

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
 Data File : BG049489.D
 Acq On : 26 Jul 2021 22:10
 Operator : CG/JU
 Sample : M3082-19
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0017

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M

Title : SVOA CALIBRATION

Signal : TIC: BG049489.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.080	3	7	16	rBV2	92514	194873	45.55%	3.182%
2	3.162	16	21	27	rVV	150099	229197	53.57%	3.742%
3	3.438	64	68	72	rBV4	4236	7342	1.72%	0.120%
4	3.861	136	140	149	rBV	32195	61902	14.47%	1.011%
5	3.967	154	158	164	rVB3	15247	24884	5.82%	0.406%
6	4.090	177	179	183	rVB3	3362	3610	0.84%	0.059%
7	4.754	289	292	296	rVB3	2611	3892	0.91%	0.064%
8	5.218	368	371	375	rBV4	3289	4864	1.14%	0.079%
9	5.289	382	383	388	rBV4	1198	1375	0.32%	0.022%
10	5.459	409	412	416	rVB3	1573	2576	0.60%	0.042%
11	5.488	416	417	421	rBV	1217	1727	0.40%	0.028%
12	6.041	510	511	514	rVB	2327	1511	0.35%	0.025%
13	6.182	532	535	538	rBV3	996	1385	0.32%	0.023%
14	6.423	573	576	579	rBV2	1237	1893	0.44%	0.031%
15	6.681	617	620	621	rBV2	1492	1709	0.40%	0.028%
16	7.204	701	709	718	rBV	73297	130160	30.42%	2.125%
17	7.368	729	737	746	rBV	50391	90087	21.06%	1.471%
18	7.480	752	756	766	rVB3	2814	6110	1.43%	0.100%
19	7.580	766	773	780	rBV	74227	132485	30.97%	2.163%
20	7.662	784	787	790	rBV3	2994	4065	0.95%	0.066%
21	7.691	790	792	794	rVB	2146	1632	0.38%	0.027%
22	8.050	846	853	863	rBV	99800	173318	40.51%	2.830%
23	8.273	885	891	897	rVB2	28254	46889	10.96%	0.766%
24	8.755	967	973	982	rBV2	71941	130464	30.49%	2.130%
25	8.884	992	995	999	rBV3	1476	1588	0.37%	0.026%
26	9.213	1045	1051	1059	rBV	76663	141943	33.18%	2.318%
27	9.272	1059	1061	1067	rVB3	2539	4320	1.01%	0.071%
28	9.366	1074	1077	1080	rBV2	1488	2219	0.52%	0.036%
29	9.947	1168	1176	1183	rBV	71095	131499	30.74%	2.147%
30	10.488	1261	1268	1277	rBV2	100914	182053	42.55%	2.972%
31	10.629	1285	1292	1304	rBV2	67537	122108	28.54%	1.994%
32	10.734	1308	1310	1314	rVB3	1909	1906	0.45%	0.031%
33	10.869	1327	1333	1339	rVB	135973	235383	55.02%	3.843%
34	11.005	1349	1356	1364	rBV3	43063	87585	20.47%	1.430%
35	11.392	1420	1422	1425	rBV3	1765	1936	0.45%	0.032%
36	11.880	1501	1505	1509	rBV2	2071	3928	0.92%	0.064%

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Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

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

37	12.221	1561	1563	1566	rBV3	2127	1926	0.45%	0.031%
38	12.268	1569	1571	1572	rBV3	2157	1687	0.39%	0.028%
39	12.461	1601	1604	1605	rVB2	2054	1355	0.32%	0.022%
40	12.532	1612	1616	1619	rVB	3721	6079	1.42%	0.099%
41	12.743	1649	1652	1656	rVB	2445	2953	0.69%	0.048%
42	12.972	1689	1691	1694	rBV	1196	1362	0.32%	0.022%
43	13.002	1694	1696	1699	rBV	1785	1422	0.33%	0.023%
44	13.061	1705	1706	1709	rBV2	2428	1691	0.40%	0.028%
45	13.184	1723	1727	1730	rBV	988	1709	0.40%	0.028%
46	13.231	1732	1735	1738	rVB2	3564	3279	0.77%	0.054%
47	13.296	1742	1746	1747	rBV	2027	2410	0.56%	0.039%
48	13.448	1769	1772	1778	rVB	1056	1855	0.43%	0.030%
49	13.513	1778	1783	1789	rBV4	5957	11093	2.59%	0.181%
50	13.578	1789	1794	1798	rBV3	7854	10384	2.43%	0.170%
51	13.848	1838	1840	1841	rBV	2714	2039	0.48%	0.033%
52	13.977	1857	1862	1864	rBV	735	1368	0.32%	0.022%
53	14.018	1864	1869	1873	rVV2	5073	8085	1.89%	0.132%
54	14.094	1873	1882	1889	rVV	197111	296264	69.25%	4.837%
55	14.165	1892	1894	1898	rVB2	2855	3538	0.83%	0.058%
56	14.253	1906	1909	1911	rBV2	1786	1787	0.42%	0.029%
57	14.335	1918	1923	1924	rBV2	2358	1845	0.43%	0.030%
58	14.394	1927	1933	1944	rVB	189611	317545	74.22%	5.185%
59	14.582	1962	1965	1969	rVB3	6979	9342	2.18%	0.153%
60	14.700	1972	1985	1991	rVV	225325	371379	86.80%	6.064%
61	14.764	1991	1996	2003	rVB2	3030	6569	1.54%	0.107%
62	14.893	2013	2018	2027	rBV3	78880	138382	32.34%	2.259%
63	15.040	2040	2043	2048	rVB2	5945	8047	1.88%	0.131%
64	15.093	2048	2052	2057	rVB5	7195	11128	2.60%	0.182%
65	15.175	2063	2066	2067	rVB2	2834	2418	0.57%	0.039%
66	15.305	2086	2088	2094	rVB3	3785	5100	1.19%	0.083%
67	15.357	2094	2097	2101	rBV2	2940	4380	1.02%	0.072%
68	15.422	2106	2108	2111	rVB2	1605	1495	0.35%	0.024%
69	15.481	2113	2118	2119	rVV3	3833	5236	1.22%	0.085%
70	15.534	2122	2127	2131	rVB2	7918	12641	2.95%	0.206%
71	15.692	2147	2154	2159	rBV	266801	407449	95.24%	6.653%
72	15.739	2159	2162	2168	rVV3	7195	10198	2.38%	0.167%
73	15.804	2168	2173	2183	rVB2	64034	99705	23.30%	1.628%
74	15.869	2183	2184	2188	rBV	2800	3326	0.78%	0.054%
75	15.927	2188	2194	2203	rVB7	7777	18286	4.27%	0.299%
76	16.057	2212	2216	2223	rVB2	46402	66364	15.51%	1.084%

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 Sampling : 1 Min Area: 0.1 % of largest Peak
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 Stop Thrs : 0 Peak Location: TOP

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

77	16.151	2227	2232	2235	rBV5	2645	5438	1.27%	0.089%
78	16.256	2248	2250	2251	rBV2	2755	1352	0.32%	0.022%
79	16.397	2270	2274	2276	rBV2	8018	12638	2.95%	0.206%
80	16.803	2339	2343	2345	rBV4	4489	6464	1.51%	0.106%
81	17.331	2429	2433	2441	rBV5	12063	20244	4.73%	0.331%
82	17.390	2441	2443	2446	rBV2	4893	5150	1.20%	0.084%
83	17.449	2447	2453	2457	rBV	283054	427834	100.00%	6.985%
84	17.490	2457	2460	2464	rVB	74340	99924	23.36%	1.632%
85	17.549	2464	2470	2475	rBV	252085	376691	88.05%	6.150%
86	18.101	2561	2564	2569	rVB4	10712	15002	3.51%	0.245%
87	18.330	2598	2603	2607	rBV3	118872	163894	38.31%	2.676%
88	19.329	2769	2773	2777	rVB2	105938	128002	29.92%	2.090%
89	19.499	2797	2802	2808	rVB	253237	348943	81.56%	5.697%
90	19.593	2816	2818	2825	rBV5	37268	50037	11.70%	0.817%
91	19.834	2854	2859	2862	rBV	257517	425476	99.45%	6.947%

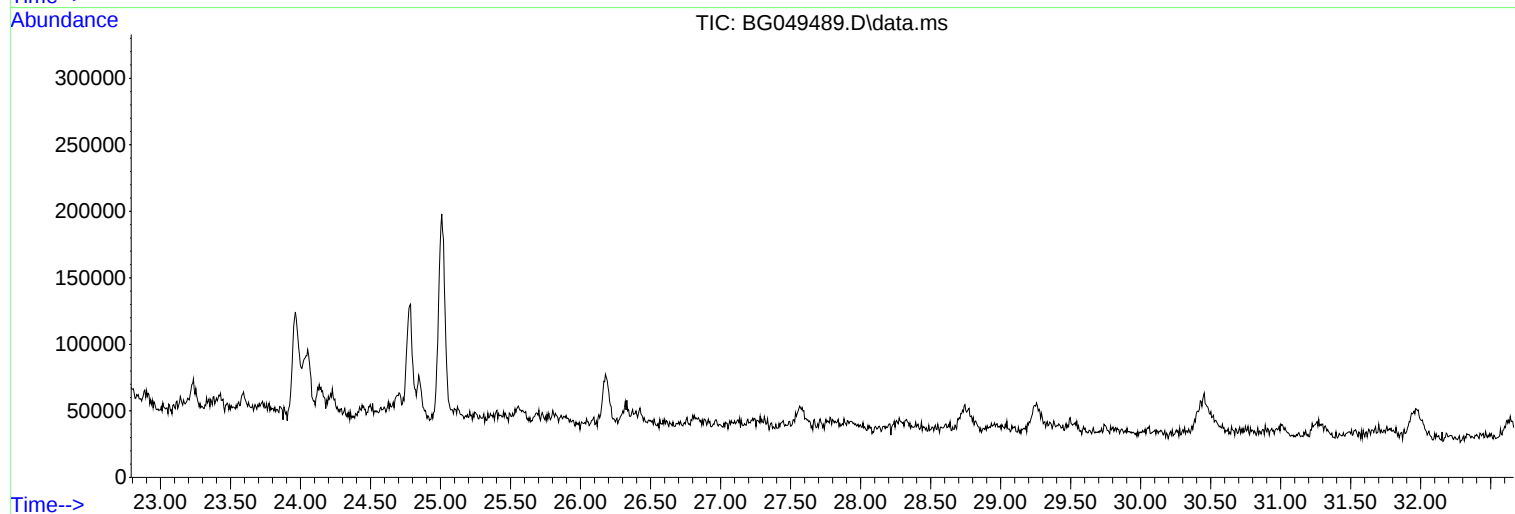
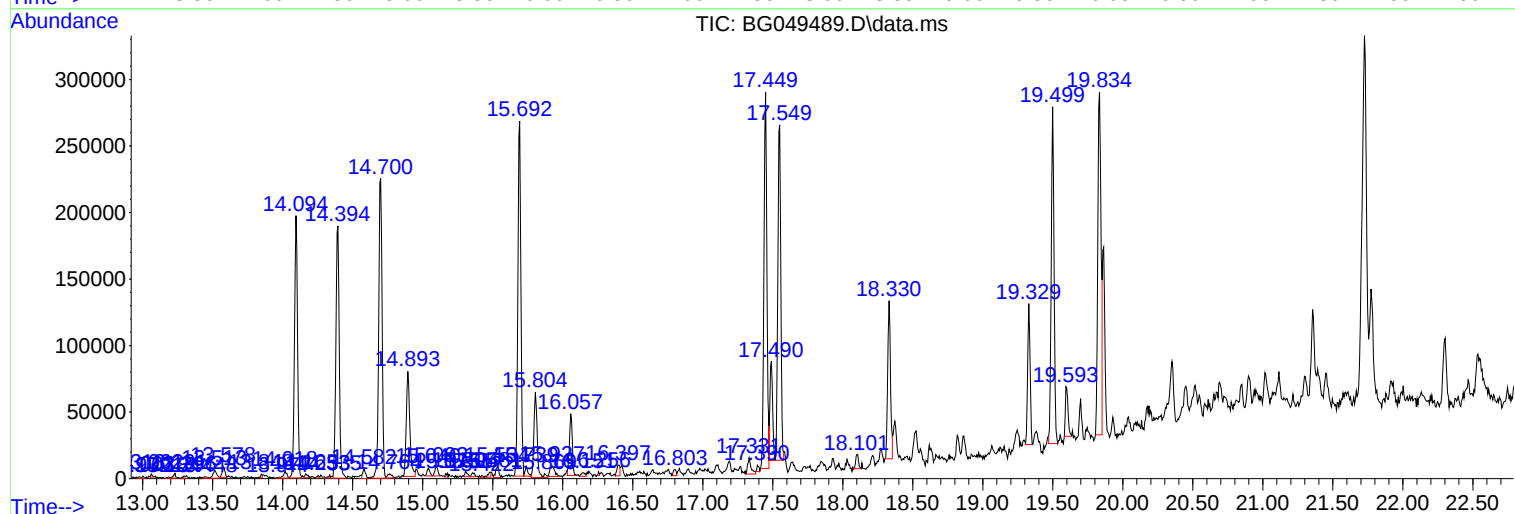
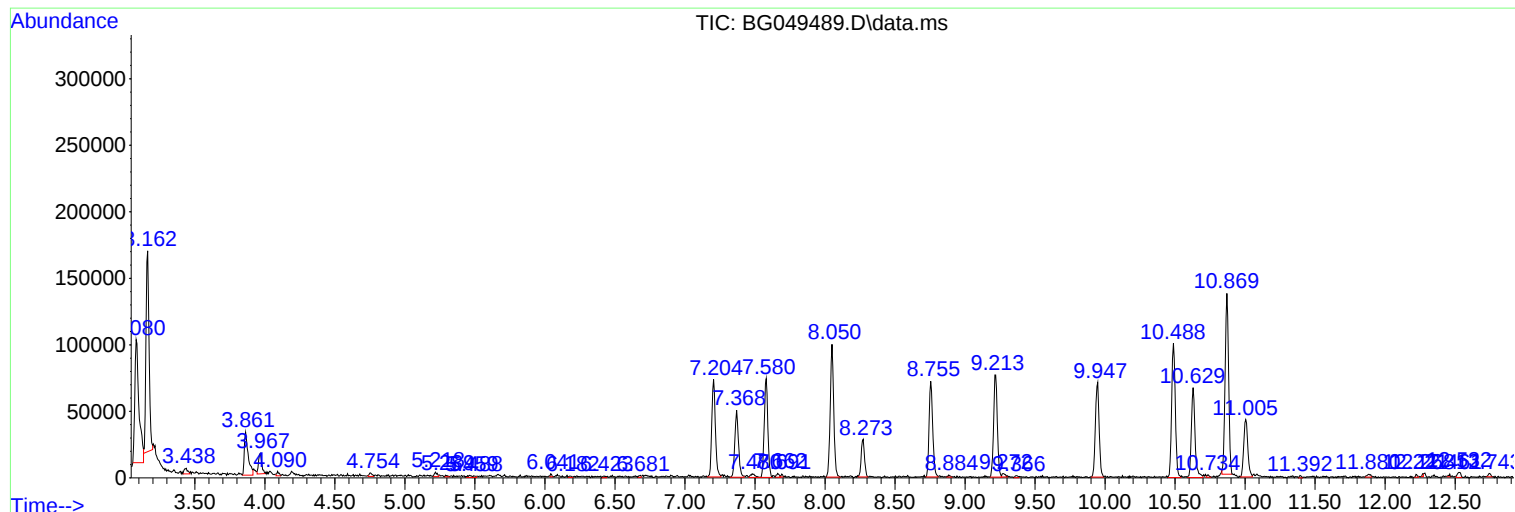
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION

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



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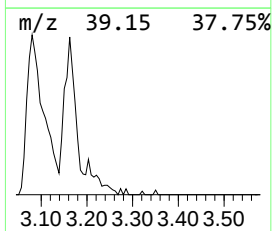
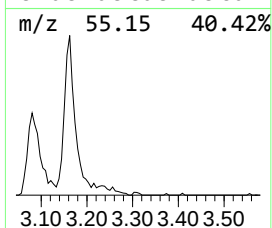
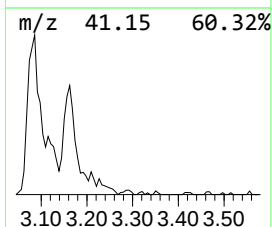
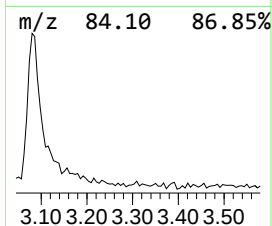
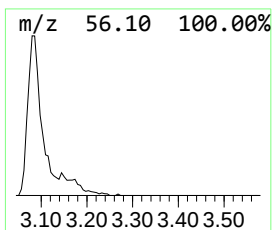
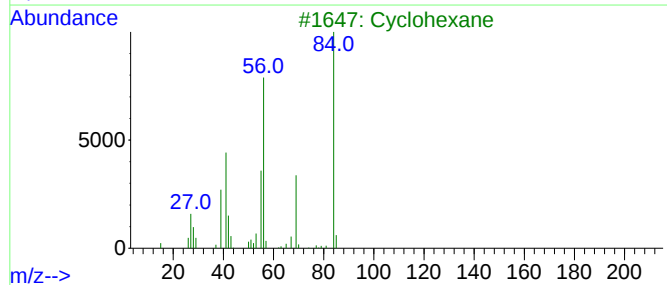
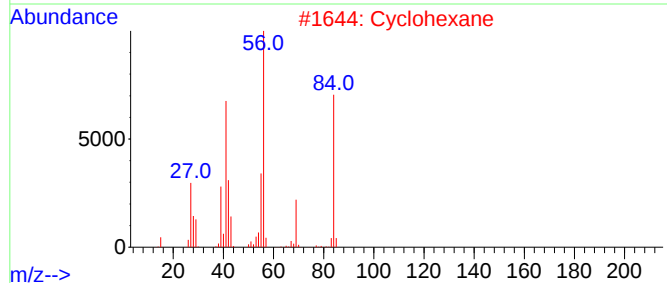
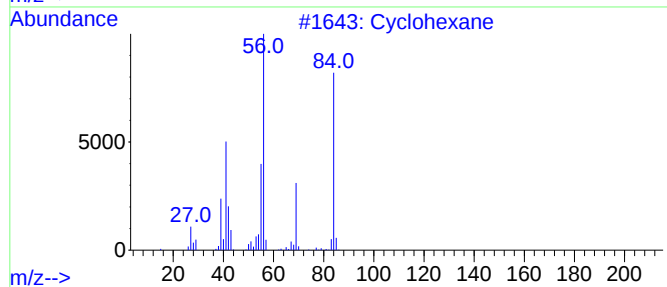
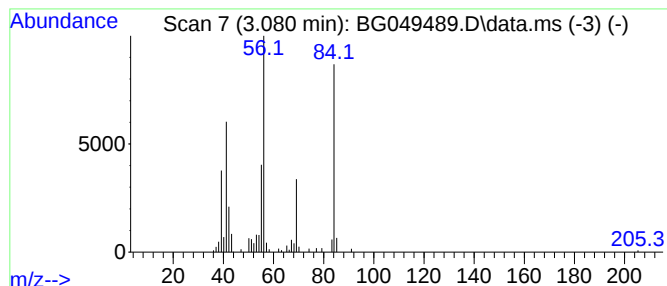
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic3.080 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.080	22.49 ng/ul	194873	1,4-Dichlorobenzene-d4	8.050

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane	84	C6H12	000110-82-7	95
2			Cyclohexane	84	C6H12	000110-82-7	91
3			Cyclohexane	84	C6H12	000110-82-7	91
4			Cyclohexane	84	C6H12	000110-82-7	91
5			Cyclohexane	84	C6H12	000110-82-7	80



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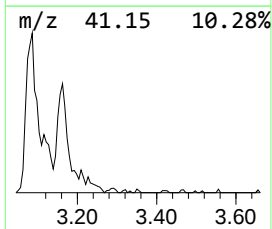
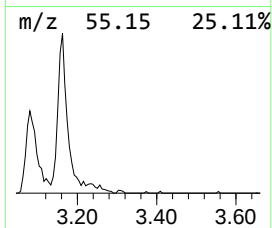
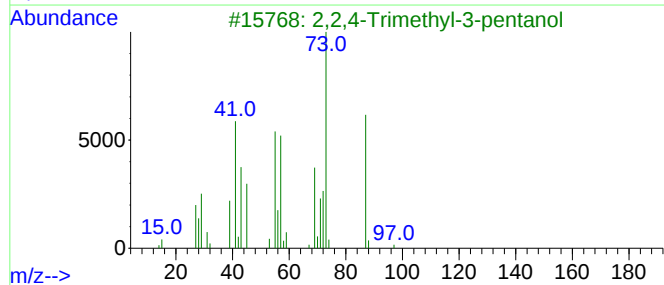
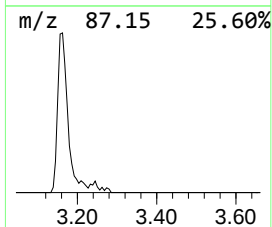
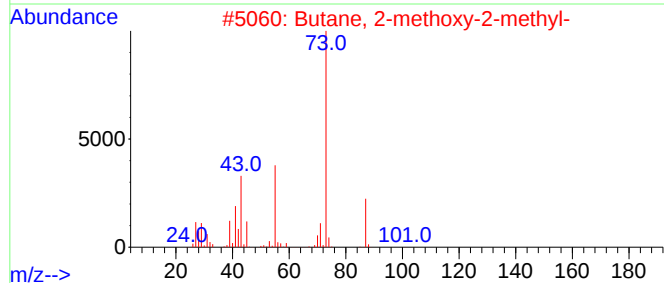
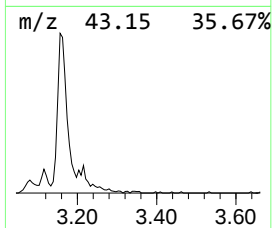
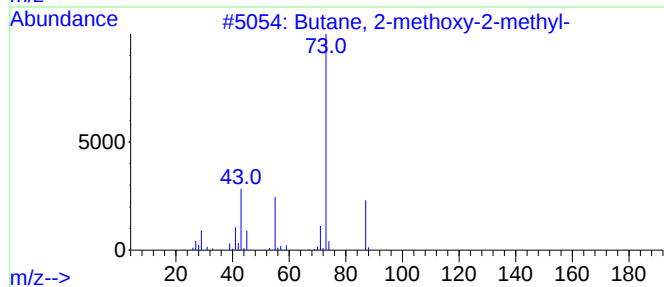
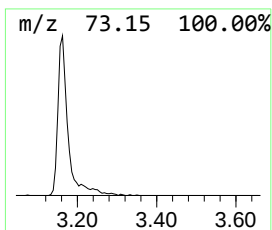
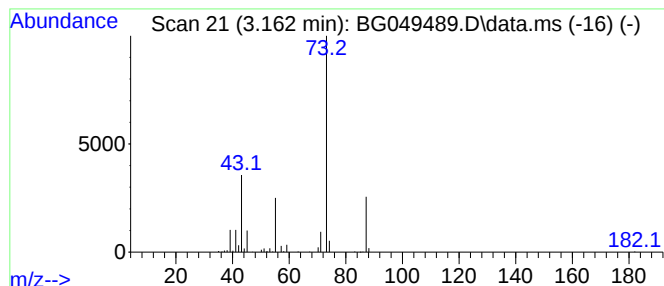
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.162	26.45 ng/ul	229197	1,4-Dichlorobenzene-d4	8.050

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	64
4			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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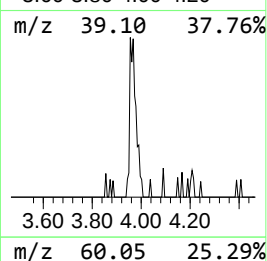
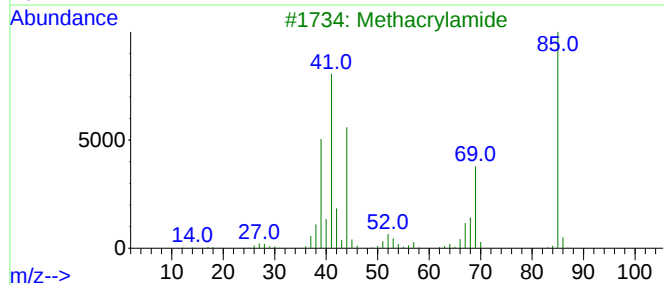
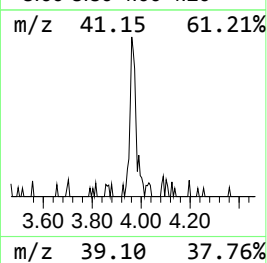
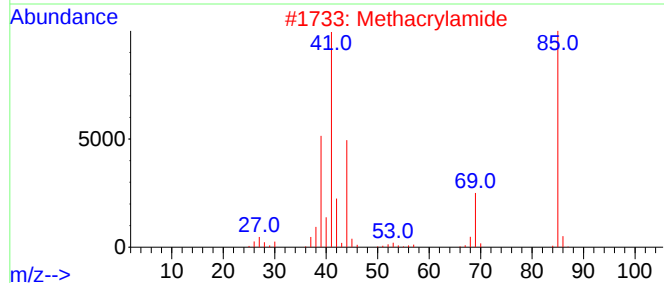
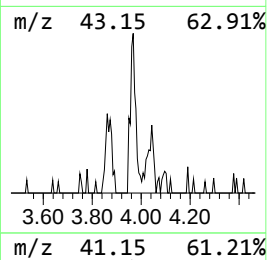
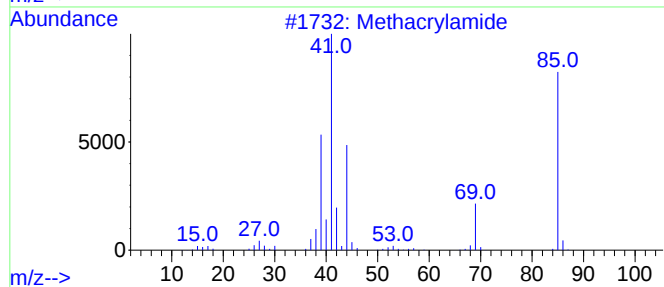
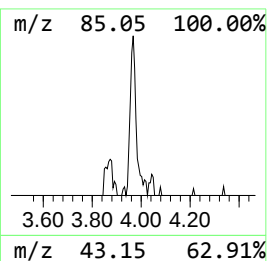
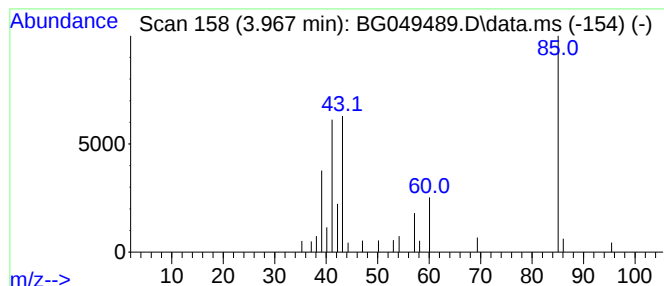
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Methacrylamide Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.967	2.87 ng/ul	24884	1,4-Dichlorobenzene-d4	8.050

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methacrylamide	85	C4H7NO	000079-39-0	50
2			Methacrylamide	85	C4H7NO	000079-39-0	50
3			Methacrylamide	85	C4H7NO	000079-39-0	49
4			2H-Pyran, 2-(4-chlorobutoxy)tetr...	192	C9H17ClO2	041302-05-0	38
5			Hexane, 2-bromo-	164	C6H13Br	003377-86-4	38



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ClientSampleId :
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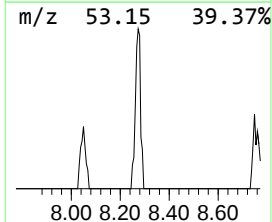
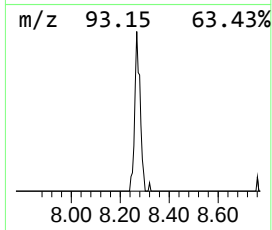
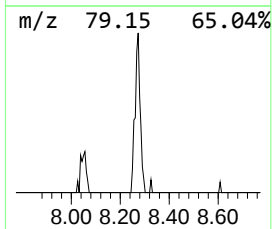
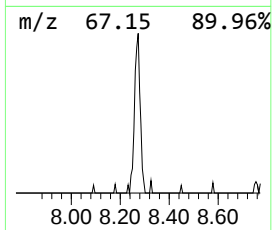
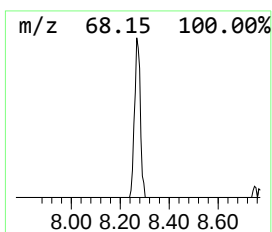
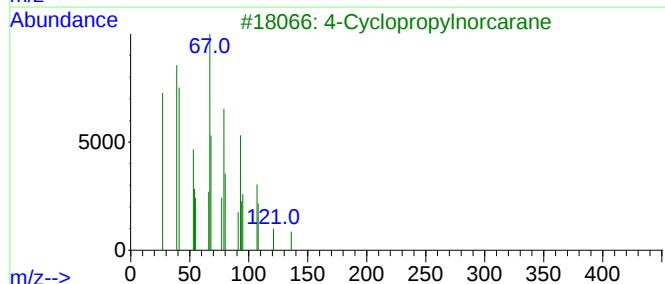
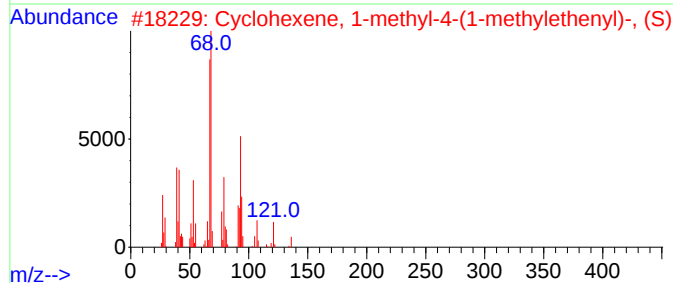
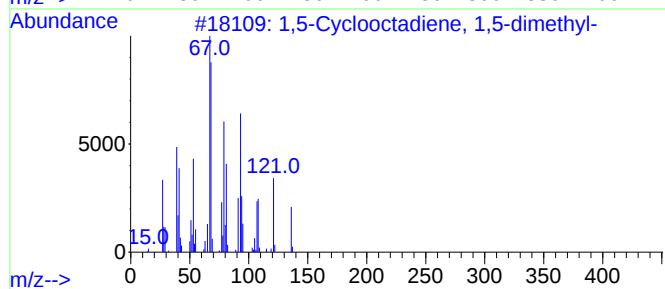
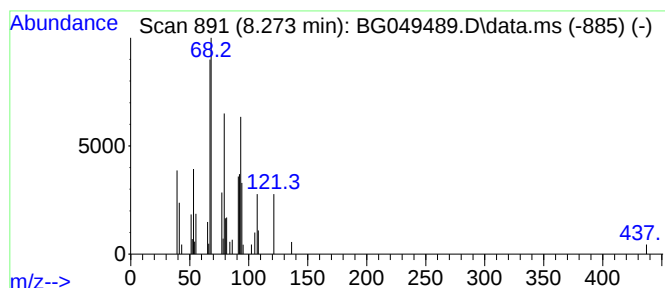
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1,5-Cyclooctadiene, 1,5-dim... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.273	5.41 ng/ul	46889	1,4-Dichlorobenzene-d4	8.050

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,5-Cyclooctadiene, 1,5-dimethyl-	136	C10H16	003760-14-3	93
2			Cyclohexene, 1-methyl-4-(1-methy...	136	C10H16	005989-54-8	83
3			4-Cyclopropylnorcarane	136	C10H16	1000222-73-9	72
4			1,5-Cyclooctadiene, 3-methyl-	122	C9H14	056564-88-6	72
5			1,7-Octadiene, 3-methylene-	122	C9H14	068695-13-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049489.D
Acq On : 26 Jul 2021 22:10
Operator : CG/JU
Sample : M3082-19
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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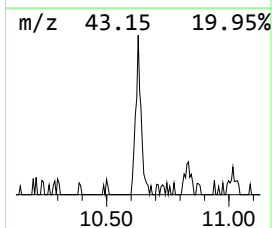
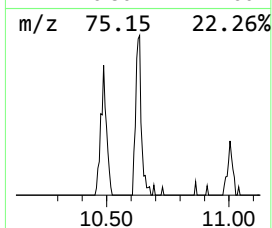
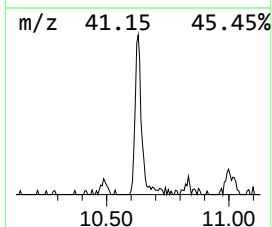
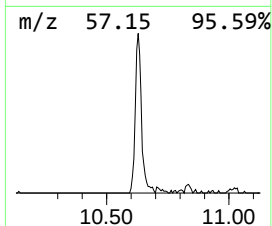
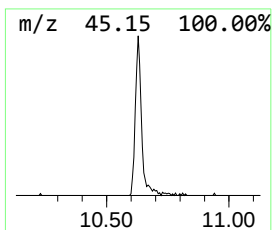
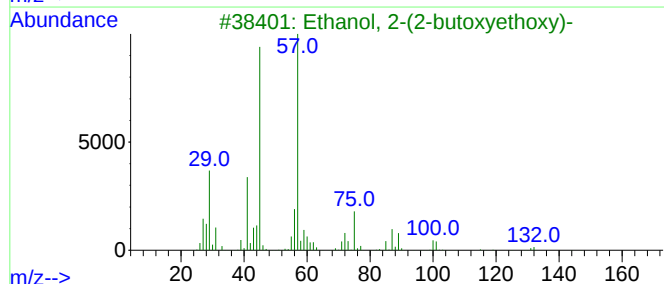
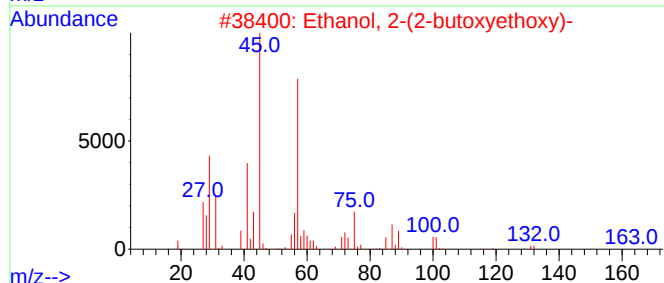
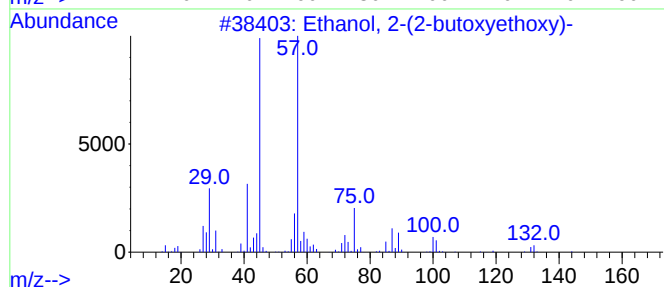
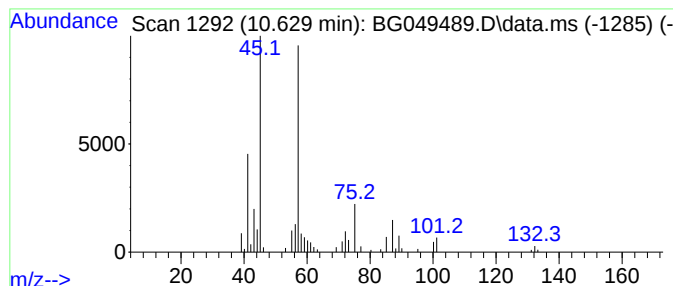
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.629	10.38 ng/ul	122108	Naphthalene-d8	10.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	86
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	74
4			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
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Operator : CG/JU
Sample : M3082-19
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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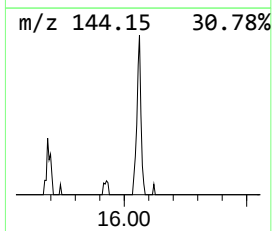
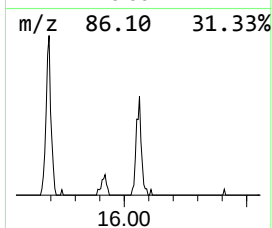
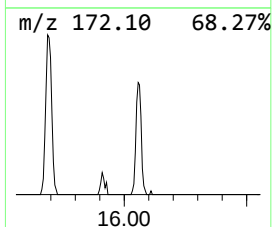
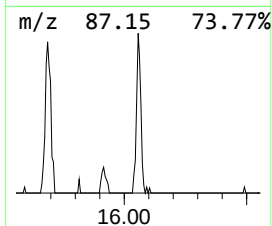
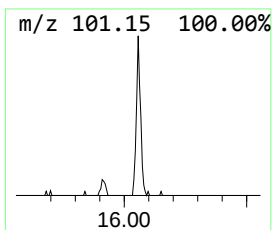
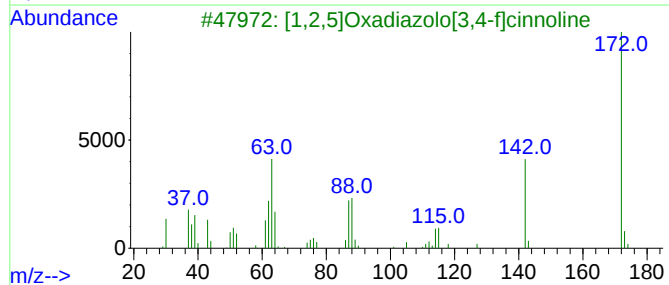
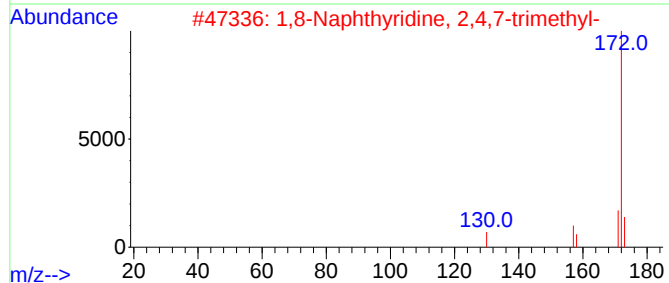
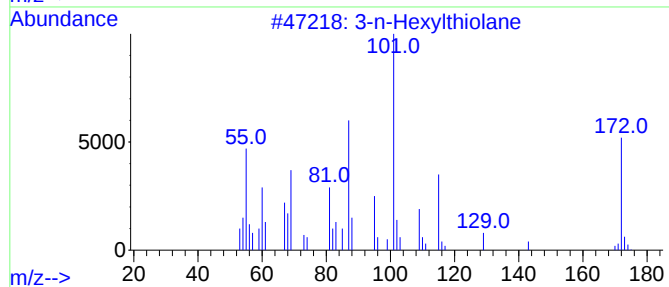
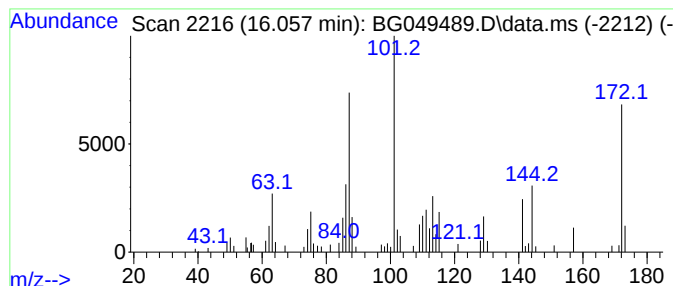
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.057	3.57 ng/ul	66364	Acenaphthene-d10	14.700

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-n-Hexylthiolane	172	C10H20S	001613-50-9	47
2			1,8-Naphthyridine, 2,4,7-trimethyl-	172	C11H12N2	014757-44-9	37
3			[1,2,5]Oxadiazolo[3,4-f]cinnoline	172	C8H4N4O	217491-04-8	15
4			Isopropyl isothiocyanate	101	C4H7NS	002253-73-8	14
5			trans-3-Methyl-2-n-propylthiophane	144	C8H16S	118068-75-0	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049489.D
Acq On : 26 Jul 2021 22:10
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Sample : M3082-19
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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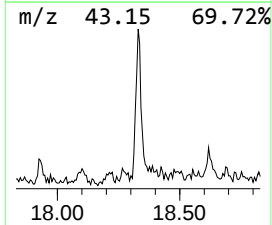
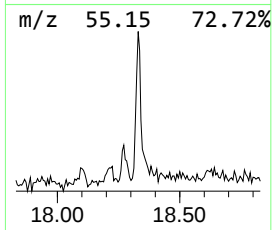
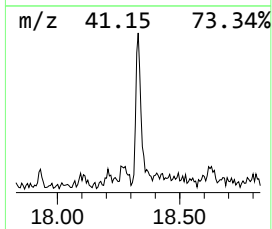
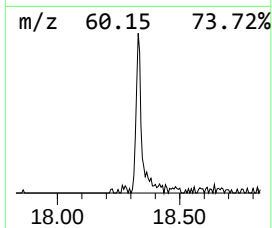
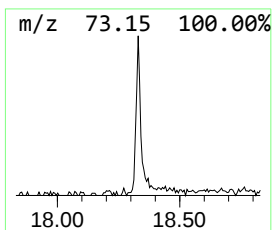
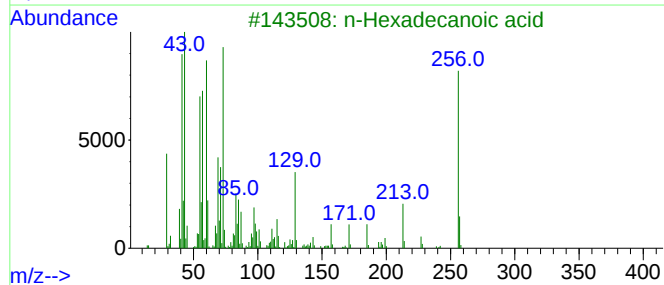
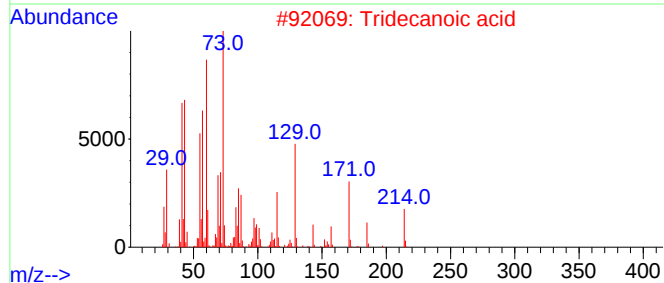
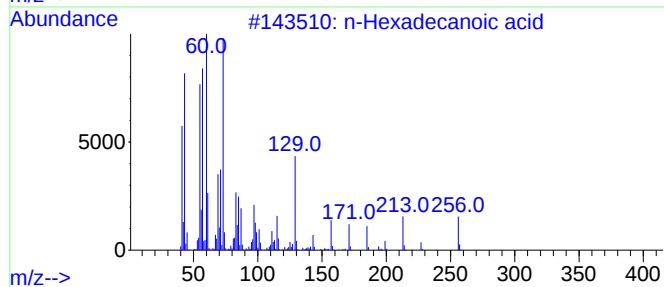
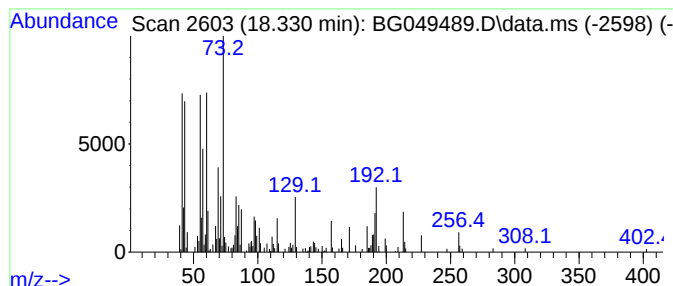
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.330	7.66 ng/ul	163894	Phenanthrene-d10	17.449

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			Tridecanoic acid	214	C13H26O2	000638-53-9	97
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
4			Tridecanoic acid	214	C13H26O2	000638-53-9	95
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049489.D
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Sample : M3082-19
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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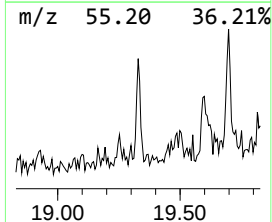
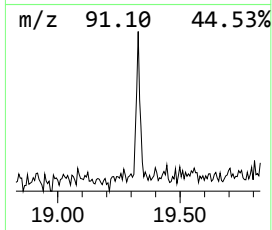
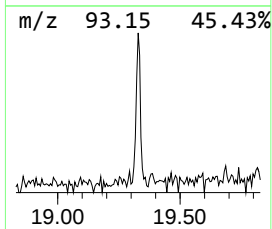
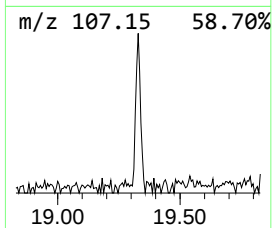
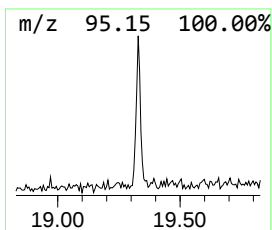
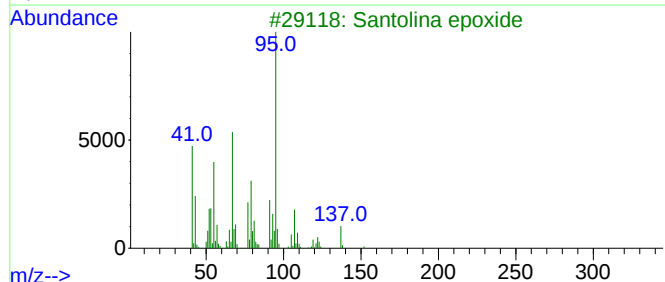
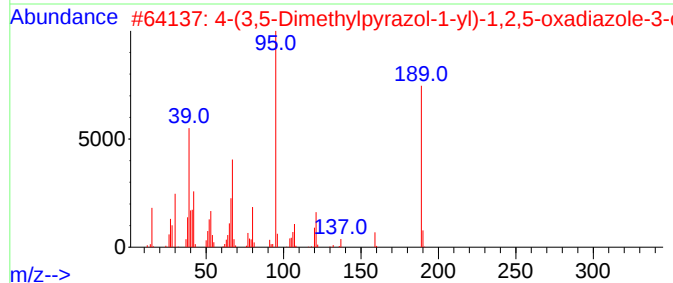
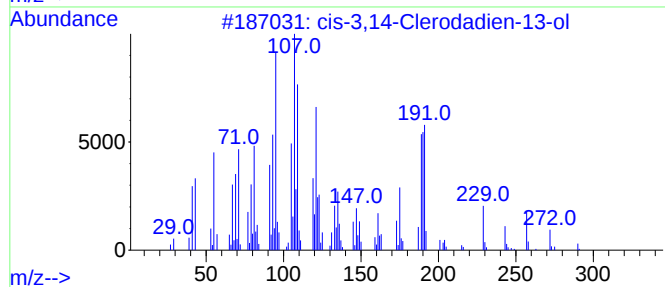
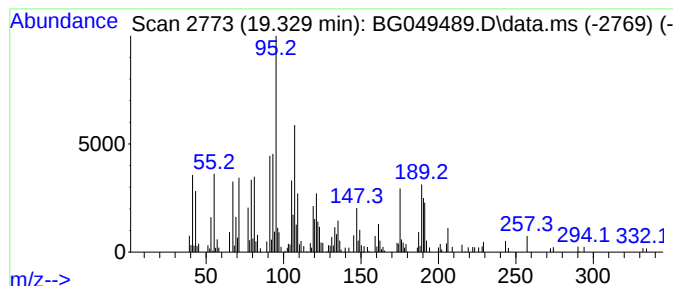
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 cis-3,14-Clerodadien-13-ol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.329	5.98 ng/ul	128002	Phenanthrene-d10	17.449

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-3,14-Clerodadien-13-ol	290	C20H34O	374925-73-2	53
2			4-(3,5-Dimethylpyrazol-1-yl)-1,2...	189	C8H7N5O	1000436-50-1	27
3			Santolina epoxide	152	C10H16O	060485-45-2	27
4			4,7-Methanobenzofuran, 2,2'-oxyb...	374	C24H38O3	087248-50-8	22
5			1,5-Cyclooctadiene, 1-t-butyl-	164	C12H20	1000152-18-6	20



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
Data File : BG049489.D
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Sample : M3082-19
Misc :
ALS Vial : 19 Sample Multiplier: 1

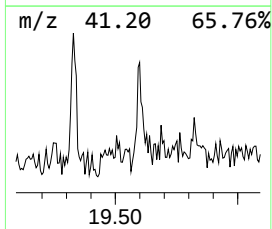
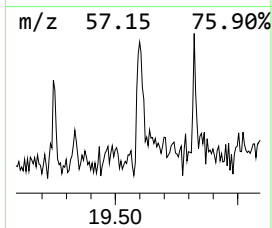
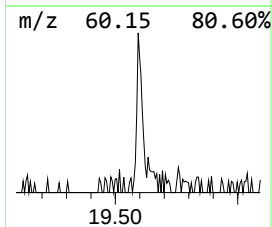
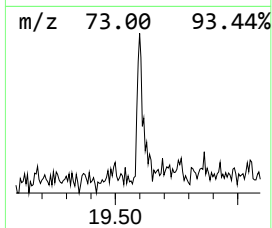
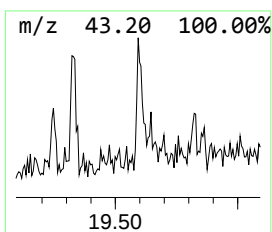
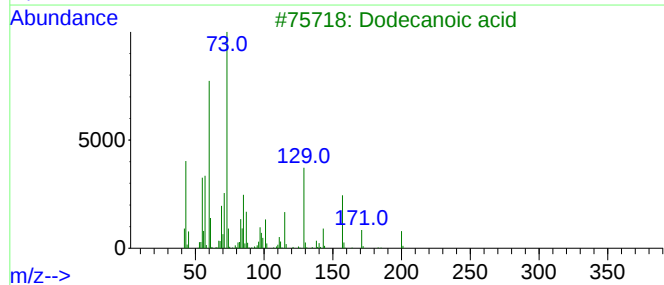
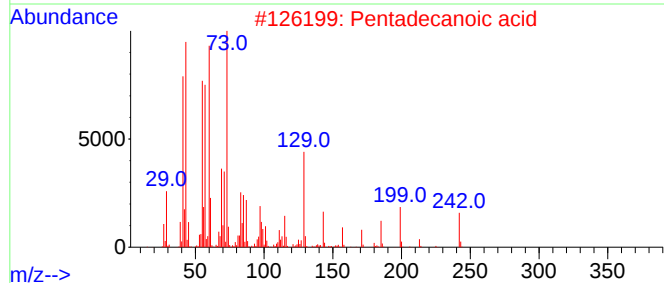
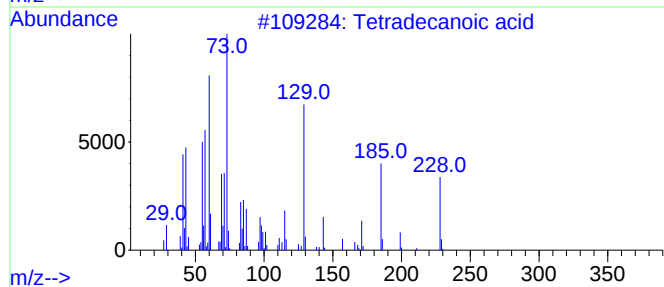
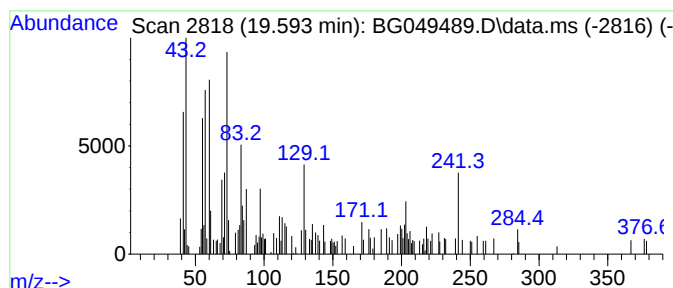
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Tetradecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.593	2.35 ng/ul	50037	Chrysene-d12	21.725	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecanoic acid	228	C14H28O2	000544-63-8	60
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	53
3	Dodecanoic acid	200	C12H24O2	000143-07-7	49
4	n-Decanoic acid	172	C10H20O2	000334-48-5	49
5	Dodecanoic acid	200	C12H24O2	000143-07-7	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
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Misc :
ALS Vial : 19 Sample Multiplier: 1

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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: C...	3.080	22.5	ng/ul	194873	1	8.050	173318	20.0
Butane, 2-metho...	3.162	26.4	ng/ul	229197	1	8.050	173318	20.0
Methacrylamide	3.967	2.9	ng/ul	24884	1	8.050	173318	20.0
1,5-Cyclooctadi...	8.273	5.4	ng/ul	46889	1	8.050	173318	20.0
Ethanol, 2-(2-b...	10.629	10.4	ng/ul	122108	2	10.869	235383	20.0
unknown-01	16.057	3.6	ng/ul	66364	3	14.700	371379	20.0
n-Hexadecanoic ...	18.330	7.7	ng/ul	163894	4	17.449	427834	20.0
cis-3,14-Clerod...	19.329	6.0	ng/ul	128002	4	17.449	427834	20.0
Tetradecanoic acid	19.593	2.4	ng/ul	50037	5	21.725	425476	20.0