

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082823\
 Data File : BP016848.D
 Acq On : 30 Aug 2023 00:11
 Operator : MA/JU
 Sample : 04148-08
 Misc :
 ALS Vial : 66 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJX2

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-BP082423.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP016848.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.334	3	8	20	rBV	3662646	6178789	39.36%	3.990%
2	3.417	20	22	37	rVB	216053	326780	2.08%	0.211%
3	3.764	77	81	87	rBV	212636	279756	1.78%	0.181%
4	3.817	87	90	106	rVB	548524	880270	5.61%	0.568%
5	4.017	119	124	129	rVB	47569	76742	0.49%	0.050%
6	4.134	140	144	150	rVB	30069	48003	0.31%	0.031%
7	4.205	152	156	162	rVB	55508	71982	0.46%	0.046%
8	4.658	226	233	250	rBV	10048658	15699311	100.00%	10.137%
9	4.805	254	258	264	rVB	21402	34868	0.22%	0.023%
10	5.069	298	303	309	rBV	19594	28989	0.18%	0.019%
11	5.293	336	341	352	rVB	198400	306403	1.95%	0.198%
12	5.993	449	460	468	rBV	343447	524048	3.34%	0.338%
13	6.463	536	540	548	rVB2	18002	32171	0.20%	0.021%
14	6.711	578	582	593	rVB	76401	114732	0.73%	0.074%
15	7.169	654	660	673	rBV	562915	894997	5.70%	0.578%
16	7.363	686	693	701	rBV	1998242	3051634	19.44%	1.971%
17	7.458	703	709	714	rVB2	31471	55146	0.35%	0.036%
18	7.546	714	724	737	rBV	1873195	2971896	18.93%	1.919%
19	8.022	798	805	814	rBV	1552540	2523118	16.07%	1.629%
20	8.375	856	865	873	rVV	350967	579891	3.69%	0.374%
21	8.728	917	925	935	rBV	1590759	2542612	16.20%	1.642%
22	8.816	936	940	941	rVV2	25235	33647	0.21%	0.022%
23	8.846	942	945	954	rVB4	28517	55427	0.35%	0.036%
24	8.946	956	962	968	rVV	263465	419831	2.67%	0.271%
25	9.140	991	995	1001	rVV2	31080	53272	0.34%	0.034%
26	9.222	1001	1009	1021	rVV	2353326	3846600	24.50%	2.484%
27	9.322	1023	1026	1031	rVV7	15838	35685	0.23%	0.023%
28	9.387	1032	1037	1042	rVV	47691	85580	0.55%	0.055%
29	9.446	1042	1047	1059	rVV5	40009	97167	0.62%	0.063%
30	9.669	1080	1085	1092	rVB2	15869	33110	0.21%	0.021%
31	9.828	1107	1112	1117	rBV5	17072	31238	0.20%	0.020%
32	9.940	1124	1131	1139	rBV	1803369	2938907	18.72%	1.898%
33	10.140	1157	1165	1172	rVB7	20488	52727	0.34%	0.034%
34	10.210	1172	1177	1186	rBV7	21622	44243	0.28%	0.029%
35	10.322	1188	1196	1208	rVB2	122117	262894	1.67%	0.170%
36	10.463	1212	1220	1230	rBV	2649922	4498659	28.66%	2.905%

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37	10.851	1274	1286	1296	rBV	2456976	4235722	26.98%	2.735%
38	11.010	1304	1313	1320	rBV	1145273	2011979	12.82%	1.299%
39	11.122	1326	1332	1341	rVB2	398197	693374	4.42%	0.448%
40	11.210	1341	1347	1360	rVB2	72972	147027	0.94%	0.095%
41	11.375	1367	1375	1380	rBV6	46889	111395	0.71%	0.072%
42	11.551	1400	1405	1409	rVB6	20052	31408	0.20%	0.020%
43	11.616	1409	1416	1417	rBV2	97188	144136	0.92%	0.093%
44	11.646	1417	1421	1429	rVV	164216	327006	2.08%	0.211%
45	11.716	1429	1433	1437	rVV2	68976	118840	0.76%	0.077%
46	11.763	1437	1441	1453	rVB3	91254	221658	1.41%	0.143%
47	11.987	1473	1479	1486	rVB	250194	366154	2.33%	0.236%
48	12.128	1499	1503	1509	rVB3	18511	29435	0.19%	0.019%
49	12.251	1518	1524	1525	rVV2	180656	265328	1.69%	0.171%
50	12.275	1525	1528	1535	rVV	285413	459273	2.93%	0.297%
51	12.375	1539	1545	1549	rVV	219452	340561	2.17%	0.220%
52	12.428	1549	1554	1563	rVB3	70868	150734	0.96%	0.097%
53	12.575	1573	1579	1585	rVB3	82245	144818	0.92%	0.094%
54	12.722	1600	1604	1615	rVB3	18096	51905	0.33%	0.034%
55	12.840	1619	1624	1629	rBV2	35771	58715	0.37%	0.038%
56	13.375	1708	1715	1721	rBV2	21114	58950	0.38%	0.038%
57	13.522	1731	1740	1746	rBV	323089	530216	3.38%	0.342%
58	13.663	1754	1764	1772	rBV	22384	78247	0.50%	0.051%
59	13.763	1772	1781	1786	rVV	376657	569997	3.63%	0.368%
60	13.810	1786	1789	1796	rVB6	41150	69607	0.44%	0.045%
61	13.975	1812	1817	1822	rBV	117684	177365	1.13%	0.115%
62	14.092	1830	1837	1843	rBV	5536848	7632957	48.62%	4.929%
63	14.228	1856	1860	1863	rBV2	34282	42181	0.27%	0.027%
64	14.381	1877	1886	1892	rBV	6007304	8924795	56.85%	5.763%
65	14.451	1892	1898	1906	rVV3	171909	408472	2.60%	0.264%
66	14.528	1907	1911	1916	rVV	113440	157412	1.00%	0.102%
67	14.592	1918	1922	1930	rVB	129205	229656	1.46%	0.148%
68	14.681	1930	1937	1943	rBV	3541110	5199204	33.12%	3.357%
69	14.887	1967	1972	1984	rBV	403137	720732	4.59%	0.465%
70	15.063	1999	2002	2007	rVB6	31684	45495	0.29%	0.029%
71	15.216	2023	2028	2035	rVB	133792	197101	1.26%	0.127%
72	15.381	2052	2056	2058	rBV3	33177	41106	0.26%	0.027%
73	15.481	2069	2073	2078	rVV2	142721	211388	1.35%	0.136%
74	15.539	2079	2083	2086	rVB3	47328	68254	0.43%	0.044%
75	15.675	2099	2106	2114	rBV	8062710	11369255	72.42%	7.341%
76	15.792	2121	2126	2136	rBV	2408837	3276813	20.87%	2.116%

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77	15.939	2147	2151	2158	rVB2	191993	294253	1.87%	0.190%
78	16.016	2160	2164	2168	rBV5	21156	38942	0.25%	0.025%
79	16.204	2191	2196	2201	rVB	205643	273465	1.74%	0.177%
80	16.257	2202	2205	2208	rVV5	24347	32303	0.21%	0.021%
81	16.298	2208	2212	2215	rVV	143615	213177	1.36%	0.138%
82	16.334	2215	2218	2223	rVV	248700	357372	2.28%	0.231%
83	16.404	2226	2230	2236	rVB	152213	215772	1.37%	0.139%
84	16.722	2276	2284	2293	rBV2	79588	203815	1.30%	0.132%
85	16.939	2317	2321	2325	rVB3	75596	103129	0.66%	0.067%
86	17.175	2357	2361	2366	rBV	248125	325142	2.07%	0.210%
87	17.286	2377	2380	2389	rVB7	33095	69958	0.45%	0.045%
88	17.445	2401	2407	2417	rBV	4428667	6152175	39.19%	3.973%
89	17.545	2418	2424	2434	rVV	8807918	12744612	81.18%	8.229%
90	18.075	2511	2514	2519	rVB4	86373	126463	0.81%	0.082%
91	18.280	2545	2549	2560	rBV	318996	506306	3.23%	0.327%
92	19.098	2685	2688	2694	rVB2	64127	91952	0.59%	0.059%
93	19.351	2727	2731	2736	rBV2	114050	142470	0.91%	0.092%
94	19.416	2739	2742	2752	rVB	111070	158209	1.01%	0.102%
95	19.516	2755	2759	2768	rBV	215811	334256	2.13%	0.216%
96	19.780	2798	2804	2810	rBV	10621116	13745612	87.56%	8.876%
97	20.063	2849	2852	2856	rBV2	64833	74777	0.48%	0.048%
98	21.539	3098	3103	3111	rBV	3813942	4998843	31.84%	3.228%
99	23.939	3503	3511	3526	rVB2	4534207	9741866	62.05%	6.291%
100	24.110	3533	3540	3551	rVB	1959022	4290706	27.33%	2.771%

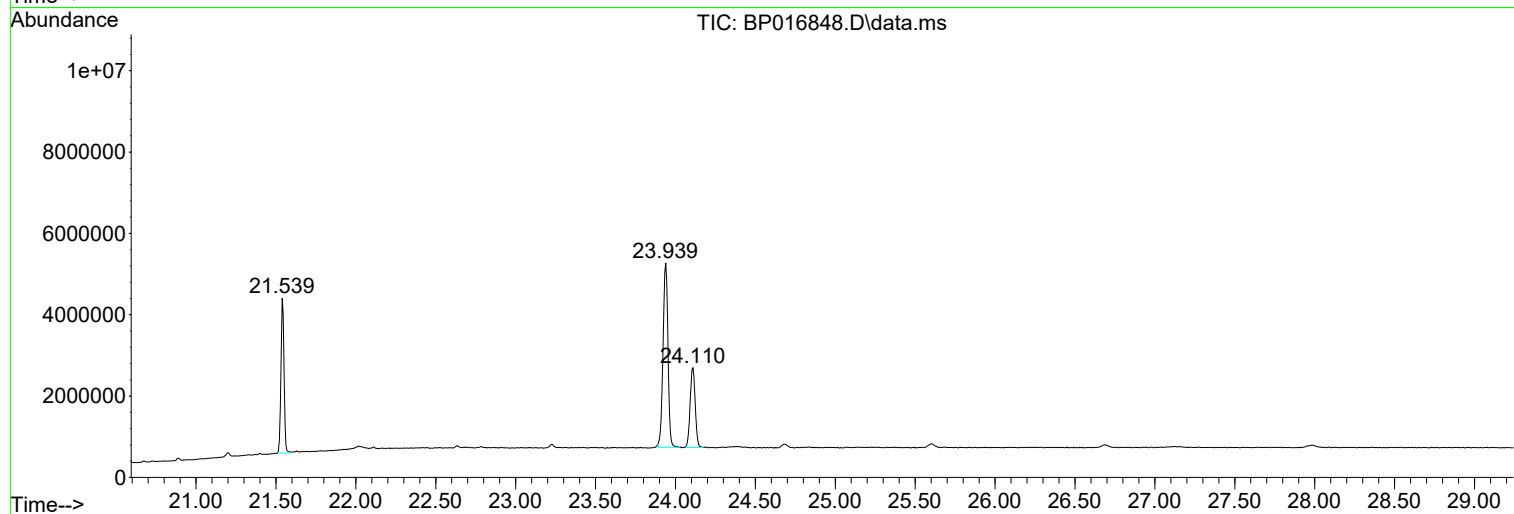
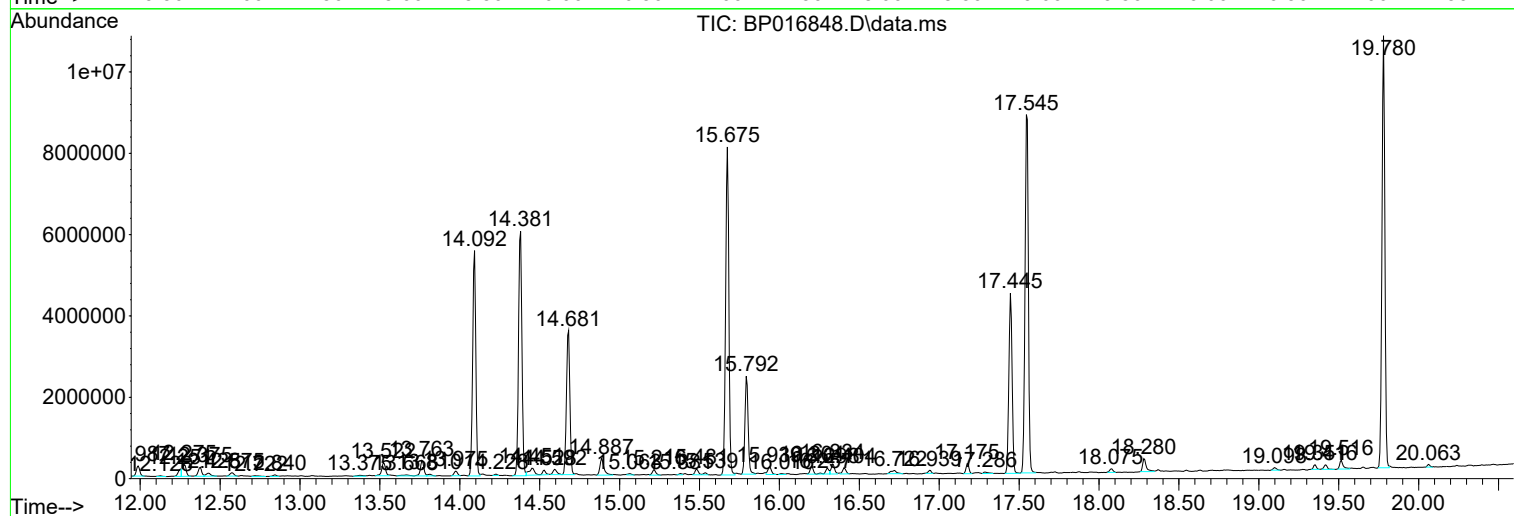
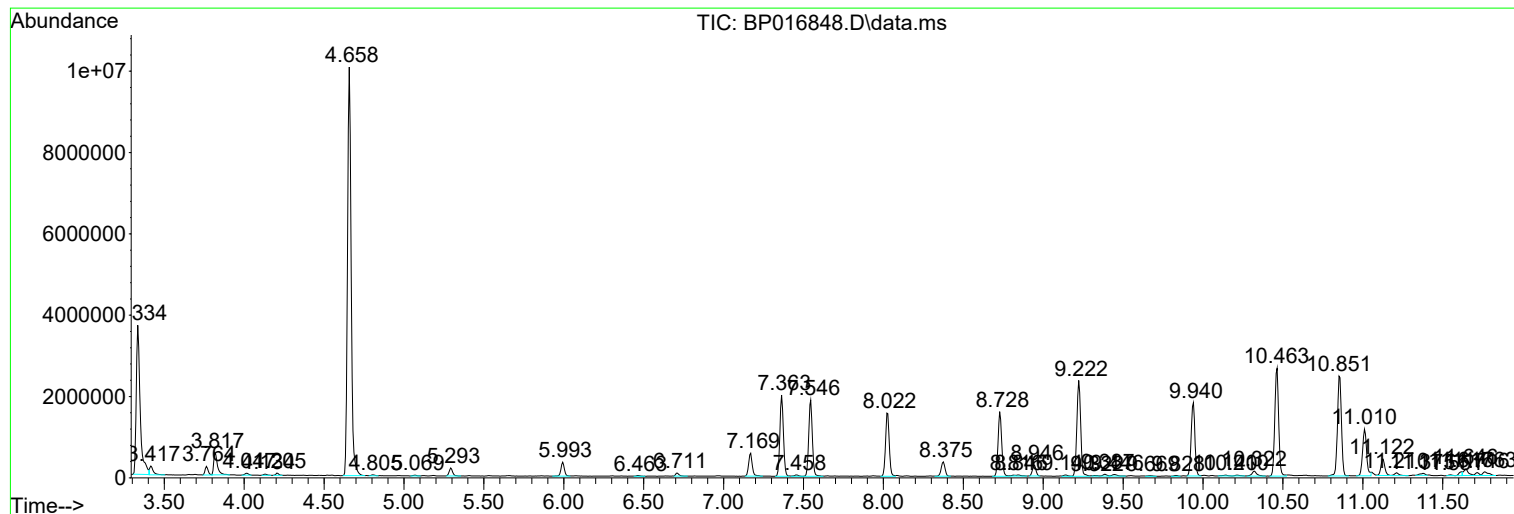
Sum of corrected areas: 154865371

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082423.MA.M
Quant Title : SVOA CALIBRATION

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



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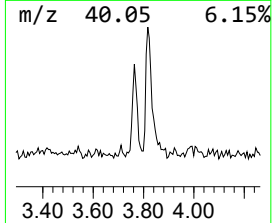
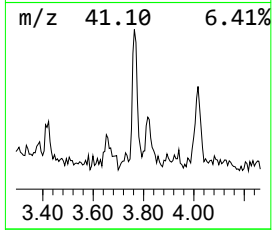
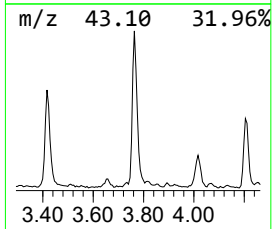
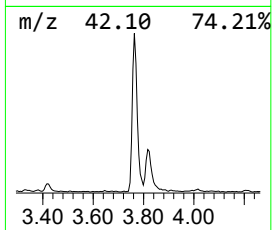
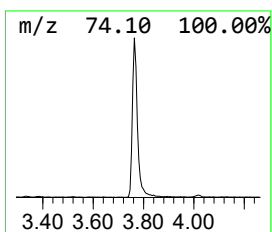
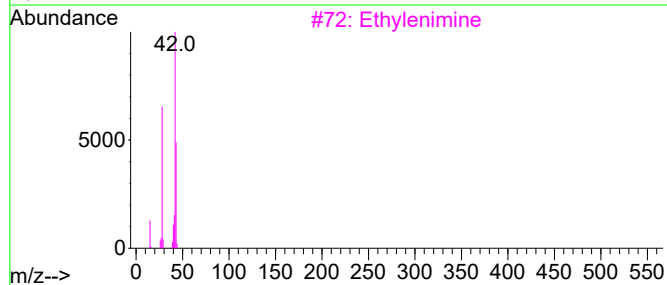
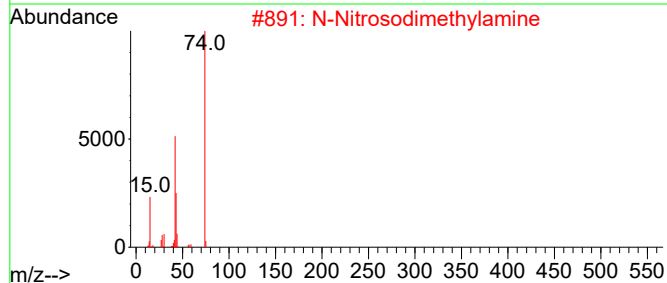
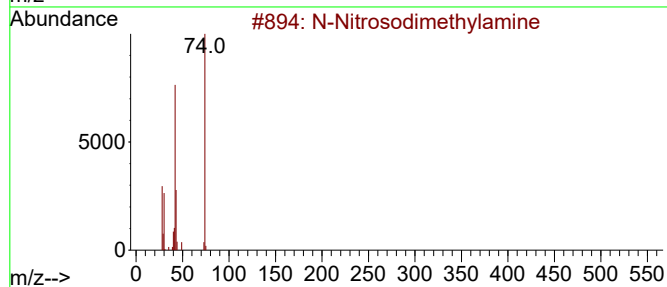
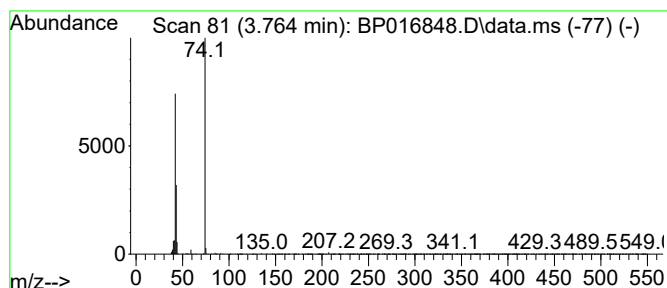
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082423.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 N-Nitrosodimethylamine Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.764	2.22 ng/ul	279756	1,4-Dichlorobenzene-d4	8.028

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N-Nitrosodimethylamine	74	C2H6N2O	000062-75-9	78
2			N-Nitrosodimethylamine	74	C2H6N2O	000062-75-9	45
3			Ethylenimine	43	C2H5N	000151-56-4	4
4			N-Nitrosodimethylamine	74	C2H6N2O	000062-75-9	4
5			Aminoguanidine	74	CH6N4	000079-17-4	4



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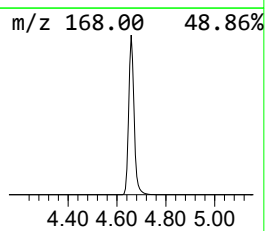
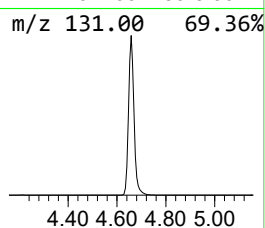
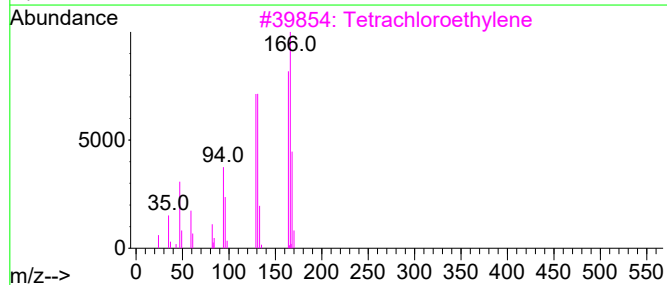
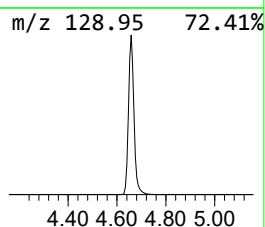
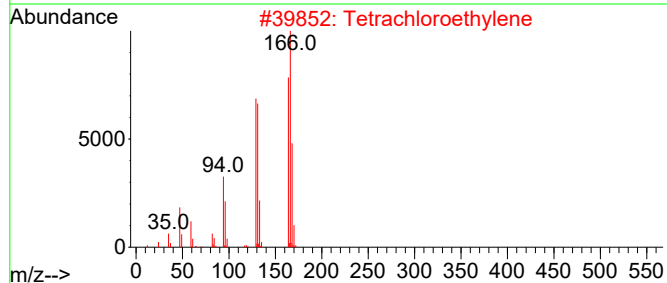
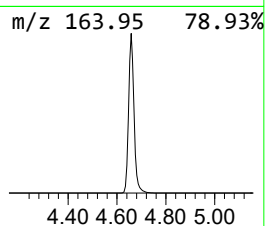
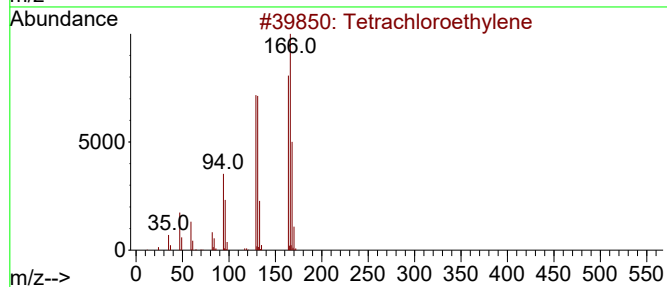
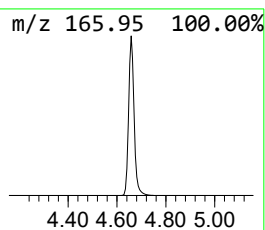
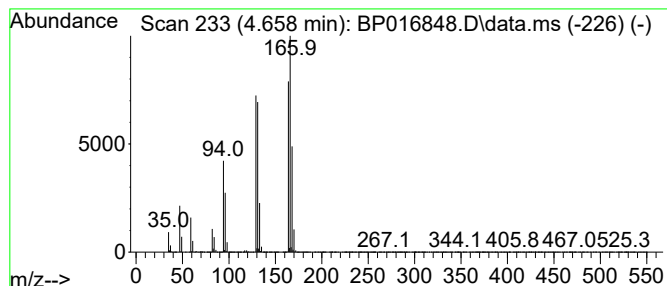
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082423.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Tetrachloroethylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.658	124.44 ng/ul	15699300	1,4-Dichlorobenzene-d4	8.028

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrachloroethylene	164	C2Cl4	000127-18-4	99
2			Tetrachloroethylene	164	C2Cl4	000127-18-4	99
3			Tetrachloroethylene	164	C2Cl4	000127-18-4	98
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	95
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HC12FN2	002927-71-1	38



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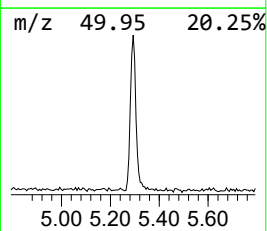
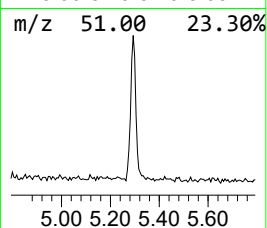
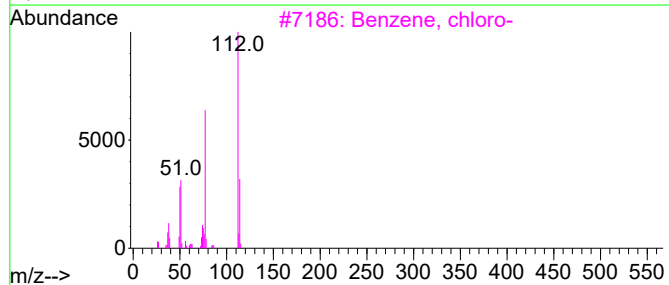
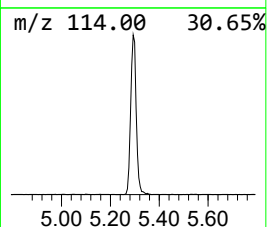
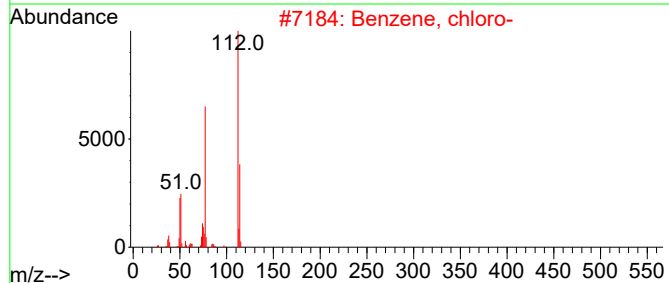
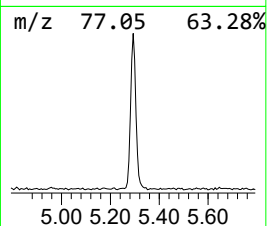
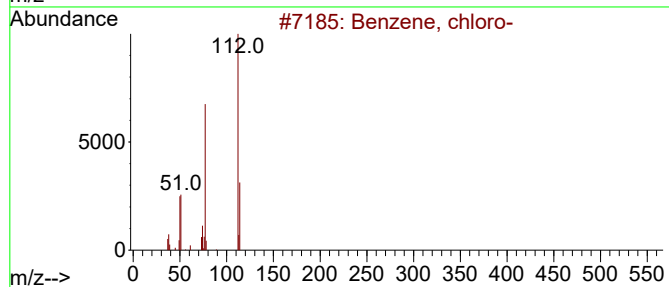
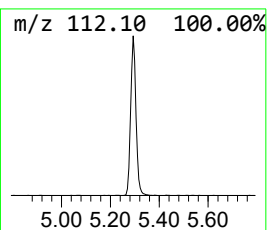
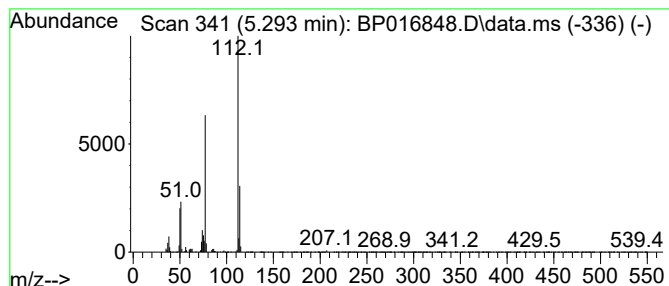
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082423.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, chloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.293	2.43 ng/ul	306403	1,4-Dichlorobenzene-d4	8.028

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, chloro-	112	C6H5Cl	000108-90-7	96
2			Benzene, chloro-	112	C6H5Cl	000108-90-7	94
3			Benzene, chloro-	112	C6H5Cl	000108-90-7	94
4			Benzene, chloro-	112	C6H5Cl	000108-90-7	94
5			Benzene, chloro-	112	C6H5Cl	000108-90-7	90



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Data File : BP016848.D
Acq On : 30 Aug 2023 00:11
Operator : MA/JU
Sample : 04148-08
Misc :
ALS Vial : 66 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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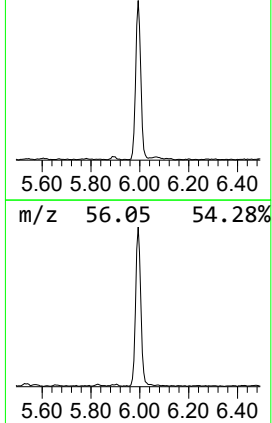
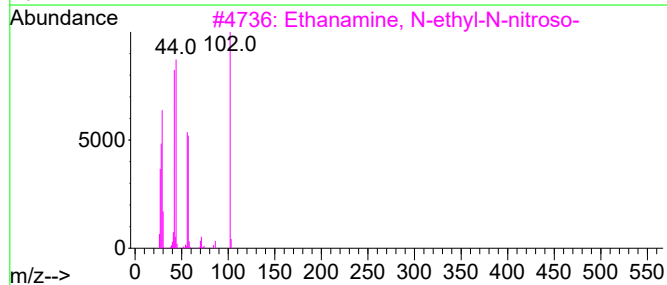
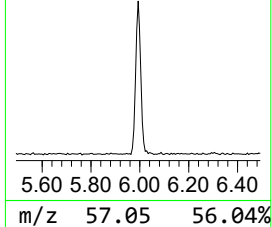
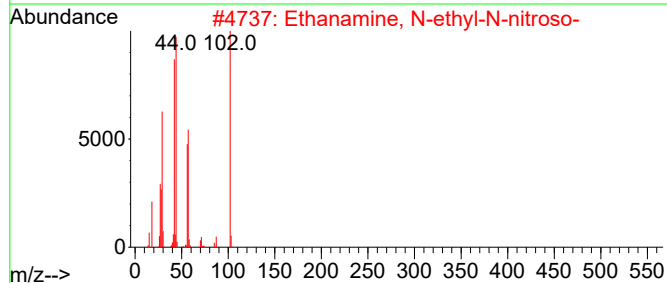
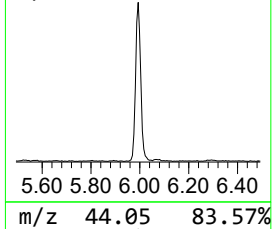
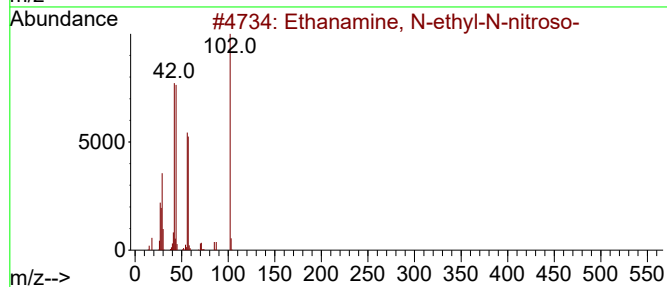
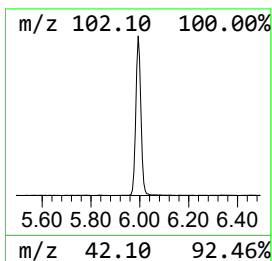
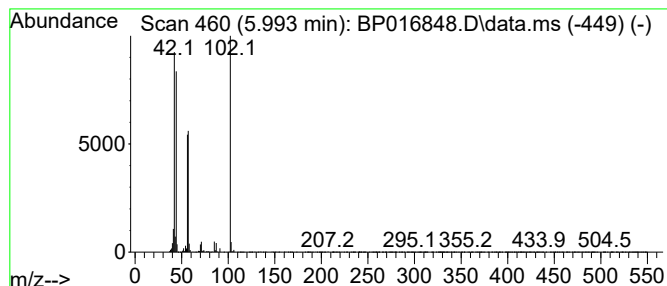
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Ethanamine, N-ethyl-N-nitroso- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.993	4.15 ng/ul	524048	1,4-Dichlorobenzene-d4	8.028

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanamine, N-ethyl-N-nitroso-	102	C4H10N2O	000055-18-5	94
2			Ethanamine, N-ethyl-N-nitroso-	102	C4H10N2O	000055-18-5	91
3			Ethanamine, N-ethyl-N-nitroso-	102	C4H10N2O	000055-18-5	91
4			Ethanamine, N-ethyl-N-nitroso-	102	C4H10N2O	000055-18-5	91
5			2-Propanamine, N-methyl-N-nitroso-	102	C4H10N2O	030533-08-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082823\
Data File : BP016848.D
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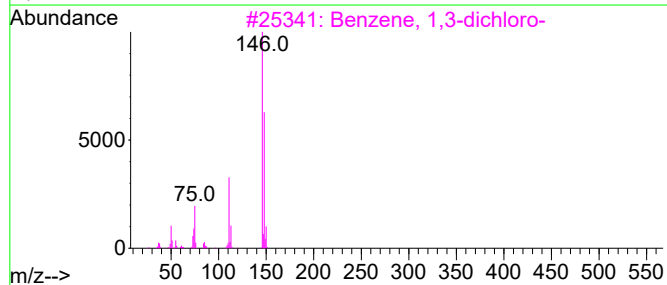
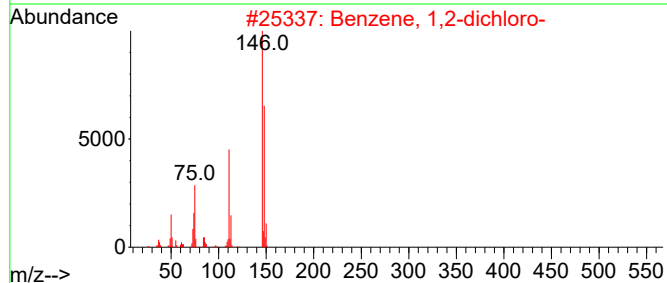
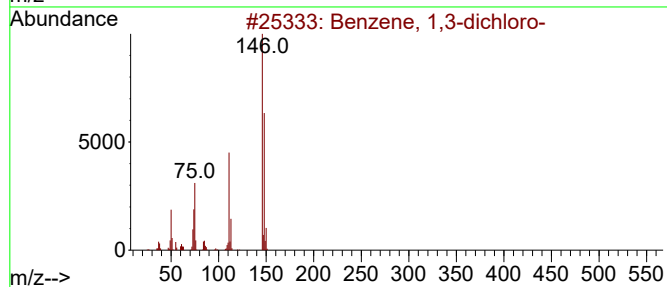
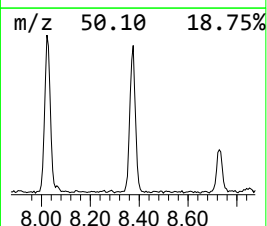
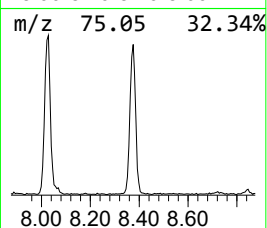
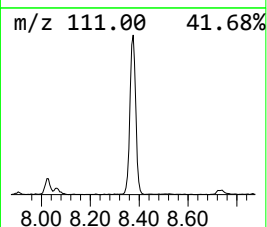
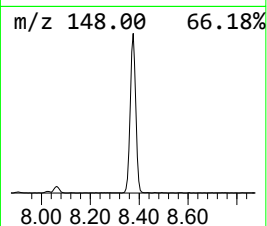
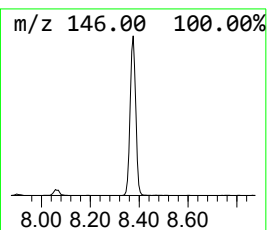
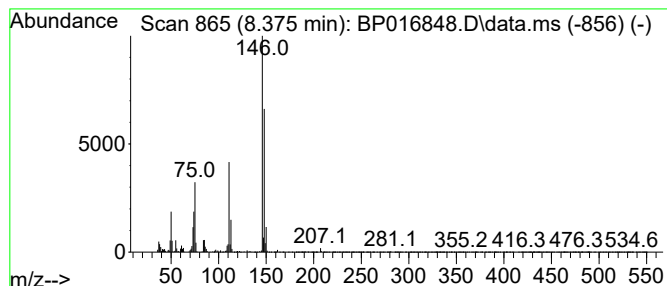
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1,3-dichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.375	4.60 ng/ul	579891	1,4-Dichlorobenzene-d4	8.028

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
2			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
3			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
4			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
5			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96



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Data File : BP016848.D
Acq On : 30 Aug 2023 00:11
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Sample : 04148-08
Misc :
ALS Vial : 66 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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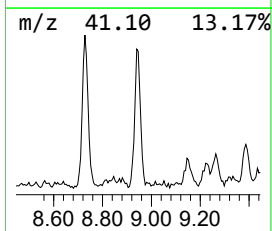
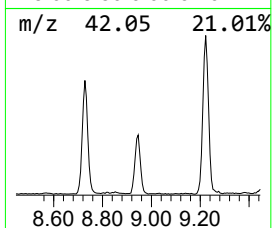
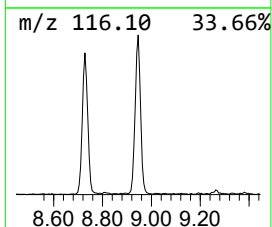
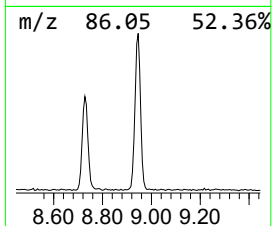
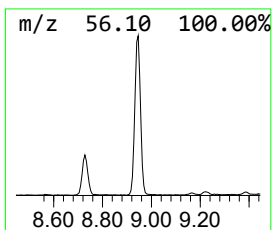
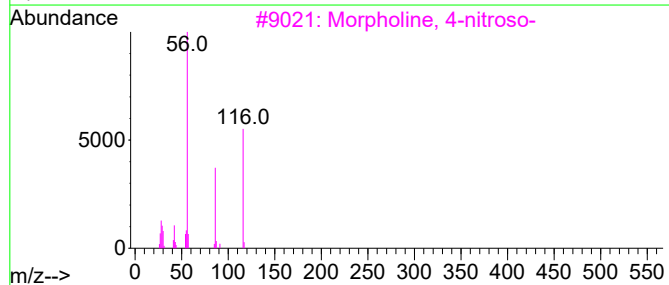
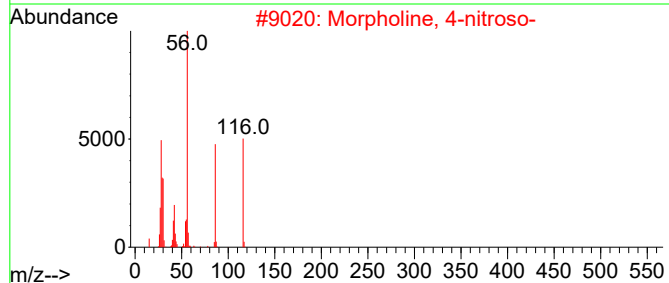
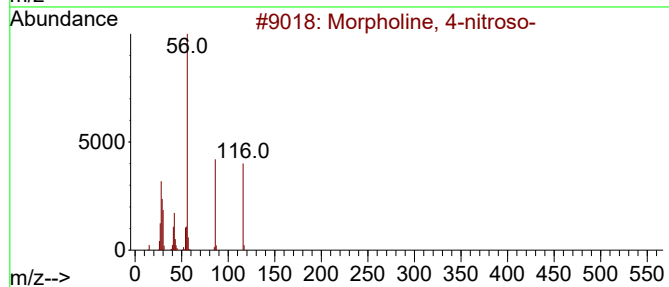
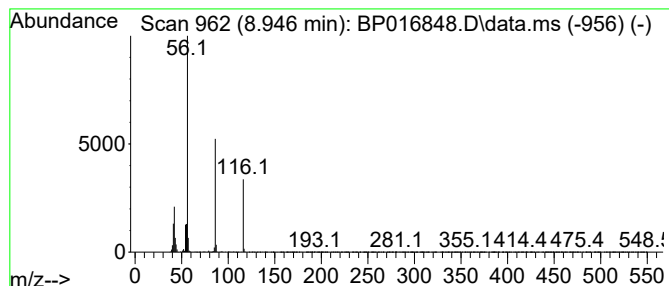
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Morpholine, 4-nitroso- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.946	3.33 ng/ul	419831	1,4-Dichlorobenzene-d4	8.028

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Morpholine, 4-nitroso-	116	C4H8N2O2	000059-89-2	91
2			Morpholine, 4-nitroso-	116	C4H8N2O2	000059-89-2	78
3			Morpholine, 4-nitroso-	116	C4H8N2O2	000059-89-2	72
4			Formic acid, hexyl ester	130	C7H14O2	000629-33-4	28
5			Bis(S-2-diethylaminoethyl) methy...	326	C13H31N2O2PS2	092030-06-3	17



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082823\
Data File : BP016848.D
Acq On : 30 Aug 2023 00:11
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Sample : 04148-08
Misc :
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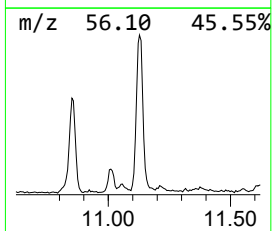
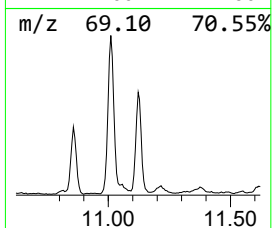
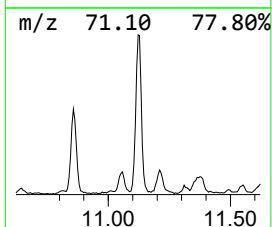
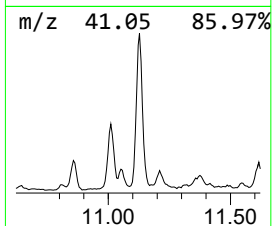
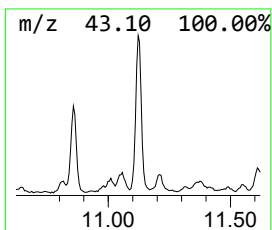
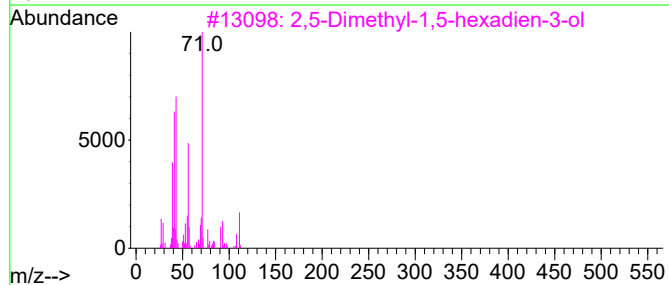
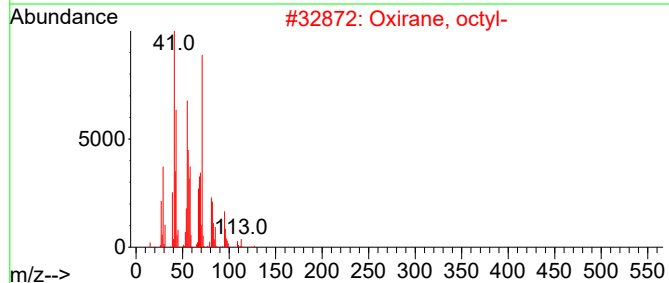
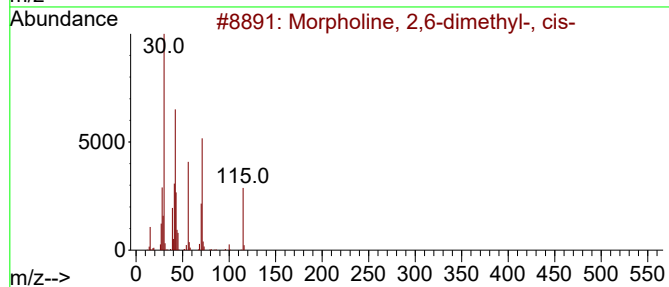
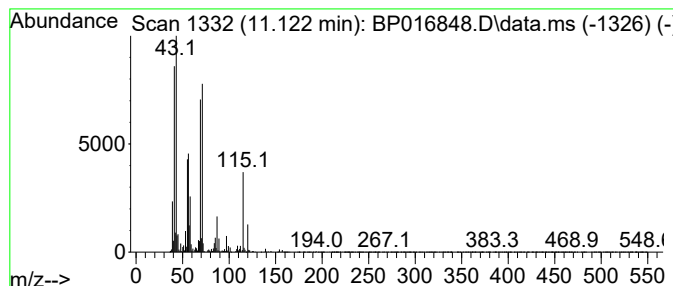
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Morpholine, 2,6-dimethyl-, ... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.122	3.27 ng/ul	693374	Naphthalene-d8	10.852

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Morpholine, 2,6-dimethyl-, cis-	115	C6H13NO	006485-55-8	53
2			Oxirane, octyl-	156	C10H20O	002404-44-6	53
3			2,5-Dimethyl-1,5-hexadien-3-ol	126	C8H14O	017123-63-6	47
4			Propanoic acid, 2-methyl-, 2-met...	144	C8H16O2	000097-85-8	47
5			Oxirane, butyl-	100	C6H12O	001436-34-6	43



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Data File : BP016848.D
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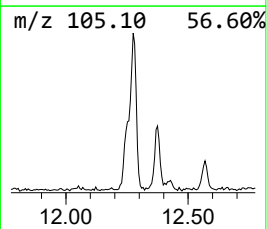
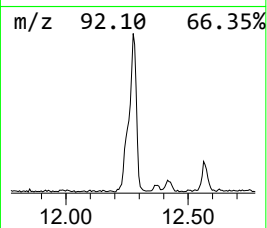
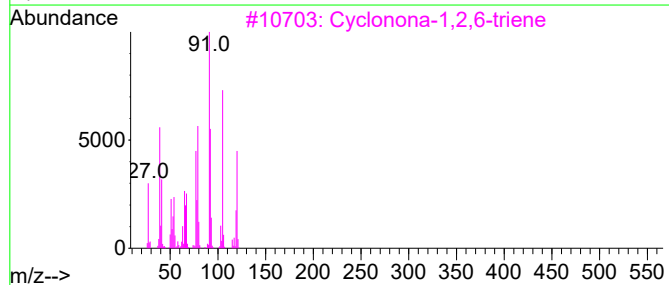
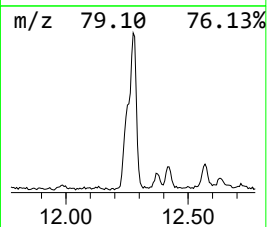
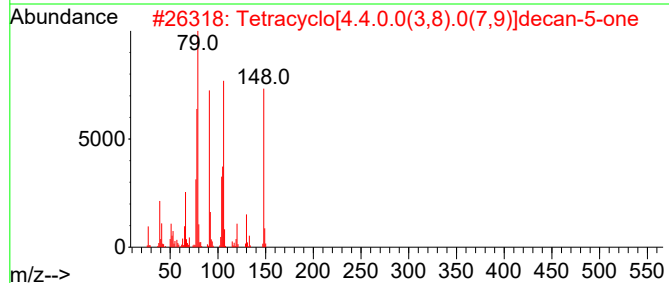
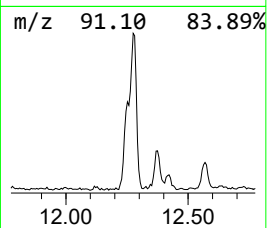
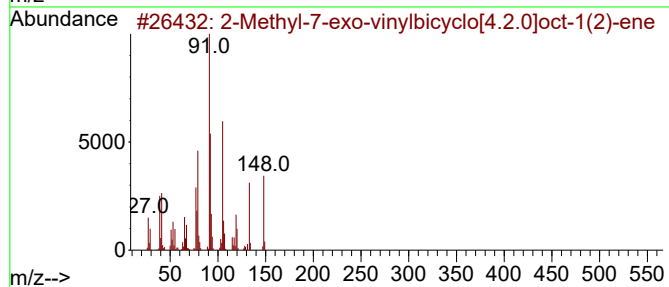
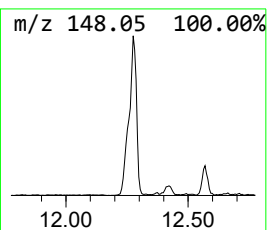
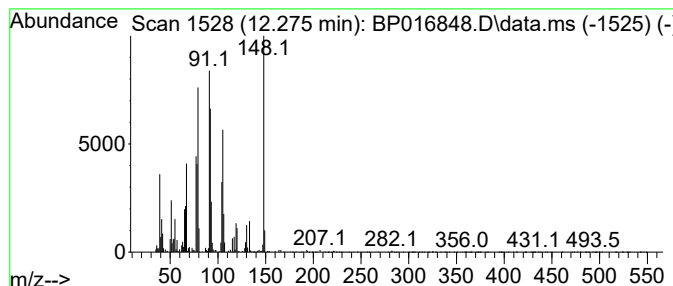
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 2-Methyl-7-exo-vinylbicyclo... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.275	2.17 ng/ul	459273	Naphthalene-d8	10.852

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Methyl-7-exo-vinylbicyclo[4.2....	148	C11H16	107914-89-6	70
2			Tetracyclo[4.4.0.0(3,8).0(7,9)]d...	148	C10H12O	1000190-20-7	64
3			Cyclonona-1,2,6-triene	120	C9H12	1000196-99-9	49
4			Tricyclo[4.2.1.1(2,5)]dec-3-en-9...	148	C10H12O	067009-89-6	47
5			3,5-Diethenylcyclopent-1-ene	120	C9H12	1000436-86-5	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082823\
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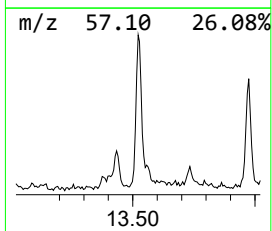
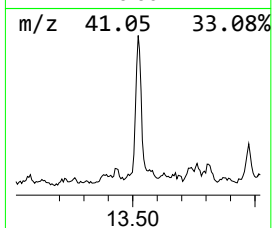
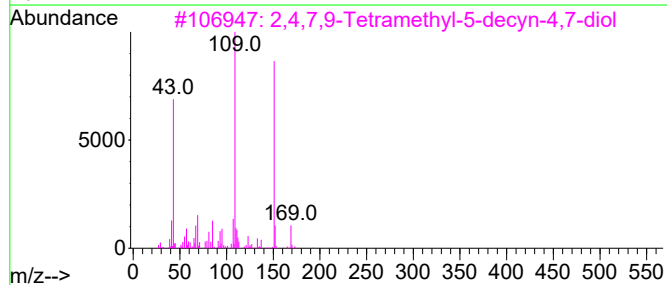
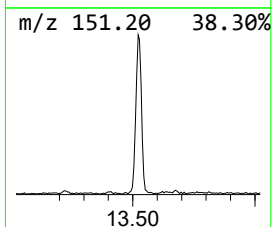
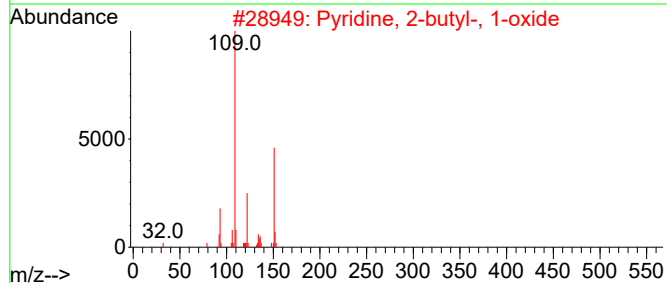
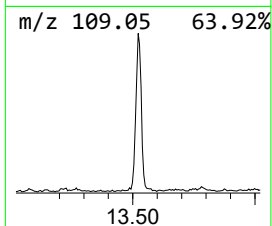
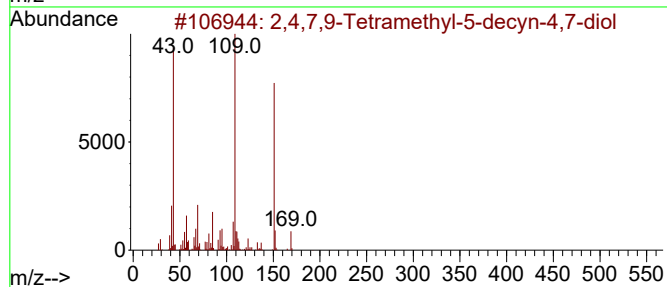
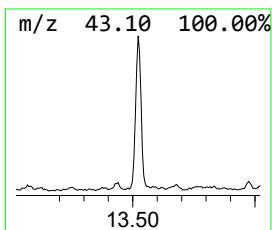
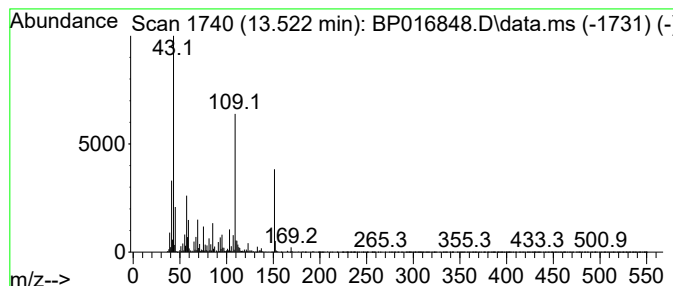
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.522	2.04 ng/ul	530216	Acenaphthene-d10	14.681	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	58
2	Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	53
3	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	50
4	3-Pyridinol, 4-methyl-, acetate ...	151	C8H9NO2	001006-96-8	47
5	Metacetamol	151	C8H9NO2	000621-42-1	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082823\
Data File : BP016848.D
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Misc :
ALS Vial : 66 Sample Multiplier: 1

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ClientSampleId :
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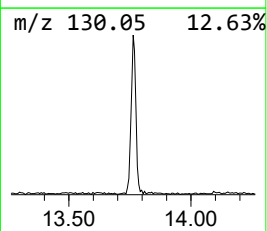
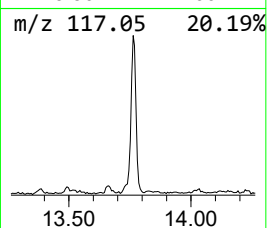
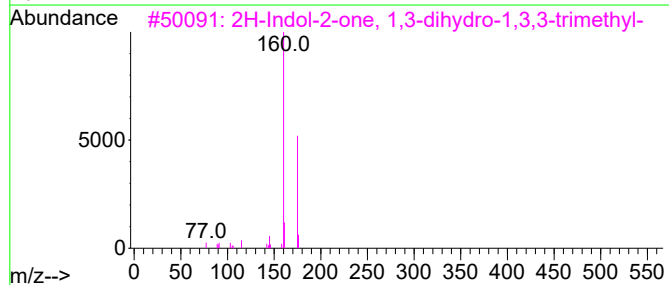
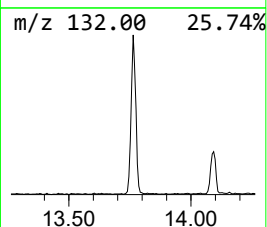
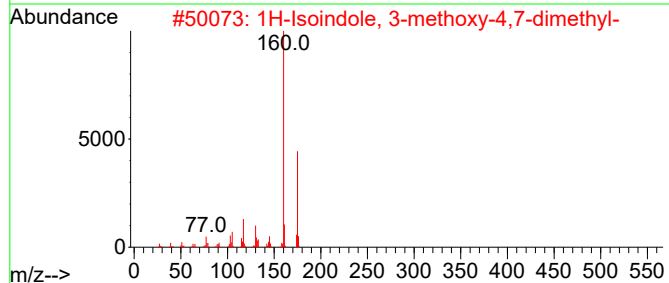
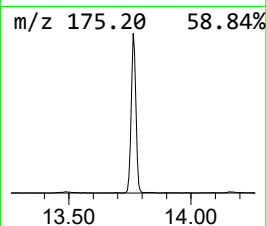
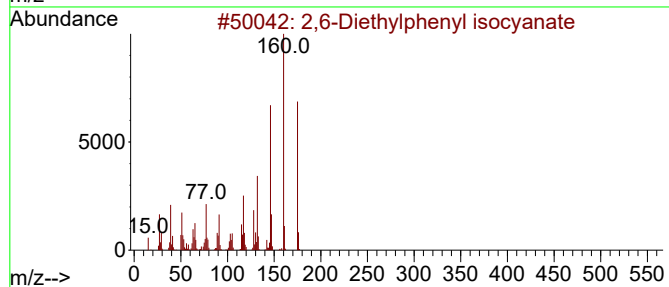
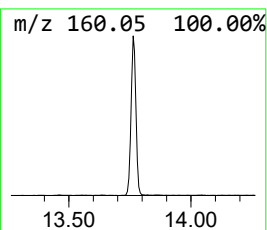
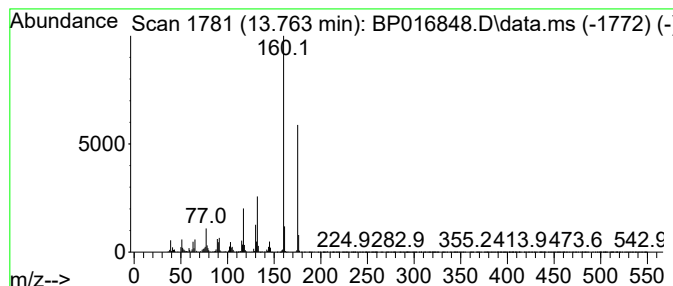
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082423.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 2,6-Diethylphenyl isocyanate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.763	2.19 ng/ul	569997	Acenaphthene-d10	14.681

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,6-Diethylphenyl isocyanate	175	C11H13NO	020458-99-5	90
2			1H-Isoindole, 3-methoxy-4,7-dime...	175	C11H13NO	100813-60-3	81
3			2H-Indol-2-one, 1,3-dihydro-1,3,...	175	C11H13NO	020200-86-6	70
4			1H-Isoindole-1,3(2H)-dione, N-et...	175	C10H9NO2	005022-29-7	53
5			1-Indanone, 5,6-dimethyl-	160	C11H12O	016440-97-4	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082823\
Data File : BP016848.D
Acq On : 30 Aug 2023 00:11
Operator : MA/JU
Sample : 04148-08
Misc :
ALS Vial : 66 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBJX2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082423.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
N-Nitrosodimeth...	3.764	2.2	ng/ul	279756	1	8.028	2523120	20.0
Tetrachloroethy...	4.658	124.4	ng/ul	15699300	1	8.028	2523120	20.0
Benzene, chloro-	5.293	2.4	ng/ul	306403	1	8.028	2523120	20.0
Ethanamine, N-e...	5.993	4.2	ng/ul	524048	1	8.028	2523120	20.0
Benzene, 1,3-di...	8.375	4.6	ng/ul	579891	1	8.028	2523120	20.0
Morpholine, 4-n...	8.946	3.3	ng/ul	419831	1	8.028	2523120	20.0
Morpholine, 2,6...	11.122	3.3	ng/ul	693374	2	10.852	4235720	20.0
2-Methyl-7-exo-...	12.275	2.2	ng/ul	459273	2	10.852	4235720	20.0
2,4,7,9-Tetrame...	13.522	2.0	ng/ul	530216	3	14.681	5199200	20.0
2,6-Diethylphen...	13.763	2.2	ng/ul	569997	3	14.681	5199200	20.0