

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012422\
 Data File : BG052126.D
 Acq On : 24 Jan 2022 14:58
 Operator : CG/JU
 Sample : N1199-16
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0Q10

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-BG010622.M
 Title : SVOA CALIBRATION

Signal : TIC: BG052126.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.546	4	10	25	rVB	179482	312478	32.19%	2.718%
2	4.004	84	88	99	rVV2	21308	37937	3.91%	0.330%
3	7.388	654	664	673	rBV2	46461	88239	9.09%	0.767%
4	7.546	685	691	701	rBV	105161	181614	18.71%	1.579%
5	7.769	722	729	736	rVB	126301	220659	22.73%	1.919%
6	8.245	803	810	817	rVV	98358	170619	17.58%	1.484%
7	8.950	923	930	940	rBV3	144363	300435	30.95%	2.613%
8	9.038	940	945	950	rBV3	5739	9575	0.99%	0.083%
9	9.185	965	970	975	rVB3	3383	5958	0.61%	0.052%
10	9.303	983	990	996	rVB3	3094	5419	0.56%	0.047%
11	9.414	1002	1009	1021	rVB	160725	287828	29.65%	2.503%
12	9.526	1021	1028	1032	rBV7	3027	6081	0.63%	0.053%
13	9.614	1037	1043	1051	rVB8	5454	9999	1.03%	0.087%
14	9.690	1051	1056	1066	rBV7	11477	23945	2.47%	0.208%
15	9.778	1066	1071	1077	rBV3	23026	41063	4.23%	0.357%
16	9.831	1077	1080	1085	rVB2	10022	13981	1.44%	0.122%
17	9.919	1090	1095	1099	rVV6	4273	8051	0.83%	0.070%
18	10.072	1117	1121	1122	rVV3	5099	7999	0.82%	0.070%
19	10.148	1126	1134	1142	rVB	138904	257470	26.52%	2.239%
20	10.260	1148	1153	1156	rBV4	5828	10279	1.06%	0.089%
21	10.360	1163	1170	1177	rVB4	24904	46726	4.81%	0.406%
22	10.436	1177	1183	1184	rBV2	8537	10027	1.03%	0.087%
23	10.695	1220	1227	1239	rBV	214738	386075	39.77%	3.358%
24	10.830	1246	1250	1258	rBV5	9417	22315	2.30%	0.194%
25	11.030	1279	1284	1287	rBV4	6679	9562	0.99%	0.083%
26	11.083	1287	1293	1298	rVV	143072	259574	26.74%	2.257%
27	11.130	1298	1301	1307	rVV6	24509	43619	4.49%	0.379%
28	11.206	1307	1314	1327	rVB	109596	238313	24.55%	2.073%
29	11.347	1336	1338	1345	rBV7	2714	4662	0.48%	0.041%
30	11.435	1349	1353	1359	rVB5	8423	16055	1.65%	0.140%
31	11.500	1359	1364	1368	rBV5	3672	6514	0.67%	0.057%
32	11.558	1368	1374	1375	rBV4	3137	4584	0.47%	0.040%
33	11.670	1385	1393	1395	rVV6	6317	13463	1.39%	0.117%
34	11.799	1413	1415	1420	rVB6	4628	6837	0.70%	0.059%
35	11.852	1420	1424	1427	rBV3	5267	6828	0.70%	0.059%
36	11.934	1433	1438	1442	rVV5	5012	8804	0.91%	0.077%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
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37	11.987	1442	1447	1451	rVB6	4523	7449	0.77%	0.065%
38	12.040	1453	1456	1464	rVB3	14993	25356	2.61%	0.221%
39	12.116	1464	1469	1472	rBV5	6399	10717	1.10%	0.093%
40	12.193	1477	1482	1485	rVB5	3191	5224	0.54%	0.045%
41	12.310	1497	1502	1504	rBV6	3092	5090	0.52%	0.044%
42	12.357	1508	1510	1516	rVB7	5196	7736	0.80%	0.067%
43	12.557	1534	1544	1548	rBV5	8875	20818	2.14%	0.181%
44	12.627	1548	1556	1562	rBV10	10344	30771	3.17%	0.268%
45	12.698	1564	1568	1574	rVB7	7925	15485	1.60%	0.135%
46	12.815	1586	1588	1593	rVB5	3362	4525	0.47%	0.039%
47	12.880	1593	1599	1601	rBV4	7747	11924	1.23%	0.104%
48	12.933	1604	1608	1614	rVV7	16061	29647	3.05%	0.258%
49	13.021	1618	1623	1627	rVV7	12867	27821	2.87%	0.242%
50	13.056	1627	1629	1634	rVB5	12182	17509	1.80%	0.152%
51	13.139	1640	1643	1649	rBV7	6663	12418	1.28%	0.108%
52	13.197	1649	1653	1655	rVB3	4492	6319	0.65%	0.055%
53	13.291	1665	1669	1675	rBV6	8313	16629	1.71%	0.145%
54	13.421	1687	1691	1698	rBV9	6785	13251	1.37%	0.115%
55	13.532	1706	1710	1713	rBV3	7980	12192	1.26%	0.106%
56	13.703	1734	1739	1740	rBV4	6320	8123	0.84%	0.071%
57	13.773	1745	1751	1756	rVV4	19115	43154	4.45%	0.375%
58	13.885	1766	1770	1778	rVV10	10465	22349	2.30%	0.194%
59	14.055	1796	1799	1801	rBV4	4692	5842	0.60%	0.051%
60	14.278	1830	1837	1844	rVB	388299	595233	61.32%	5.177%
61	14.413	1856	1860	1866	rVB9	5118	11791	1.21%	0.103%
62	14.525	1877	1879	1882	rVV4	8617	10641	1.10%	0.093%
63	14.584	1883	1889	1899	rVB	453987	725544	74.74%	6.310%
64	14.766	1915	1920	1921	rBV5	5428	7659	0.79%	0.067%
65	14.889	1932	1941	1947	rBV	234256	390002	40.17%	3.392%
66	15.077	1967	1973	1983	rVV2	37827	76913	7.92%	0.669%
67	15.336	2015	2017	2022	rVB6	3800	4937	0.51%	0.043%
68	15.424	2028	2032	2036	rVV6	12470	19441	2.00%	0.169%
69	15.482	2039	2042	2043	rVV3	8049	8353	0.86%	0.073%
70	15.512	2043	2047	2054	rVB9	16116	26630	2.74%	0.232%
71	15.618	2061	2065	2067	rBV5	5116	5969	0.61%	0.052%
72	15.876	2103	2109	2121	rVB	553892	857940	88.38%	7.461%
73	15.988	2122	2128	2135	rBV3	179421	308318	31.76%	2.681%
74	16.228	2167	2169	2173	rVB5	8677	10487	1.08%	0.091%
75	16.557	2221	2225	2228	rBV6	6339	11069	1.14%	0.096%
76	16.751	2254	2258	2259	rBV4	7165	8486	0.87%	0.074%

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77	16.910	2283	2285	2286	rBV2	6030	5316	0.55%	0.046%
78	16.980	2295	2297	2300	rBV4	5375	6258	0.64%	0.054%
79	17.098	2316	2317	2321	rVB3	9400	8628	0.89%	0.075%
80	17.286	2344	2349	2357	rBV	222865	355938	36.67%	3.095%
81	17.638	2403	2409	2417	rBV	300831	453223	46.69%	3.942%
82	17.738	2419	2426	2431	rVV2	574532	873003	89.93%	7.592%
83	18.285	2515	2519	2522	rVB	31776	39269	4.05%	0.342%
84	18.508	2554	2557	2560	rBV5	16242	19925	2.05%	0.173%
85	19.442	2712	2716	2721	rBV2	54945	74989	7.72%	0.652%
86	19.876	2787	2790	2794	rBV2	34674	46455	4.79%	0.404%
87	20.017	2809	2814	2822	rVB	671538	957751	98.66%	8.329%
88	21.956	3139	3144	3151	rVB	254473	441944	45.53%	3.843%
89	23.530	3408	3412	3421	rVB7	28691	63571	6.55%	0.553%
90	24.212	3525	3528	3529	rBV3	7171	8250	0.85%	0.072%
91	24.412	3560	3562	3563	rBV2	5937	4708	0.48%	0.041%
92	24.441	3563	3567	3573	rVB7	15768	29442	3.03%	0.256%
93	24.823	3631	3632	3636	rBV4	5051	4767	0.49%	0.041%
94	25.210	3687	3698	3712	rVB2	323785	970760	100.00%	8.442%
95	25.445	3729	3738	3751	rVB2	155183	499816	51.49%	4.347%
96	26.726	3950	3956	3966	rVB8	14450	43972	4.53%	0.382%
97	26.920	3987	3989	3993	rBV5	4504	7030	0.72%	0.061%
98	28.224	4204	4211	4220	rVV5	12745	42453	4.37%	0.369%
99	30.022	4512	4517	4519	rBV5	6877	12760	1.31%	0.111%
100	30.961	4674	4677	4679	rBV4	4209	5112	0.53%	0.044%

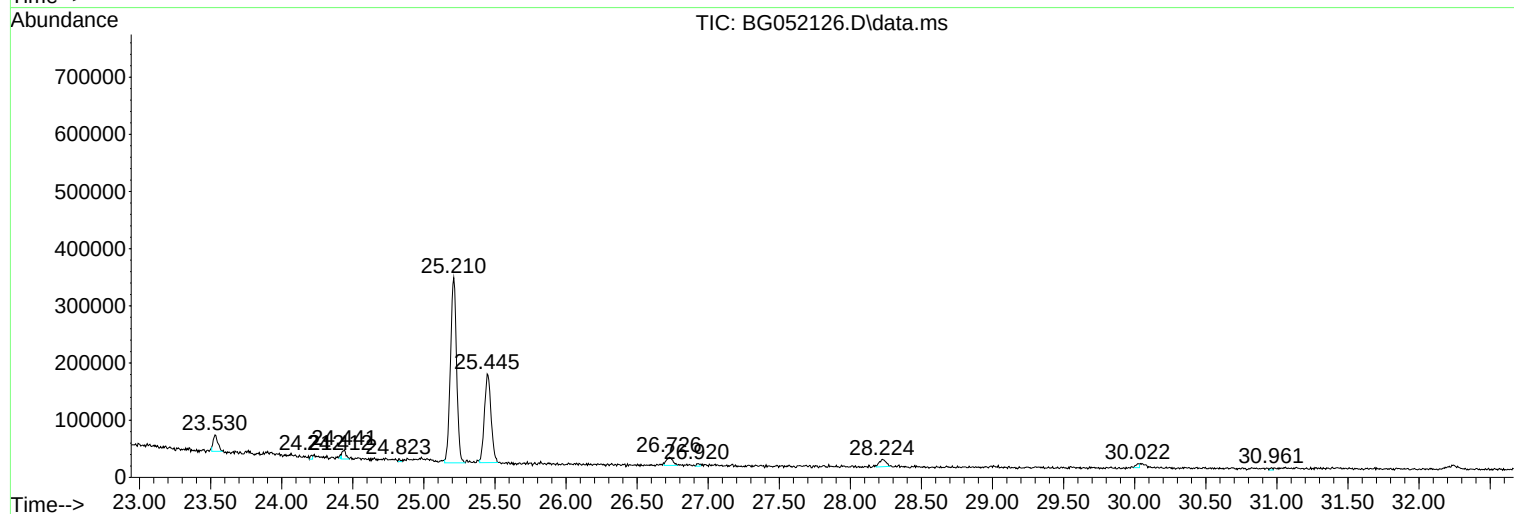
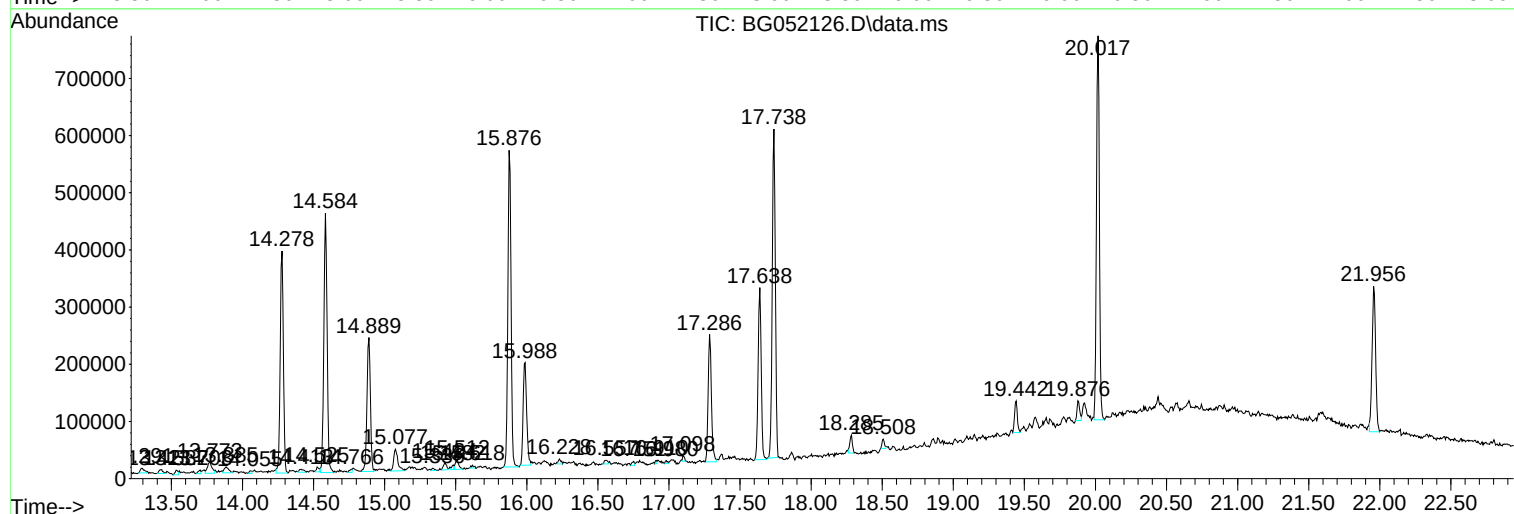
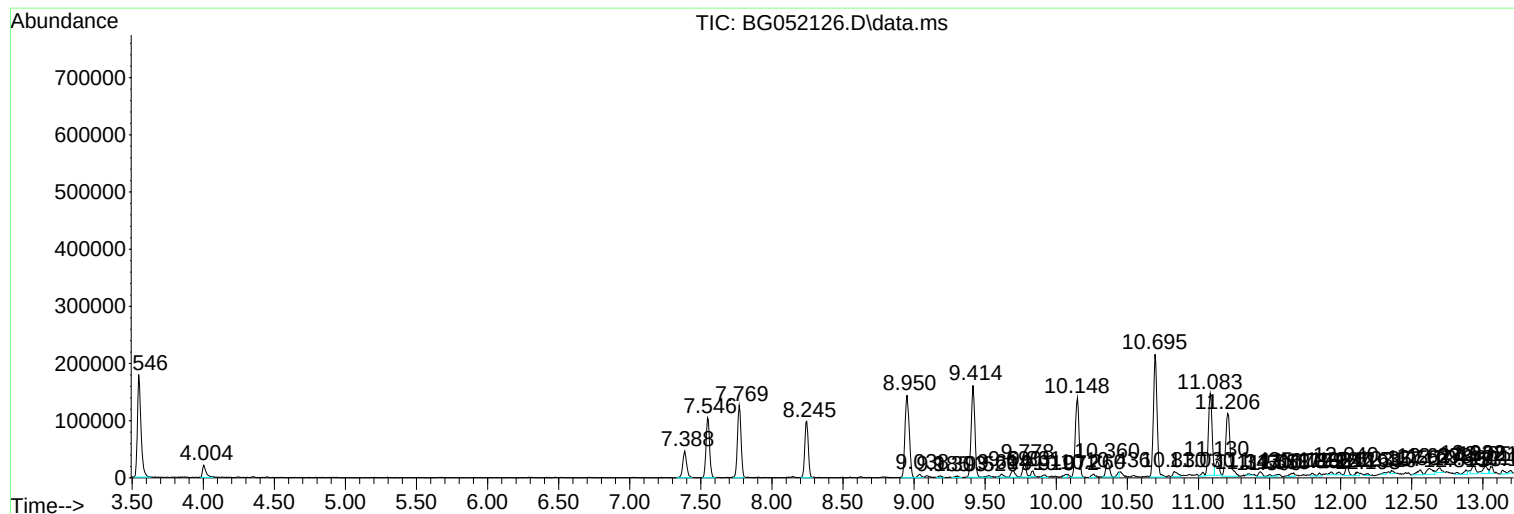
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
Quant Title : SVOA CALIBRATION

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



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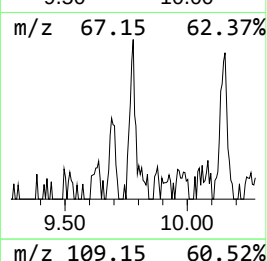
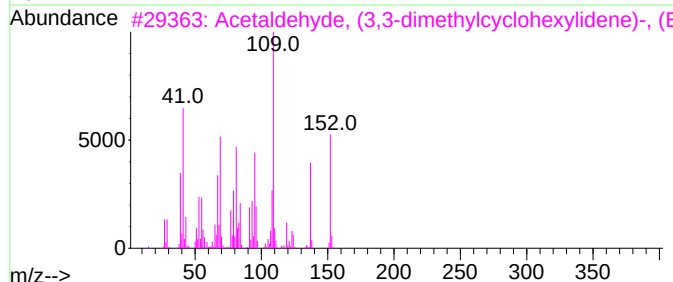
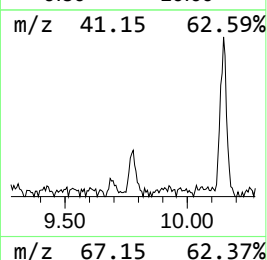
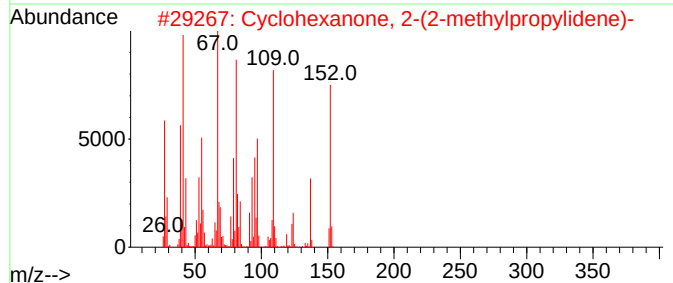
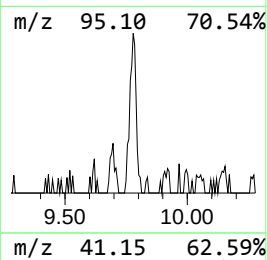
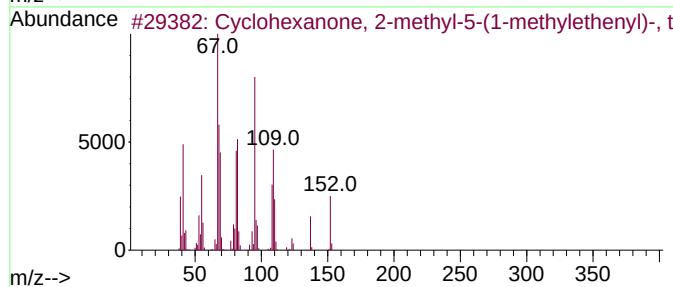
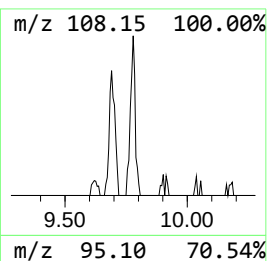
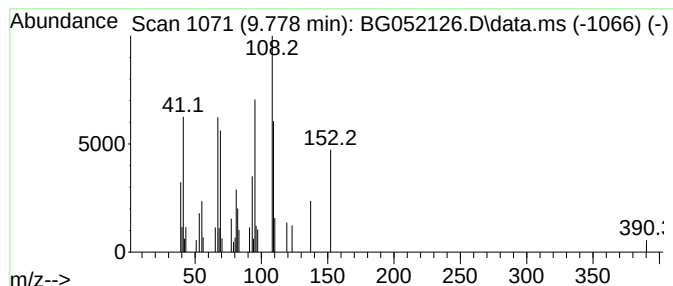
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Cyclohexanone, 2-methyl-5-(... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.778	3.16 ng/ul	41063	Naphthalene-d8	11.083

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanone, 2-methyl-5-(1-met...	152	C10H16O	005948-04-9	62
2			Cyclohexanone, 2-(2-methylpropyl...	152	C10H16O	043108-69-6	47
3			Acetaldehyde, (3,3-dimethylcyclo...	152	C10H16O	026532-25-2	46
4			1-Methyldecahydronaphthalene	152	C11H20	002958-75-0	43
5			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	43



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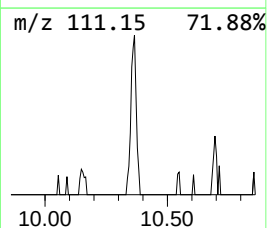
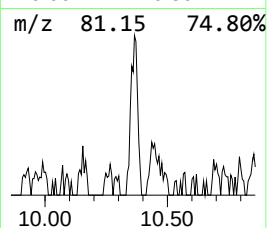
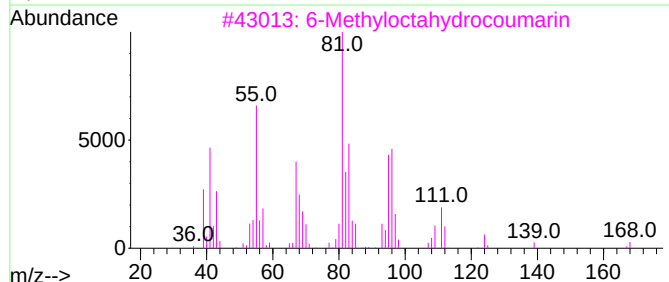
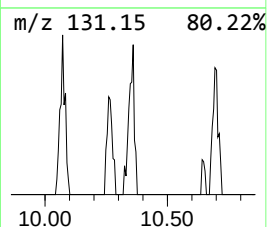
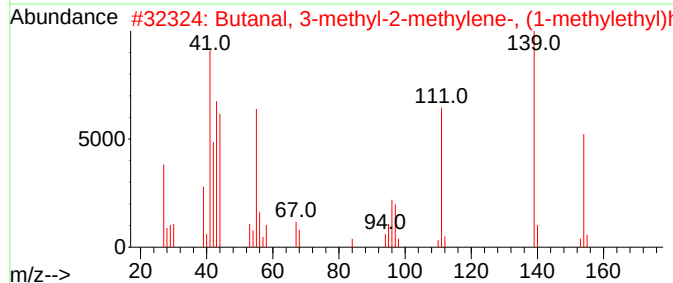
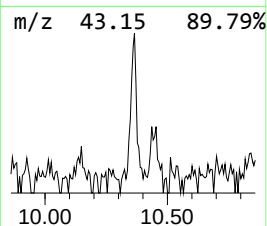
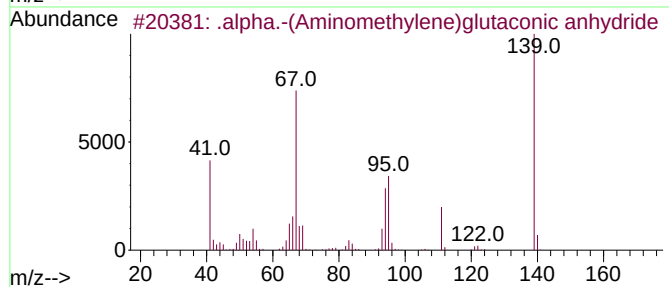
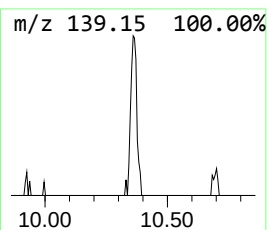
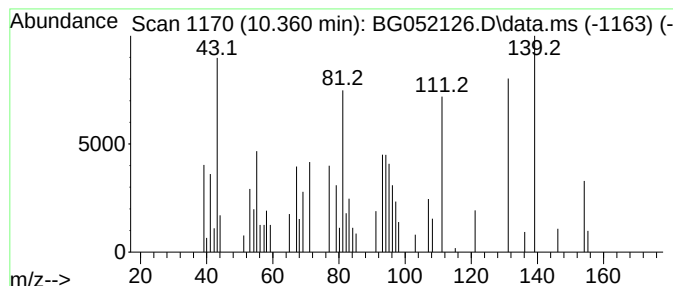
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.360	3.60 ng/ul	46726	Naphthalene-d8	11.083

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.		.alpha.-(Aminomethylene)glutacon...	139	C6H5NO3	067598-07-6	30
2	Butanal, 3-methyl-2-methylene-, ...	154	C9H18N2			033063-81-9	27
3	6-Methyloctahydrocoumarin	168	C10H16O2			080648-29-9	22
4	2(3H)-Benzofuranone, hexahydro-7...	154	C9H14O2			039163-24-1	22
5	2(3H)-Benzofuranone, hexahydro-7...	154	C9H14O2			039163-24-1	22



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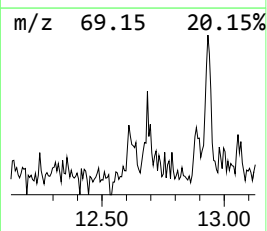
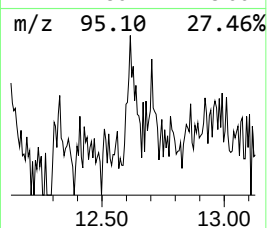
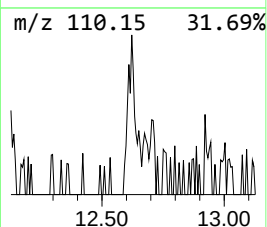
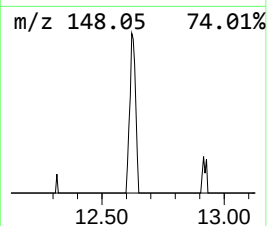
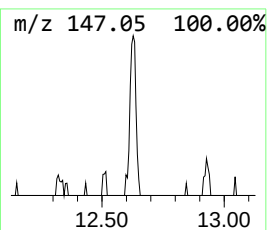
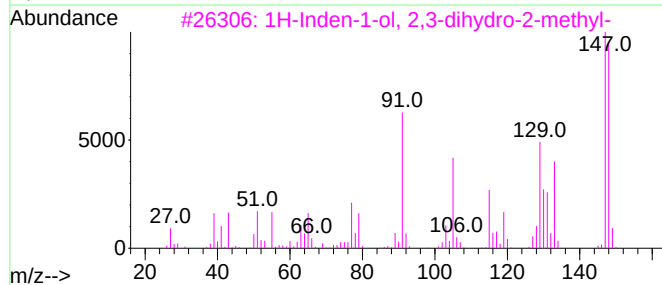
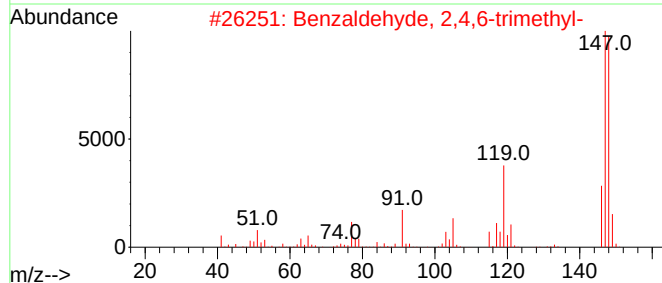
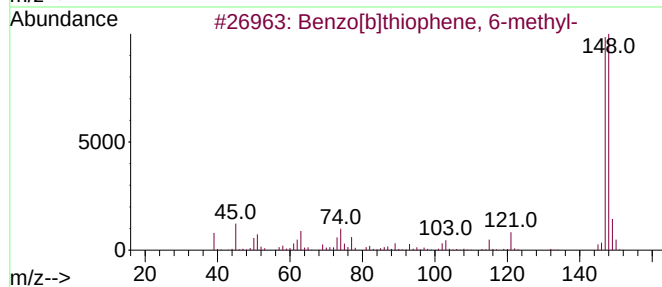
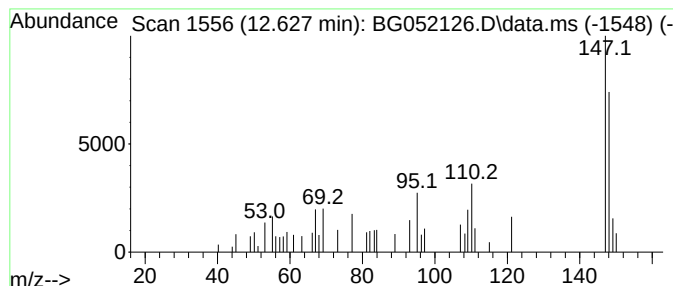
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzo[b]thiophene, 6-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.628	2.37 ng/ul	30771	Naphthalene-d8	11.083

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]thiophene, 6-methyl-	148	C9H8S	016587-47-6	52
2			Benzaldehyde, 2,4,6-trimethyl-	148	C10H12O	000487-68-3	50
3			1H-Inden-1-ol, 2,3-dihydro-2-met...	148	C10H12O	017496-18-3	47
4			Benzaldehyde, 2,4,6-trimethyl-	148	C10H12O	000487-68-3	47
5			Benzene, 1-ethynyl-2-(methylthio)-	148	C9H8S	078905-08-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012422\
Data File : BG052126.D
Acq On : 24 Jan 2022 14:58
Operator : CG/JU
Sample : N1199-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0Q10

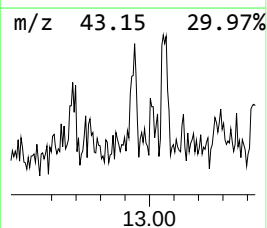
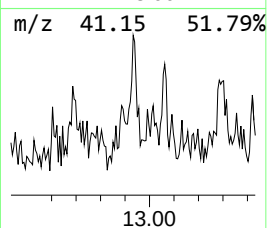
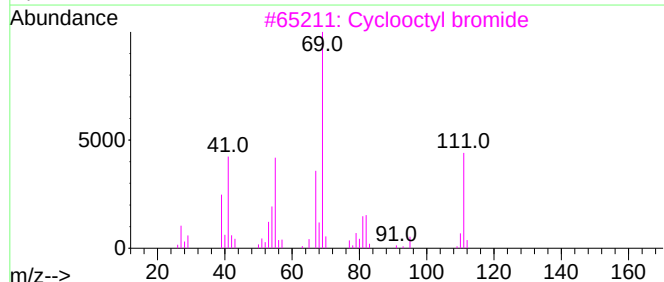
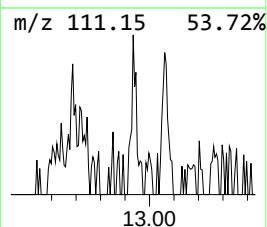
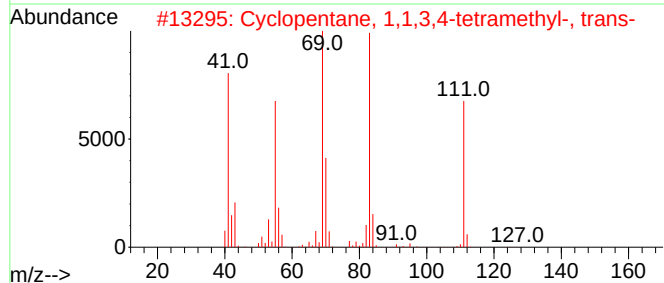
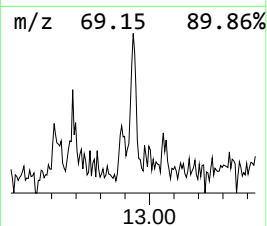
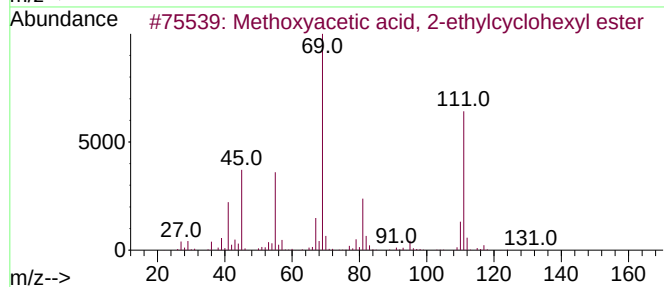
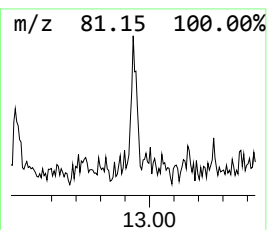
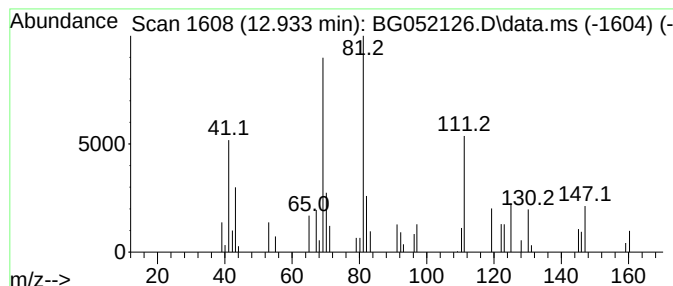
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-02 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.933	2.28 ng/ul	29647	Naphthalene-d8	11.083

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methoxyacetic acid, 2-ethylcyclo...	200	C11H20O3	1000282-69-7	12
2			Cyclopentane, 1,1,3,4-tetramethy...	126	C9H18	020309-77-7	12
3			Cyclooctyl bromide	190	C8H15Br	001556-09-8	12
4			Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	12
5			Bicyclo[2.2.2]octan-1-amine	125	C8H15N	001193-42-6	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012422\
Data File : BG052126.D
Acq On : 24 Jan 2022 14:58
Operator : CG/JU
Sample : N1199-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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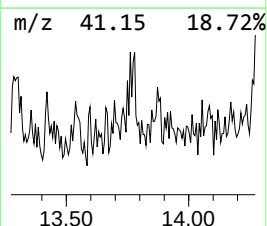
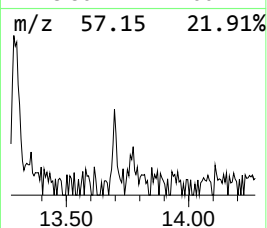
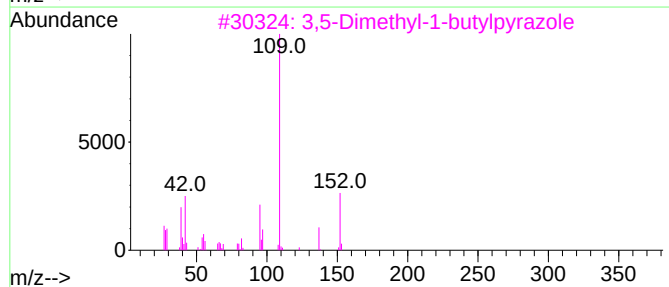
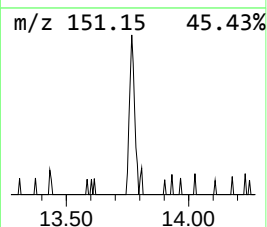
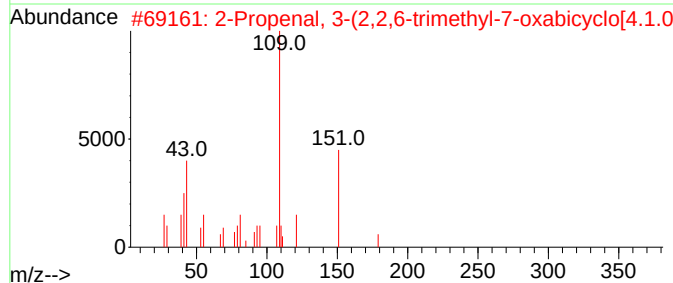
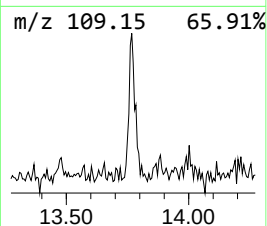
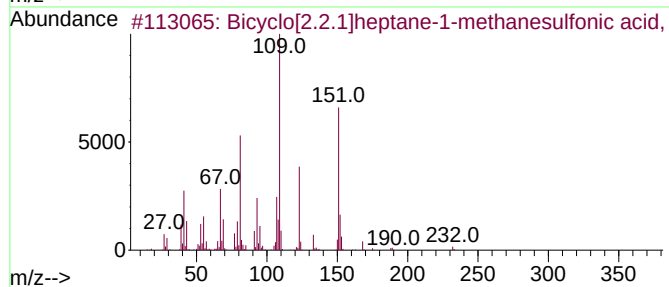
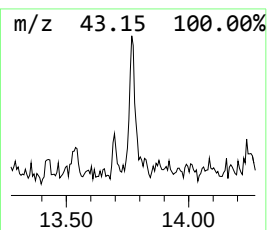
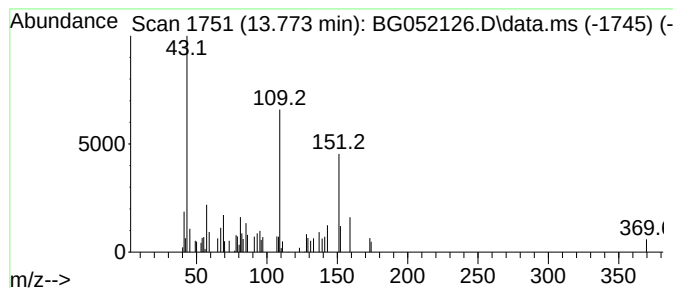
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-03 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.773	2.21 ng/ul	43154	Acenaphthene-d10	14.889

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[2.2.1]heptane-1-methanes...	232	C10H16O4S	005872-08-2	40
2			2-Propenal, 3-(2,2,6-trimethyl-7...	194	C12H18O2	055759-91-6	38
3			3,5-Dimethyl-1-butylpyrazole	152	C9H16N2	002655-37-0	35
4			Benzene, 1-fluoro-2-(2-methoxyet...	152	C9H9FO	054533-36-7	35
5			trans-2-Caren-4-ol	152	C10H16O	004017-82-7	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012422\
Data File : BG052126.D
Acq On : 24 Jan 2022 14:58
Operator : CG/JU
Sample : N1199-16
Misc :
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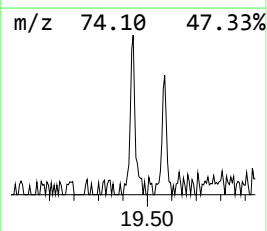
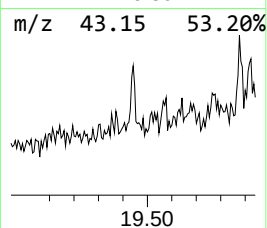
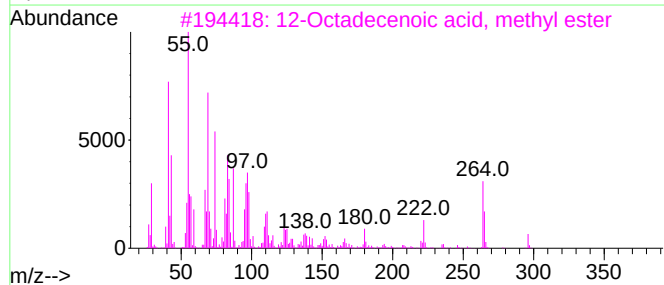
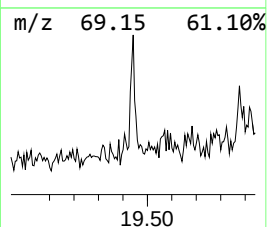
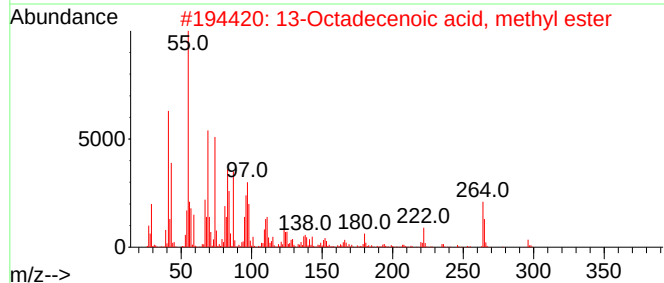
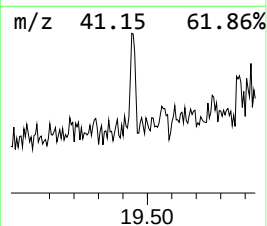
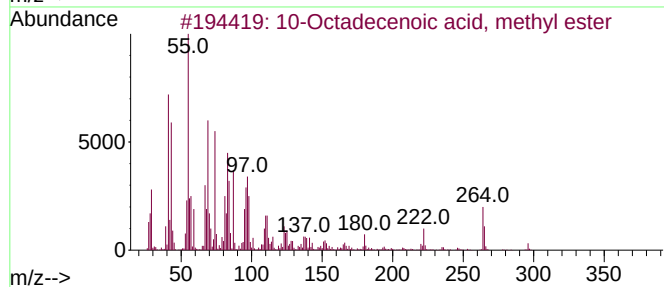
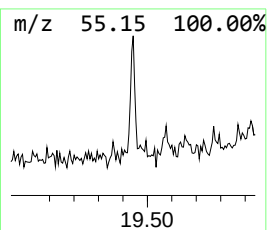
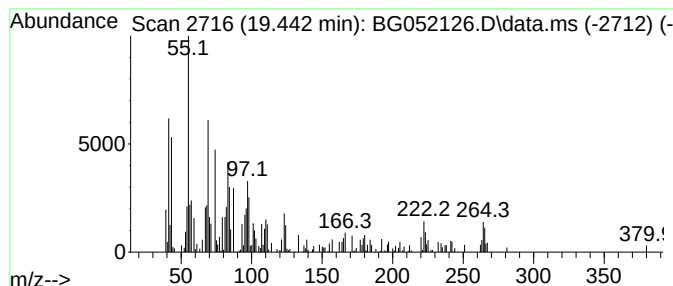
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 10-Octadecenoic acid, methyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.442	3.31 ng/ul	74989	Phenanthrene-d10	17.638	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	95
2	13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	94
3	12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	93
4	14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	91
5	11-Octadecenoic acid, methyl est...	296	C19H36O2	001937-63-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012422\
Data File : BG052126.D
Acq On : 24 Jan 2022 14:58
Operator : CG/JU
Sample : N1199-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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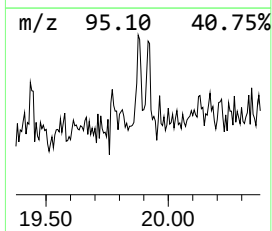
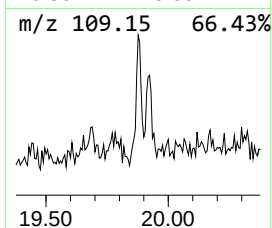
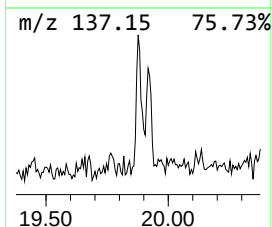
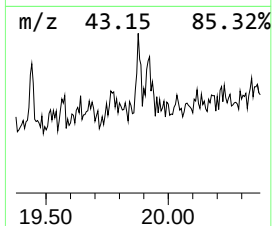
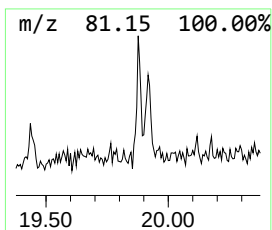
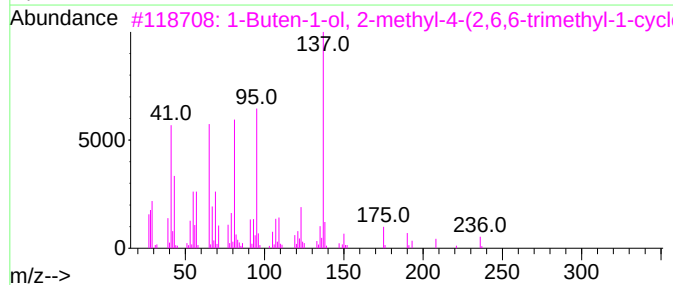
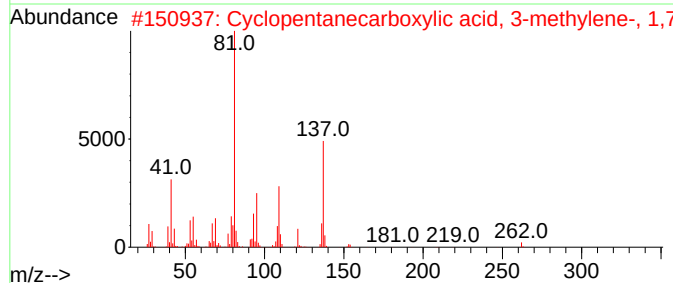
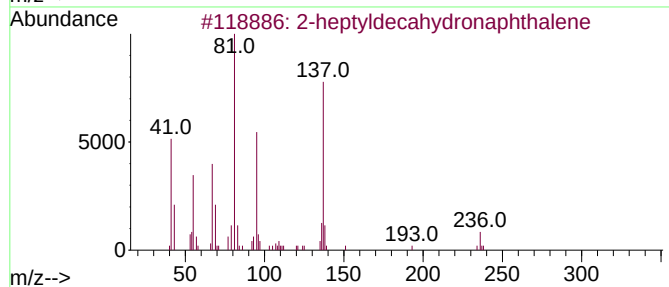
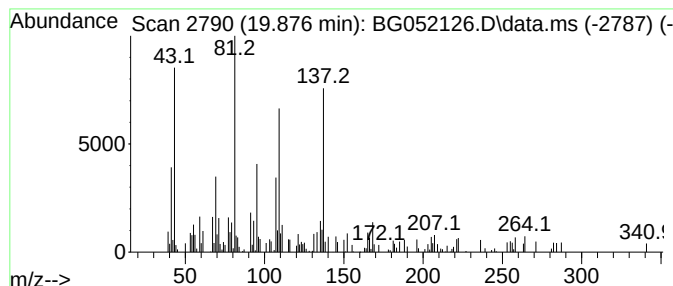
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 2-heptyldecahydronaphthalene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.877	2.10 ng/ul	46455	Chrysene-d12	21.956

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-heptyldecahydronaphthalene	236	C17H32	1000441-03-2	58
2			Cyclopentanecarboxylic acid, 3-m...	262	C17H26O2	074793-59-2	43
3			1-Buten-1-ol, 2-methyl-4-(2,6,6-...	236	C15H24O2	021730-91-6	38
4			Artemiseole	152	C10H16O	060485-46-3	35
5			3,4-Dimethylthiophenol, S-pentaf...	284	C11H9F5OS	1000470-52-3	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012422\
Data File : BG052126.D
Acq On : 24 Jan 2022 14:58
Operator : CG/JU
Sample : N1199-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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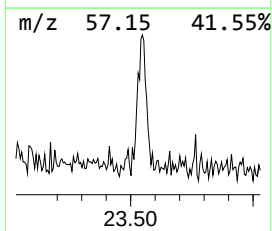
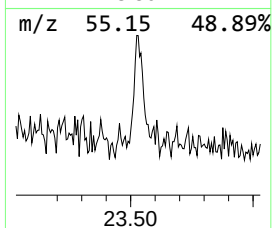
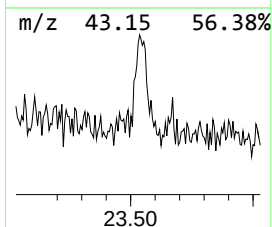
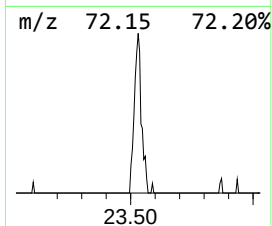
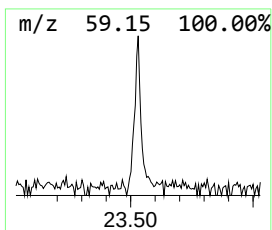
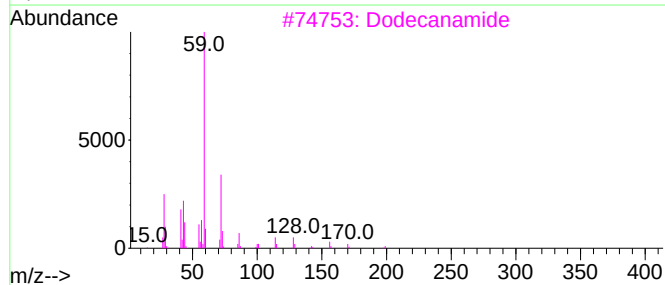
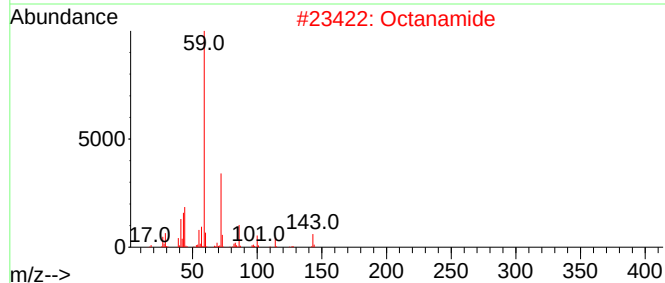
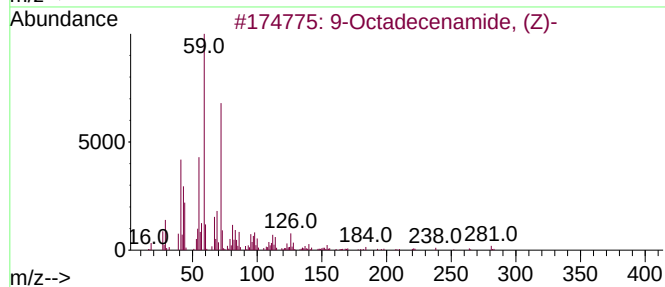
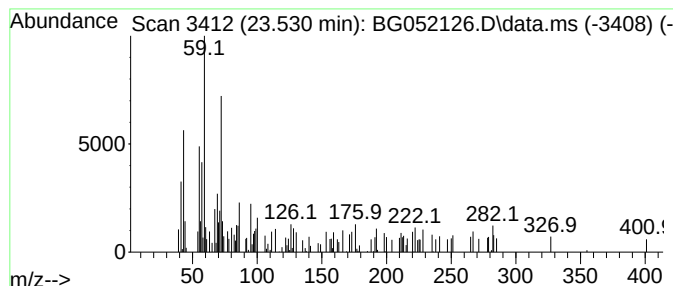
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 9-Octadecenamide, (Z)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.530	2.88 ng/ul	63571	Chrysene-d12	21.956

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	64
2	Octanamide			143	C8H17NO	000629-01-6	53
3	Dodecanamide			199	C12H25NO	001120-16-7	50
4	Pentadecanamide, 15-bromo-			319	C15H30BrNO	1000163-86-1	50
5	Palmitoleamide			253	C16H31NO	106010-22-4	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012422\
Data File : BG052126.D
Acq On : 24 Jan 2022 14:58
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Sample : N1199-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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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
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unknown-01	10.360	3.6	ng/ul	46726	2	11.083	259574	20.0
Benzo[b]thiophe...	12.628	2.4	ng/ul	30771	2	11.083	259574	20.0
unknown-02	12.933	2.3	ng/ul	29647	2	11.083	259574	20.0
unknown-03	13.773	2.2	ng/ul	43154	3	14.889	390002	20.0
10-Octadecenoic...	19.442	3.3	ng/ul	74989	4	17.638	453223	20.0
2-heptyldecahyd...	19.877	2.1	ng/ul	46455	5	21.956	441944	20.0
9-Octadecenamid...	23.530	2.9	ng/ul	63571	5	21.956	441944	20.0