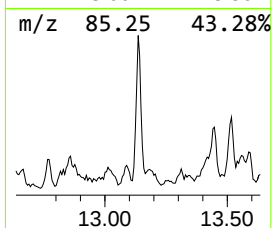
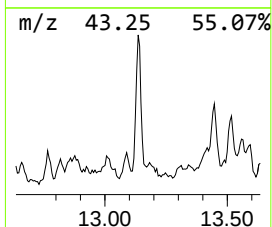
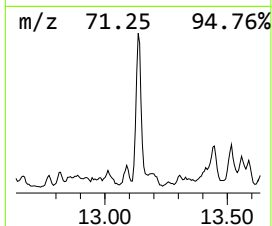
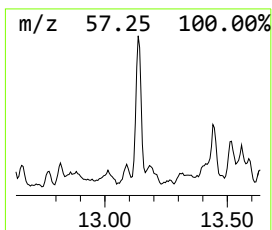
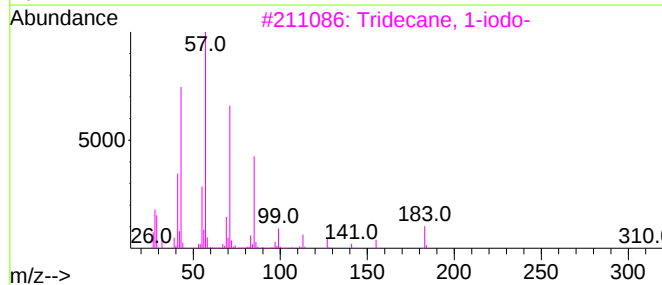
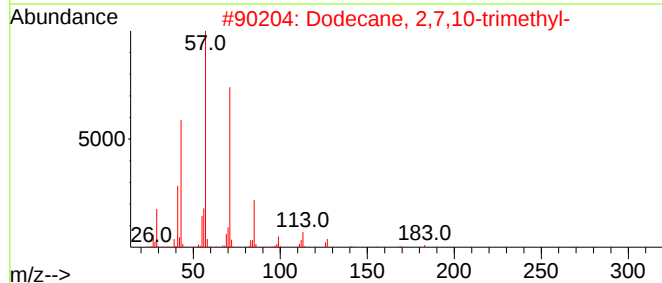
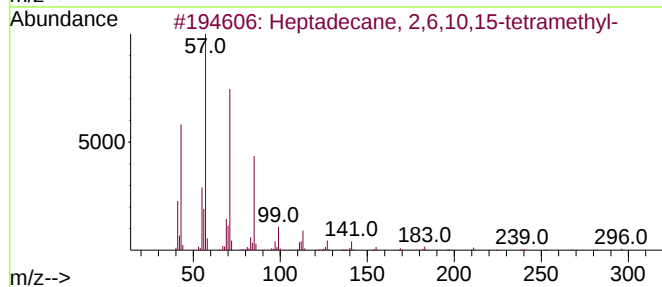
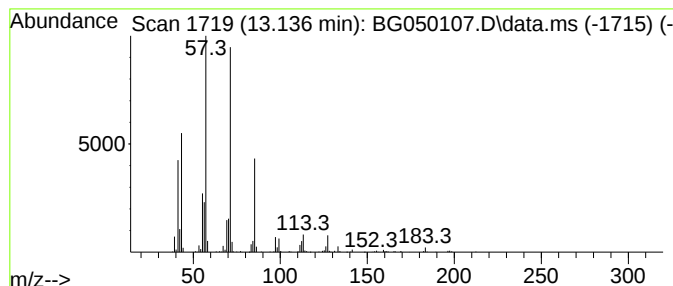
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

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

Peak Number 52 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.136	45.29 ng/ul	748474	Acenaphthene-d10	14.622	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	90
2	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	80
3	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	72
4	Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	72
5	Sulfurous acid, decyl pentyl ester	292	C15H32O3S	1000309-14-3	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW7C4

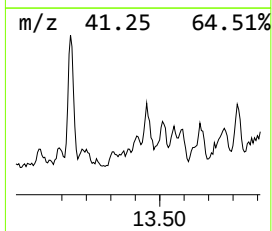
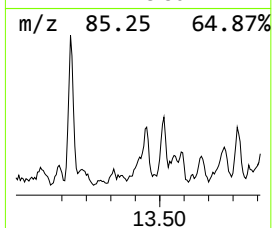
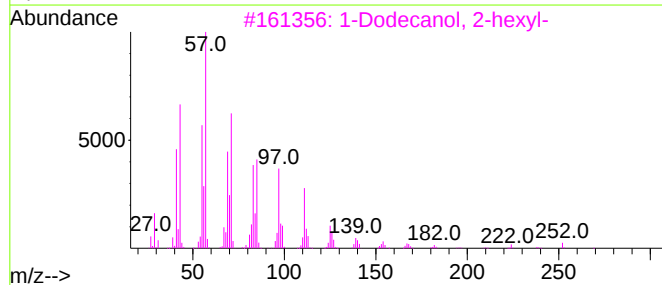
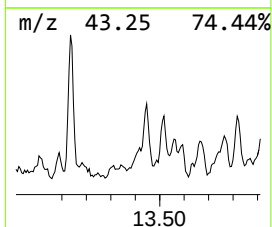
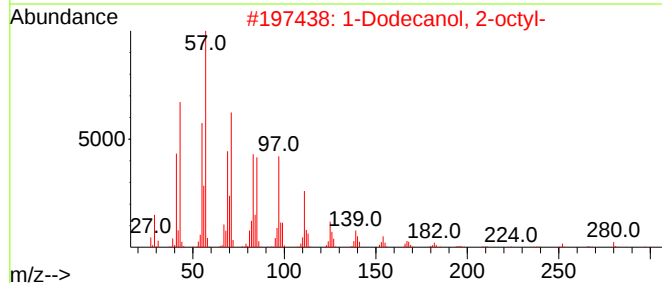
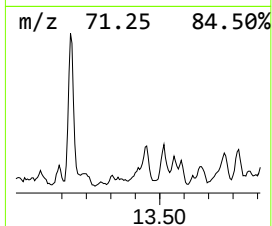
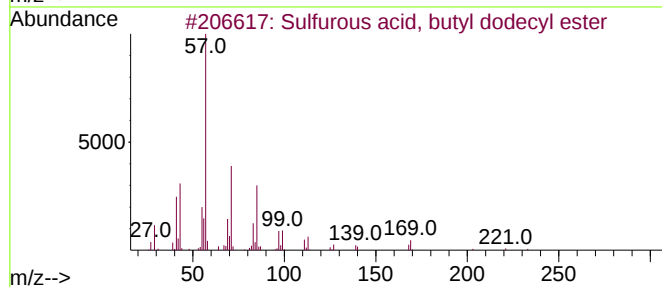
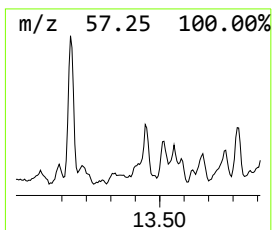
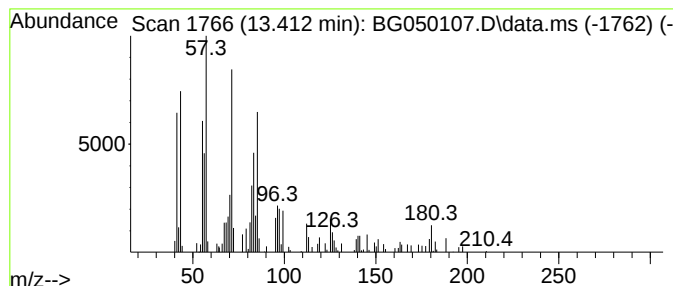
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Quant Title : SVOA CALIBRATION

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

Peak Number 53 Sulfurous acid, butyl dodec... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.412	7.32 ng/ul	120971	Acenaphthene-d10	14.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	59
2			1-Dodecanol, 2-octyl-	298	C20H42O	005333-42-6	59
3			1-Dodecanol, 2-hexyl-	270	C18H38O	110225-00-8	58
4			Cyclopentane, 1-methyl-2-propyl-	126	C9H18	003728-57-2	38
5			Nonane, 3-methyl-	142	C10H22	005911-04-6	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

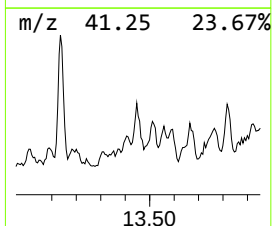
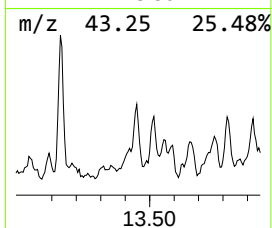
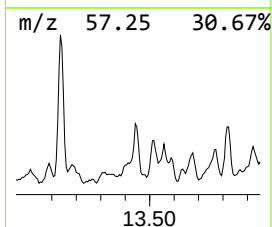
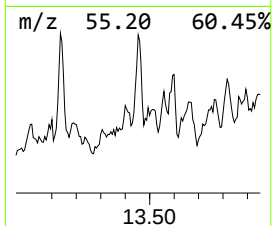
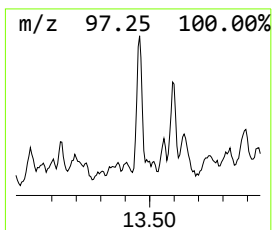
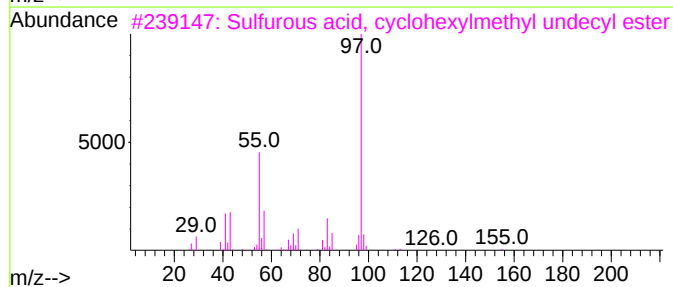
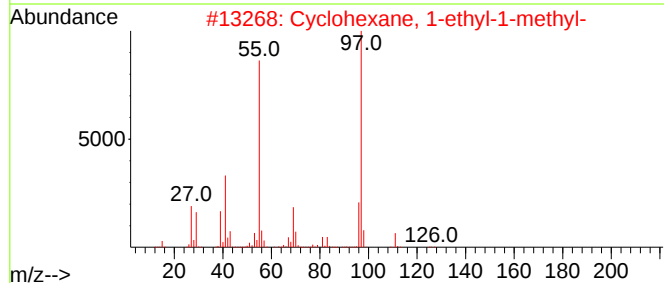
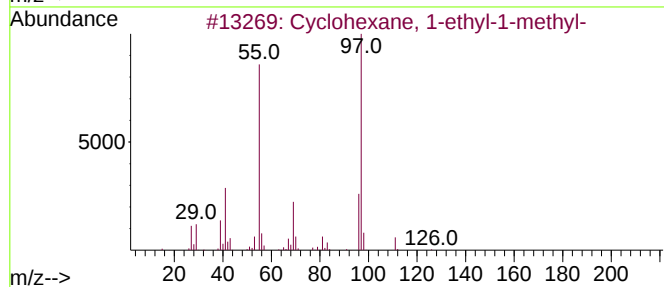
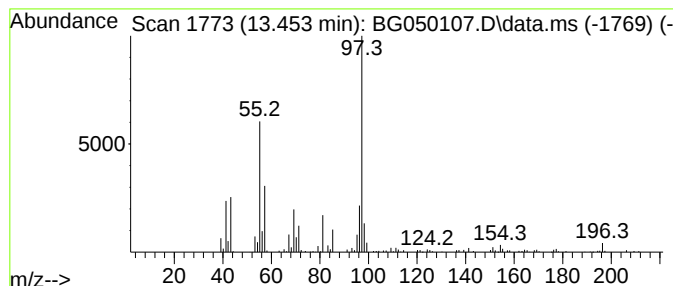
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

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

Peak Number 54 (DEL) Alkane: Cyclic13.453 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.453	18.24 ng/ul	301505	Acenaphthene-d10	14.622	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1-ethyl-1-methyl-	126	C9H18	004926-90-3	70
2	Cyclohexane, 1-ethyl-1-methyl-	126	C9H18	004926-90-3	68
3	Sulfurous acid, cyclohexylmethyl...	332	C18H36O3S	1000309-21-9	64
4	Sulfurous acid, cyclohexylmethyl...	304	C16H32O3S	1000309-21-8	64
5	Sulfurous acid, cyclohexylmethyl...	374	C21H42O3S	1000309-22-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

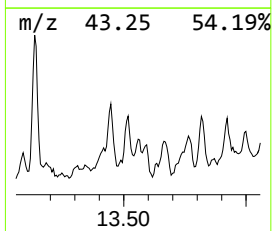
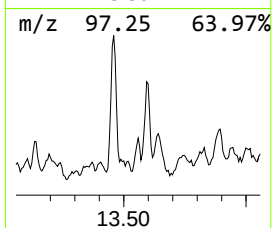
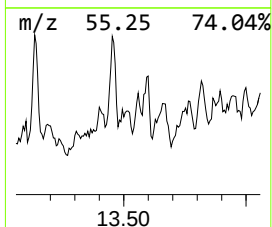
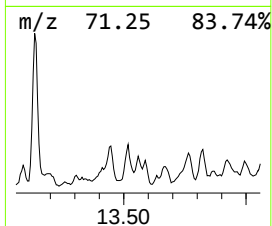
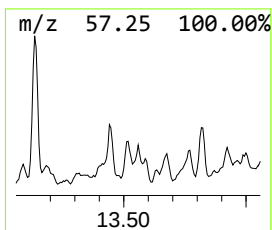
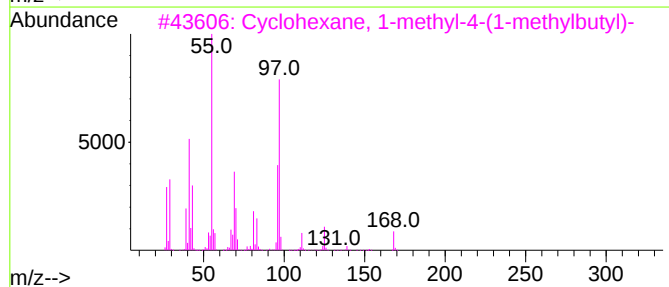
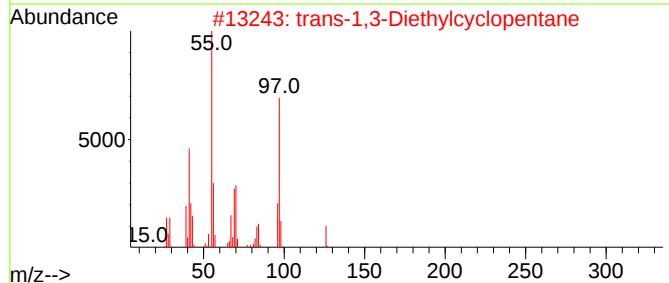
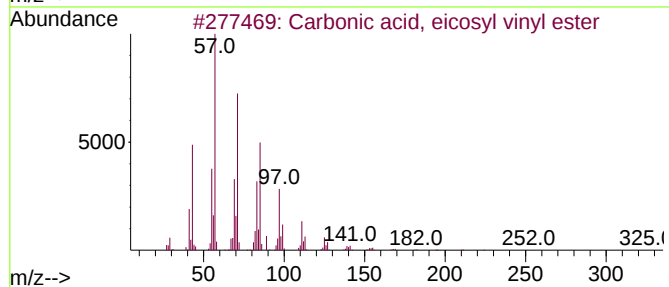
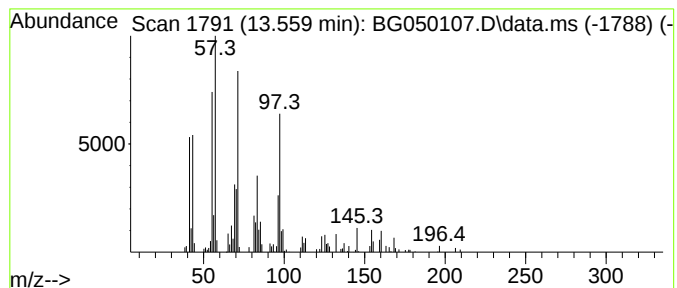
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BNA_G
ClientSampleId :
EW7C4

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Quant Title : SVOA CALIBRATION

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

Peak Number 55 Carbonic acid, eicosyl vinyl... Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.559	5.52 ng/ul	91263	Acenaphthene-d10			14.622
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Carbonic acid, eicosyl vinyl ester		368	C23H44O3	1000382-54-3	52
2	trans-1,3-Diethylcyclopentane		126	C9H18	1000113-87-1	38
3	Cyclohexane, 1-methyl-4-(1-methy...		168	C12H24	054411-00-6	35
4	Cyclopentane, 2-isopropyl-1,3-di...		140	C10H20	032281-85-9	30
5	Oxalic acid, cyclohexylmethyl is...		270	C15H26O4	1000309-68-3	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW7C4

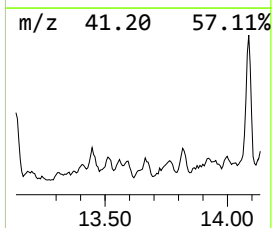
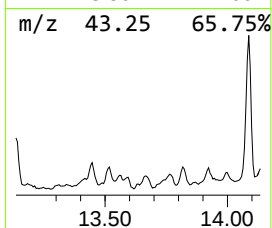
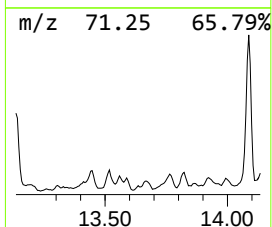
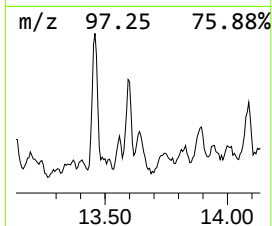
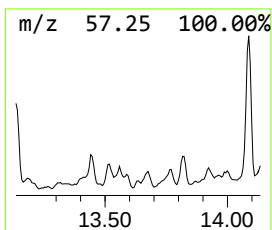
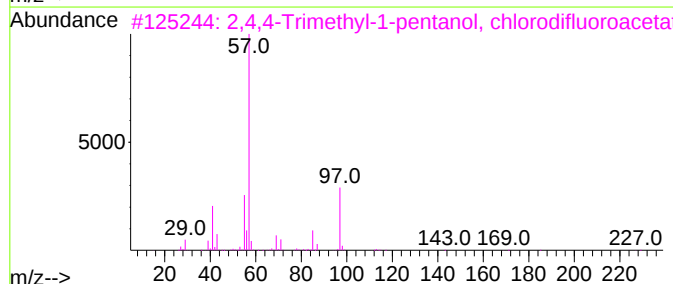
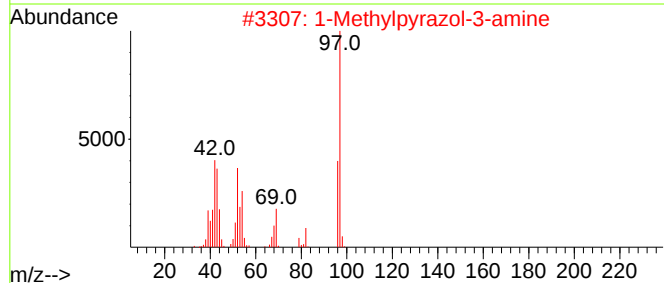
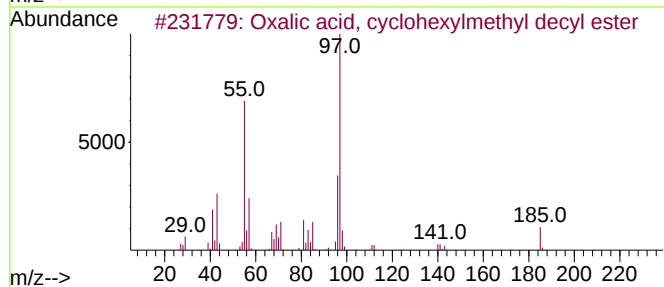
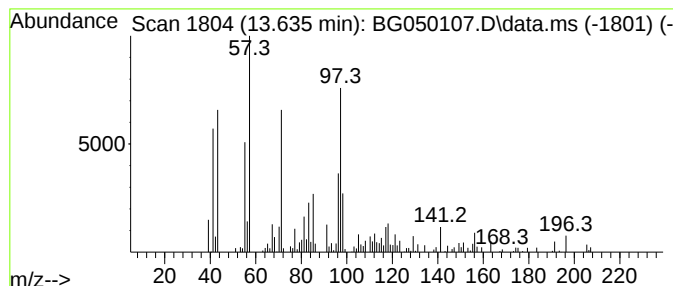
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Quant Title : SVOA CALIBRATION

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

Peak Number 56 unknown-08 Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.635	4.55 ng/ul	75240	Acenaphthene-d10	14.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, cyclohexylmethyl de...	326	C19H34O4	1000309-68-7	43
2			1-Methylpyrazol-3-amine	97	C4H7N3	001904-31-0	38
3			2,4,4-Trimethyl-1-pentanol, chlo...	242	C10H17ClF2O2	1000447-27-2	38
4			Oxalic acid, cyclohexylmethyl un...	340	C20H36O4	1000309-68-8	38
5			Oxalic acid, cyclohexylmethyl oc...	298	C17H30O4	1000309-68-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW7C4

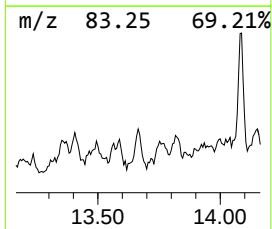
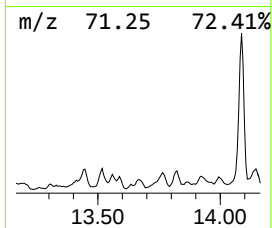
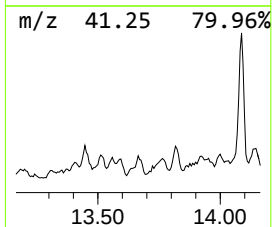
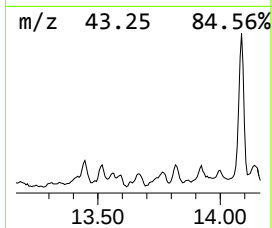
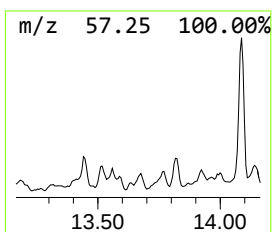
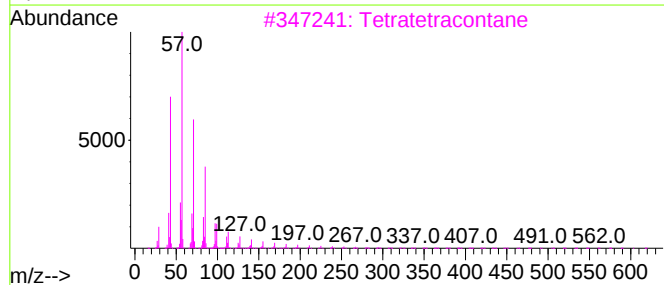
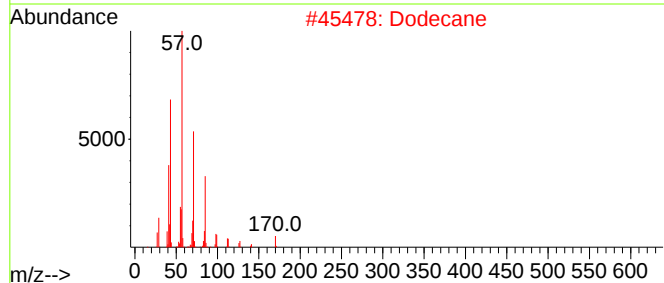
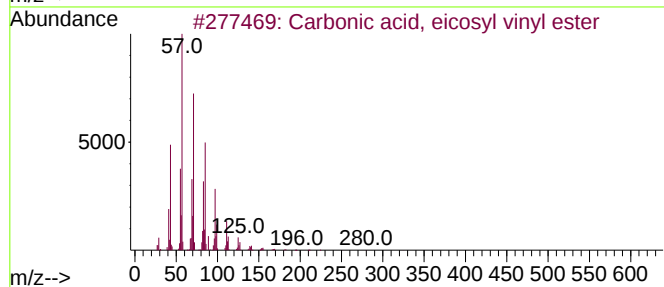
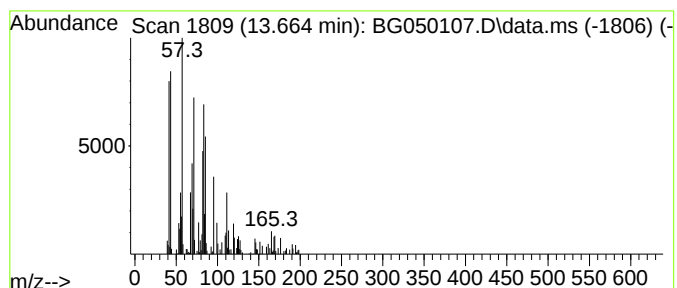
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Quant Title : SVOA CALIBRATION

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

Peak Number 57 unknown-09 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.664	16.00 ng/ul	264364	Acenaphthene-d10	14.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	38
2			Dodecane	170	C12H26	000112-40-3	38
3			Tetratetracontane	619	C44H90	007098-22-8	35
4			Tetradecanal	212	C14H28O	000124-25-4	35
5			Cyclopentane, 1-methyl-3-(1-meth...	126	C9H18	053771-88-3	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW7C4

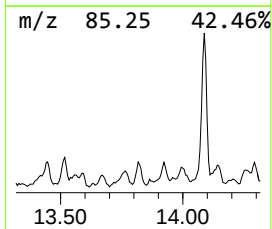
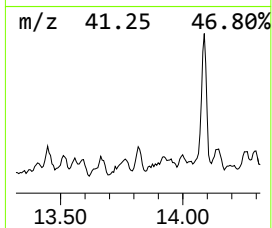
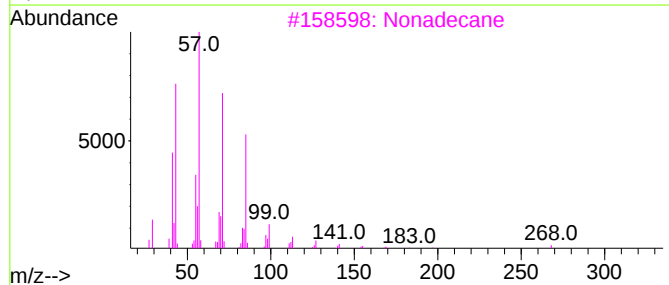
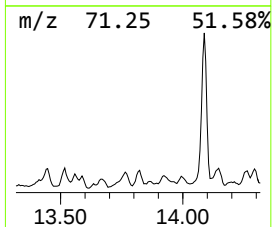
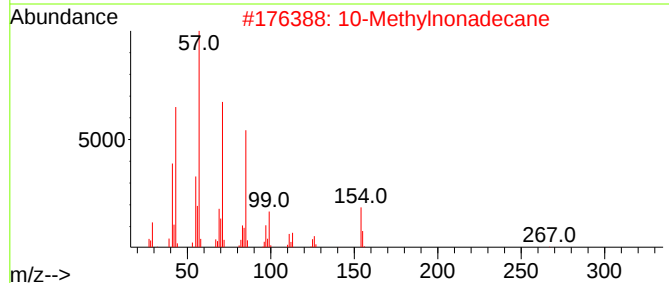
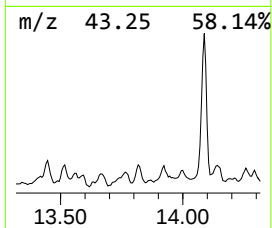
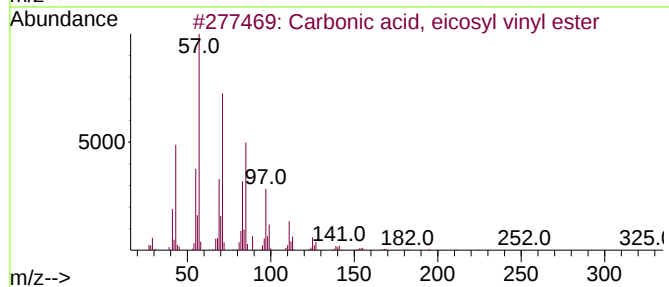
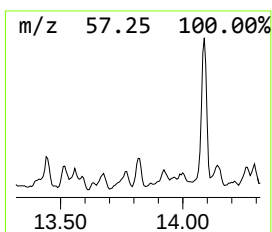
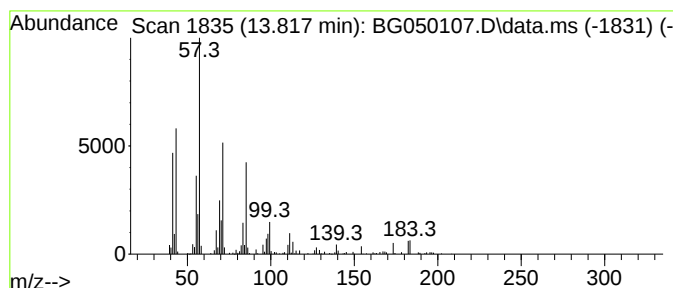
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Quant Title : SVOA CALIBRATION

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

Peak Number 58 10-Methylnonadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.817	19.35 ng/ul	319794	Acenaphthene-d10	14.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	72
2			10-Methylnonadecane	282	C20H42	056862-62-5	72
3			Nonadecane	268	C19H40	000629-92-5	64
4			2-Bromotetradecane	276	C14H29Br	074036-95-6	59
5			Octacosane	394	C28H58	000630-02-4	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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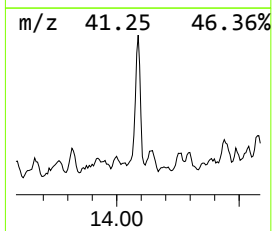
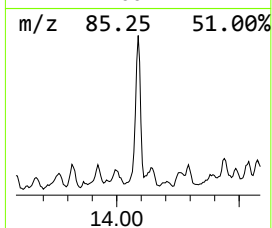
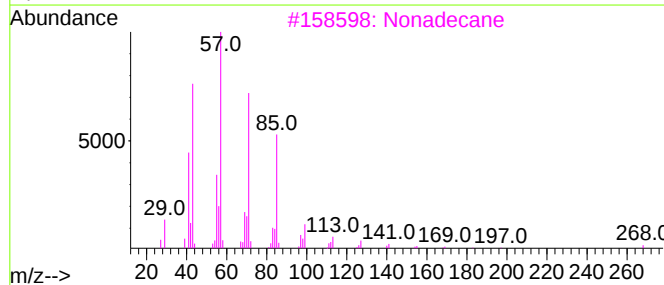
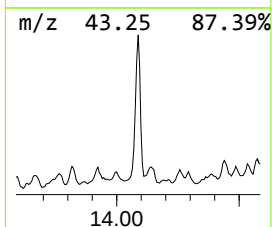
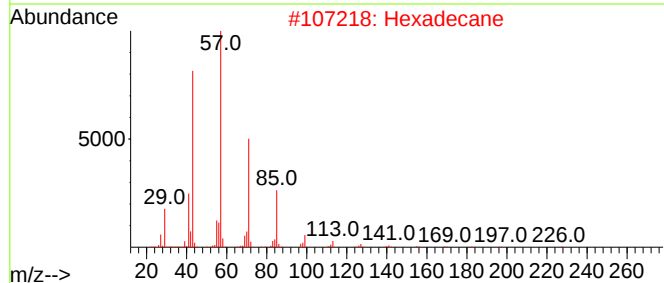
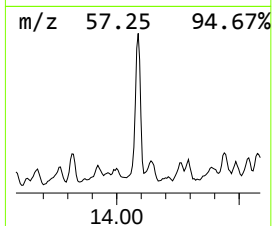
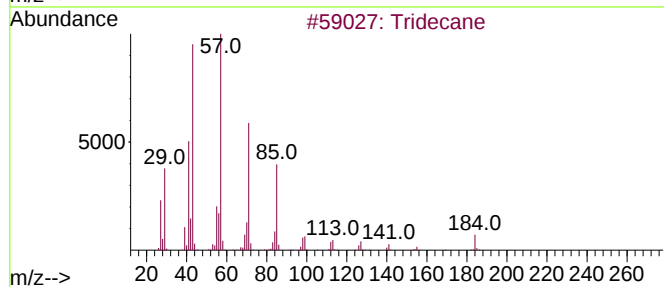
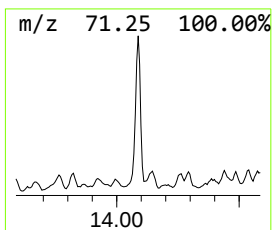
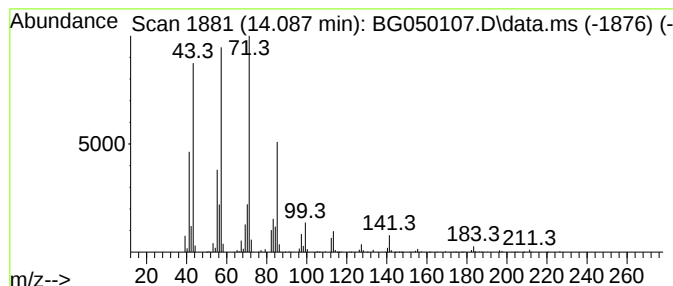
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Quant Title : SVOA CALIBRATION

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

Peak Number 59 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.087	101.60 ng/ul	1679050	Acenaphthene-d10	14.622

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane	184	C13H28	000629-50-5	90
2		Hexadecane	226	C16H34	000544-76-3	87
3		Nonadecane	268	C19H40	000629-92-5	86
4		Pentadecane, 7-methyl-	226	C16H34	006165-40-8	80
5		Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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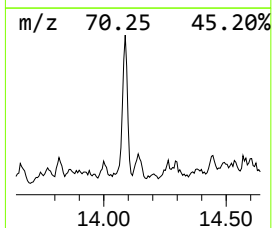
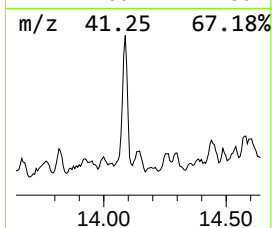
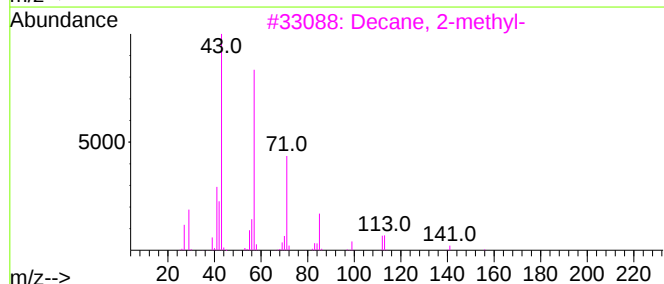
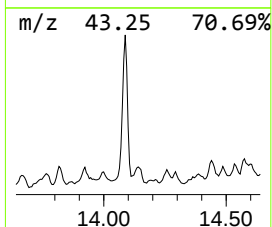
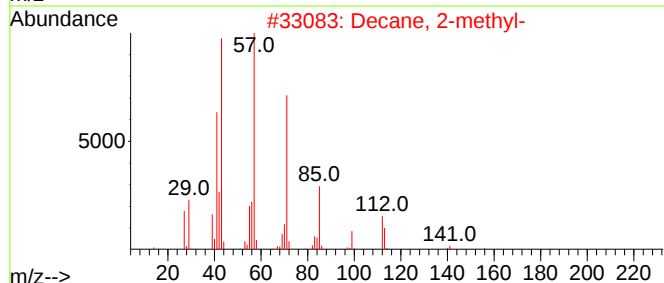
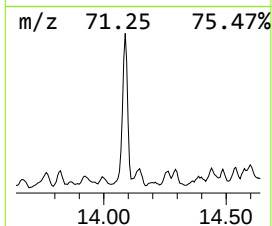
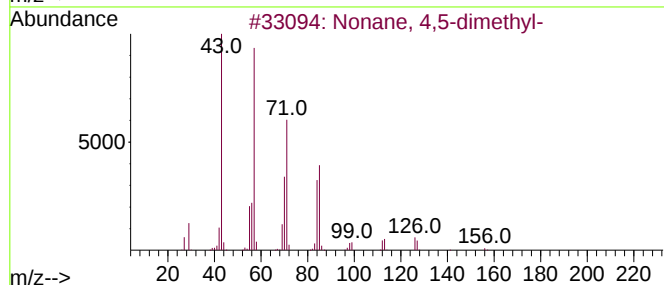
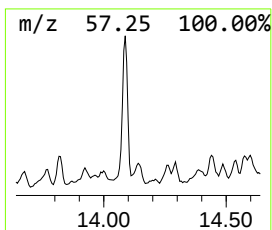
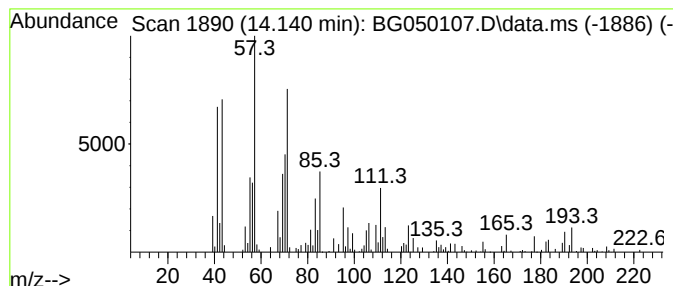
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Quant Title : SVOA CALIBRATION

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

Peak Number 60 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.140	21.42 ng/ul	353986	Acenaphthene-d10	14.622

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonane, 4,5-dimethyl-	156	C11H24	017302-23-7	49
2		Decane, 2-methyl-	156	C11H24	006975-98-0	45
3		Decane, 2-methyl-	156	C11H24	006975-98-0	43
4		Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	38
5		Hexadecane, 1-chloro-	260	C16H33Cl	004860-03-1	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
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Misc :
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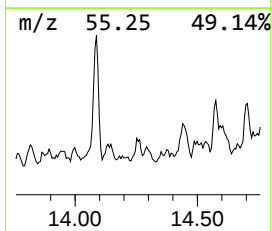
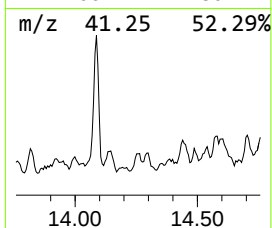
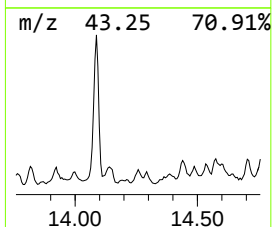
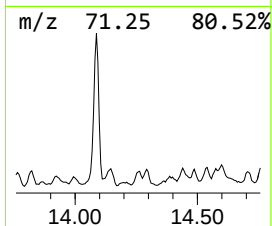
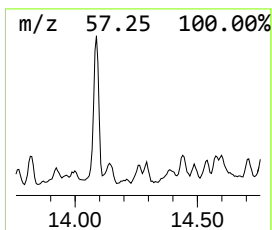
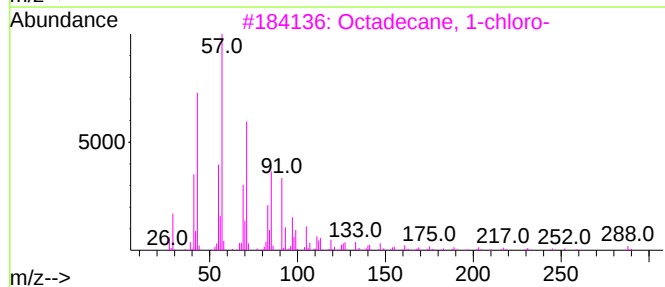
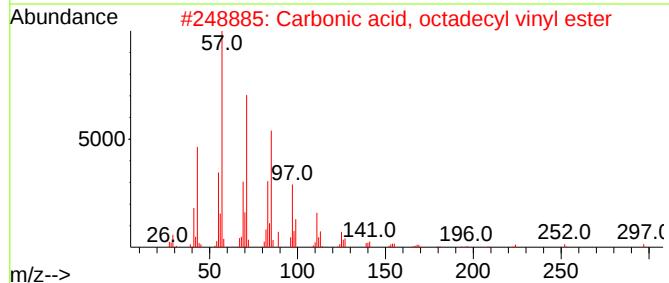
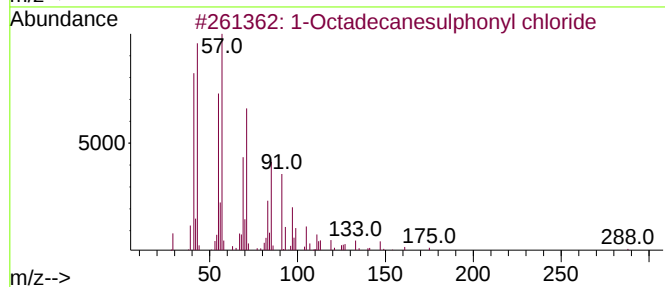
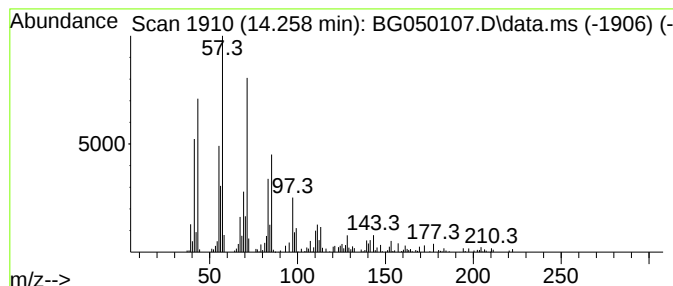
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BNA_G
ClientSampleId :
EW7C4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION

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

Peak Number 61 1-Octadecanesulphonyl chloride Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.258	14.11 ng/ul	233175	Acenaphthene-d10	14.622	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	91
2	Carbonic acid, octadecyl vinyl e...	340	C21H40O3	1000382-54-4	90
3	Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	90
4	Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	90
5	Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EW7C4

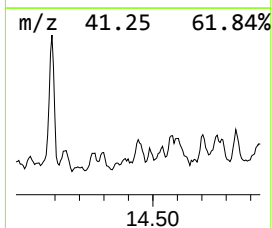
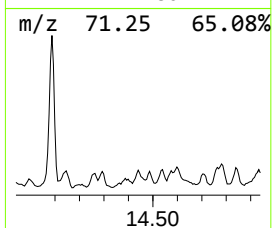
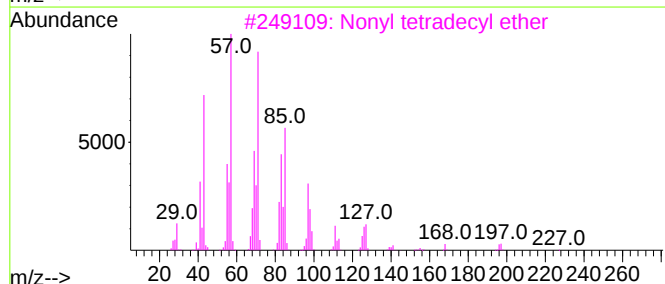
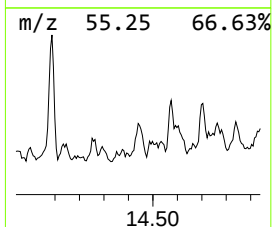
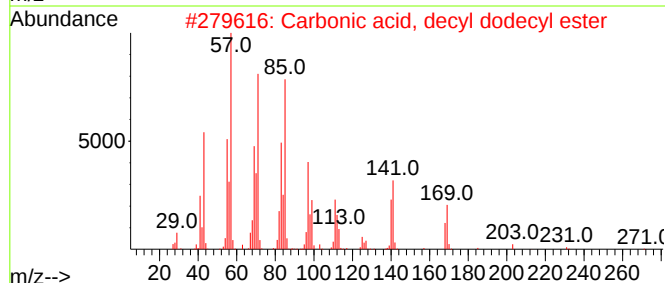
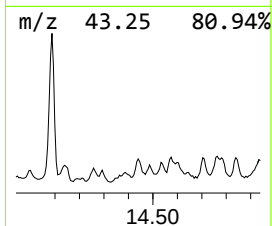
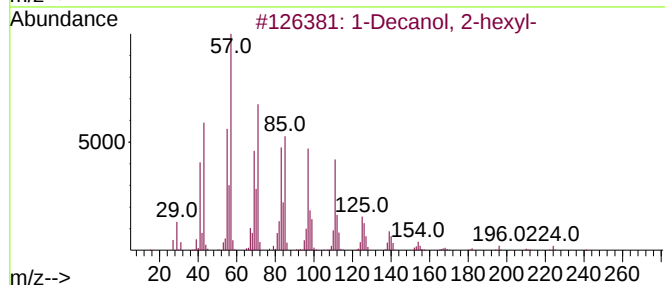
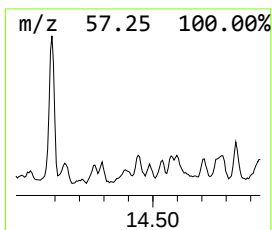
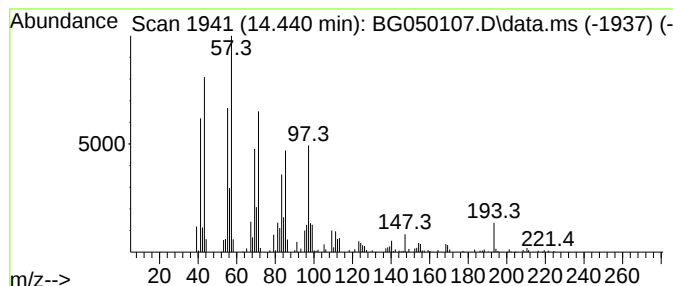
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Quant Title : SVOA CALIBRATION

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

Peak Number 62 1-Decanol, 2-hexyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.440	24.17 ng/ul	399410	Acenaphthene-d10	14.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Decanol, 2-hexyl-			242	C16H34O	002425-77-6	68
2	Carbonic acid, decyl dodecyl ester			370	C23H46O3	1000383-16-1	58
3	Nonyl tetradecyl ether			340	C23H48O	1000406-37-6	58
4	Carbonic acid, decyl undecyl ester			356	C22H44O3	1000383-16-0	52
5	Carbonic acid, decyl tridecyl ester			384	C24H48O3	1000383-16-2	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW7C4

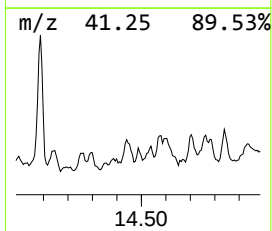
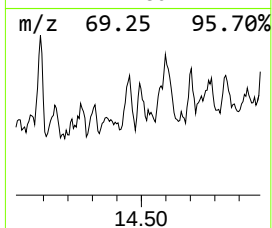
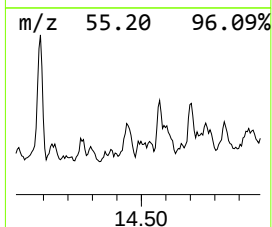
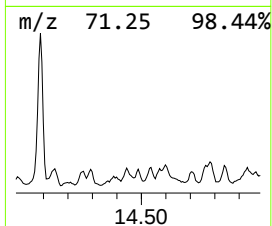
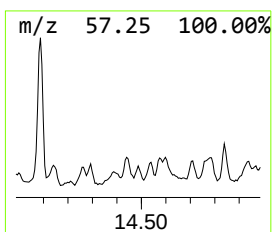
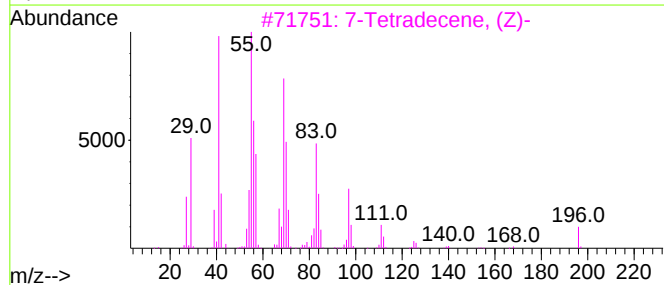
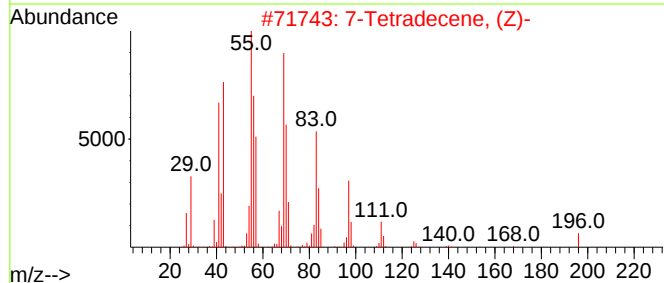
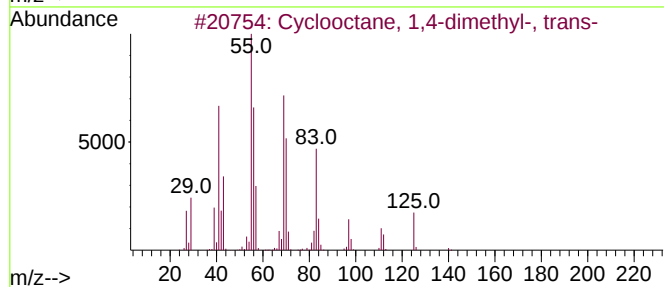
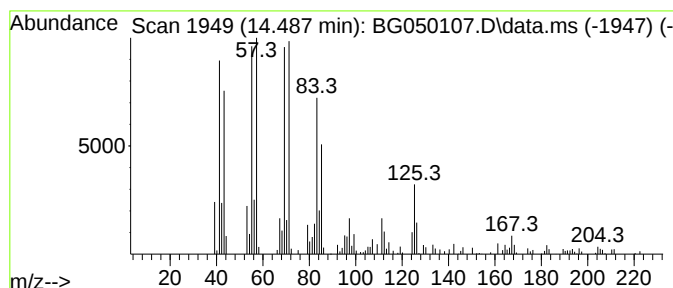
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Quant Title : SVOA CALIBRATION

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

Peak Number 63 (DEL) Alkane: Cyclic14.487 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.487	9.43 ng/ul	155898	Acenaphthene-d10	14.622

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclooctane, 1,4-dimethyl-, trans-	140	C10H20	013151-98-9	46
2		7-Tetradecene, (Z)-	196	C14H28	041446-60-0	43
3		7-Tetradecene, (Z)-	196	C14H28	041446-60-0	41
4		7-Tetradecene, (Z)-	196	C14H28	041446-60-0	41
5		3-Hexene, 2,3-dimethyl-	112	C8H16	007145-23-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW7C4

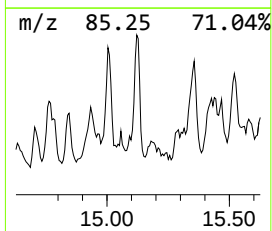
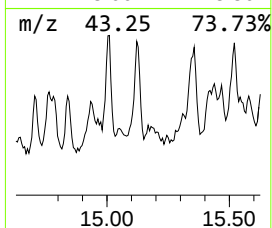
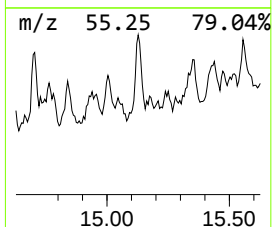
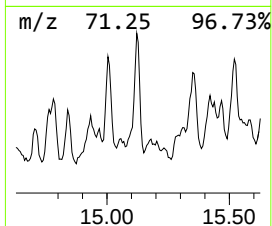
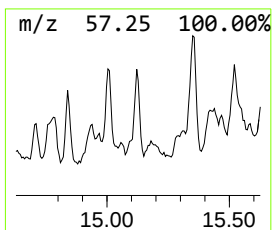
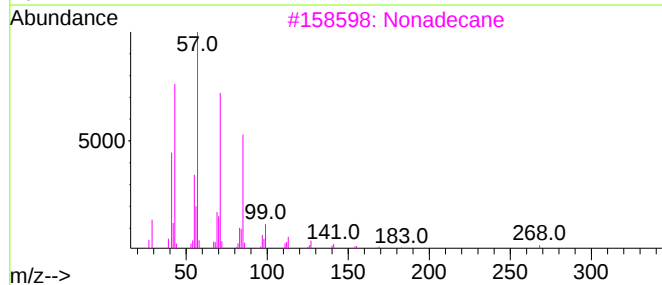
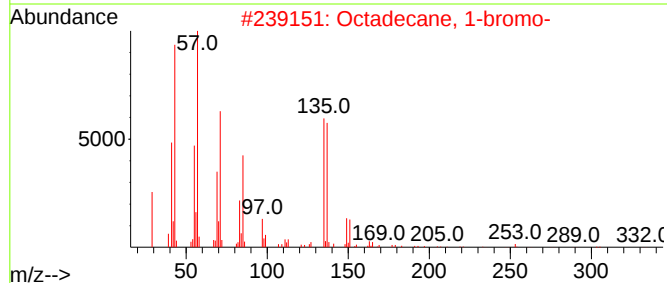
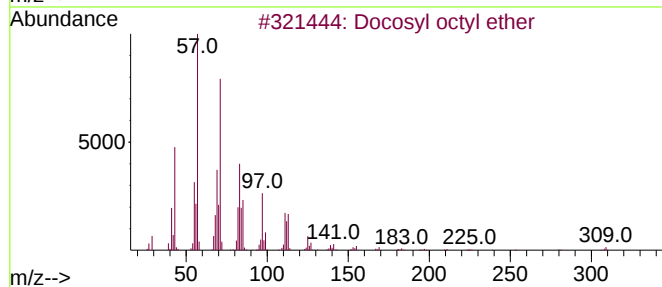
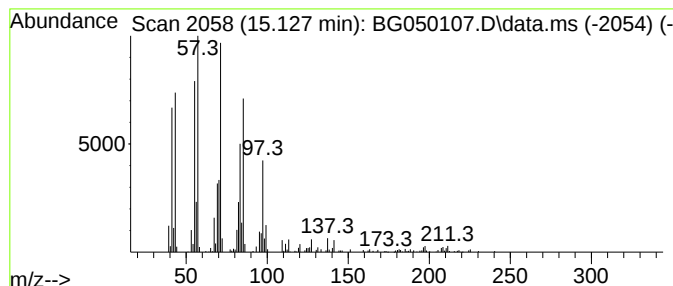
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Quant Title : SVOA CALIBRATION

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

Peak Number 64 Docosyl octyl ether Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.127	30.08 ng/ul	497124	Acenaphthene-d10	14.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Docosyl octyl ether	438	C30H62O	1000406-38-9	64
2			Octadecane, 1-bromo-	332	C18H37Br	000112-89-0	50
3			Nonadecane	268	C19H40	000629-92-5	49
4			Decane, 2,3,8-trimethyl-	184	C13H28	062238-14-6	47
5			Succinic acid, dodec-2-en-1-yl t...	368	C21H36O5	1000390-72-8	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

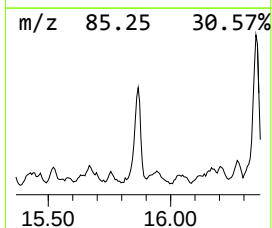
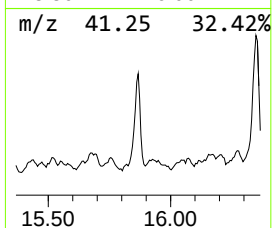
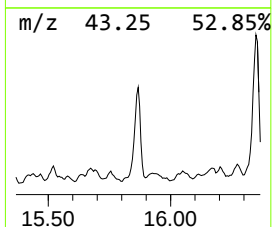
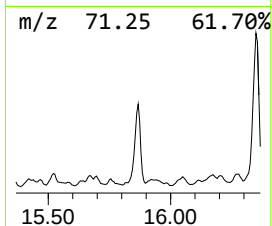
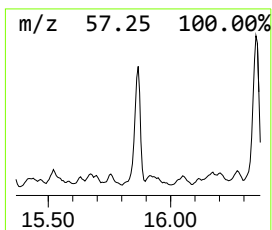
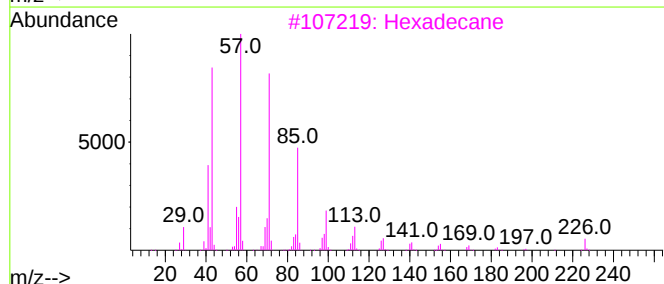
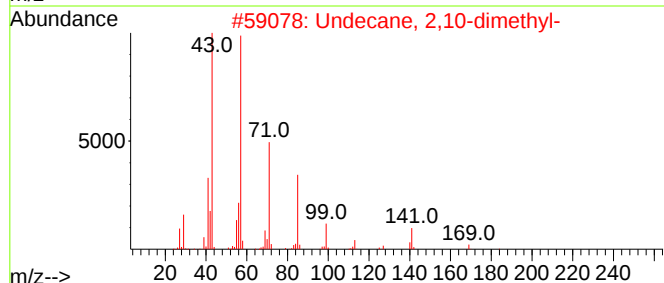
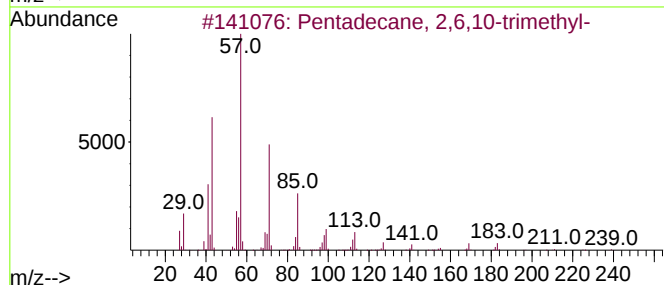
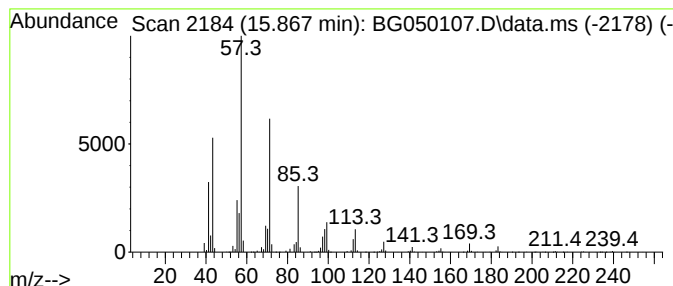
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 65 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD			R.T.
15.867	165.14 ng/ul	2729230	Acenaphthene-d10			14.622
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10-trimethyl-		254	C18H38	003892-00-0	94
2	Undecane, 2,10-dimethyl-		184	C13H28	017301-27-8	80
3	Hexadecane		226	C16H34	000544-76-3	80
4	Hexadecane, 2,6,10,14-tetramethyl-		282	C20H42	000638-36-8	80
5	Heneicosane		296	C21H44	000629-94-7	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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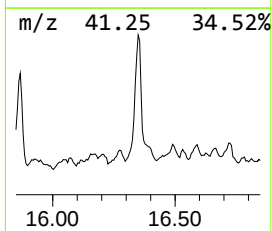
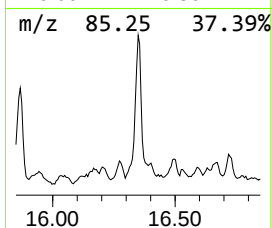
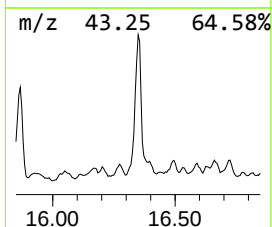
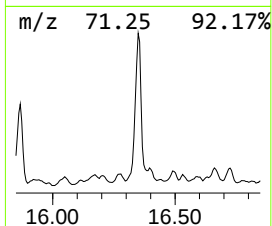
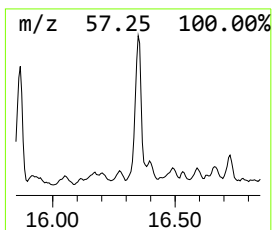
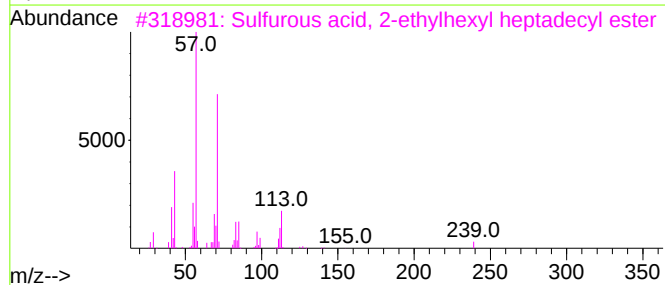
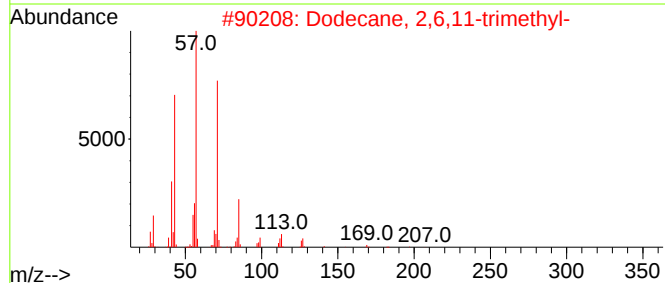
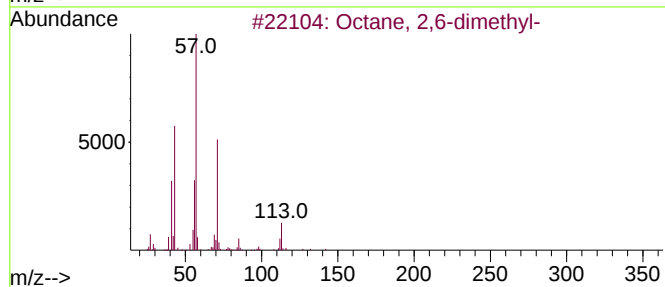
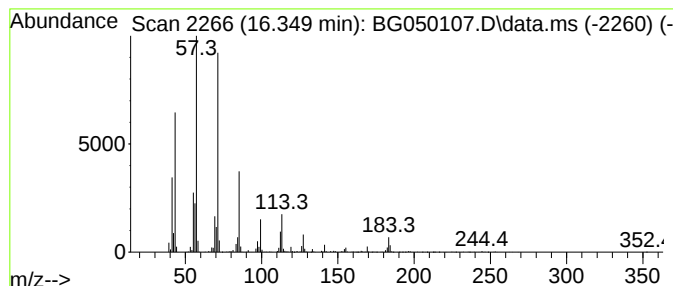
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Quant Title : SVOA CALIBRATION

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

Peak Number 66 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.349	259.14 ng/ul	5134670	Phenanthrene-d10	17.377

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-		142	C10H22		002051-30-1	72
2	Dodecane, 2,6,11-trimethyl-		212	C15H32		031295-56-4	72
3	Sulfurous acid, 2-ethylhexyl hep...		432	C25H52O3S		1000309-20-0	72
4	Sulfurous acid, decyl 2-ethylhex...		334	C18H38O3S		1000309-19-3	72
5	Sulfurous acid, 2-ethylhexyl non...		320	C17H36O3S		1010309-19-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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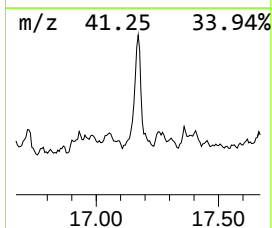
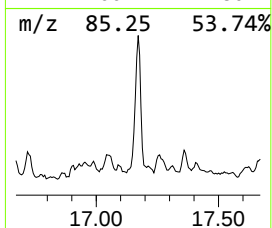
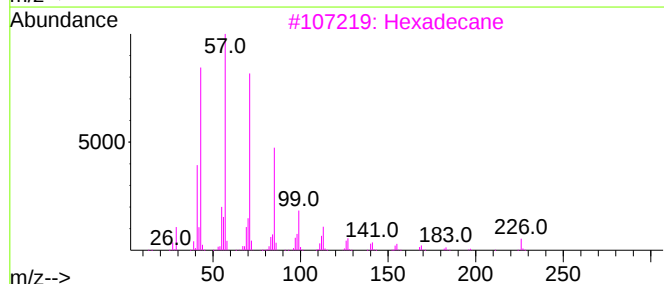
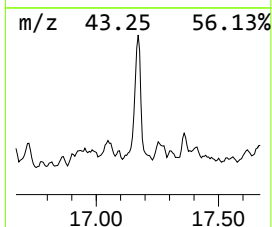
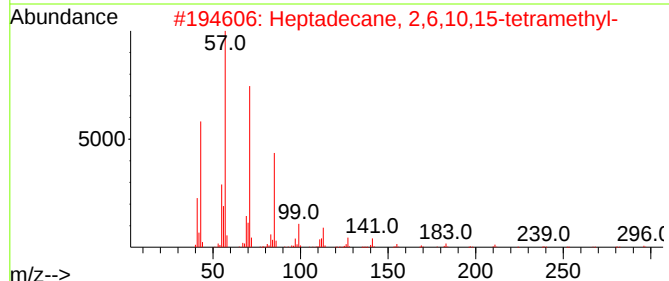
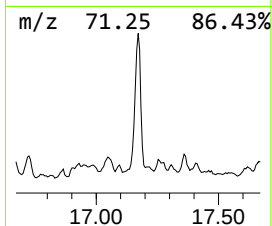
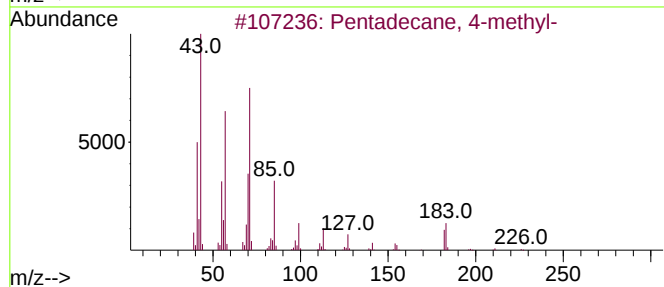
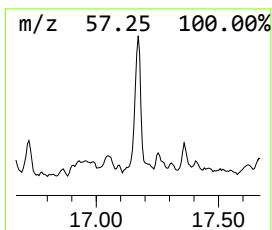
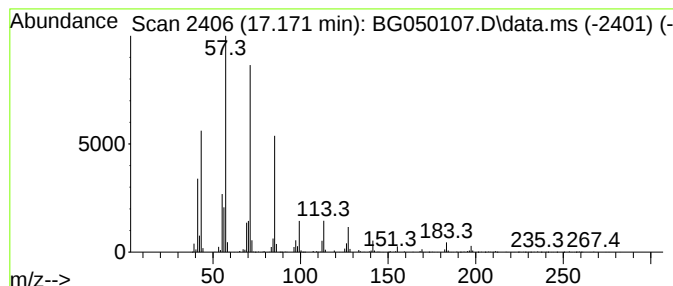
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Quant Title : SVOA CALIBRATION

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

Peak Number 67 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.171	168.51 ng/ul	3338970	Phenanthrene-d10	17.377

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane, 4-methyl-	226	C16H34	002801-87-8	90
2			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	87
3			Hexadecane	226	C16H34	000544-76-3	87
4			Tridecane, 4-methyl-	198	C14H30	026730-12-1	87
5			Heneicosane	296	C21H44	000629-94-7	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050107.D
Acq On : 2 Sep 2021 21:53
Operator : CG/JU
Sample : M3526-14
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW7C4

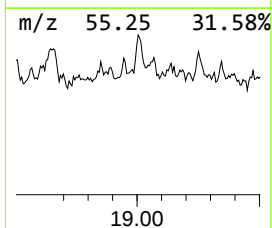
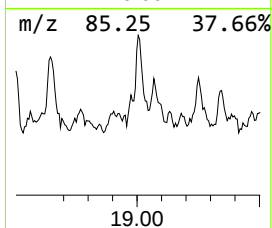
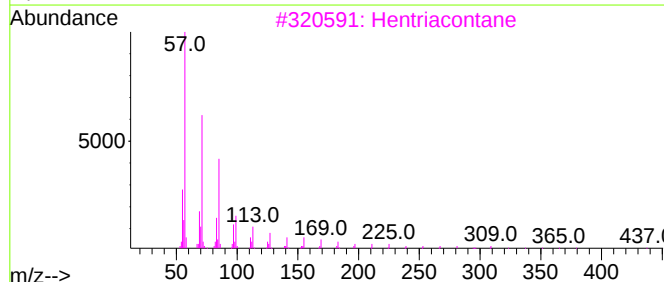
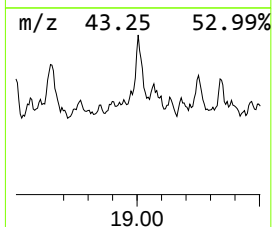
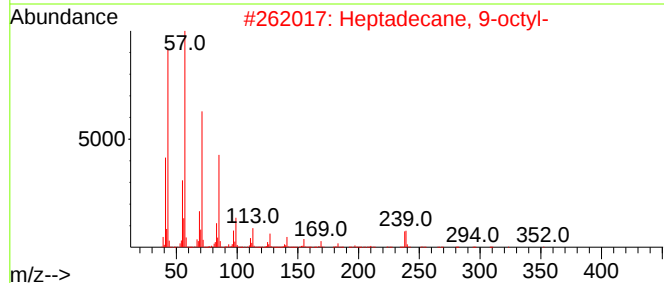
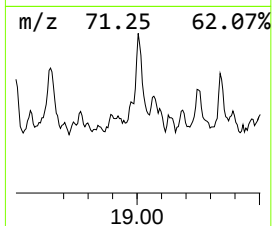
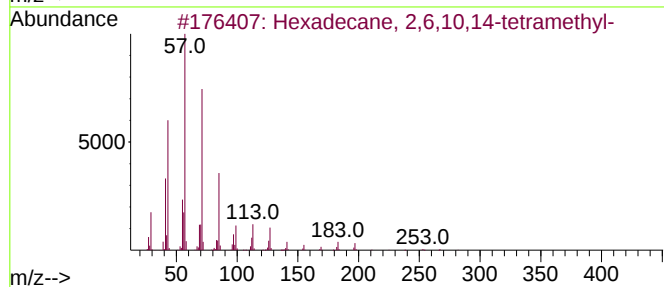
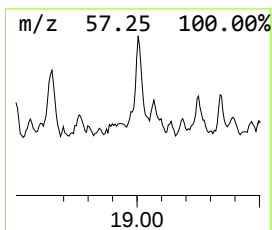
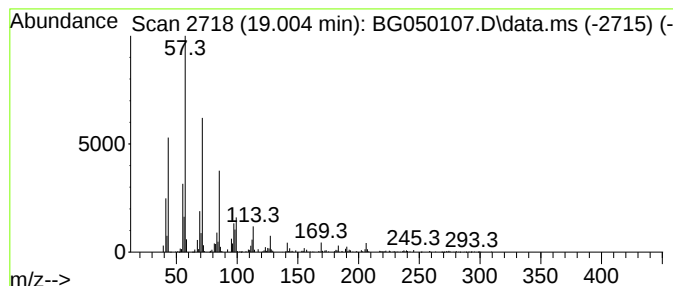
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Quant Title : SVOA CALIBRATION

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

Peak Number 68 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.004	30.43 ng/ul	602941	Phenanthrene-d10	17.377

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	90
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	80
3			Hentriacontane	437	C31H64	000630-04-6	80
4			Sulfurous acid, butyl tetradecyl...	334	C18H38O3S	1010309-18-1	80
5			Hexacosane	366	C26H54	000630-01-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
 Data File : BG050107.D
 Acq On : 2 Sep 2021 21:53
 Operator : CG/JU
 Sample : M3526-14
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW7C4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.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
Divinyl sulfide	3.848	4.6	ng/ul	38438	1	7.949	167072	20.0
(DEL) Alkane: S...	6.920	2.5	ng/ul	21145	1	7.949	167072	20.0
(DEL) Alkane: C...	7.432	2.3	ng/ul	19425	1	7.949	167072	20.0
(DEL) Alkane: S...	8.131	4.6	ng/ul	38735	1	7.949	167072	20.0
(DEL) Alkane: C...	8.225	3.0	ng/ul	25299	1	7.949	167072	20.0
Naphthalene, de...	8.789	5.5	ng/ul	45834	1	7.949	167072	20.0
Oxalic acid, cy...	9.059	11.3	ng/ul	94634	1	7.949	167072	20.0
unknown-01	9.212	3.6	ng/ul	29885	1	7.949	167072	20.0
Sulfurous acid,...	9.270	5.5	ng/ul	45771	1	7.949	167072	20.0
(DEL) Alkane: C...	9.311	9.7	ng/ul	80811	1	7.949	167072	20.0
Oxirane, [(dode...	9.423	6.1	ng/ul	86372	2	10.774	281868	20.0
(DEL) Alkane: S...	9.529	3.0	ng/ul	41599	2	10.774	281868	20.0
(DEL) Alkane: S...	9.587	5.9	ng/ul	82653	2	10.774	281868	20.0
(DEL) Alkane: C...	9.646	2.8	ng/ul	39012	2	10.774	281868	20.0
trans-Decalin, ...	9.681	6.3	ng/ul	88230	2	10.774	281868	20.0
Naphthalene, de...	9.940	7.8	ng/ul	110302	2	10.774	281868	20.0
(DEL) Alkane: C...	10.075	2.1	ng/ul	29044	2	10.774	281868	20.0
unknown-02	10.175	3.0	ng/ul	42490	2	10.774	281868	20.0
Trifluoroacetox...	10.339	7.0	ng/ul	99304	2	10.774	281868	20.0
unknown-03	10.539	4.0	ng/ul	56610	2	10.774	281868	20.0
(DEL) Alkane: C...	10.657	10.8	ng/ul	151938	2	10.774	281868	20.0
Decyl isobutyl ...	10.980	4.5	ng/ul	62869	2	10.774	281868	20.0
Octadecane, 1-c...	11.138	8.4	ng/ul	119044	2	10.774	281868	20.0
Tetrapentaconta...	11.250	3.9	ng/ul	54776	2	10.774	281868	20.0
(DEL) Alkane: S...	11.315	6.0	ng/ul	83975	2	10.774	281868	20.0
(DEL) Alkane: S...	11.420	4.8	ng/ul	66967	2	10.774	281868	20.0
(DEL) Alkane: C...	11.467	10.6	ng/ul	148735	2	10.774	281868	20.0
unknown-04	11.532	4.9	ng/ul	69578	2	10.774	281868	20.0
(DEL) Alkane: C...	11.602	10.7	ng/ul	150782	2	10.774	281868	20.0
2-Decanol, trif...	11.696	10.9	ng/ul	153209	2	10.774	281868	20.0
(DEL) Alkane: S...	11.785	35.0	ng/ul	493554	2	10.774	281868	20.0
unknown-05	11.826	8.5	ng/ul	119378	2	10.774	281868	20.0
(DEL) Alkane: C...	12.090	5.1	ng/ul	71706	2	10.774	281868	20.0
(DEL) Alkane: C...	12.172	2.6	ng/ul	37355	2	10.774	281868	20.0
2,4-Dimethyl-1,...	12.319	7.3	ng/ul	102780	2	10.774	281868	20.0
unknown-06	12.501	6.5	ng/ul	91607	2	10.774	281868	20.0
Sulfurous acid,...	12.765	8.0	ng/ul	132328	3	14.622	330537	20.0
unknown-07	12.818	10.3	ng/ul	170092	3	14.622	330537	20.0
(DEL) Alkane: C...	12.877	17.8	ng/ul	294890	3	14.622	330537	20.0
(DEL) Alkane: S...	13.136	45.3	ng/ul	748474	3	14.622	330537	20.0
Sulfurous acid,...	13.412	7.3	ng/ul	120971	3	14.622	330537	20.0
(DEL) Alkane: C...	13.453	18.2	ng/ul	301505	3	14.622	330537	20.0
Carbonic acid, ...	13.559	5.5	ng/ul	91263	3	14.622	330537	20.0
unknown-08	13.635	4.5	ng/ul	75240	3	14.622	330537	20.0
unknown-09	13.664	16.0	ng/ul	264364	3	14.622	330537	20.0
10-Methylnonade...	13.817	19.4	ng/ul	319794	3	14.622	330537	20.0
(DEL) Alkane: S...	14.087	101.6	ng/ul	1679050	3	14.622	330537	20.0
(DEL) Alkane: S...	14.140	21.4	ng/ul	353986	3	14.622	330537	20.0
1-Octadecanesul...	14.258	14.1	ng/ul	233175	3	14.622	330537	20.0
1-Decanol, 2-he...	14.440	24.2	ng/ul	399410	3	14.622	330537	20.0
(DEL) Alkane: C...	14.487	9.4	ng/ul	155898	3	14.622	330537	20.0
Docosyl octyl e...	15.127	30.1	ng/ul	497124	3	14.622	330537	20.0

(DEL) Alkane: S...	15.867	165.1	ng/ul	2729230	3	14.622	330537	20.0
(DEL) Alkane: S...	16.349	259.1	ng/ul	5134670	4	17.377	396289	20.0
(DEL) Alkane: S...	17.171	168.5	ng/ul	3338970	4	17.377	396289	20.0
(DEL) Alkane: S...	19.004	30.4	ng/ul	602941	4	17.377	396289	20.0

Instrument :
BNA_G
ClientSampleId :
EW7C4