

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012924\
 Data File : BG060467.D
 Acq On : 29 Jan 2024 18:20
 Operator : MA/JU
 Sample : P1239-12
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0FC2

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-BG012624.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG060467.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.364	44	47	49	rBV2	26594	33580	0.43%	0.044%
2	3.399	49	53	60	rVB	58864	81173	1.04%	0.107%
3	3.634	89	93	98	rVB4	27147	46436	0.59%	0.061%
4	3.775	112	117	127	rBV	193364	343183	4.38%	0.453%
5	4.110	168	174	184	rBV	126112	201672	2.57%	0.266%
6	4.222	184	193	200	rVB4	21748	43304	0.55%	0.057%
7	4.710	271	276	283	rBV7	15253	30661	0.39%	0.040%
8	5.127	342	347	355	rVB3	35495	63506	0.81%	0.084%
9	5.362	381	387	391	rBV	54763	76280	0.97%	0.101%
10	5.432	395	399	402	rBV2	20366	29394	0.37%	0.039%
11	5.585	416	425	429	rBV2	15185	32284	0.41%	0.043%
12	5.902	475	479	481	rBV3	11816	19583	0.25%	0.026%
13	5.938	481	485	489	rVV4	14208	22139	0.28%	0.029%
14	6.002	492	496	503	rVV3	46559	76362	0.97%	0.101%
15	6.672	606	610	617	rVB6	14240	19881	0.25%	0.026%
16	7.013	663	668	671	rBV4	12905	19218	0.25%	0.025%
17	7.089	673	681	693	rBV2	154822	375297	4.79%	0.495%
18	7.183	693	697	702	rVB4	21657	35531	0.45%	0.047%
19	7.248	702	708	717	rBV	647631	1113493	14.20%	1.469%
20	7.453	737	743	752	rVV	544115	970046	12.37%	1.279%
21	7.571	757	763	771	rVB10	15669	42354	0.54%	0.056%
22	7.923	811	823	835	rBV	661233	1185722	15.12%	1.564%
23	8.029	840	841	848	rVB5	13107	18807	0.24%	0.025%
24	8.282	878	884	890	rBV5	11969	28766	0.37%	0.038%
25	8.388	894	902	908	rVB9	16003	36078	0.46%	0.048%
26	8.458	910	914	917	rBV6	11590	19421	0.25%	0.026%
27	8.570	928	933	934	rBV5	15784	20265	0.26%	0.027%
28	8.629	937	943	953	rVV	516463	908949	11.59%	1.199%
29	8.805	968	973	978	rBV6	20245	36244	0.46%	0.048%
30	8.952	990	998	1000	rBV	71866	134777	1.72%	0.178%
31	9.022	1007	1010	1014	rVB4	29651	43175	0.55%	0.057%
32	9.087	1014	1021	1031	rBV	748027	1357038	17.31%	1.790%
33	9.193	1035	1039	1044	rVB2	50441	78174	1.00%	0.103%
34	9.357	1063	1067	1077	rVV3	109200	305703	3.90%	0.403%
35	9.492	1086	1090	1096	rVB2	40336	64917	0.83%	0.086%
36	9.657	1109	1118	1127	rBV7	52733	123488	1.57%	0.163%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG012624.MA.M
 Title : SVOA CALIBRATION

37	9.745	1127	1133	1138	rBV	86939	160075	2.04%	0.211%
38	9.810	1138	1144	1152	rVB	567545	1040055	13.26%	1.372%
39	10.003	1173	1177	1186	rVB2	45030	110571	1.41%	0.146%
40	10.133	1194	1199	1206	rBV3	116014	206421	2.63%	0.272%
41	10.197	1206	1210	1215	rBV7	29692	52056	0.66%	0.069%
42	10.268	1217	1222	1229	rBV4	82500	160831	2.05%	0.212%
43	10.344	1229	1235	1245	rBV	1034590	1924632	24.54%	2.538%
44	10.509	1255	1263	1270	rBV5	65286	161668	2.06%	0.213%
45	10.579	1270	1275	1281	rVB2	55169	97367	1.24%	0.128%
46	10.650	1281	1287	1291	rBV9	25463	55766	0.71%	0.074%
47	10.726	1293	1300	1310	rBV	1059595	1877063	23.94%	2.476%
48	10.861	1315	1323	1337	rVB	665775	1365450	17.41%	1.801%
49	11.079	1357	1360	1366	rVB6	24786	42652	0.54%	0.056%
50	11.249	1385	1389	1395	rVB3	57369	95676	1.22%	0.126%
51	11.308	1395	1399	1403	rBV2	44344	76329	0.97%	0.101%
52	11.531	1434	1437	1441	rVV	26073	38625	0.49%	0.051%
53	11.584	1443	1446	1451	rVB6	30793	59387	0.76%	0.078%
54	11.719	1465	1469	1472	rBV4	40172	76095	0.97%	0.100%
55	11.831	1482	1488	1503	rVB7	72533	201019	2.56%	0.265%
56	11.972	1505	1512	1519	rBV	225294	440670	5.62%	0.581%
57	12.066	1524	1528	1537	rVB3	85358	168632	2.15%	0.222%
58	12.177	1542	1547	1554	rBV2	140576	267034	3.41%	0.352%
59	12.283	1560	1565	1567	rBV5	38241	58277	0.74%	0.077%
60	12.395	1580	1584	1587	rBV6	17639	24278	0.31%	0.032%
61	12.647	1623	1627	1633	rBV7	45466	85236	1.09%	0.112%
62	12.906	1666	1671	1676	rBV9	31518	66695	0.85%	0.088%
63	13.018	1687	1690	1691	rBV3	42887	44154	0.56%	0.058%
64	13.082	1697	1701	1707	rVB3	91054	148462	1.89%	0.196%
65	13.258	1723	1731	1738	rBV3	57527	198863	2.54%	0.262%
66	13.576	1781	1785	1788	rBV7	78252	129897	1.66%	0.171%
67	13.770	1816	1818	1825	rVB5	42719	55510	0.71%	0.073%
68	13.893	1834	1839	1843	rBV5	82363	156790	2.00%	0.207%
69	13.963	1843	1851	1863	rVB	2451510	3973994	50.68%	5.241%
70	14.169	1882	1886	1888	rBV4	28030	41489	0.53%	0.055%
71	14.257	1894	1901	1906	rBV	2938284	4599161	58.65%	6.066%
72	14.304	1906	1909	1914	rVB3	256897	360839	4.60%	0.476%
73	14.375	1919	1921	1928	rVB4	60430	114894	1.47%	0.152%
74	14.486	1936	1940	1943	rBV6	43107	69940	0.89%	0.092%
75	14.522	1943	1946	1947	rVV3	57225	62721	0.80%	0.083%
76	14.563	1947	1953	1960	rVB	1932011	3012979	38.42%	3.974%

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 Title : SVOA CALIBRATION

77	14.663	1967	1970	1977	rBV2	67199	123911	1.58%	0.163%
78	14.774	1986	1989	1997	rBV3	94811	191651	2.44%	0.253%
79	15.033	2031	2033	2037	rVB4	30573	36104	0.46%	0.048%
80	15.086	2037	2042	2047	rBV	249182	345278	4.40%	0.455%
81	15.215	2061	2064	2069	rVB3	117216	141532	1.80%	0.187%
82	15.297	2074	2078	2085	rVB	75884	127807	1.63%	0.169%
83	15.385	2091	2093	2096	rVB3	41936	39308	0.50%	0.052%
84	15.556	2116	2122	2132	rBV	4187545	6100763	77.80%	8.046%
85	15.673	2135	2142	2151	rVB	943830	1483672	18.92%	1.957%
86	17.312	2415	2421	2430	rVB	2550756	3752281	47.85%	4.949%
87	17.412	2432	2438	2446	rVV	4256438	6415594	81.81%	8.461%
88	17.671	2478	2482	2487	rBV	99275	178633	2.28%	0.236%
89	18.200	2569	2572	2579	rVB2	277699	395601	5.04%	0.522%
90	18.417	2606	2609	2616	rBV	118283	184108	2.35%	0.243%
91	19.704	2822	2828	2833	rBV	5443492	7729640	98.57%	10.194%
92	21.220	3082	3086	3092	rBV2	252853	391714	5.00%	0.517%
93	21.572	3141	3146	3152	rBV	2410383	3835593	48.91%	5.059%
94	23.023	3388	3393	3400	rVB5	147437	290156	3.70%	0.383%
95	23.817	3523	3528	3534	rVB2	112484	234523	2.99%	0.309%
96	24.522	3637	3648	3662	rBV2	2860685	7841874	100.00%	10.342%
97	24.739	3675	3685	3699	rVB	1637136	4490253	57.26%	5.922%
98	25.844	3866	3873	3883	rVB2	135843	351943	4.49%	0.464%
99	27.166	4089	4098	4108	rBV5	113772	392359	5.00%	0.517%
100	31.267	4784	4796	4810	rVB5	176705	829276	10.57%	1.094%

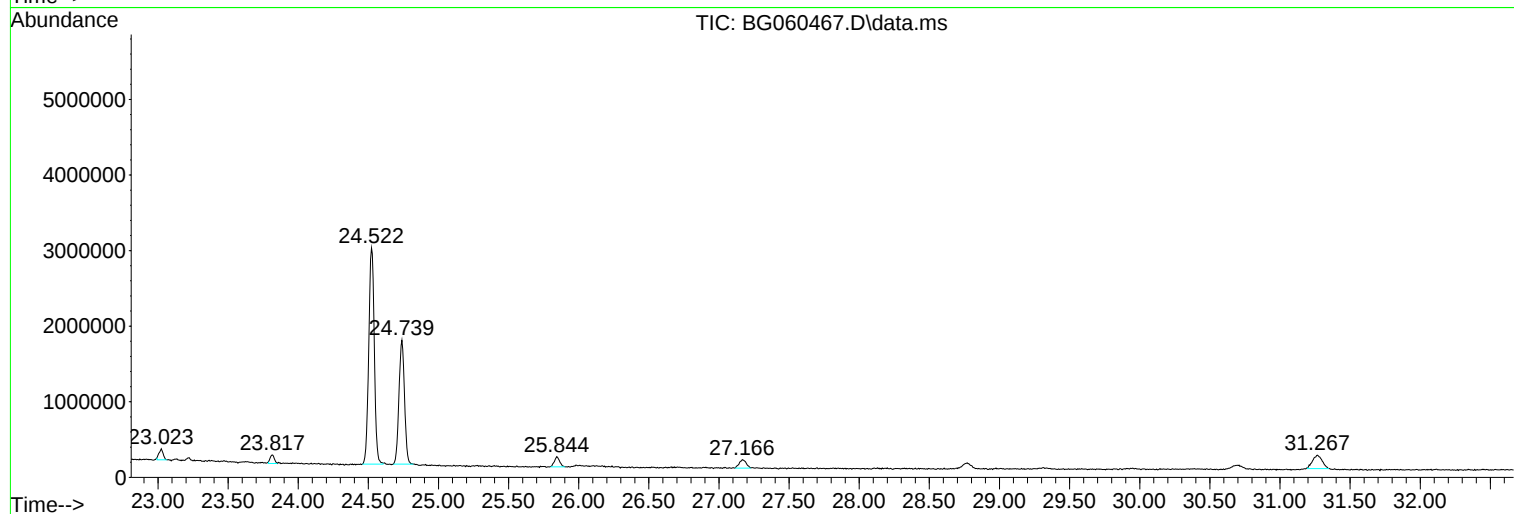
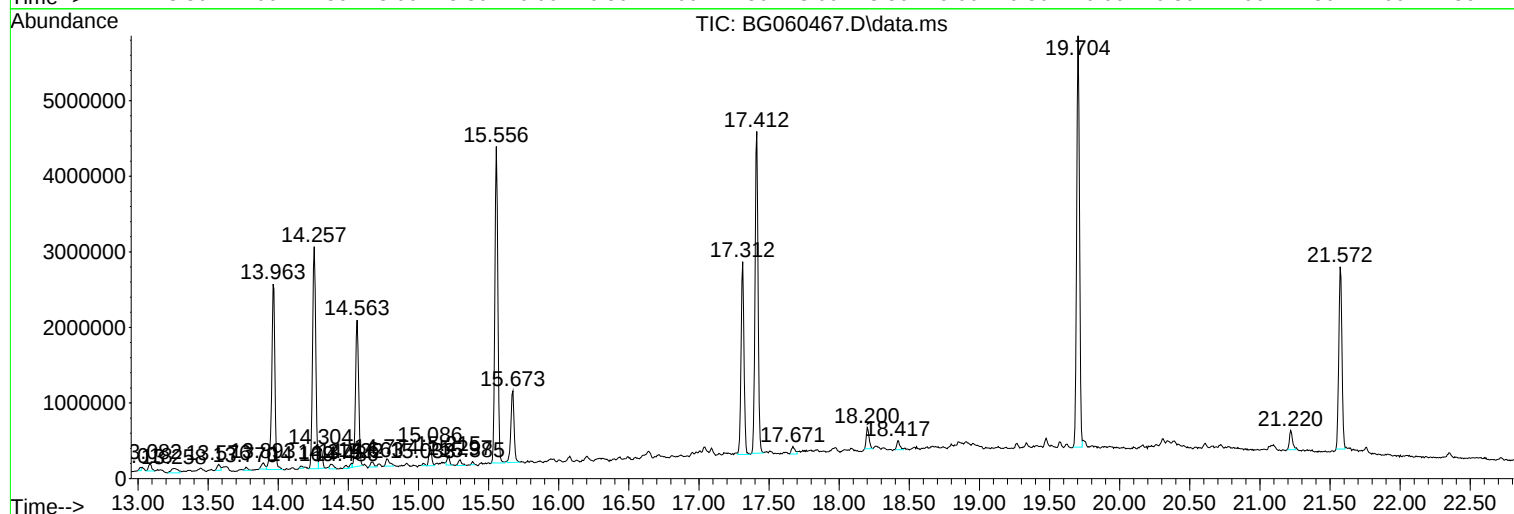
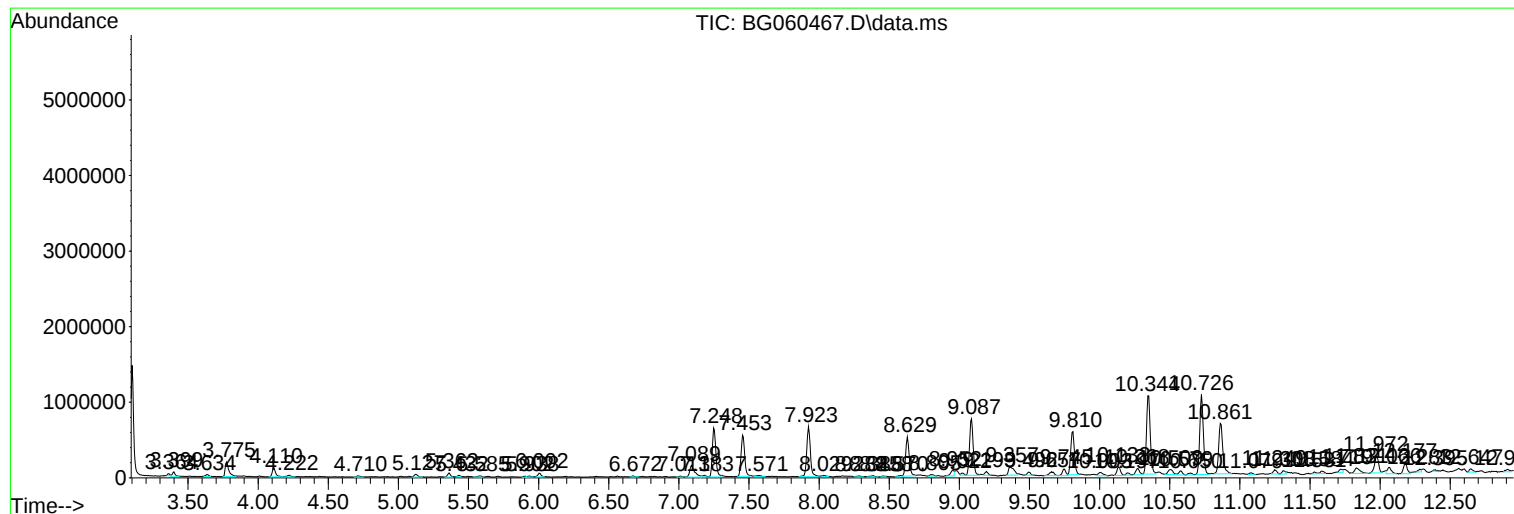
Sum of corrected areas: 75822728

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG012624.MA.M
Quant Title : SVOA CALIBRATION

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



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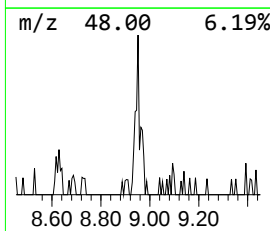
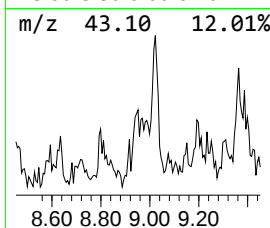
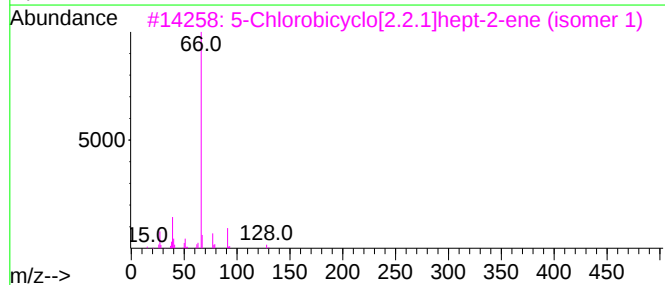
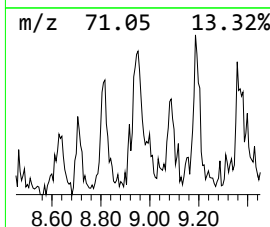
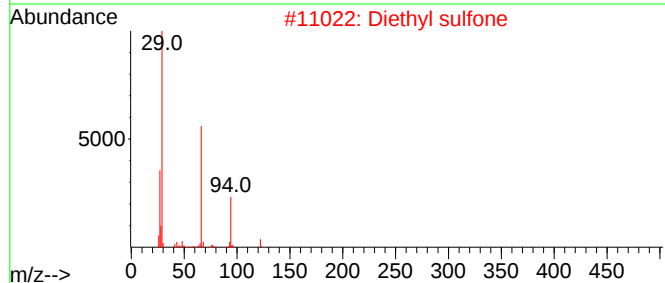
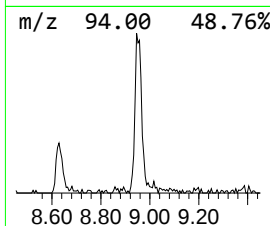
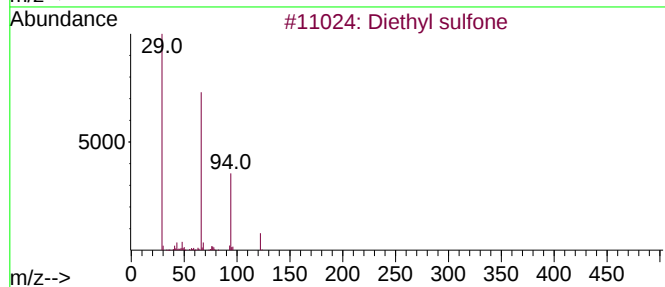
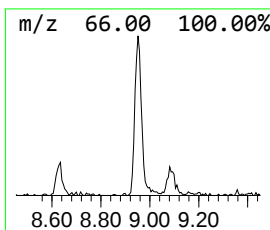
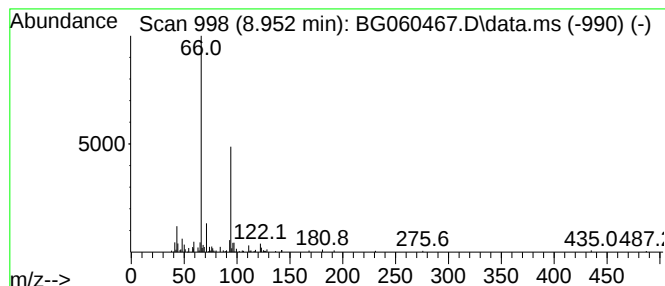
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG012624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Diethyl sulfone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.952	2.27 ng/ul	134777	1,4-Dichlorobenzene-d4	7.923

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diethyl sulfone	122	C4H10O2S	000597-35-3	78
2			Diethyl sulfone	122	C4H10O2S	000597-35-3	78
3			5-Chlorobicyclo[2.2.1]hept-2-ene...	128	C7H9Cl	1000436-27-6	9
4			Propanedinitrile	66	C3H2N2	000109-77-3	9
5			2-Amino-bicyclo[2.2.1]hept-5-ene...	134	C8H10N2	1000190-89-6	9



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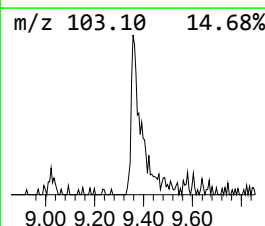
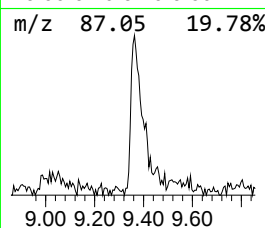
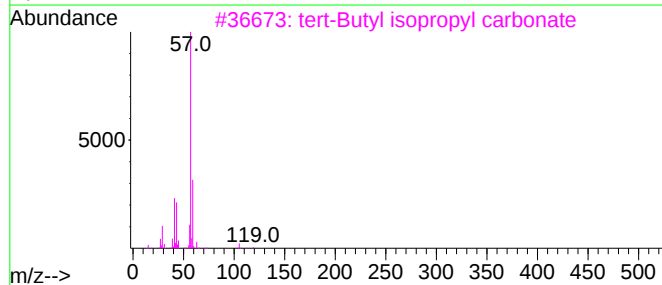
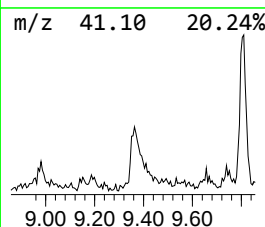
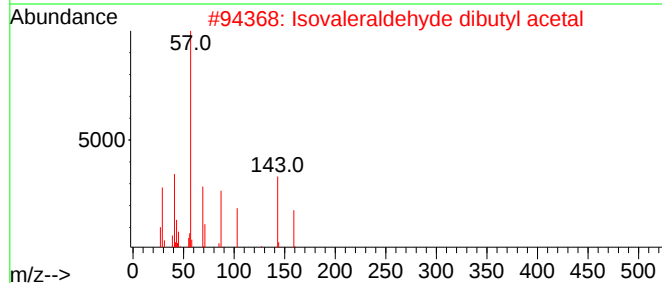
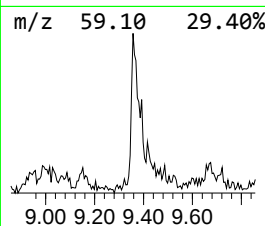
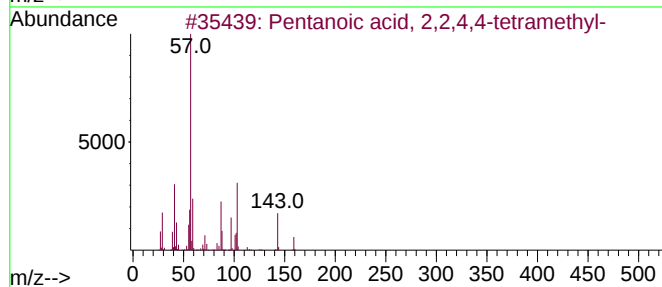
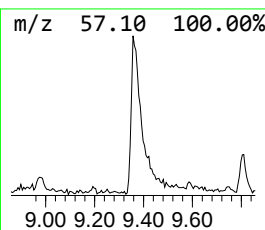
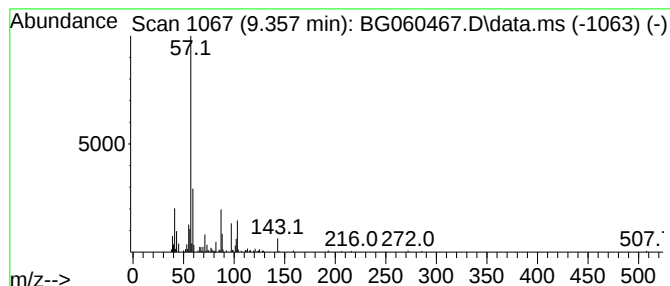
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG012624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.357	3.26 ng/ul	305703	Naphthalene-d8	10.726

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentanoic acid, 2,2,4,4-tetramet...	158	C9H18O2	003302-12-3	42
2			Isovaleraldehyde dibutyl acetal	216	C13H28O2	1000431-42-2	32
3			tert-Butyl isopropyl carbonate	160	C8H16O3	030221-21-7	16
4			Carbonic acid, monoamide, tert-b...	201	C11H23NO2	1000420-59-4	16
5			Propanoic acid, 2,2-dimethyl-, p...	144	C8H16O2	005129-35-1	12



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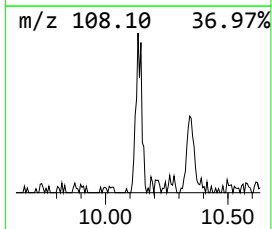
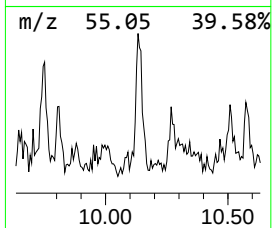
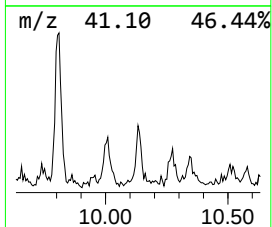
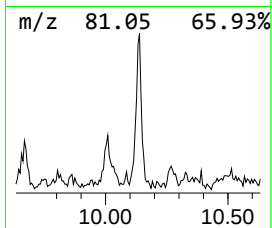
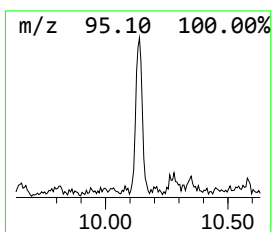
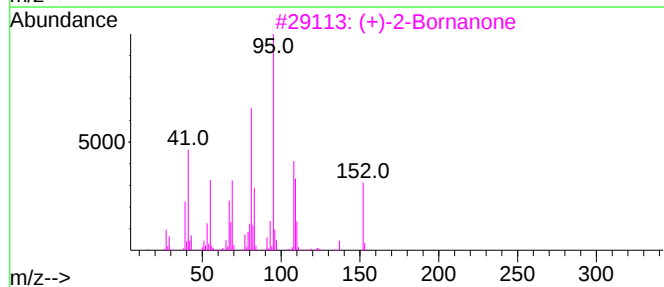
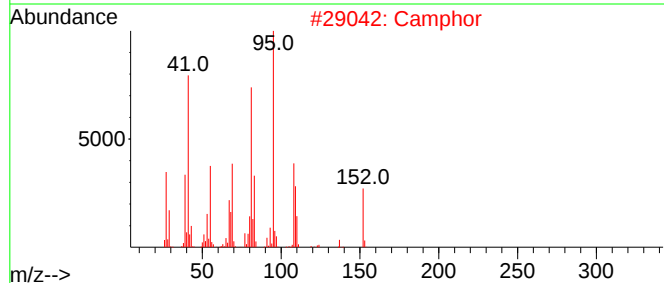
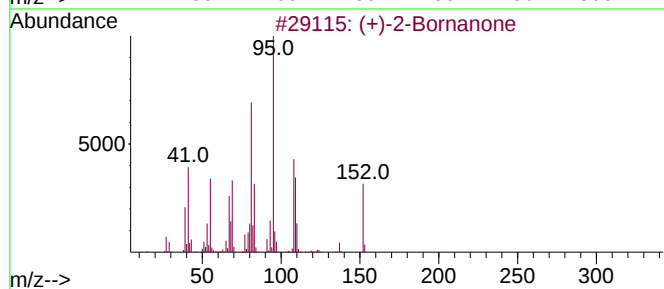
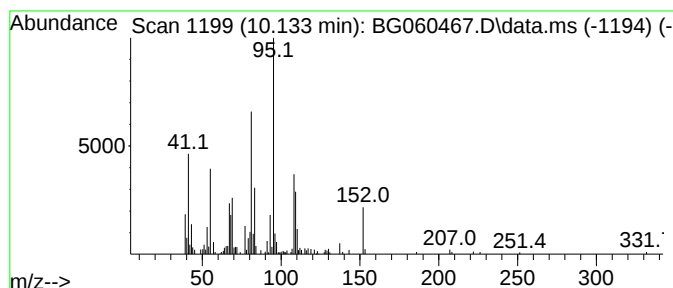
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (+)-2-Bornanone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.133	2.20 ng/ul	206421	Naphthalene-d8	10.726	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(+)-2-Bornanone	152	C10H16O	000464-49-3	97
2	Camphor	152	C10H16O	000076-22-2	97
3	(+)-2-Bornanone	152	C10H16O	000464-49-3	97
4	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	96
5	(+)-2-Bornanone	152	C10H16O	000464-49-3	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012924\
Data File : BG060467.D
Acq On : 29 Jan 2024 18:20
Operator : MA/JU
Sample : P1239-12
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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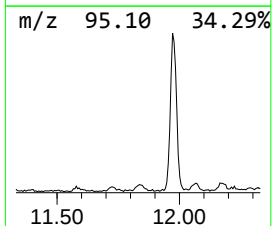
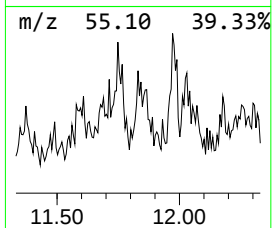
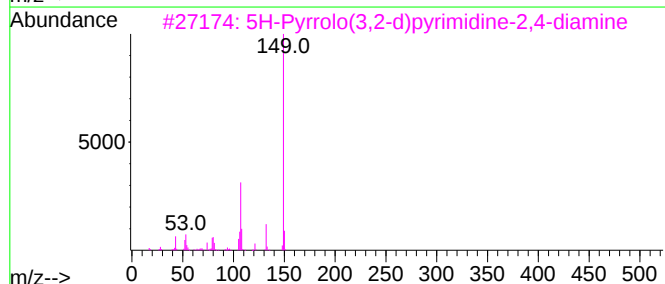
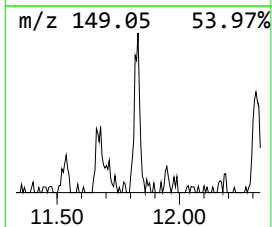
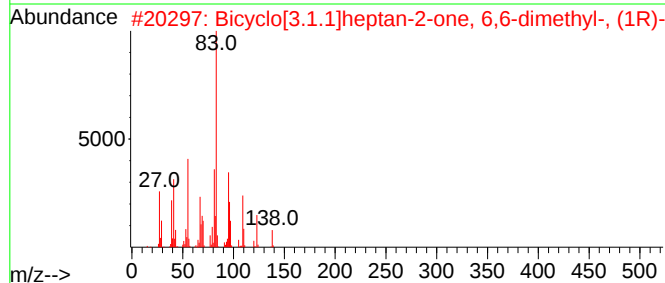
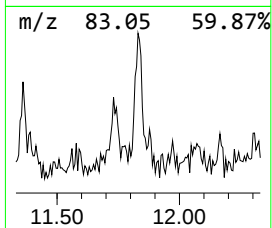
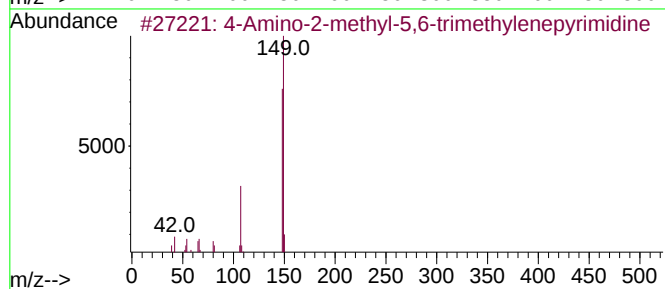
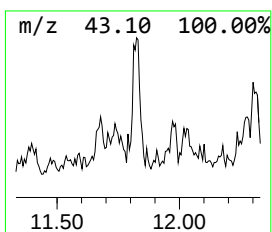
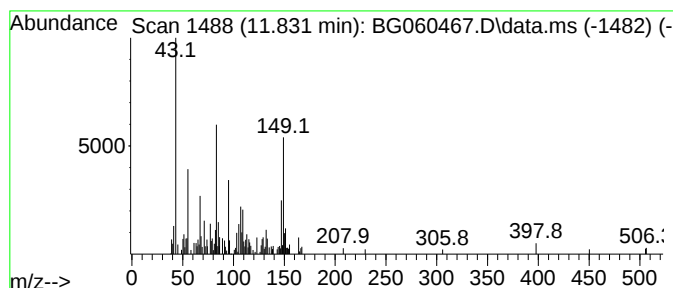
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-02 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.831	2.14 ng/ul	201019	Naphthalene-d8	10.726

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Amino-2-methyl-5,6-trimethylen...	149	C8H11N3	076881-49-7	22
2			Bicyclo[3.1.1]heptan-2-one, 6,6-...	138	C9H14O	038651-65-9	22
3			5H-Pyrrolo(3,2-d)pyrimidine-2,4-...	149	C6H7N5	1000244-21-4	22
4			Phosphonoselenoic difluoride	150	F2HPSe	026144-62-7	22
5			Phosphonoselenoic difluoride	150	F2HPSe	041916-58-9	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012924\
Data File : BG060467.D
Acq On : 29 Jan 2024 18:20
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Sample : P1239-12
Misc :
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ClientSampleId :
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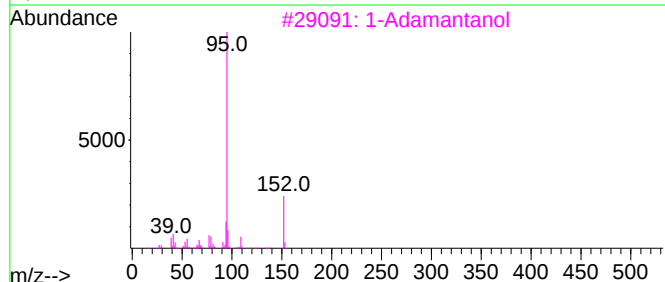
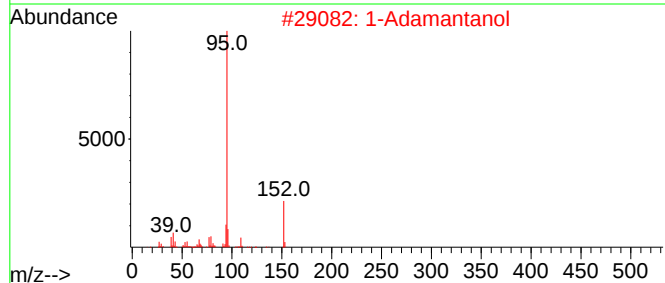
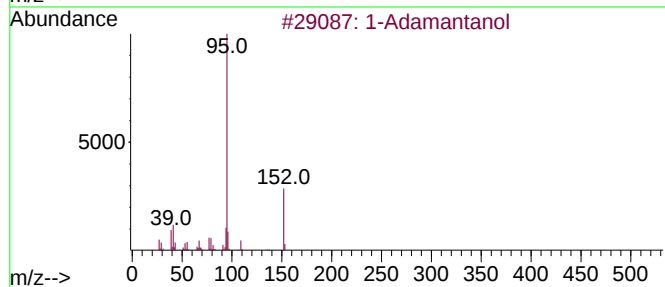
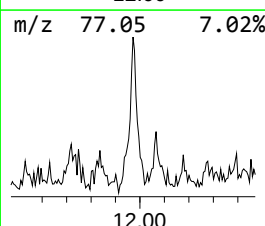
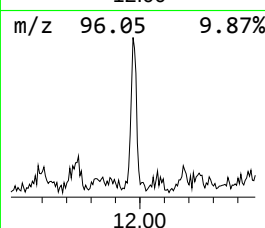
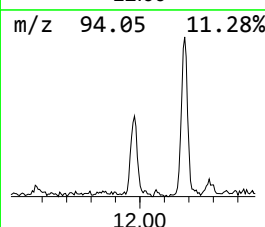
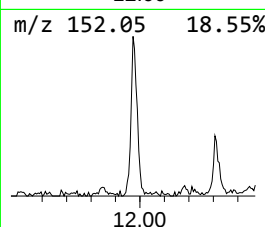
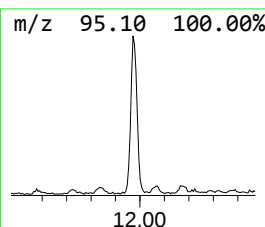
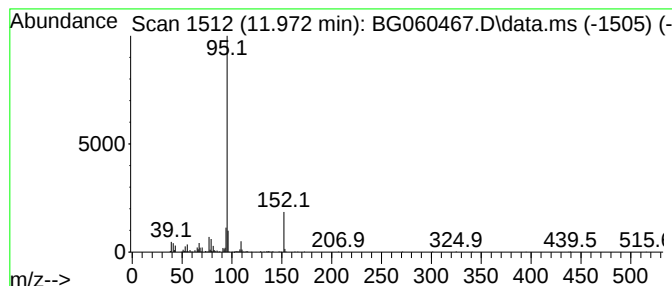
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1-Adamantanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.972	4.70 ng/ul	440670	Naphthalene-d8	10.726

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Adamantanol			152	C10H16O	000768-95-6	91
2	1-Adamantanol			152	C10H16O	000768-95-6	91
3	1-Adamantanol			152	C10H16O	000768-95-6	60
4	2-Cyclopentene-1-carboxylic acid...			168	C10H16O2	005809-03-0	50
5	1-Adamantanol			152	C10H16O	000768-95-6	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012924\
Data File : BG060467.D
Acq On : 29 Jan 2024 18:20
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Sample : P1239-12
Misc :
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Instrument :
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ClientSampleId :
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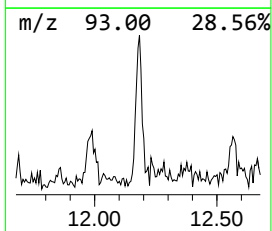
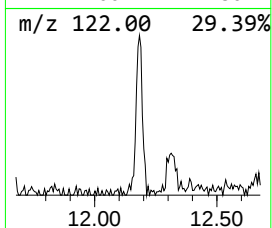
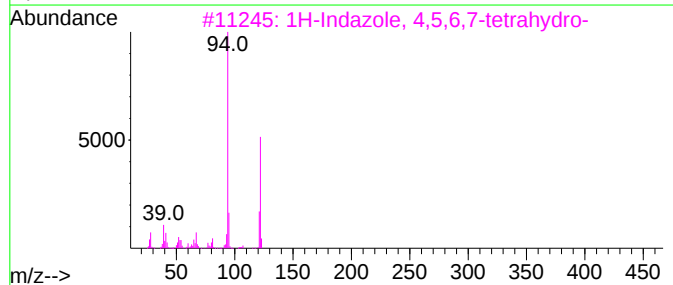
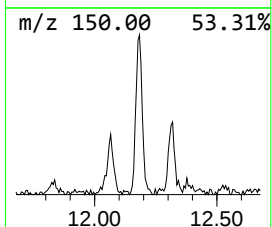
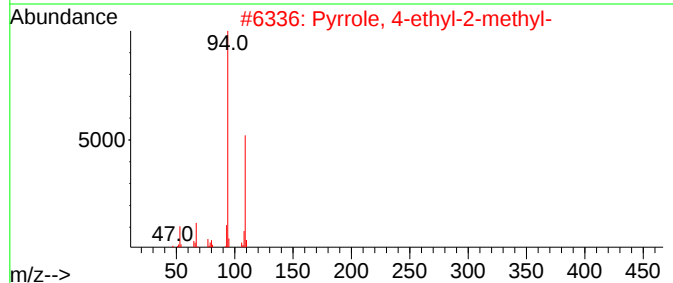
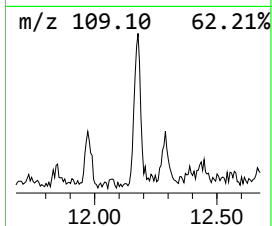
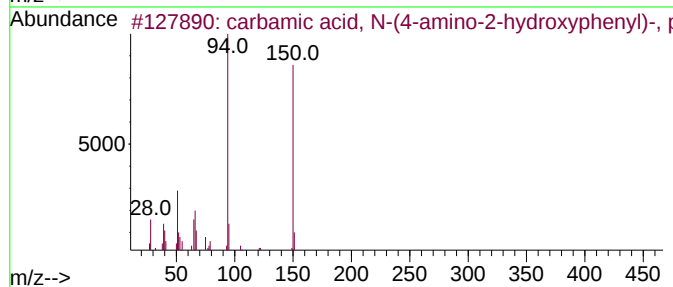
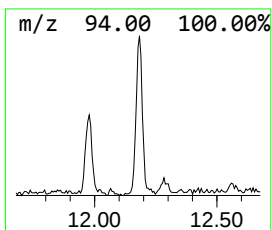
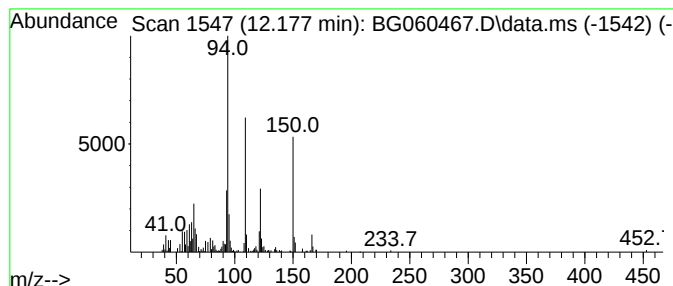
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-03 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.177	2.85 ng/ul	267034	Naphthalene-d8	10.726

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			carbamic acid, N-(4-amino-2-hydr...	244	C13H12N2O3	1000396-55-9	47
2			Pyrrole, 4-ethyl-2-methyl-	109	C7H11N	005690-96-0	46
3			1H-Indazole, 4,5,6,7-tetrahydro-	122	C7H10N2	002305-79-5	43
4			1H-Pyrrole, 2-ethyl-4-methyl-	109	C7H11N	069687-77-0	43
5			Triethylphosphine sulfide	150	C6H15PS	000597-51-3	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012924\
Data File : BG060467.D
Acq On : 29 Jan 2024 18:20
Operator : MA/JU
Sample : P1239-12
Misc :
ALS Vial : 13 Sample Multiplier: 1

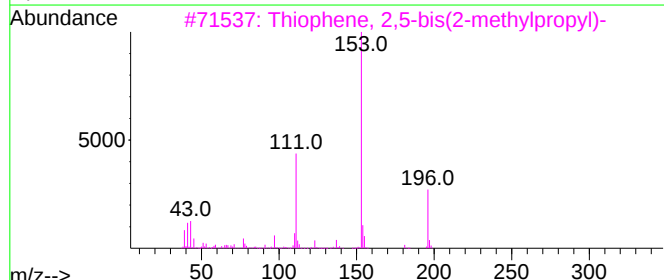
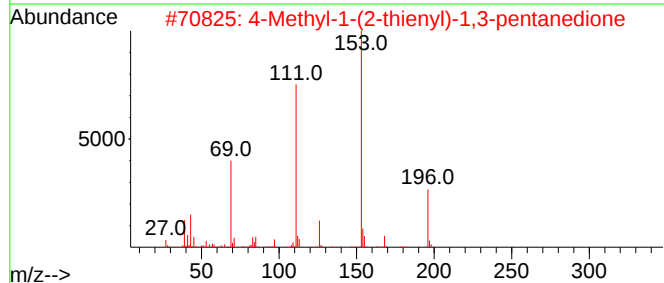
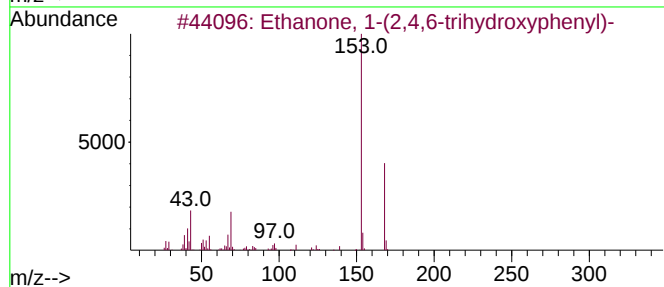
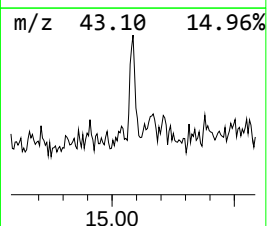
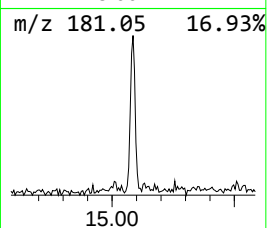
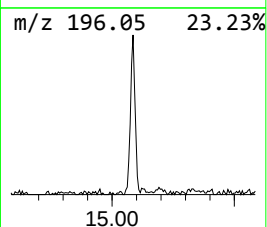
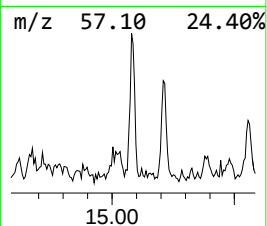
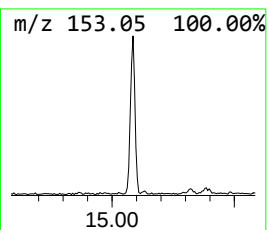
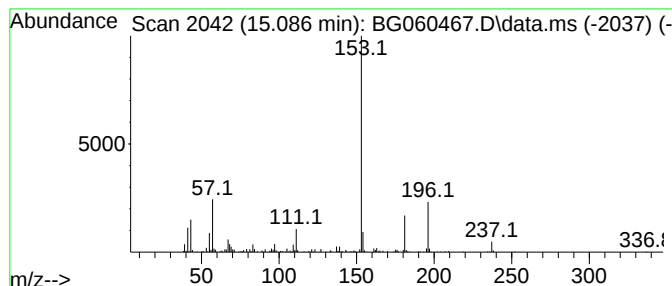
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Ethanone, 1-(2,4,6-trihydro... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.086	2.29 ng/ul	345278	Acenaphthene-d10	14.563	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanone, 1-(2,4,6-trihydroxyphe...	168	C8H8O4	000480-66-0	50
2	4-Methyl-1-(2-thienyl)-1,3-penta...	196	C10H12O2S	030984-27-1	50
3	Thiophene, 2,5-bis(2-methylpropyl)-	196	C12H20S	054845-33-9	50
4	Thiophene, 2-butyl-5-(2-methylpr...	196	C12H20S	054845-35-1	50
5	4-(1-Pyrrolidinyl)-2(5H)-furanone	153	C8H11NO2	1000427-46-1	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012924\
Data File : BG060467.D
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Operator : MA/JU
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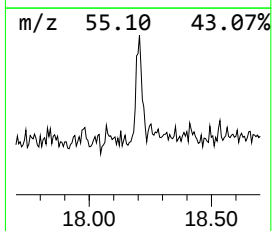
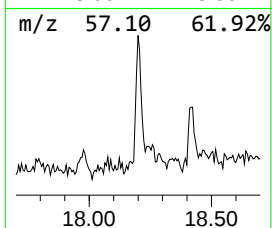
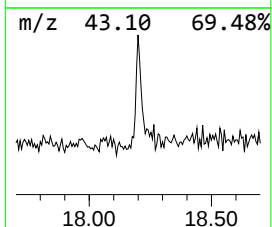
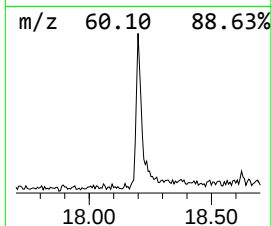
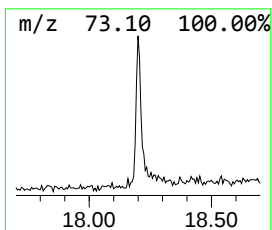
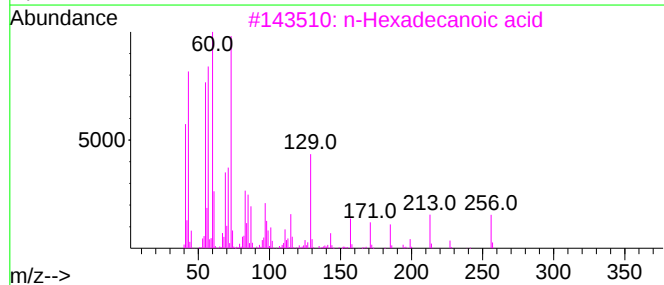
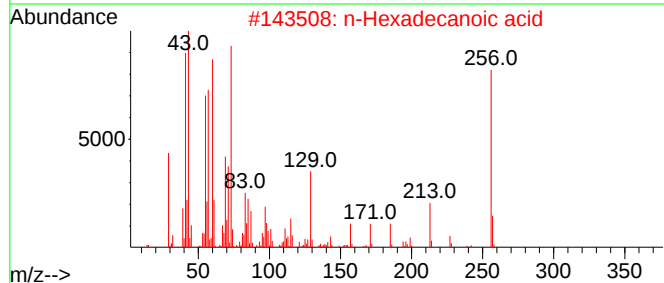
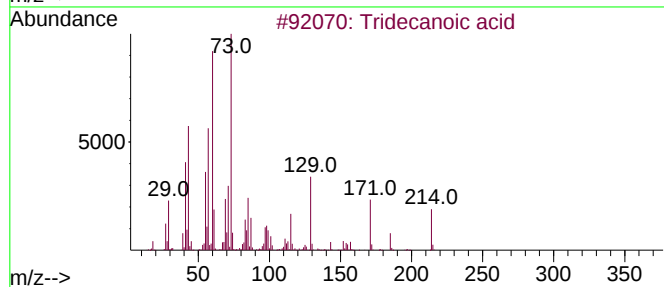
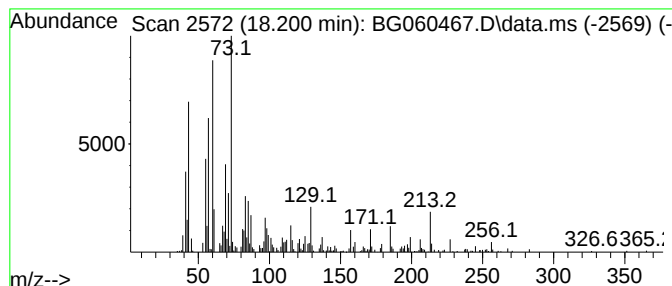
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG012624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 Tridecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.200	2.11 ng/ul	395601	Phenanthrene-d10		17.313	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tridecanoic acid		214	C13H26O2	000638-53-9	95
2	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	83
3	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	81
4	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	72
5	Dodecanoic acid, 1-methylethyl e...		242	C15H30O2	010233-13-3	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012924\
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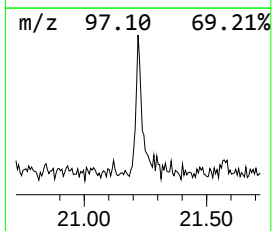
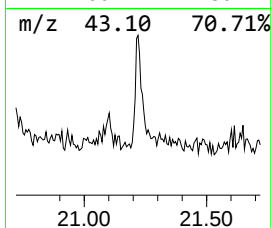
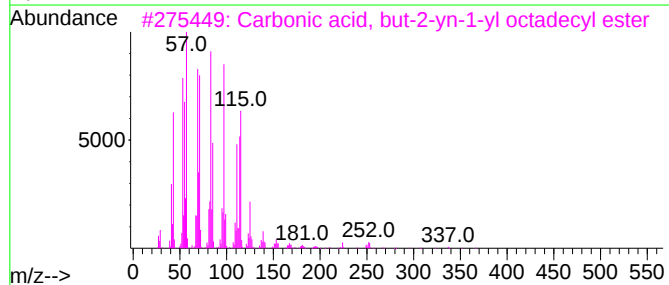
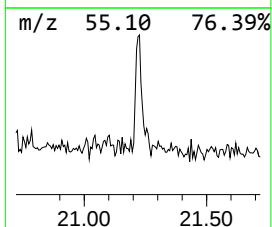
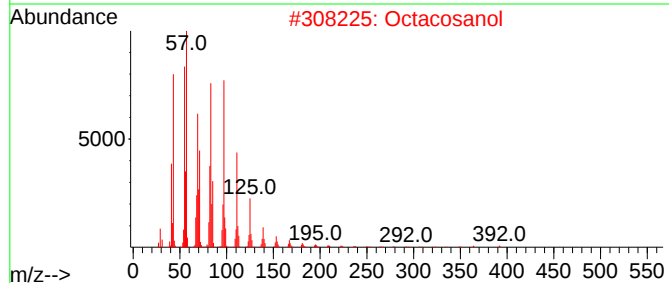
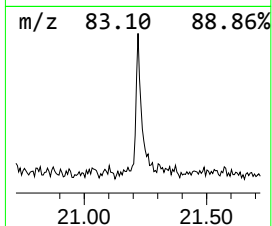
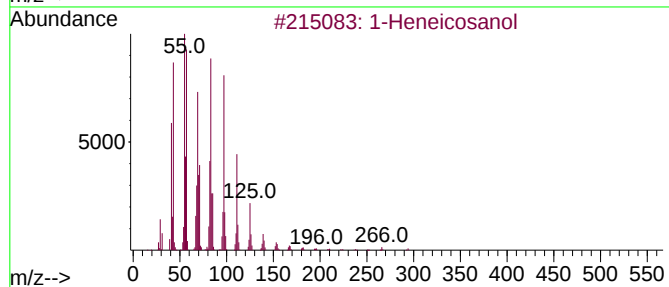
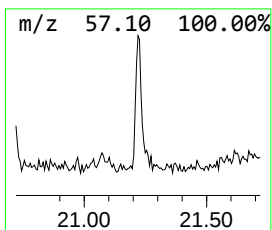
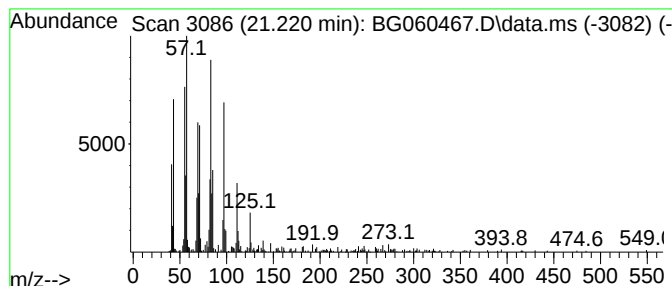
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 1-Heneicosanol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.220	2.04 ng/ul	391714	Chrysene-d12	21.572	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	83
2	Octacosanol	410	C28H58O	000557-61-9	74
3	Carbonic acid, but-2-yn-1-yl oct...	366	C23H42O3	1000383-21-3	68
4	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	68
5	Nonacos-1-ene	406	C29H58	018835-35-3	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012924\
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ALS Vial : 13 Sample Multiplier: 1

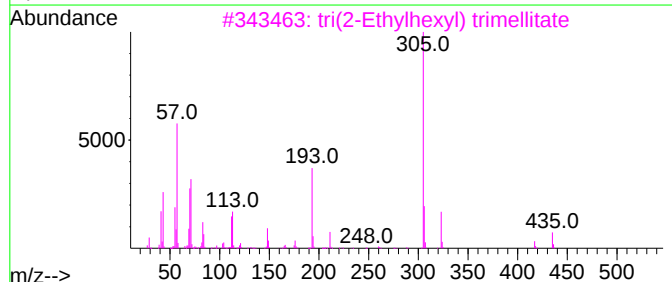
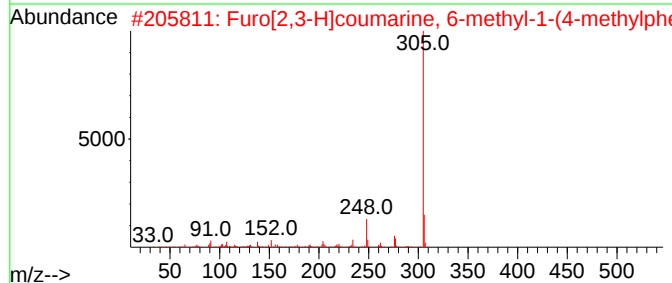
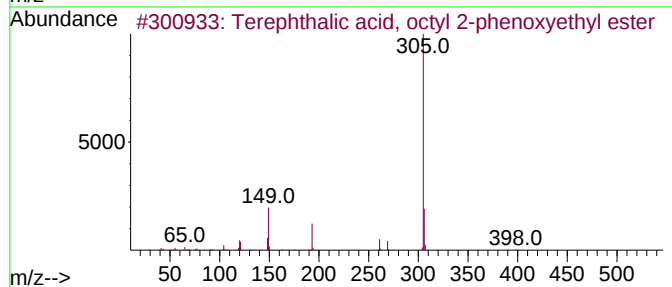
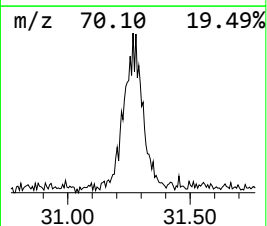
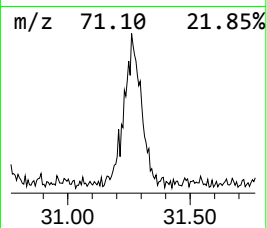
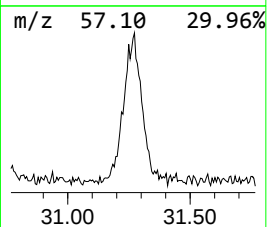
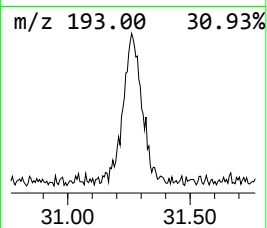
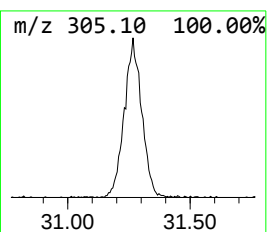
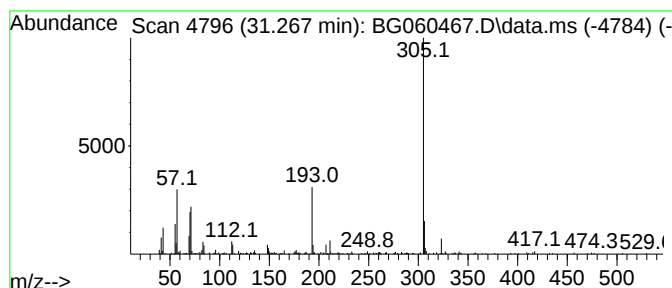
Instrument :
BNA_G
ClientSampleId :
C0FC2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG012624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-04 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
31.267	3.69 ng/ul	829276	Perylene-d12	24.739	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Terephthalic acid, octyl 2-pheno...	398	C24H30O5	1000416-04-7	43
2	Furo[2,3-H]coumarine, 6-methyl-1...	305	C19H15NO3	298684-60-3	38
3	tri(2-Ethylhexyl) trimellitate	546	C33H54O6	003319-31-1	37
4	Phthalic acid, 2-(4-bromophenoxy...	476	C24H29BrO5	1010382-89-7	33
5	1H-Purine-2,6-dione, 3,7-dihydro...	340	C16H25C1N4O2	309742-23-2	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012924\
Data File : BG060467.D
Acq On : 29 Jan 2024 18:20
Operator : MA/JU
Sample : P1239-12
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0FC2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG012624.MA.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
Diethyl sulfone	8.952	2.3	ng/ul	134777	1	7.923	1185720	20.0
unknown-01	9.357	3.3	ng/ul	305703	2	10.726	1877060	20.0
(+)-2-Bornanone	10.133	2.2	ng/ul	206421	2	10.726	1877060	20.0
unknown-02	11.831	2.1	ng/ul	201019	2	10.726	1877060	20.0
1-Adamantanol	11.972	4.7	ng/ul	440670	2	10.726	1877060	20.0
unknown-03	12.177	2.9	ng/ul	267034	2	10.726	1877060	20.0
Ethanone, 1-(2,...	15.086	2.3	ng/ul	345278	3	14.563	3012980	20.0
Tridecanoic acid	18.200	2.1	ng/ul	395601	4	17.313	3752280	20.0
1-Heneicosanol	21.220	2.0	ng/ul	391714	5	21.572	3835590	20.0
unknown-04	31.267	3.7	ng/ul	829276	6	24.739	4490250	20.0