

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
 Data File : BP011410.D
 Acq On : 12 Aug 2022 04:51
 Operator : CG/JU
 Sample : N4129-02
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGBW2

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_P\Methods\SFAM-EPA-BP080822.M
 Title : SVOA CALIBRATION

Signal : TIC: BP011410.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.081	13	16	28	rVB	31207	44714	1.62%	0.172%
2	3.175	29	32	36	rVB2	8722	10515	0.38%	0.041%
3	3.252	41	45	52	rVB3	19970	26950	0.98%	0.104%
4	3.358	60	63	68	rVB4	9120	11368	0.41%	0.044%
5	3.499	85	87	97	rBV	94933	147773	5.37%	0.570%
6	3.605	103	105	109	rVB3	7262	8144	0.30%	0.031%
7	3.705	117	122	127	rBV2	17363	26722	0.97%	0.103%
8	3.763	129	132	134	rVV3	7212	8928	0.32%	0.034%
9	3.793	135	137	140	rVB3	9024	10377	0.38%	0.040%
10	4.234	208	212	215	rBV	20772	27644	1.00%	0.107%
11	4.263	215	217	218	rVV	14983	14423	0.52%	0.056%
12	4.299	218	223	236	rVB	2035993	2753065	100.00%	10.616%
13	4.546	261	265	266	rBV	15151	15663	0.57%	0.060%
14	4.581	266	271	283	rVB	741347	1027085	37.31%	3.960%
15	4.669	283	286	290	rVB4	4380	5718	0.21%	0.022%
16	4.793	302	307	318	rVB	77294	115537	4.20%	0.446%
17	5.463	417	421	431	rVB3	7132	14100	0.51%	0.054%
18	5.634	444	450	457	rVB	11553	20195	0.73%	0.078%
19	5.799	473	478	481	rBV7	3077	5579	0.20%	0.022%
20	5.846	481	486	494	rVB4	29247	47400	1.72%	0.183%
21	6.710	626	633	638	rBV2	11344	19459	0.71%	0.075%
22	6.763	638	642	650	rVB	11428	18509	0.67%	0.071%
23	6.922	663	669	676	rBV	370860	572708	20.80%	2.208%
24	6.987	676	680	689	rVB	58164	94496	3.43%	0.364%
25	7.110	694	701	710	rVV	204524	317374	11.53%	1.224%
26	7.199	713	716	722	rVV7	5426	7870	0.29%	0.030%
27	7.287	725	731	739	rVB	34553	53685	1.95%	0.207%
28	7.575	772	780	786	rBV	377026	579553	21.05%	2.235%
29	7.640	786	791	798	rVB2	30172	53118	1.93%	0.205%
30	7.840	820	825	829	rBV7	3436	5342	0.19%	0.021%
31	7.946	836	843	850	rVB10	7126	13178	0.48%	0.051%
32	8.063	858	863	868	rVB8	5790	9264	0.34%	0.036%
33	8.187	877	884	889	rBV5	8298	16755	0.61%	0.065%
34	8.299	897	903	914	rBV2	53728	98555	3.58%	0.380%
35	8.440	920	927	935	rVB2	18163	33715	1.22%	0.130%
36	8.740	969	978	986	rBV	481642	780412	28.35%	3.009%

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37	8.846	991	996	1007	rVB	55572	94059	3.42%	0.363%
38	9.146	1042	1047	1055	rBV	12244	21770	0.79%	0.084%
39	9.457	1093	1100	1118	rBV	312487	512372	18.61%	1.976%
40	9.993	1181	1191	1208	rBV	390066	675946	24.55%	2.606%
41	10.151	1211	1218	1229	rVV2	78863	152622	5.54%	0.589%
42	10.275	1233	1239	1247	rVV2	13023	28340	1.03%	0.109%
43	10.363	1247	1254	1266	rVB	500000	797850	28.98%	3.077%
44	10.528	1274	1282	1289	rBV2	22221	43397	1.58%	0.167%
45	11.198	1390	1396	1402	rBV2	3243	6797	0.25%	0.026%
46	11.304	1407	1414	1425	rVB	30722	56012	2.03%	0.216%
47	11.410	1425	1432	1443	rBV	86613	170318	6.19%	0.657%
48	11.651	1463	1473	1482	rBV6	24178	61719	2.24%	0.238%
49	11.845	1500	1506	1514	rVB2	35538	58542	2.13%	0.226%
50	11.975	1519	1528	1534	rVB5	6242	13933	0.51%	0.054%
51	12.087	1541	1547	1554	rBV3	26694	42665	1.55%	0.165%
52	12.351	1587	1592	1596	rBV7	5843	10918	0.40%	0.042%
53	12.645	1630	1642	1653	rBV4	79423	197376	7.17%	0.761%
54	12.922	1685	1689	1693	rBV6	6203	10596	0.38%	0.041%
55	13.039	1702	1709	1716	rBV5	7964	15057	0.55%	0.058%
56	13.228	1735	1741	1748	rBV	50128	81385	2.96%	0.314%
57	13.675	1803	1817	1825	rBV	1089052	1472958	53.50%	5.680%
58	13.751	1825	1830	1835	rVV	72782	121236	4.40%	0.467%
59	13.810	1836	1840	1850	rVB3	28747	52258	1.90%	0.202%
60	13.939	1855	1862	1872	rBV	483149	705939	25.64%	2.722%
61	14.181	1897	1903	1908	rBV7	11617	20289	0.74%	0.078%
62	14.251	1908	1915	1923	rBV	693612	1013136	36.80%	3.907%
63	14.381	1933	1937	1942	rVB8	4572	6506	0.24%	0.025%
64	14.492	1948	1956	1966	rBV2	27328	76729	2.79%	0.296%
65	14.581	1969	1971	1974	rVV3	6313	8247	0.30%	0.032%
66	14.616	1974	1977	1985	rVB10	5166	10488	0.38%	0.040%
67	14.751	1994	2000	2005	rBV7	15347	25277	0.92%	0.097%
68	14.810	2005	2010	2017	rVB2	48762	70468	2.56%	0.272%
69	14.922	2023	2029	2031	rBV7	4502	7650	0.28%	0.029%
70	15.104	2056	2060	2071	rVB5	23745	42947	1.56%	0.166%
71	15.251	2078	2085	2102	rBV	1123110	1661383	60.35%	6.406%
72	15.386	2102	2108	2122	rVV	383992	555730	20.19%	2.143%
73	15.492	2123	2126	2131	rVB3	9936	13508	0.49%	0.052%
74	15.739	2160	2168	2173	rBV7	10431	21932	0.80%	0.085%
75	15.951	2198	2204	2215	rBV4	30491	79424	2.88%	0.306%
76	16.045	2218	2220	2226	rVV6	4799	7248	0.26%	0.028%

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77	16.139	2229	2236	2240	rVV10	4408	10027	0.36%	0.039%
78	16.504	2295	2298	2304	rVB3	5673	7501	0.27%	0.029%
79	16.675	2323	2327	2335	rVB2	22278	36904	1.34%	0.142%
80	16.904	2362	2366	2373	rBV2	17354	23770	0.86%	0.092%
81	17.022	2380	2386	2396	rBV	843465	1166636	42.38%	4.499%
82	17.122	2396	2403	2414	rVV3	1150402	1723834	62.62%	6.647%
83	17.210	2415	2418	2423	rVV3	19689	38782	1.41%	0.150%
84	17.275	2423	2429	2447	rVB2	83853	172627	6.27%	0.666%
85	17.486	2462	2465	2473	rVB8	6708	10935	0.40%	0.042%
86	17.586	2479	2482	2489	rVB3	7711	10173	0.37%	0.039%
87	17.845	2523	2526	2530	rVV4	6844	10257	0.37%	0.040%
88	17.951	2536	2544	2548	rBV	61173	102986	3.74%	0.397%
89	18.251	2590	2595	2599	rBV5	10815	21558	0.78%	0.083%
90	18.404	2616	2621	2627	rBV2	47836	74198	2.70%	0.286%
91	18.539	2642	2644	2652	rVB9	4917	8287	0.30%	0.032%
92	18.663	2662	2665	2667	rBV4	5712	7448	0.27%	0.029%
93	18.716	2670	2674	2680	rVB7	22766	37162	1.35%	0.143%
94	18.957	2712	2715	2722	rVB3	14945	19764	0.72%	0.076%
95	19.033	2723	2728	2737	rBV5	37936	89729	3.26%	0.346%
96	19.386	2782	2788	2797	rBV	1525869	1968005	71.48%	7.589%
97	20.580	2987	2991	2997	rBV3	54235	75003	2.72%	0.289%
98	21.151	3083	3088	3099	rBV	1008088	1327537	48.22%	5.119%
99	23.309	3449	3455	3463	rBV	920398	1669952	60.66%	6.439%
100	23.457	3474	3480	3490	rVB2	664244	1279609	46.48%	4.934%

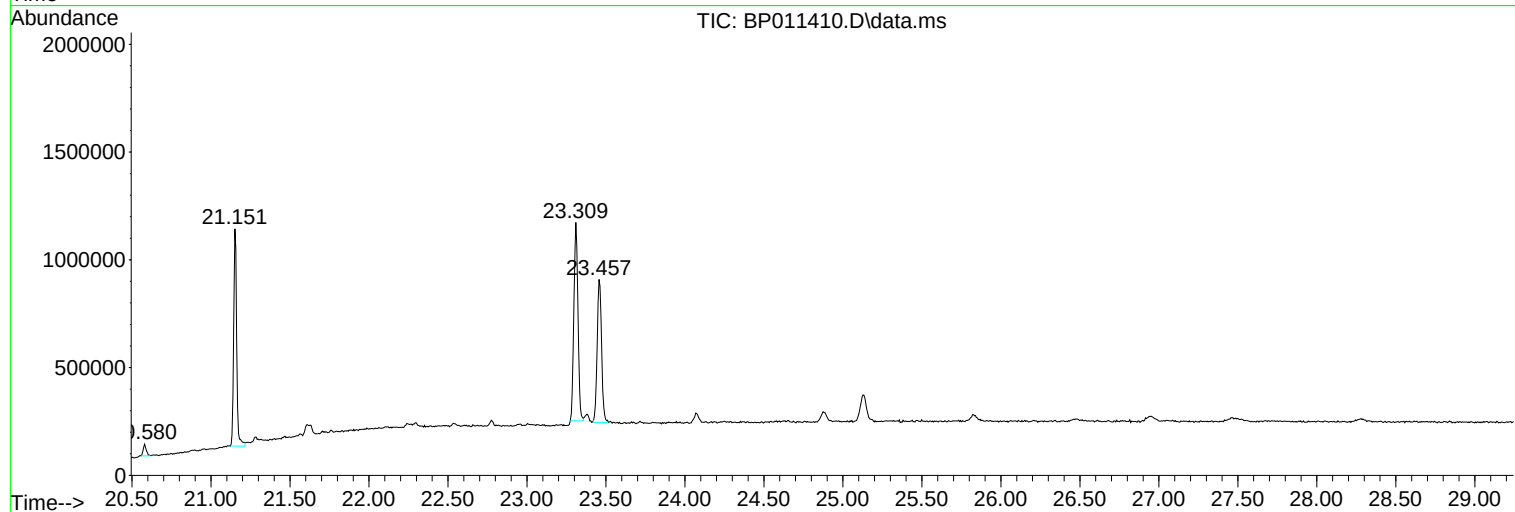
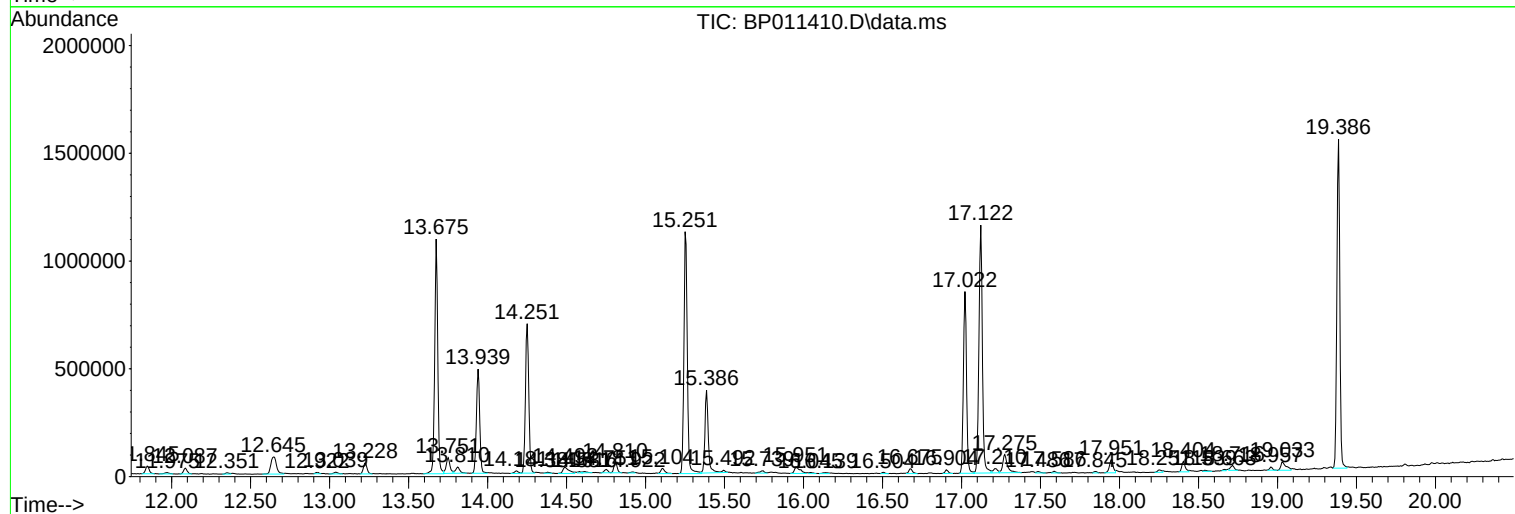
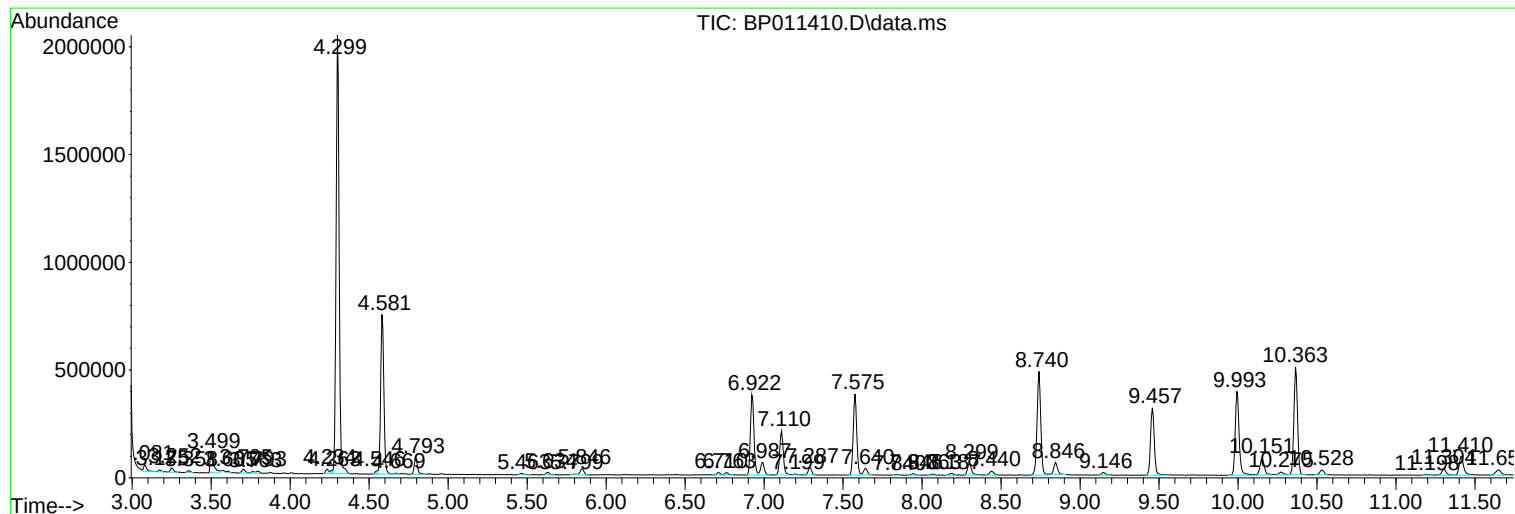
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

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



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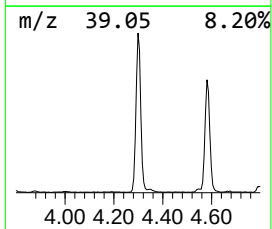
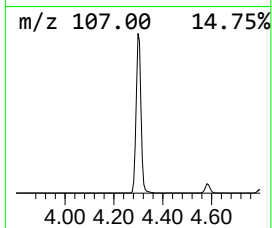
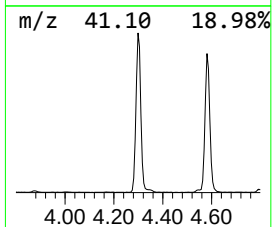
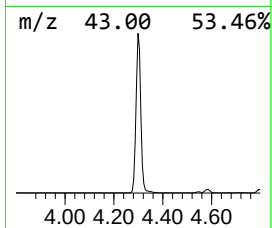
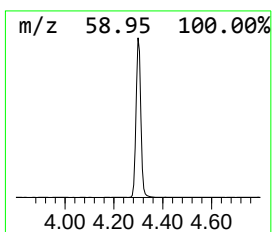
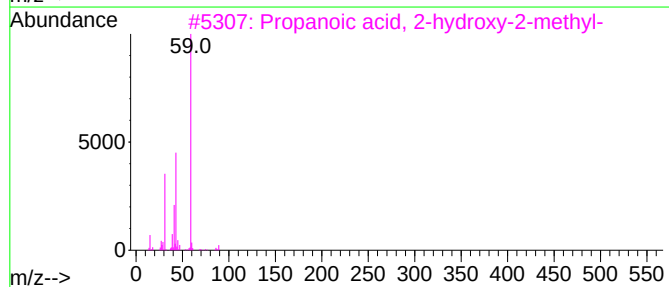
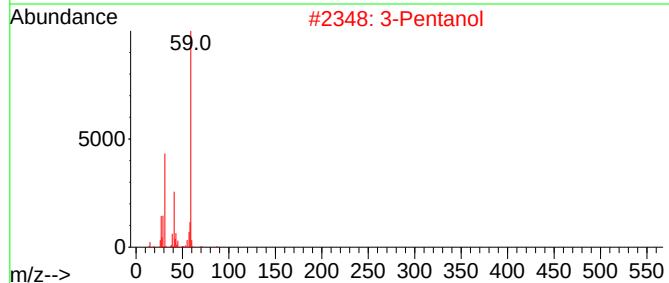
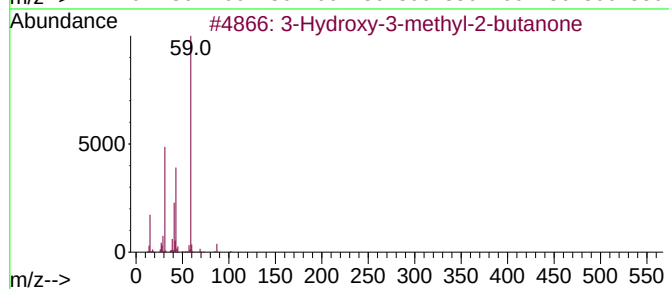
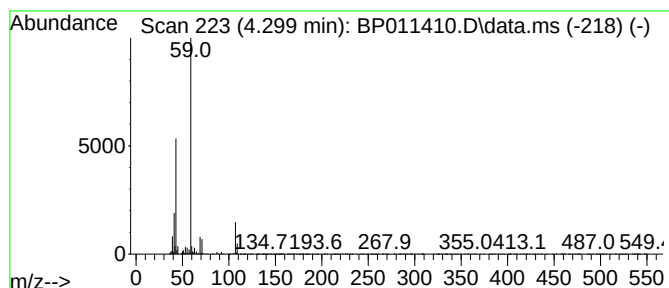
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.299	95.01 ng/ul	2753070	1,4-Dichlorobenzene-d4	7.575

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	45		
2	3-Pentanol	88	C5H12O	000584-02-1	42		
3	Propanoic acid, 2-hydroxy-2-methyl-	104	C4H8O3	000594-61-6	42		
4	2-Propanol, 2-methyl-	74	C4H10O	000075-65-0	40		
5	3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	38		



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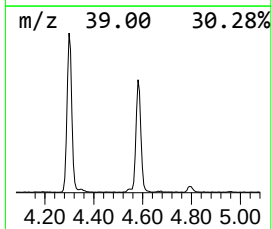
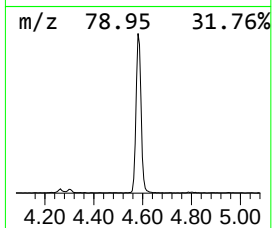
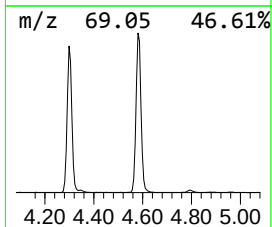
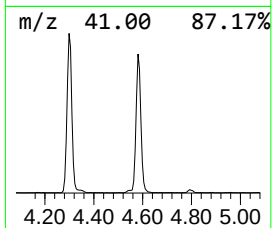
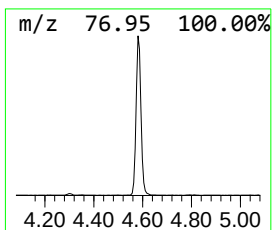
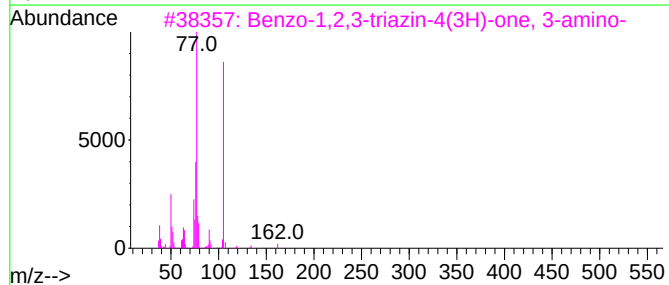
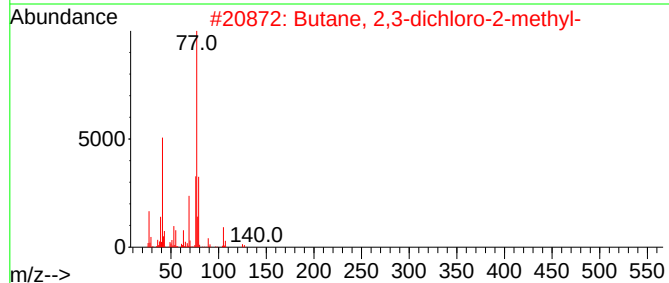
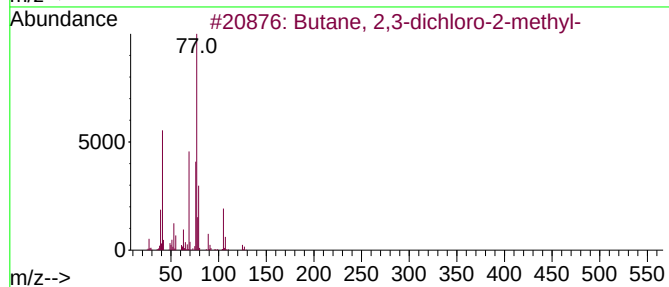
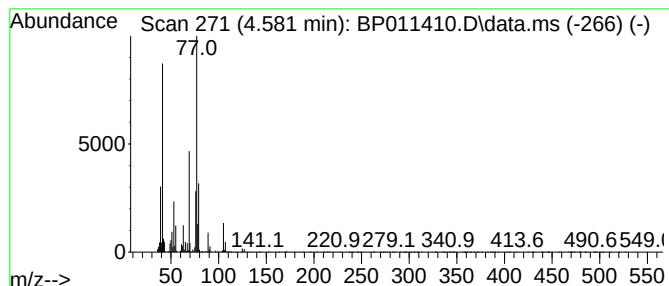
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butane, 2,3-dichloro-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.581	35.44 ng/ul	1027090	1,4-Dichlorobenzene-d4	7.575

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,3-dichloro-2-methyl-	140	C5H10Cl2	000507-45-9	86
2			Butane, 2,3-dichloro-2-methyl-	140	C5H10Cl2	000507-45-9	78
3			Benzo-1,2,3-triazin-4(3H)-one, 3...	162	C7H6N4O	041225-81-4	40
4			Propane, 2-bromo-1-chloro-	156	C3H6BrCl	003017-95-6	37
5			Propane, 1,1,2-trichloro-2-methyl-	160	C4H7Cl3	029559-52-2	33



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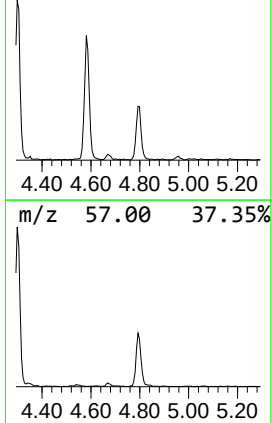
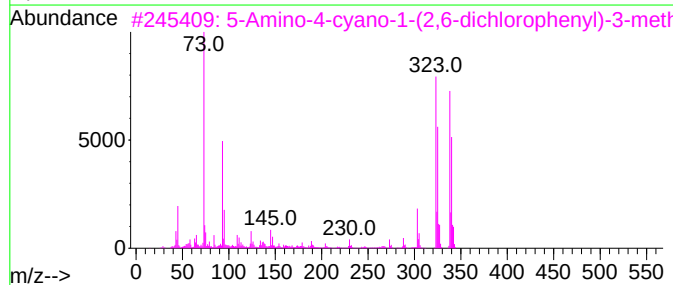
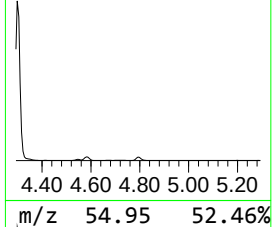
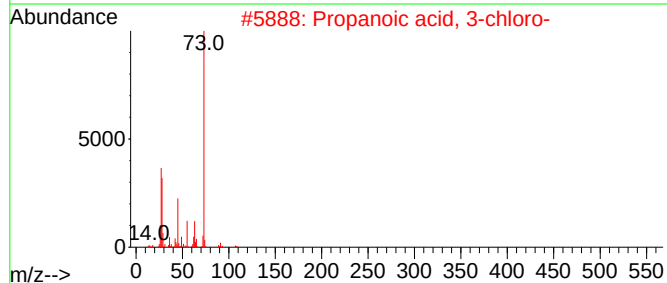
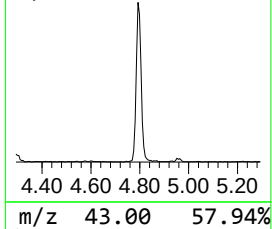
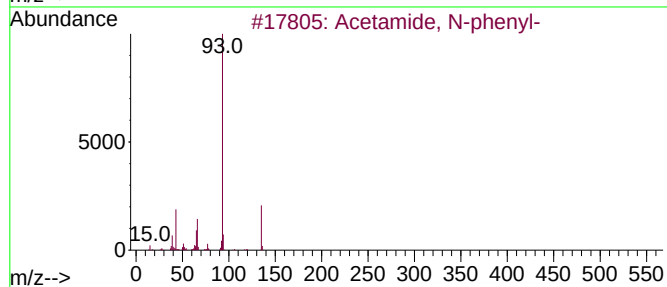
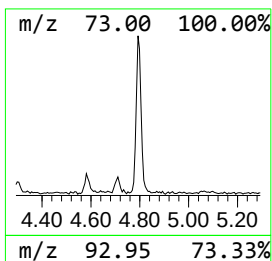
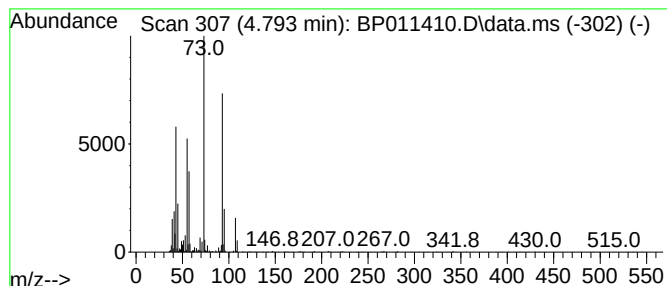
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-02 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.793	3.99 ng/ul	115537	1,4-Dichlorobenzene-d4	7.575

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetamide, N-phenyl-	135	C8H9NO	000103-84-4	17
2			Propanoic acid, 3-chloro-	108	C3H5ClO2	000107-94-8	9
3			5-Amino-4-cyano-1-(2,6-dichlorop...	338	C14H16Cl2N4Si	1000505-14-6	9
4			2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9
5			1,3-Dioxolane, 2-(chloromethyl)-	122	C4H7ClO2	002568-30-1	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
Data File : BP011410.D
Acq On : 12 Aug 2022 04:51
Operator : CG/JU
Sample : N4129-02
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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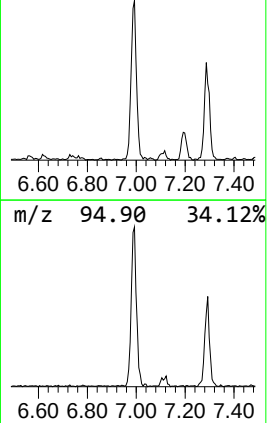
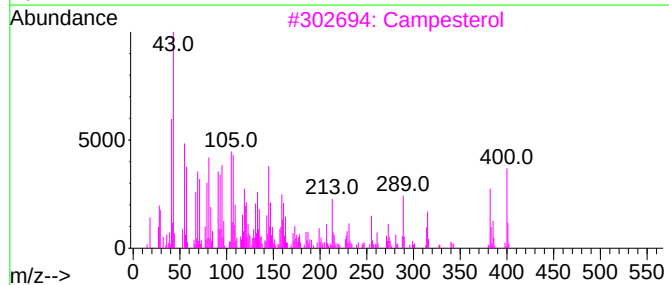
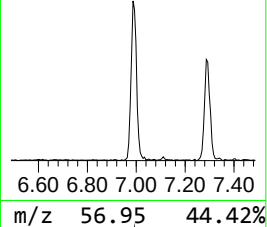
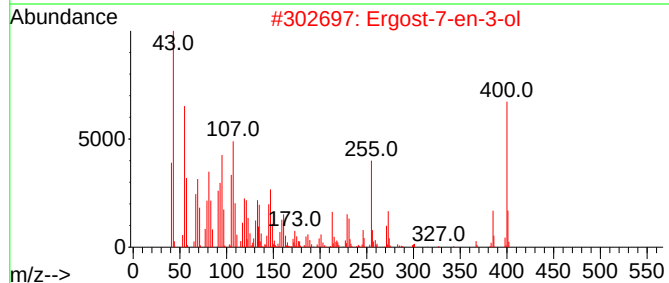
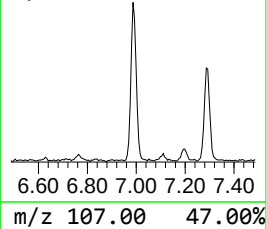
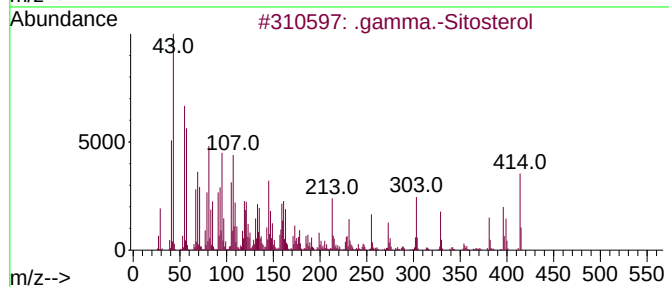
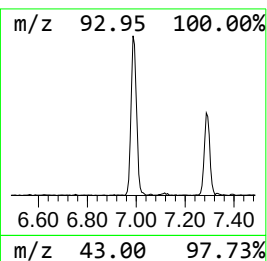
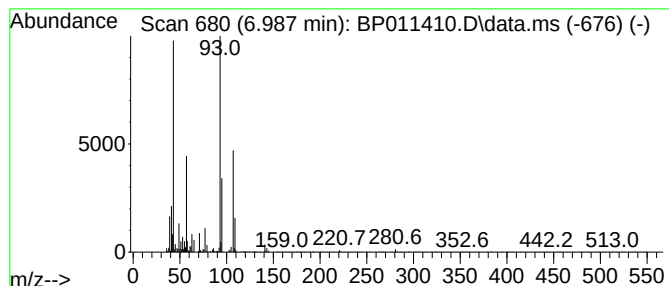
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-03 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.987	3.26 ng/ul	94496	1,4-Dichlorobenzene-d4	7.575

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	gamma.-Sitosterol	414	C29H50O	000083-47-6	10	
2	Ergost-7-en-3-ol	400	C28H48O	116179-22-7	10		
3	Campesterol	400	C28H48O	000474-62-4	10		
4	Stigmasterol, acetate	458	C31H54O2	002364-21-8	10		
5	(1S)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-26-4	9		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
Data File : BP011410.D
Acq On : 12 Aug 2022 04:51
Operator : CG/JU
Sample : N4129-02
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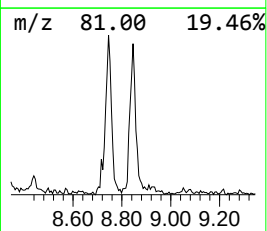
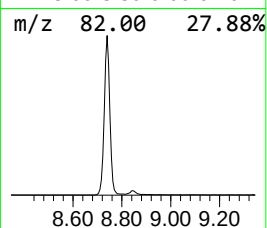
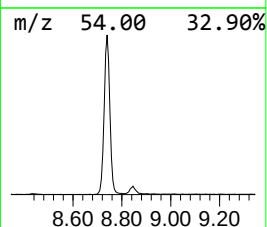
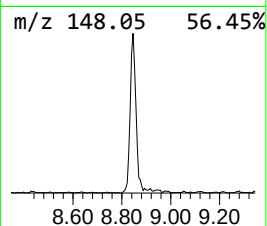
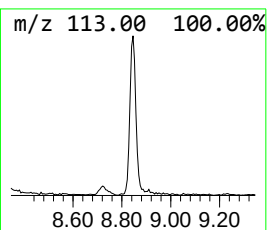
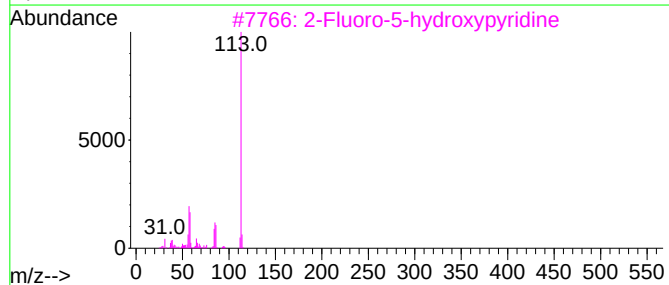
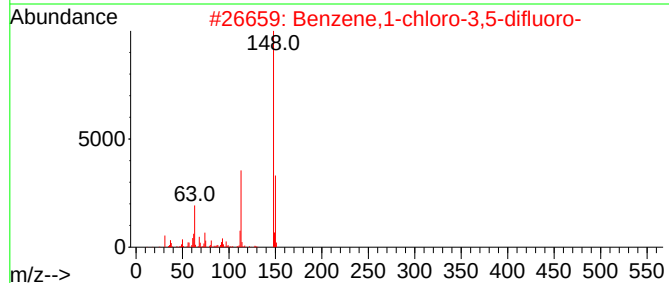
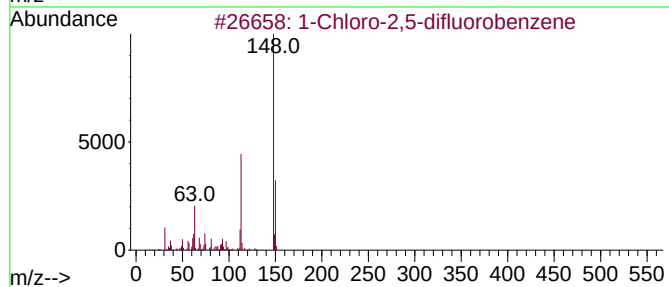
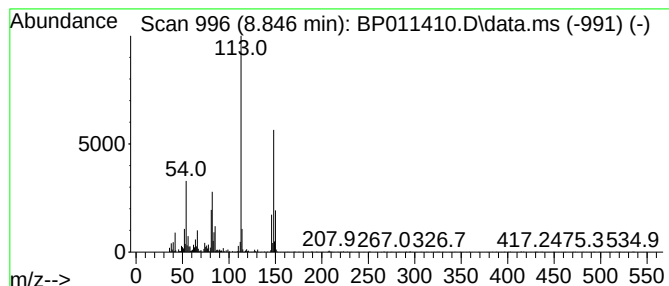
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-04 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.846	3.25 ng/ul	94059	1,4-Dichlorobenzene-d4	7.575

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Chloro-2,5-difluorobenzene	148	C6H3ClF2	002367-91-1	40
2			Benzene,1-chloro-3,5-difluoro-	148	C6H3ClF2	001435-43-4	38
3			2-Fluoro-5-hydroxypyridine	113	C5H4FNO	055758-32-2	38
4			3-Fluoro-4-pyridinol	113	C5H4FNO	022282-73-1	38
5			1-Chloro-2,4-difluorobenzene	148	C6H3ClF2	001435-44-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
Data File : BP011410.D
Acq On : 12 Aug 2022 04:51
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Sample : N4129-02
Misc :
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Instrument :
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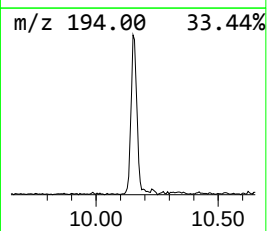
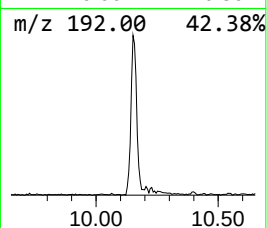
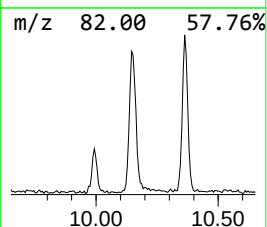
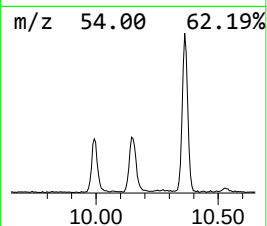
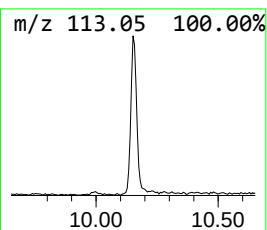
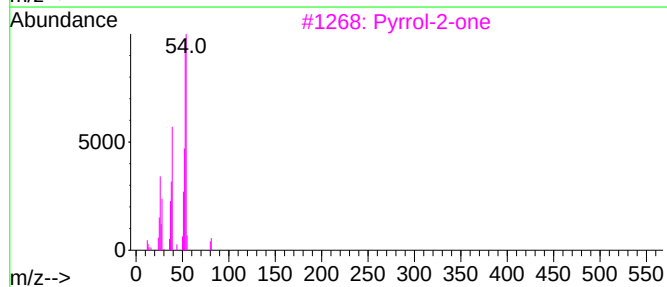
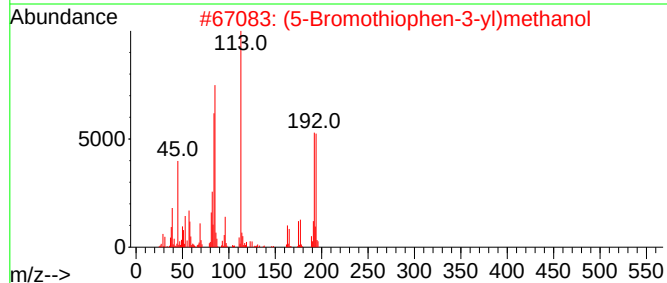
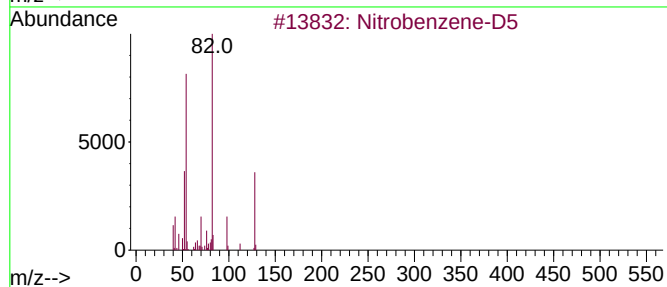
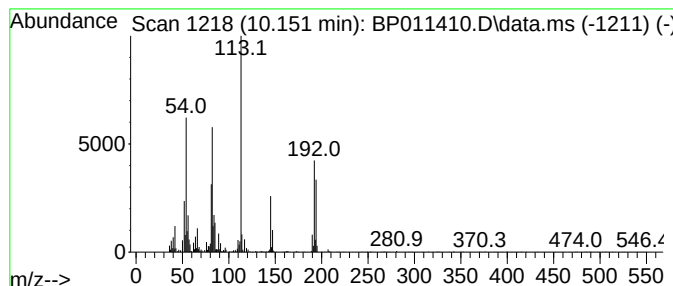
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-05 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.151	3.83 ng/ul	152622	Naphthalene-d8	10.363

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nitrobenzene-D5	128	C6D5NO2	004165-60-0	25
2		(5-Bromothiophen-3-yl)methanol	192	C5H5BrOS	073919-88-7	25
3		Pyrrol-2-one	81	C4H3NO	1010436-46-3	22
4		But-2-enedioic acid, dimethyl ester	144	C6H8O4	023055-10-9	14
5		4-Methylamino-2(5H)-furanone	113	C5H7NO2	1000427-45-5	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
Data File : BP011410.D
Acq On : 12 Aug 2022 04:51
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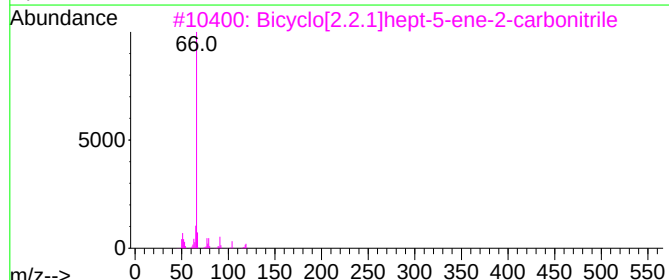
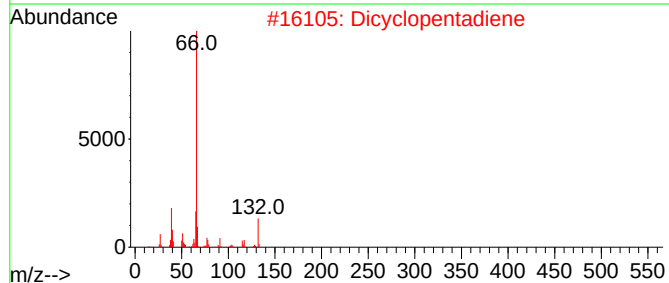
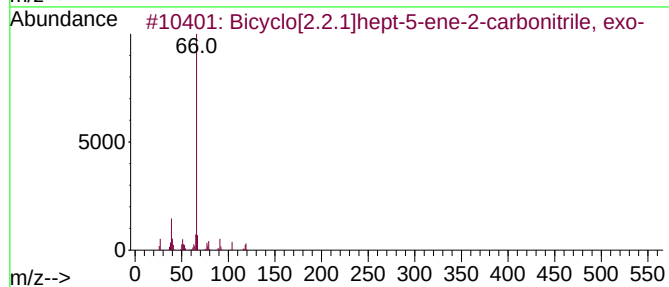
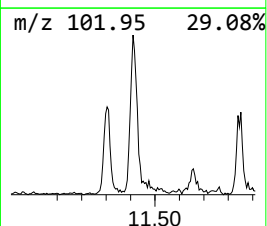
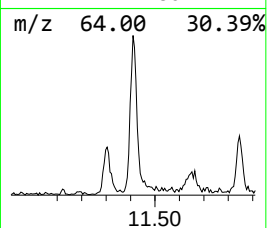
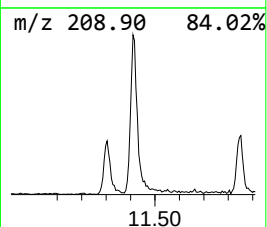
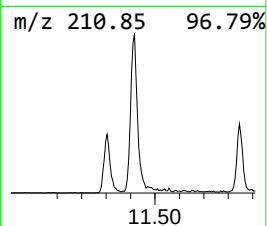
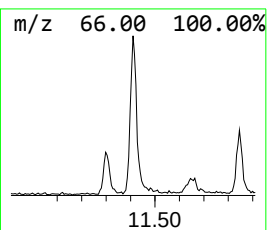
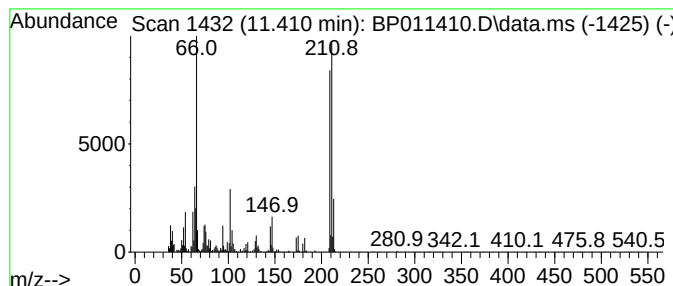
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-06 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.410	4.27 ng/ul	170318	Naphthalene-d8	10.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[2.2.1]hept-5-ene-2-carbo...	119	C8H9N	002890-96-2	46
2			Dicyclopentadiene	132	C10H12	000077-73-6	43
3			Bicyclo[2.2.1]hept-5-ene-2-carbo...	119	C8H9N	000095-11-4	43
4			4,7-Methano-1H-inden-1-one, 2,3,...	148	C10H12O	086105-40-0	43
5			2-Norbornene	94	C7H10	000498-66-8	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
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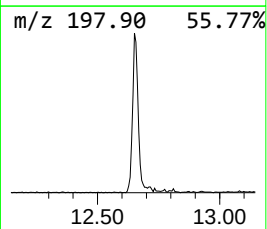
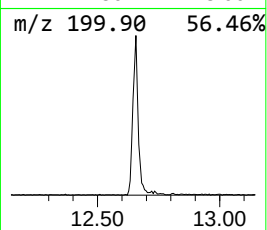
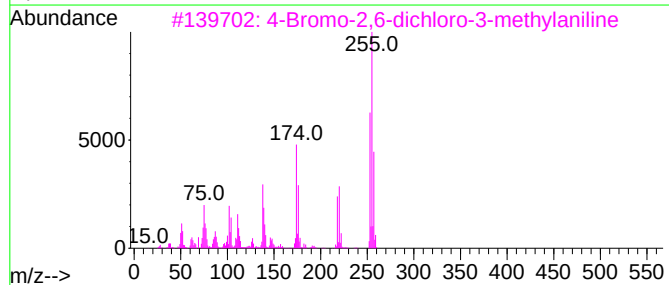
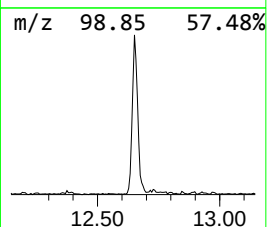
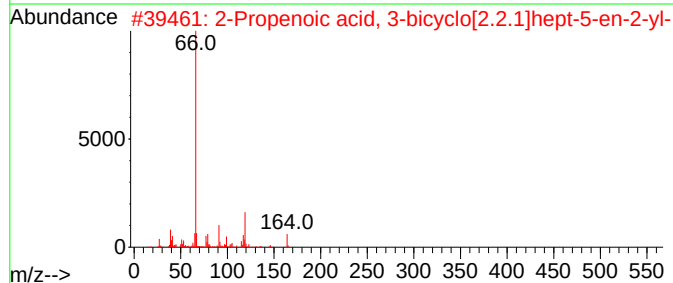
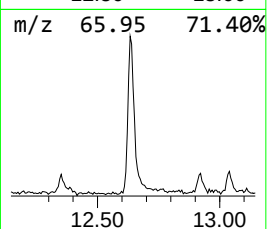
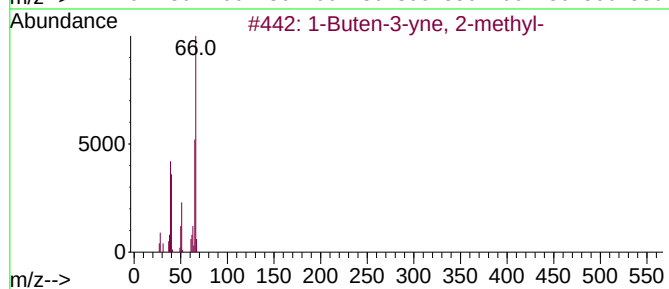
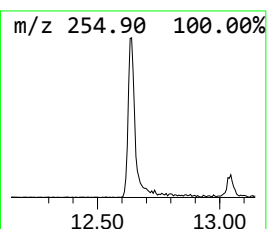
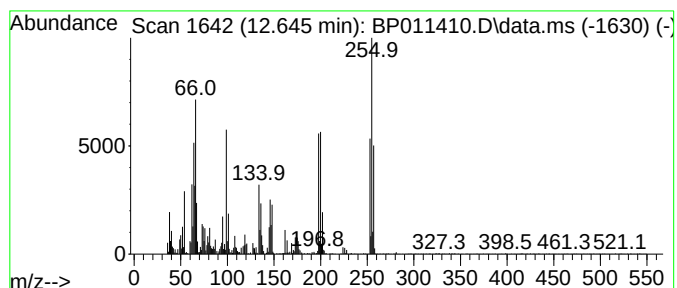
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-07 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.645	3.90 ng/ul	197376	Acenaphthene-d10	14.251	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Buten-3-yne, 2-methyl-	66	C5H6	000078-80-8	11
2	2-Propenoic acid, 3-bicyclo[2.2.1]hept-5-en-2-yl-	164	C10H12O2	015222-64-7	11
3	4-Bromo-2,6-dichloro-3-methylaniline	253	C7H6BrCl2N	062406-68-2	11
4	3-Butenenitrile, 3-chloro-	101	C4H4ClN	021031-46-9	11
5	3,4-Dibromo-5-fluoropyridine	253	C5H2Br2FN	1260843-59-1	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
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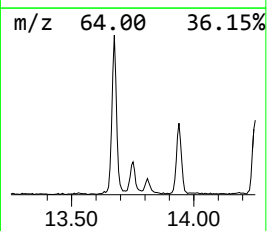
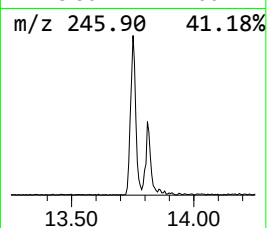
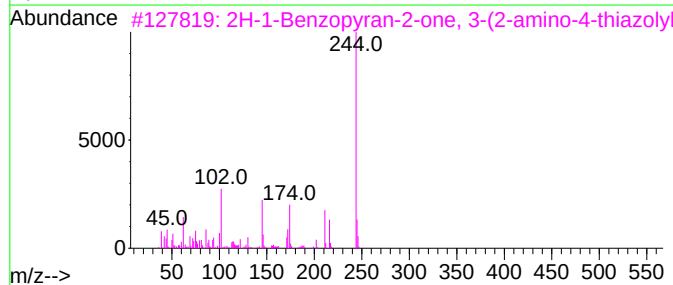
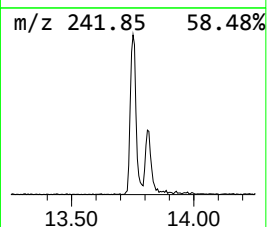
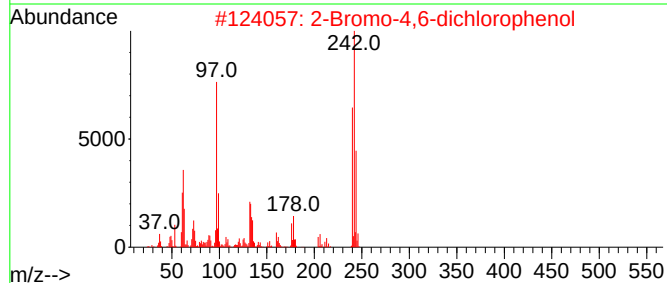
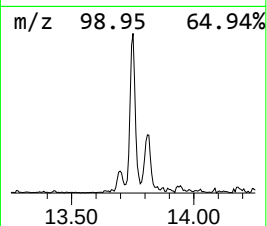
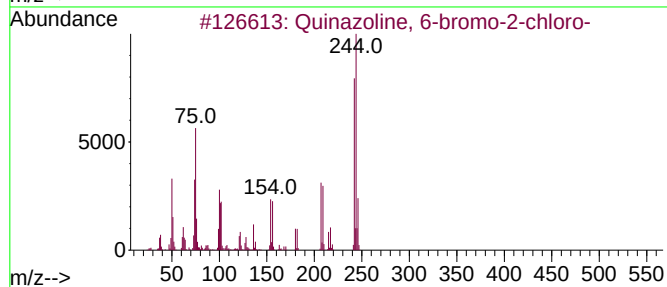
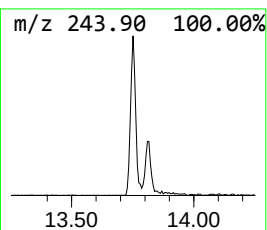
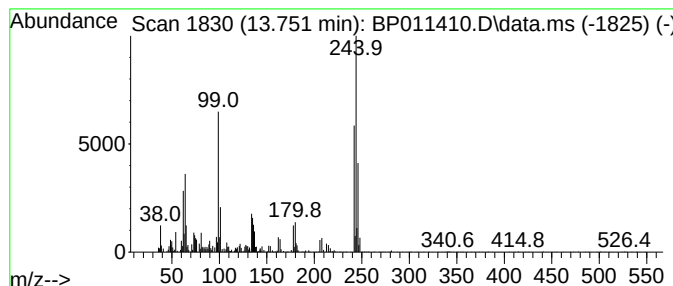
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-08 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.751	2.39 ng/ul	121236	Acenaphthene-d10	14.251

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Quinazoline, 6-bromo-2-chloro-	242	C8H4BrClN2	882672-05-1	43		
2	2-Bromo-4,6-dichlorophenol	240	C6H3BrCl2O	004524-77-0	38		
3	2H-1-Benzopyran-2-one, 3-(2-amin...	244	C12H8N2O2S	061636-28-0	27		
4	Phenazine, 2-chloro-8-methoxy-	244	C13H9ClN2O	018450-08-3	12		
5	Hexanoic acid, 2-chloroethyl ester	178	C8H15ClO2	1000330-92-7	12		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
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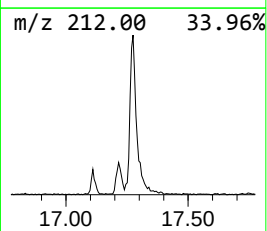
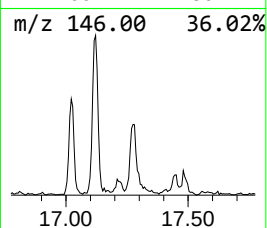
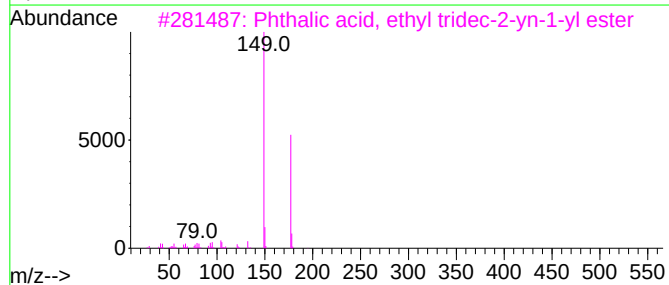
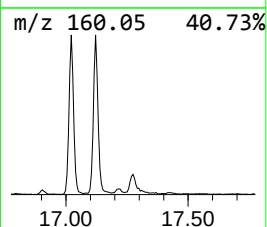
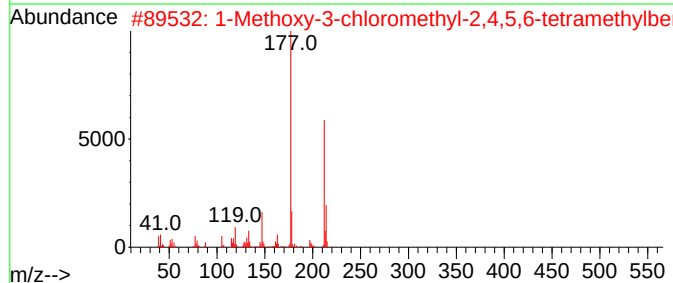
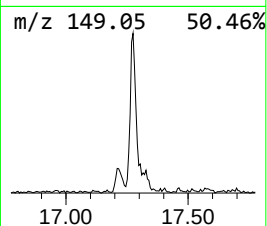
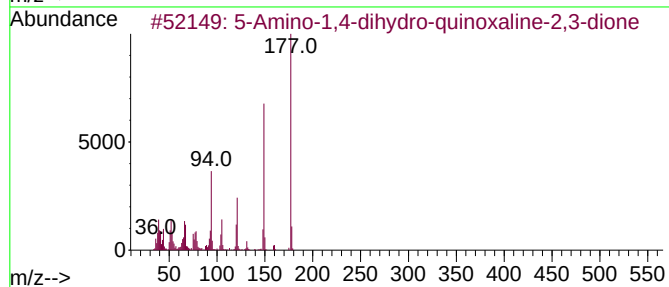
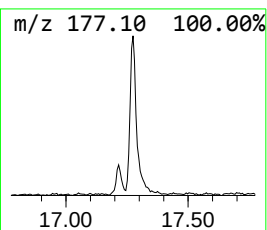
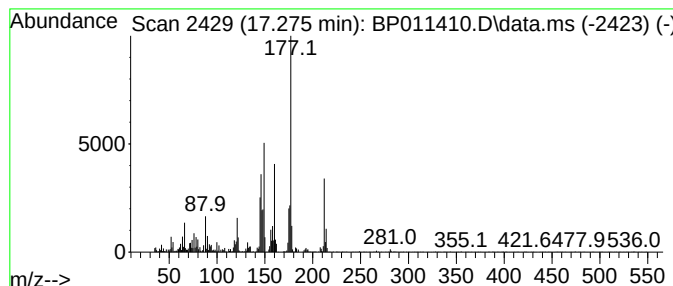
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-09 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.275	2.96 ng/ul	172627	Phenanthrene-d10	17.022

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		5-Amino-1,4-dihydro-quinoxaline-...	177	C8H7N3O2	076097-87-5	38
2		1-Methoxy-3-chloromethyl-2,4,5,6...	212	C12H17ClO	073942-80-0	27
3		Phthalic acid, ethyl tridec-2-yn...	372	C23H32O4	1000315-43-5	27
4		Benzofuran-2-one, 4-amino-2,3-di...	177	C10H11NO2	1000129-51-7	27
5		4-Pyrimidinamine, 5-(2-thienyl)-	177	C8H7N3S	058758-95-5	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
Data File : BP011410.D
Acq On : 12 Aug 2022 04:51
Operator : CG/JU
Sample : N4129-02
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGBW2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.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
unknown-01	4.299	95.0	ng/u1	2753070	1	7.575	579553	20.0
Butane, 2,3-dic...	4.581	35.4	ng/u1	1027090	1	7.575	579553	20.0
unknown-02	4.793	4.0	ng/u1	115537	1	7.575	579553	20.0
unknown-03	6.987	3.3	ng/u1	94496	1	7.575	579553	20.0
unknown-04	8.846	3.3	ng/u1	94059	1	7.575	579553	20.0
unknown-05	10.151	3.8	ng/u1	152622	2	10.363	797850	20.0
unknown-06	11.410	4.3	ng/u1	170318	2	10.363	797850	20.0
unknown-07	12.645	3.9	ng/u1	197376	3	14.251	1013140	20.0
unknown-08	13.751	2.4	ng/u1	121236	3	14.251	1013140	20.0
unknown-09	17.275	3.0	ng/u1	172627	4	17.022	1166640	20.0