

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
 Data File : BM042912.D
 Acq On : 19 Nov 2023 19:13
 Operator : MA/JU
 Sample : 05369-06
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GCDN3

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_M\Methods\SFAM-EPA-BM111323.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM042912.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.169	28	31	38	rVB5	10444	13243	1.13%	0.114%
2	3.616	103	107	119	rBV	47867	94739	8.05%	0.816%
3	3.793	136	137	144	rVB4	6726	8551	0.73%	0.074%
4	4.069	182	184	186	rBV2	1687	1199	0.10%	0.010%
5	4.434	243	246	252	rVB6	9833	15341	1.30%	0.132%
6	4.645	280	282	286	rVB3	2184	1397	0.12%	0.012%
7	4.804	307	309	312	rVB3	1502	1246	0.11%	0.011%
8	5.369	403	405	407	rBV2	2111	1938	0.16%	0.017%
9	5.392	407	409	413	rVB4	1218	1473	0.13%	0.013%
10	6.057	518	522	526	rBV5	1295	2399	0.20%	0.021%
11	6.986	676	680	692	rBV4	20007	50284	4.27%	0.433%
12	7.075	692	695	703	rVB4	3875	7604	0.65%	0.065%
13	7.157	703	709	720	rBV	133043	238417	20.26%	2.054%
14	7.334	733	739	750	rBV	138821	249951	21.24%	2.153%
15	7.686	797	799	806	rVB5	1857	2500	0.21%	0.022%
16	7.810	814	820	833	rBV	122096	217441	18.48%	1.873%
17	8.322	904	907	910	rVB3	1076	1433	0.12%	0.012%
18	8.539	938	944	953	rBV	73263	153141	13.01%	1.319%
19	9.010	1018	1024	1041	rBV	163909	326241	27.72%	2.810%
20	9.186	1053	1054	1061	rVB5	1455	2173	0.18%	0.019%
21	9.727	1139	1146	1163	rBV	136210	267947	22.77%	2.308%
22	10.251	1229	1235	1252	rBV	146671	328572	27.92%	2.830%
23	10.398	1259	1260	1265	rVB3	2066	1400	0.12%	0.012%
24	10.622	1292	1298	1312	rBV	165578	282844	24.03%	2.436%
25	10.798	1322	1328	1350	rBV	118687	325601	27.67%	2.805%
26	10.939	1350	1352	1354	rVB3	2440	2211	0.19%	0.019%
27	11.116	1378	1382	1384	rBV4	1424	1776	0.15%	0.015%
28	11.157	1384	1389	1391	rVV4	5325	8294	0.70%	0.071%
29	11.180	1391	1393	1398	rVV6	5383	7923	0.67%	0.068%
30	11.227	1398	1401	1405	rVB5	2296	3230	0.27%	0.028%
31	11.433	1426	1436	1445	rBV5	10826	35222	2.99%	0.303%
32	11.504	1445	1448	1451	rVV3	3477	5219	0.44%	0.045%
33	11.545	1451	1455	1465	rVB7	5024	11171	0.95%	0.096%
34	11.798	1496	1498	1501	rBV4	1334	1704	0.14%	0.015%
35	12.245	1572	1574	1581	rVB3	1240	2590	0.22%	0.022%
36	13.310	1750	1755	1761	rBV2	10324	17577	1.49%	0.151%

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Integrator: RTE

Smoothing : OFF

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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Peak separation: 5

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37	13.698	1811	1821	1828	rBV	529416	1110520	94.36%	9.566%
38	13.780	1828	1835	1848	rVV2	69563	256692	21.81%	2.211%
39	13.892	1848	1854	1866	rVB	396975	569254	48.37%	4.903%
40	14.168	1894	1901	1916	rBV2	438382	630420	53.57%	5.430%
41	14.468	1946	1952	1963	rBV	239335	353013	30.00%	3.041%
42	14.798	2001	2008	2010	rBV4	2025	3356	0.29%	0.029%
43	14.845	2015	2016	2020	rVB2	1158	1237	0.11%	0.011%
44	15.015	2042	2045	2050	rBV4	1196	2353	0.20%	0.020%
45	15.257	2079	2086	2087	rBV4	797	1340	0.11%	0.012%
46	15.462	2115	2121	2131	rBV	515475	760334	64.60%	6.549%
47	15.615	2141	2147	2160	rBV	133011	228795	19.44%	1.971%
48	15.704	2160	2162	2167	rVB4	3597	5174	0.44%	0.045%
49	16.180	2238	2243	2246	rBV3	2162	2291	0.19%	0.020%
50	16.492	2291	2296	2300	rVB5	2013	3141	0.27%	0.027%
51	16.533	2300	2303	2304	rBV3	1379	1755	0.15%	0.015%
52	16.592	2307	2313	2319	rVB6	2859	5843	0.50%	0.050%
53	17.145	2403	2407	2410	rBV3	1175	1376	0.12%	0.012%
54	17.221	2414	2420	2431	rBV	278951	404866	34.40%	3.487%
55	17.321	2431	2437	2454	rVV	561124	842839	71.62%	7.260%
56	17.668	2493	2496	2502	rVB3	1436	2077	0.18%	0.018%
57	17.721	2502	2505	2507	rBV4	1316	1345	0.11%	0.012%
58	17.750	2507	2510	2513	rVB3	1173	1228	0.10%	0.011%
59	17.845	2523	2526	2532	rBV4	2022	3709	0.32%	0.032%
60	18.015	2553	2555	2558	rBV4	887	1181	0.10%	0.010%
61	18.080	2562	2566	2574	rVV6	7846	13680	1.16%	0.118%
62	18.145	2574	2577	2578	rVV3	1460	1431	0.12%	0.012%
63	18.174	2580	2582	2585	rVB3	1534	1597	0.14%	0.014%
64	18.303	2598	2604	2619	rBV	243007	387414	32.92%	3.337%
65	18.539	2642	2644	2647	rVB2	1935	1918	0.16%	0.017%
66	19.003	2719	2723	2724	rBV3	1349	1766	0.15%	0.015%
67	19.180	2749	2753	2757	rBV3	6958	8948	0.76%	0.077%
68	19.256	2763	2766	2771	rVB4	4864	6188	0.53%	0.053%
69	19.362	2778	2784	2788	rBV6	4397	9527	0.81%	0.082%
70	19.492	2804	2806	2811	rBV5	1532	1817	0.15%	0.016%
71	19.621	2822	2828	2840	rBV	815602	1097660	93.27%	9.455%
72	20.350	2950	2952	2953	rBV2	2367	1735	0.15%	0.015%
73	21.409	3127	3132	3144	rBV	306594	469187	39.87%	4.041%
74	23.609	3499	3506	3521	rBV	587435	1176901	100.00%	10.138%
75	23.768	3526	3533	3544	rVB	257255	541451	46.01%	4.664%

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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_M\Methods\SFAM-EPA-BM111323.MA.M
Title : SVOA CALIBRATION

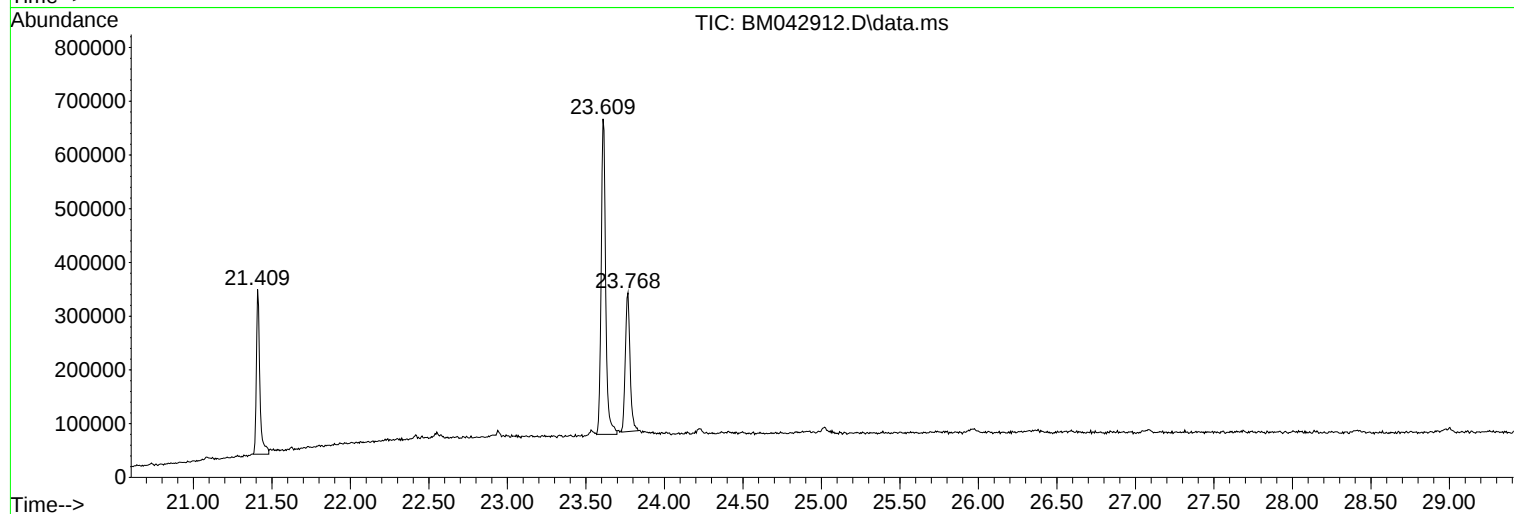
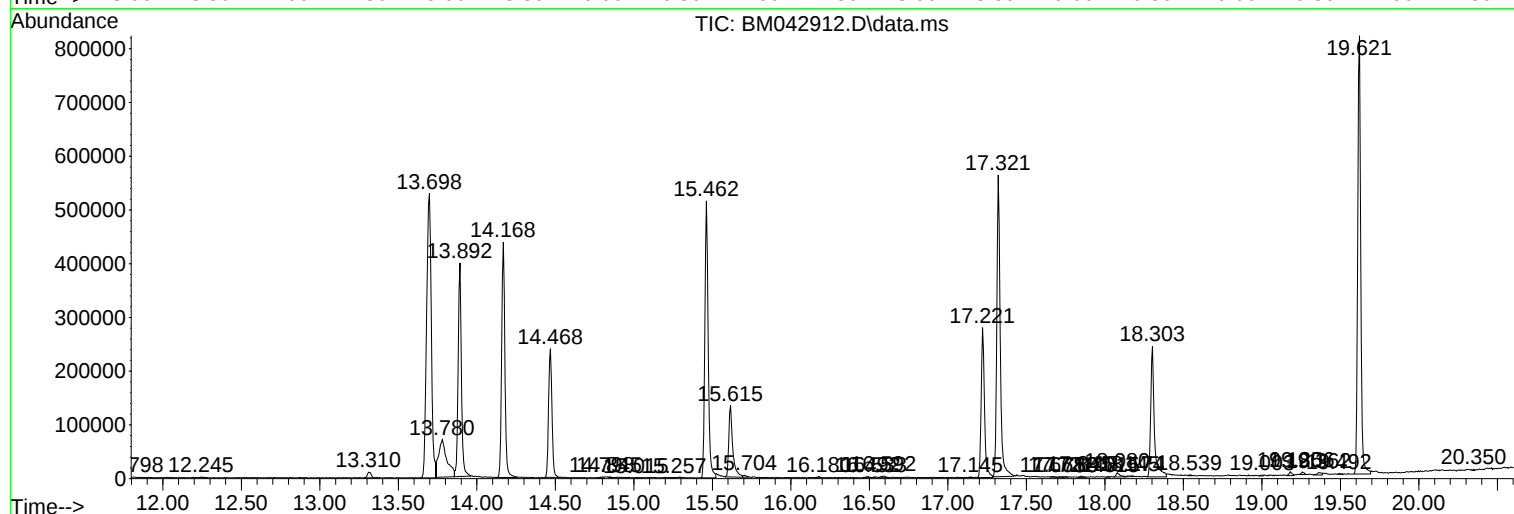
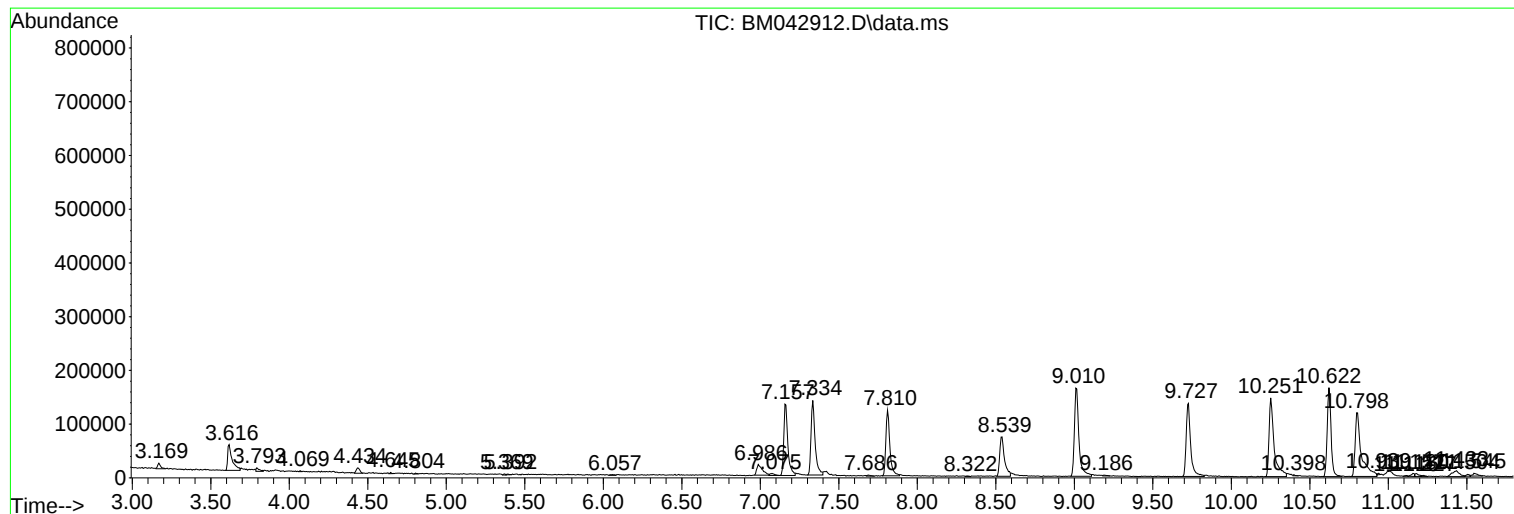
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Instrument :
BNA_M
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
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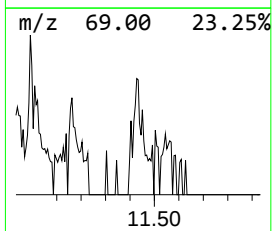
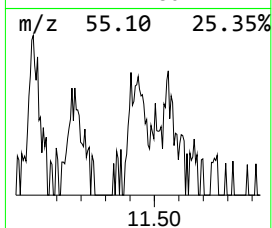
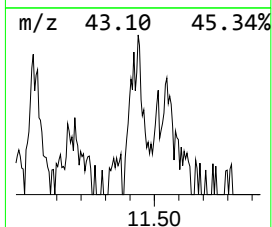
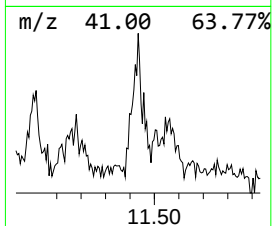
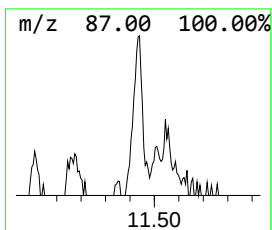
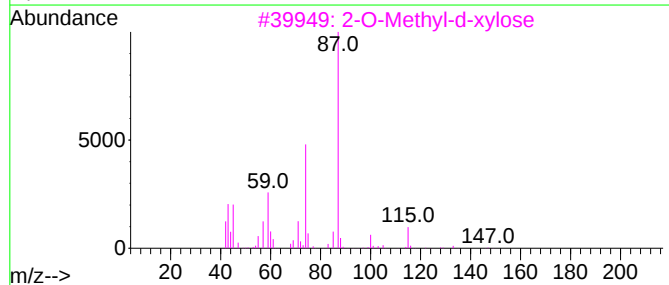
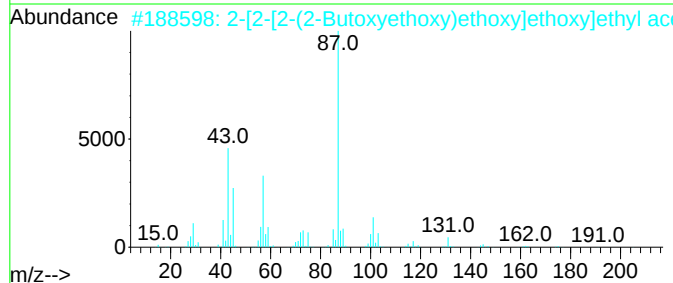
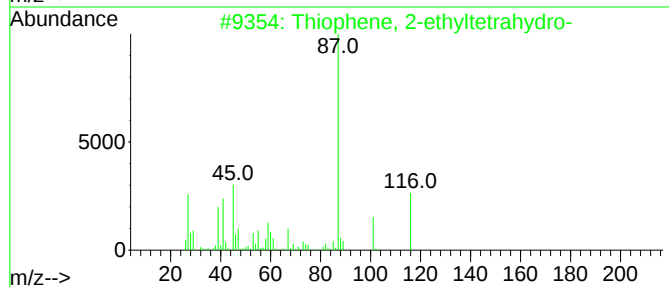
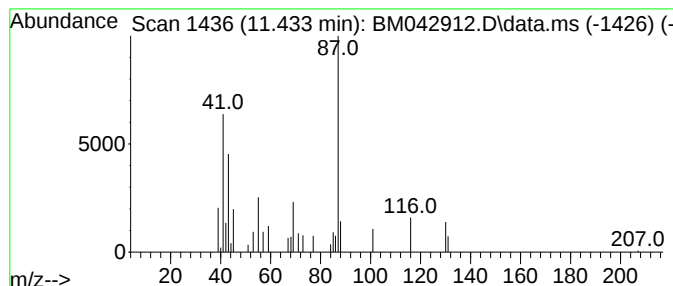
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Thiophene, 2-ethyltetrahydro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.433	2.49 ng/ul	35222	Naphthalene-d8	10.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Thiophene, 2-ethyltetrahydro-	116	C6H12S	001551-32-2	64
2			2-[2-[2-(2-Butoxyethoxy)ethoxy]e...	292	C14H28O6	1000351-93-9	64
3			2-O-Methyl-d-xylose	164	C6H12O5	1000130-01-3	59
4			Thiophane, propyl-	130	C7H14S	001551-34-4	58
5			Trimethylsilyl 23-acetoxy-3,6,9,...	498	C21H42O11Si	1000366-84-4	56



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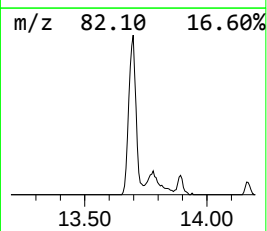
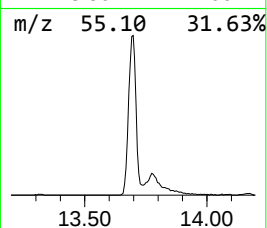
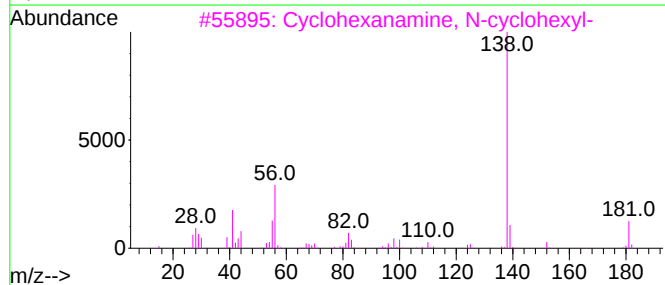
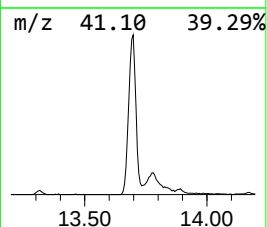
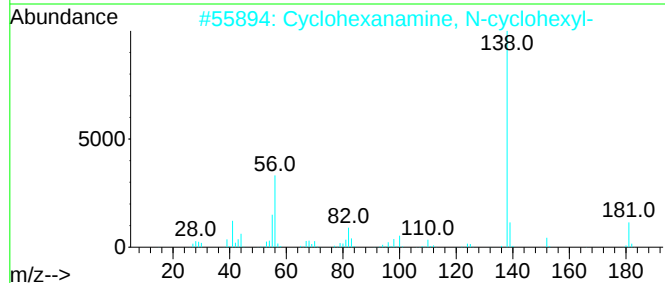
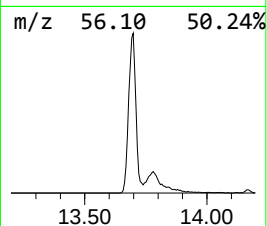
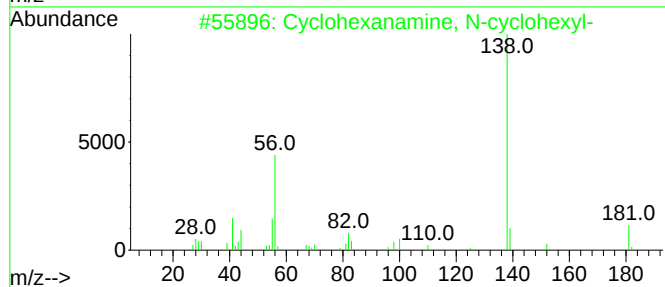
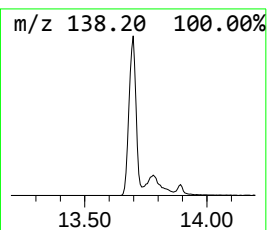
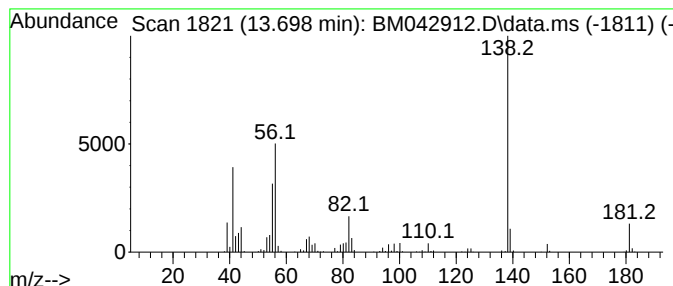
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Cyclohexanamine, N-cyclohexyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.698	62.92 ng/ul	1110520	Acenaphthene-d10	14.468

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	94
2			Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	93
3			Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	93
4			(Bicyclohexyl)-2-amine	181	C12H23N	006283-14-3	91
5			Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	87



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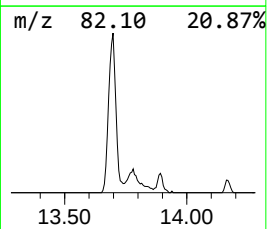
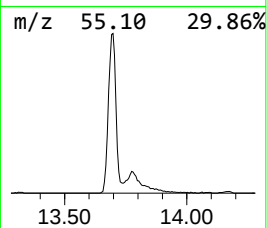
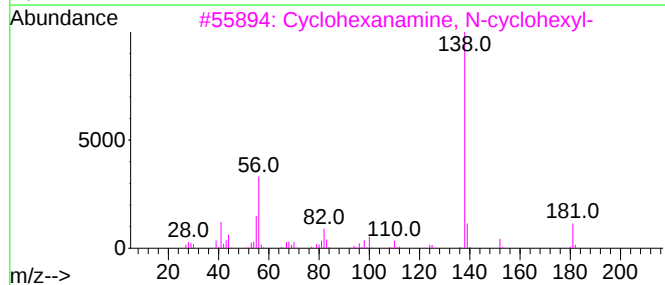
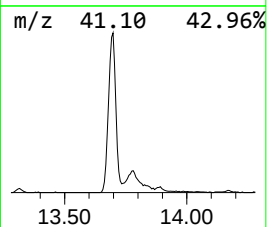
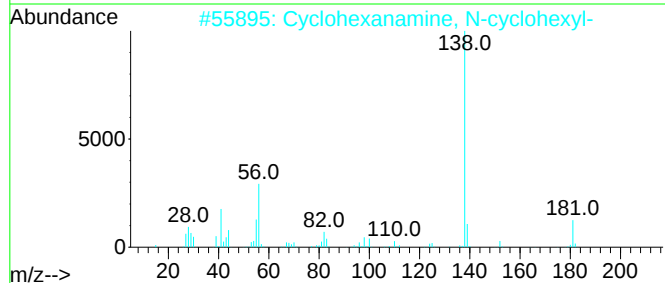
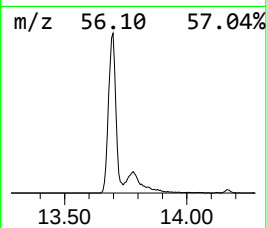
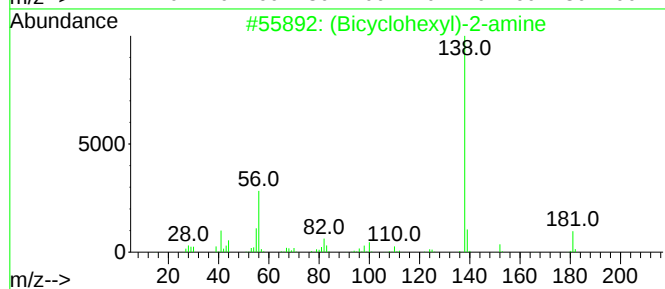
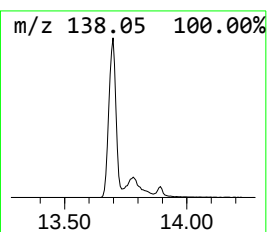
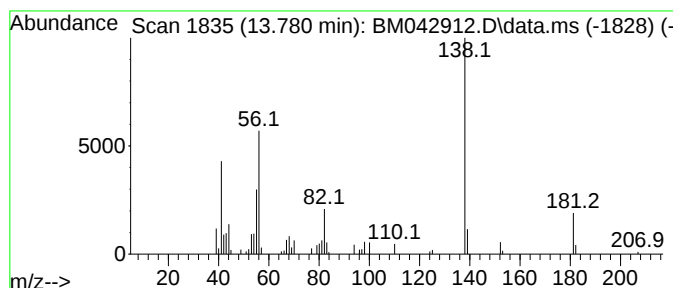
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (Bicyclohexyl)-2-amine Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.780	14.54 ng/ul	256692	Acenaphthene-d10	14.468

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(Bicyclohexyl)-2-amine	181	C12H23N	006283-14-3	93
2			Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	90
3			Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	90
4			Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	90
5			Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	90



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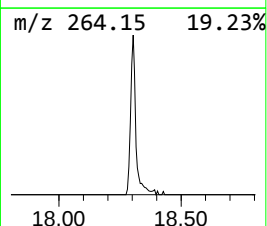
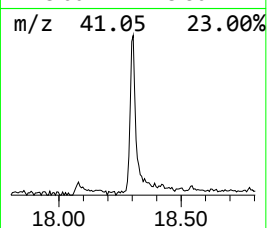
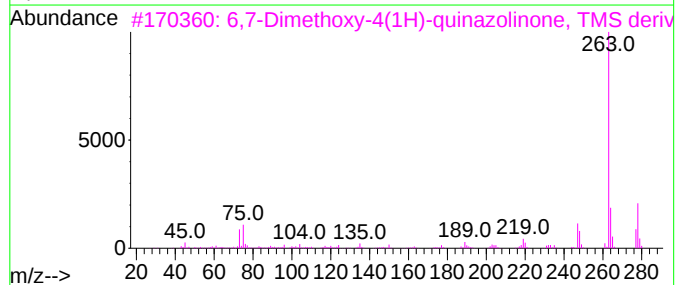
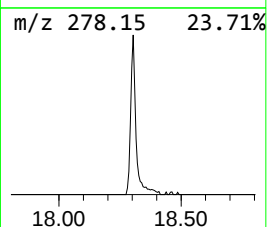
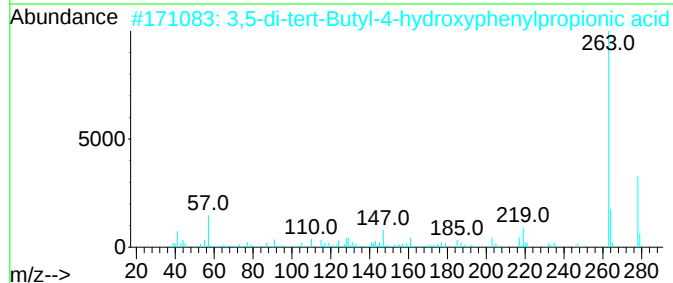
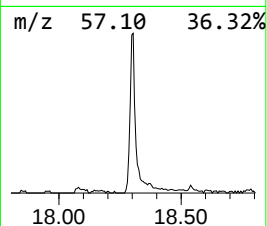
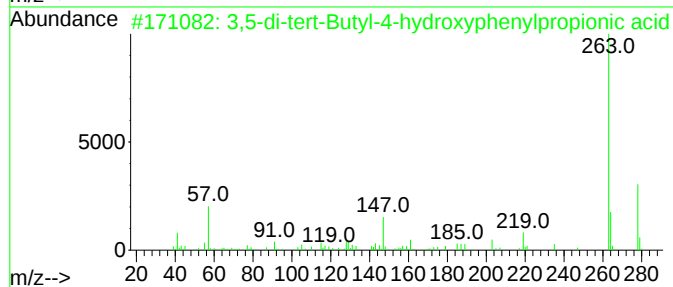
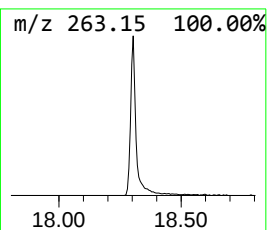
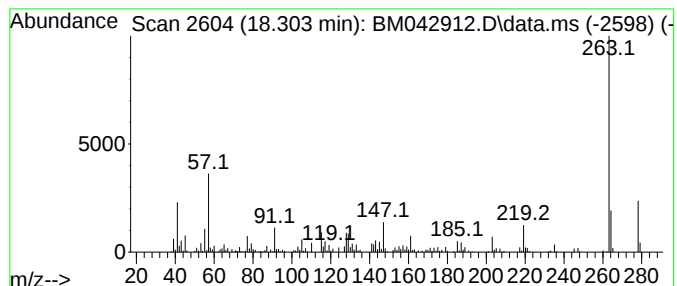
Instrument :
BNA_M
ClientSampleId :
GCDN3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.303	19.14 ng/ul	387414	Phenanthrene-d10	17.221	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	98
2	3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	91
3	6,7-Dimethoxy-4(1H)-quinazolinon...	278	C13H18N2O3Si	1000479-65-5	80
4	Cyanic acid, (1-methylethylidene...	278	C17H14N2O2	001156-51-0	74
5	5-(2,3-Difluorophenyl)-2-pyridin...	278	C14H16F2N2Si	1000504-49-4	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
Data File : BM042912.D
Acq On : 19 Nov 2023 19:13
Operator : MA/JU
Sample : 05369-06
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCDN3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Thiophene, 2-et...	11.433	2.5	ng/u1	35222	2	10.621	282844	20.0
Cyclohexanamine...	13.698	62.9	ng/u1	1110520	3	14.468	353013	20.0
(Bicyclohexyl)-...	13.780	14.5	ng/u1	256692	3	14.468	353013	20.0
3,5-di-tert-But...	18.303	19.1	ng/u1	387414	4	17.221	404866	20.0