

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
 Data File : BM034687.D
 Acq On : 15 Apr 2022 03:03
 Operator : CG/JU
 Sample : N2316-07
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ETNQ6

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-BM040922.M
 Title : SVOA CALIBRATION

Signal : TIC: BM034687.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.243	32	35	42	rVB2	35356	47792	1.60%	0.134%
2	3.667	104	107	130	rVB	240119	404852	13.60%	1.139%
3	4.678	276	279	286	rVB4	9857	15270	0.51%	0.043%
4	4.943	321	324	326	rBV3	5214	6306	0.21%	0.018%
5	5.149	358	359	365	rVB4	10361	12734	0.43%	0.036%
6	5.308	383	386	393	rBV4	9410	19162	0.64%	0.054%
7	5.425	403	406	413	rVB6	9799	14532	0.49%	0.041%
8	6.090	514	519	528	rVB2	25692	46918	1.58%	0.132%
9	6.837	642	646	652	rVB	10598	18745	0.63%	0.053%
10	6.955	661	666	670	rBV4	7724	13325	0.45%	0.037%
11	7.025	671	678	693	rBV	417368	760134	25.53%	2.138%
12	7.190	700	706	719	rBV	345171	605151	20.32%	1.702%
13	7.390	733	740	752	rBV	461657	792812	26.62%	2.230%
14	7.490	755	757	762	rVB5	6520	7796	0.26%	0.022%
15	7.702	790	793	797	rBV6	3420	4842	0.16%	0.014%
16	7.860	813	820	829	rBV	497354	796640	26.75%	2.241%
17	7.937	829	833	838	rVB3	10314	17496	0.59%	0.049%
18	8.078	852	857	867	rBV2	24700	45918	1.54%	0.129%
19	8.160	867	871	879	rVB3	15750	28580	0.96%	0.080%
20	8.319	892	898	908	rVB	133439	213327	7.16%	0.600%
21	8.484	921	926	933	rBV2	59788	102121	3.43%	0.287%
22	8.572	935	941	955	rVV	538160	972970	32.67%	2.737%
23	8.696	957	962	970	rVB3	16938	34730	1.17%	0.098%
24	8.878	984	993	1009	rVB5	34180	132580	4.45%	0.373%
25	9.025	1011	1018	1031	rVB	402052	701283	23.55%	1.973%
26	9.360	1070	1075	1082	rBV3	18298	30883	1.04%	0.087%
27	9.484	1089	1096	1107	rBV5	19182	66174	2.22%	0.186%
28	9.572	1109	1111	1117	rVB8	4337	8357	0.28%	0.024%
29	9.749	1134	1141	1152	rBV	334259	611532	20.54%	1.720%
30	9.972	1176	1179	1185	rVB6	4078	7837	0.26%	0.022%
31	10.078	1192	1197	1200	rBV3	5964	10421	0.35%	0.029%
32	10.296	1227	1234	1252	rBV	457717	884167	29.69%	2.487%
33	10.607	1282	1287	1291	rBV3	16557	27939	0.94%	0.079%
34	10.672	1291	1298	1309	rVB	651071	1120385	37.62%	3.152%
35	10.807	1313	1321	1332	rBV	347867	716368	24.06%	2.015%
36	11.096	1363	1370	1372	rBV8	8733	17554	0.59%	0.049%

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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-BM040922.M
 Title : SVOA CALIBRATION

37	11.178	1378	1384	1397	rBV	96011	238314	8.00%	0.670%
38	11.648	1459	1464	1471	rBV8	15359	35510	1.19%	0.100%
39	11.896	1503	1506	1509	rVB5	3491	4444	0.15%	0.013%
40	12.101	1536	1541	1547	rBV4	8962	17075	0.57%	0.048%

41	12.207	1551	1559	1563	rBV8	6731	18503	0.62%	0.052%
42	12.272	1566	1570	1576	rVB6	7802	14986	0.50%	0.042%
43	12.560	1614	1619	1623	rBV4	5115	8415	0.28%	0.024%
44	12.672	1633	1638	1639	rBV4	5582	6952	0.23%	0.020%
45	12.884	1670	1674	1677	rBV3	10625	17046	0.57%	0.048%

46	13.025	1692	1698	1702	rBV	20661	34116	1.15%	0.096%
47	13.178	1722	1724	1729	rVV2	4012	5047	0.17%	0.014%
48	13.348	1748	1753	1762	rVB3	29525	50286	1.69%	0.141%
49	13.448	1765	1770	1773	rBV7	5248	9505	0.32%	0.027%
50	13.731	1812	1818	1822	rBV5	8168	13421	0.45%	0.038%

51	13.919	1844	1850	1859	rVV	1028239	1447706	48.62%	4.073%
52	13.995	1860	1863	1872	rVV3	14533	30352	1.02%	0.085%
53	14.054	1872	1873	1878	rVB5	3243	4256	0.14%	0.012%
54	14.201	1888	1898	1910	rBV	1106606	1620627	54.42%	4.559%
55	14.278	1910	1911	1913	rVV2	4161	4114	0.14%	0.012%

56	14.301	1913	1915	1921	rVB7	7454	8950	0.30%	0.025%
57	14.425	1930	1936	1942	rBV	40637	76837	2.58%	0.216%
58	14.507	1944	1950	1961	rVV	974497	1445483	48.54%	4.066%
59	14.595	1961	1965	1975	rVB	13761	39584	1.33%	0.111%
60	14.748	1983	1991	1992	rBV6	11571	21039	0.71%	0.059%

61	14.789	1994	1998	2002	rVV3	6240	15871	0.53%	0.045%
62	15.160	2055	2061	2068	rVB7	7690	19770	0.66%	0.056%
63	15.236	2068	2074	2077	rBV6	3444	6199	0.21%	0.017%
64	15.336	2082	2091	2113	rBV	991733	2853403	95.82%	8.027%
65	15.501	2113	2119	2134	rVV2	1456212	2152083	72.27%	6.054%

66	15.619	2134	2139	2150	rVV	262156	451316	15.16%	1.270%
67	15.725	2152	2157	2176	rVB2	56123	141867	4.76%	0.399%
68	15.948	2192	2195	2197	rBV4	5065	5889	0.20%	0.017%
69	16.131	2222	2226	2231	rVB3	10357	17376	0.58%	0.049%
70	16.183	2231	2235	2238	rBV3	10085	17093	0.57%	0.048%

71	16.225	2238	2242	2244	rVB5	4249	5359	0.18%	0.015%
72	16.289	2249	2253	2257	rVB5	7379	12122	0.41%	0.034%
73	16.348	2257	2263	2266	rBV	29683	47861	1.61%	0.135%
74	16.378	2266	2268	2274	rVB2	22422	31152	1.05%	0.088%
75	16.489	2282	2287	2293	rBV6	4751	10269	0.34%	0.029%

76	16.654	2309	2315	2319	rBV8	6348	11061	0.37%	0.031%
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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-BM040922.M
 Title : SVOA CALIBRATION

77	16.919	2356	2360	2361	rBV4	3309	4087	0.14%	0.011%
78	16.936	2361	2363	2367	rVB4	3635	4549	0.15%	0.013%
79	17.025	2373	2378	2379	rBV5	6230	7837	0.26%	0.022%
80	17.178	2399	2404	2411	rBV	147892	244516	8.21%	0.688%
81	17.254	2411	2417	2425	rVV	1066718	1571123	52.76%	4.420%
82	17.348	2426	2433	2448	rVV2	1637035	2456911	82.51%	6.912%
83	17.930	2526	2532	2537	rVV7	10848	20019	0.67%	0.056%
84	18.077	2551	2557	2566	rVV	384478	531091	17.83%	1.494%
85	18.154	2566	2570	2575	rVV6	20431	33938	1.14%	0.095%
86	18.225	2579	2582	2584	rVV3	4721	4913	0.16%	0.014%
87	18.789	2675	2678	2681	rVB5	7449	8275	0.28%	0.023%
88	18.836	2681	2686	2691	rBV	42231	59534	2.00%	0.167%
89	18.907	2693	2698	2703	rBV	199665	249854	8.39%	0.703%
90	19.207	2745	2749	2755	rBV2	23229	38705	1.30%	0.109%
91	19.277	2758	2761	2765	rVB	19604	27309	0.92%	0.077%
92	19.601	2811	2816	2818	rBV	102202	126109	4.23%	0.355%
93	19.642	2818	2823	2829	rVV	2220096	2977807	100.00%	8.377%
94	19.689	2829	2831	2840	rVB3	55365	84931	2.85%	0.239%
95	19.783	2845	2847	2852	rVB6	17235	20266	0.68%	0.057%
96	20.289	2929	2933	2937	rBV	125419	152135	5.11%	0.428%
97	21.142	3074	3078	3086	rBV	424451	610673	20.51%	1.718%
98	21.430	3122	3127	3140	rBV	927332	1561257	52.43%	4.392%
99	23.648	3497	3504	3521	rBV2	1238103	2798841	93.99%	7.874%
100	23.801	3523	3530	3542	rVV2	764772	1663486	55.86%	4.680%

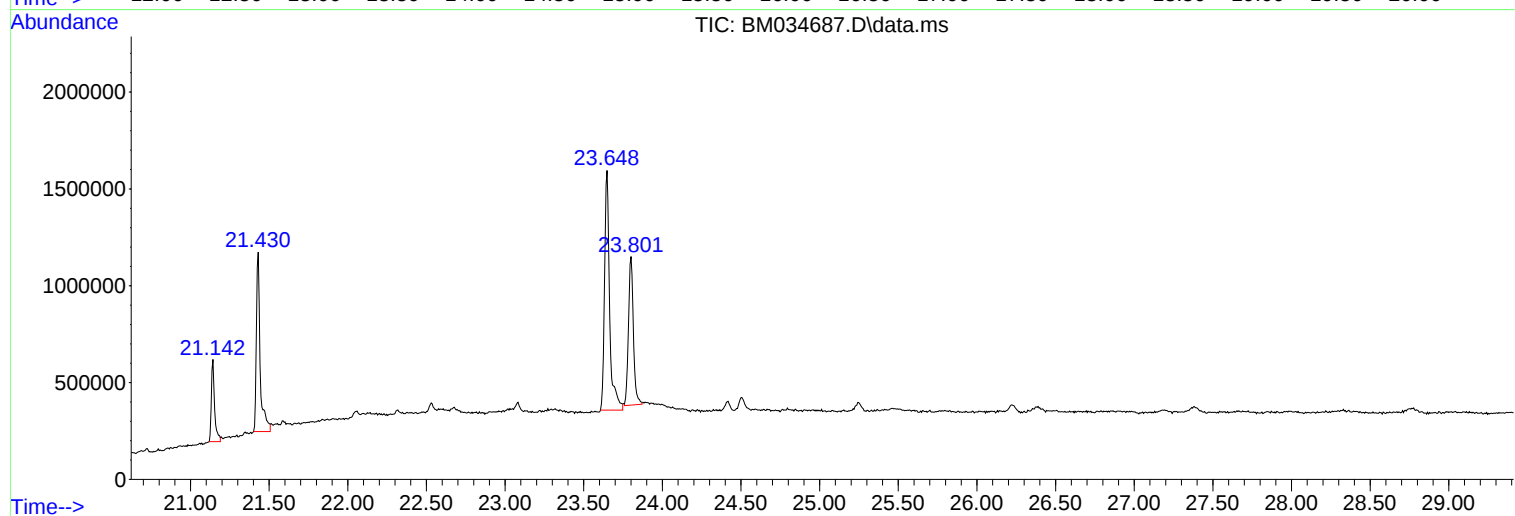
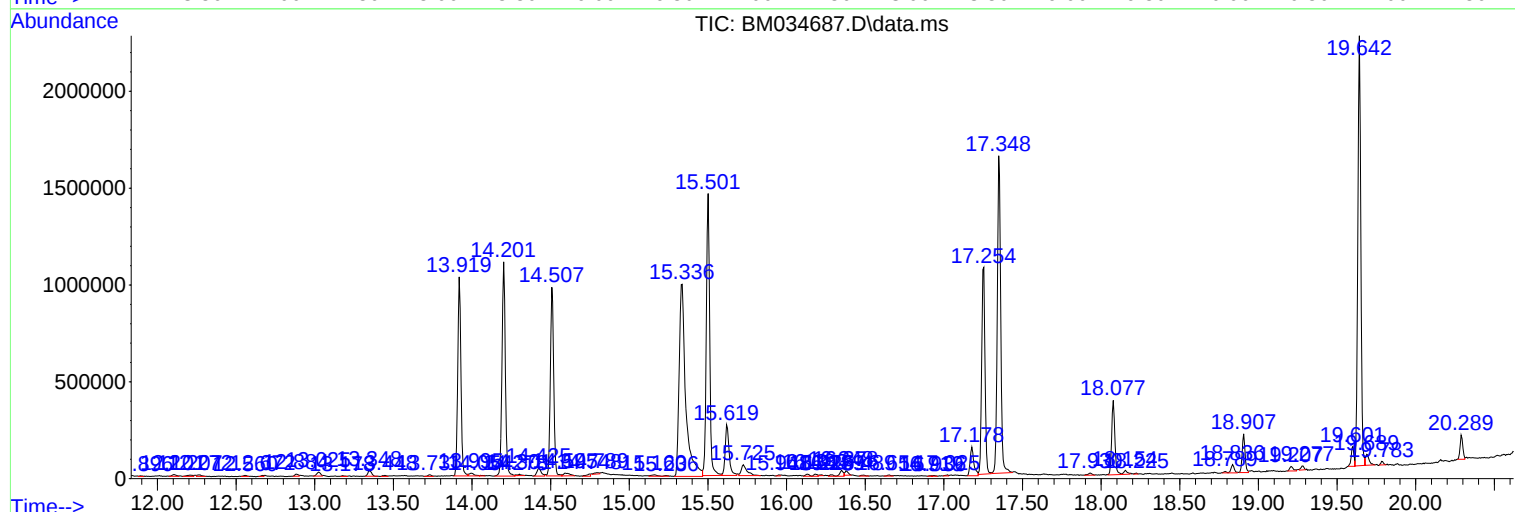
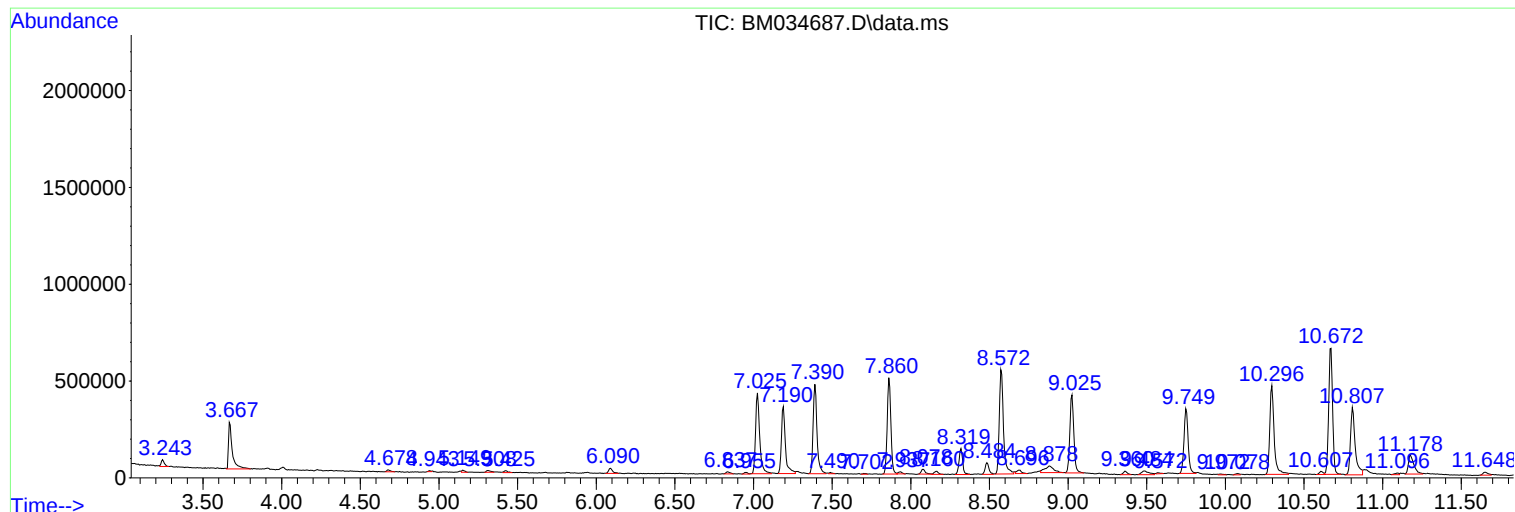
Sum of corrected areas: 35547058

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ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_M
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

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



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ALS Vial : 29 Sample Multiplier: 1

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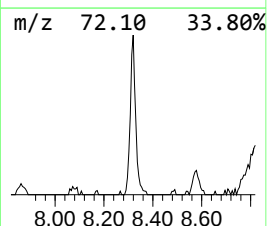
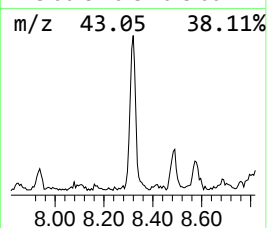
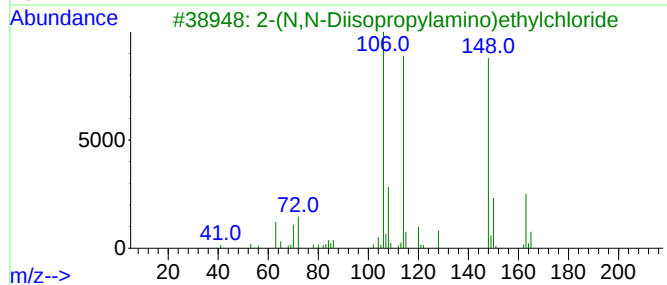
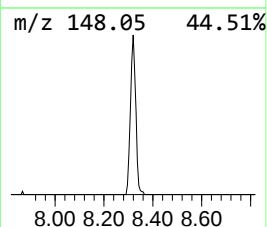
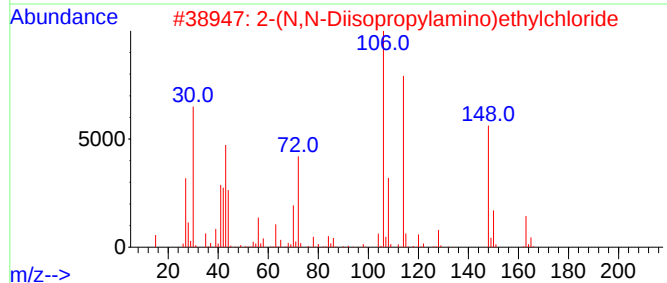
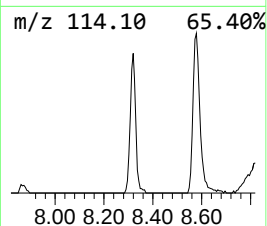
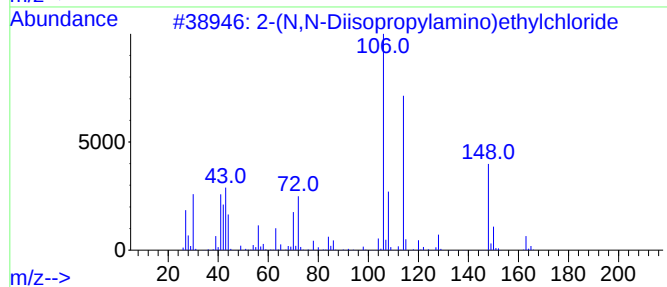
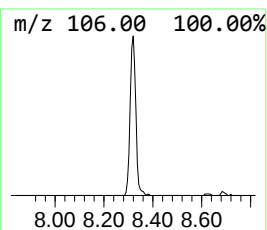
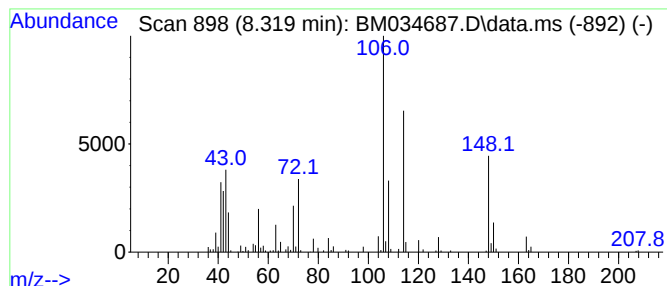
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-(N,N-Diisopropylamino)eth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.319	5.36 ng/ul	213327	1,4-Dichlorobenzene-d4	7.860

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(N,N-Diisopropylamino)ethylchl...	163	C8H18ClN	000096-79-7	93		
2	2-(N,N-Diisopropylamino)ethylchl...	163	C8H18ClN	000096-79-7	81		
3	2-(N,N-Diisopropylamino)ethylchl...	163	C8H18ClN	000096-79-7	32		
4	Bis(2-chloroethyl)methylamine	155	C5H11Cl2N	000051-75-2	27		
5	2-(N,N-Diisopropylamino)ethylchl...	163	C8H18ClN	000096-79-7	27		



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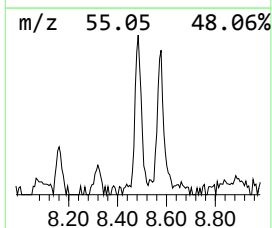
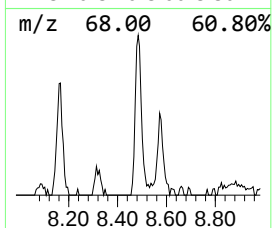
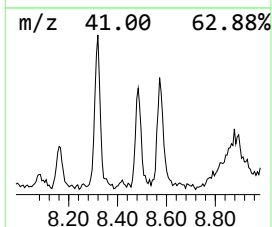
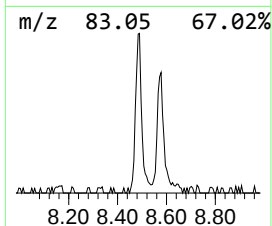
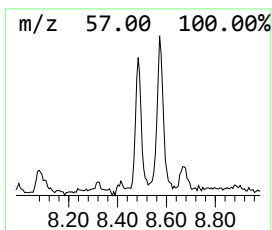
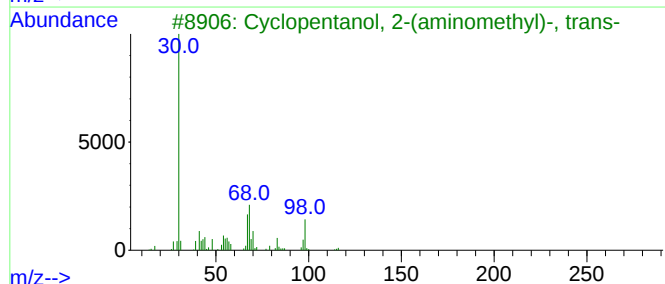
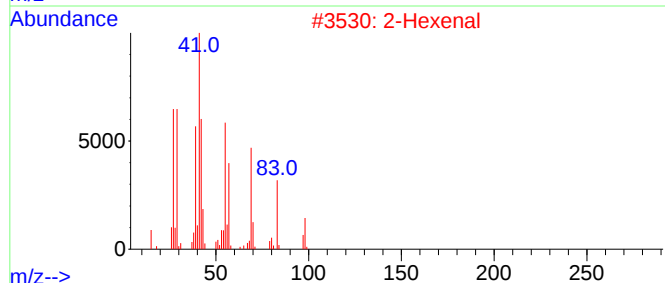
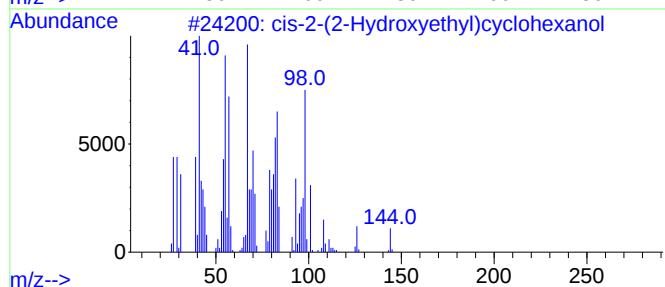
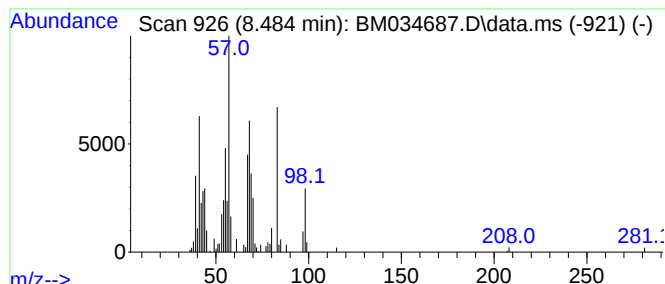
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.484	2.56 ng/ul	102121	1,4-Dichlorobenzene-d4	7.860

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-2-(2-Hydroxyethyl)cyclohexanol	144	C8H16O2	040894-17-5	43
2			2-Hexenal	98	C6H10O	000505-57-7	43
3			Cyclopentanol, 2-(aminomethyl)-,...	115	C6H13NO	040482-02-8	43
4			2-Hexenal, (E)-	98	C6H10O	006728-26-3	43
5			Furan, 2,5-dihydro-2,5-dimethyl-	98	C6H10O	059242-27-2	38



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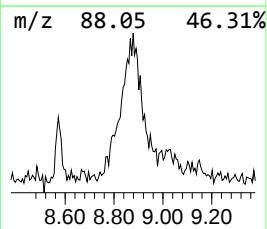
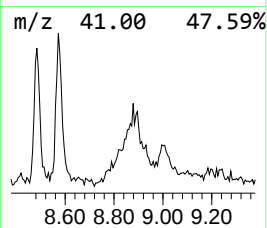
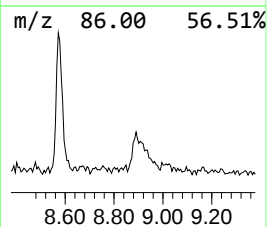
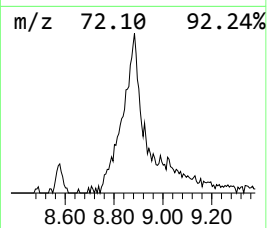
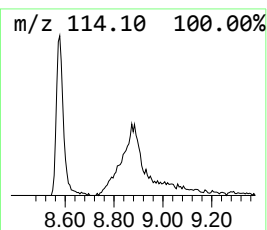
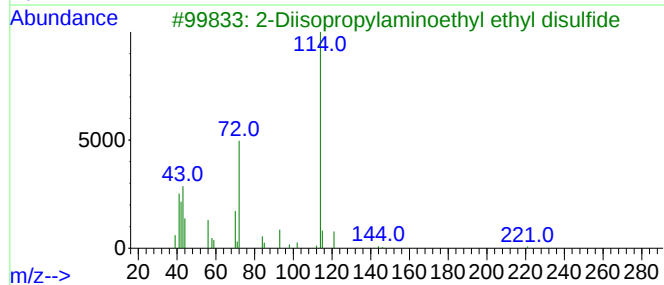
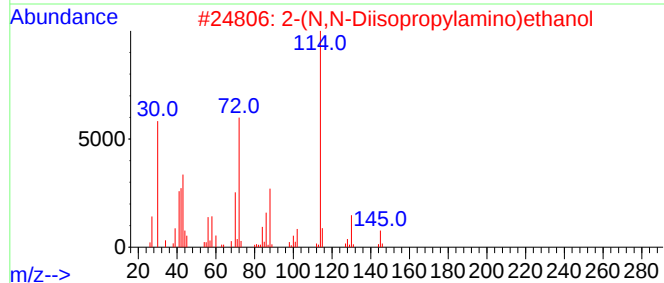
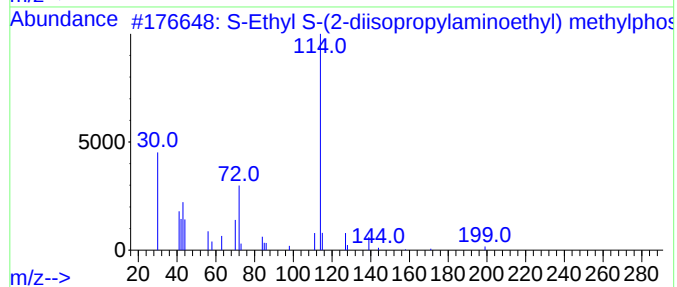
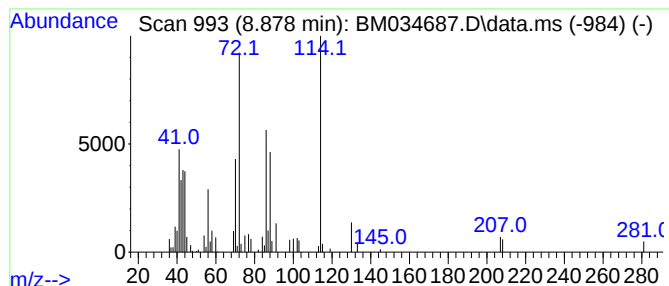
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 S-Ethyl S-(2-diisopropylami... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.878	3.33 ng/ul	132580	1,4-Dichlorobenzene-d4	7.860

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			S-Ethyl S-(2-diisopropylaminoeth...	283	C11H26NOPS2	110501-55-8	50
2			2-(N,N-Diisopropylamino)ethanol	145	C8H19NO	000096-80-0	50
3			2-Diisopropylaminoethyl ethyl di...	221	C10H23NS2	310882-85-0	50
4			2-(N,N-Diisopropylamino)ethanol	145	C8H19NO	000096-80-0	50
5			Ethanol, 2-[(1-methylethyl)amino]-	103	C5H13NO	000109-56-8	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034687.D
Acq On : 15 Apr 2022 03:03
Operator : CG/JU
Sample : N2316-07
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNQ6

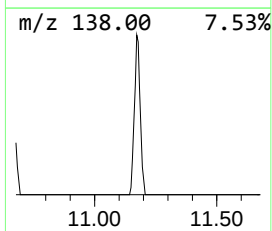
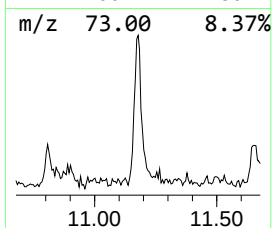
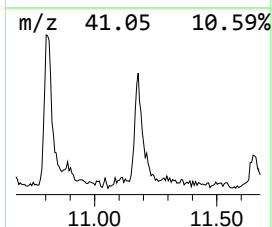
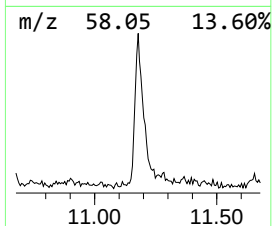
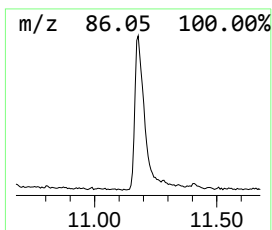
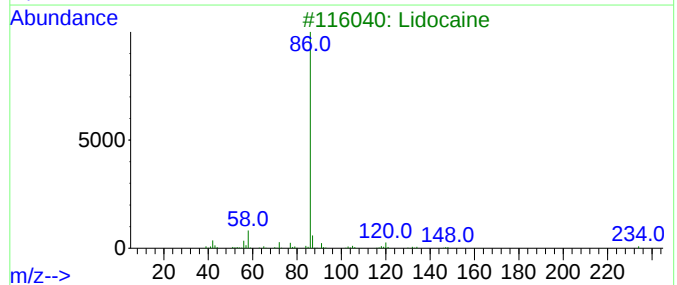
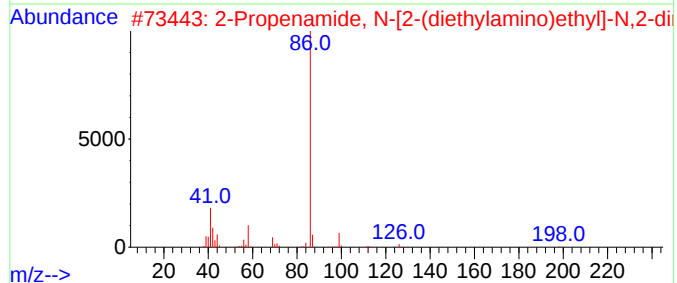
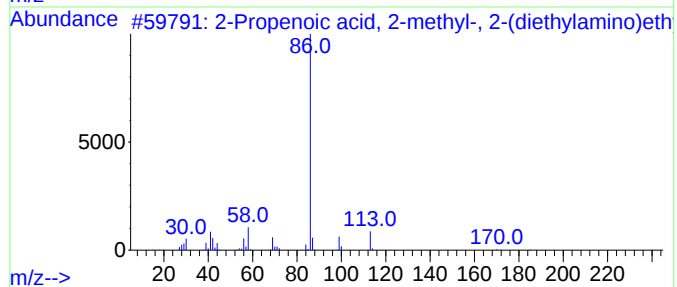
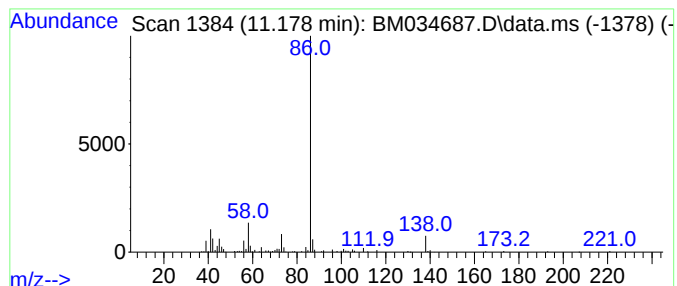
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 2-Propenoic acid, 2-methyl-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.178	4.25 ng/ul	238314	Naphthalene-d8	10.672

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propenoic acid, 2-methyl-, 2-(...	185	C10H19NO2	000105-16-8	64
2			2-Propenamide, N-[2-(diethylamin...	198	C11H22N2O	013173-19-8	64
3			Lidocaine	234	C14H22N2O	000137-58-6	64
4			1,2-Ethanediamine, N,N,N',N'-tet...	172	C10H24N2	000150-77-6	64
5			3-Fluorobenzoic acid, 2-(diethyl...	239	C13H18FNO2	002994-22-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034687.D
Acq On : 15 Apr 2022 03:03
Operator : CG/JU
Sample : N2316-07
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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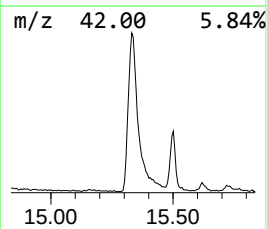
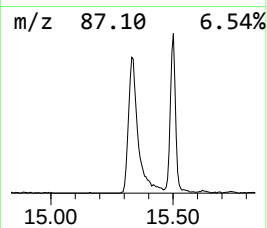
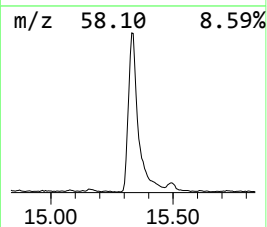
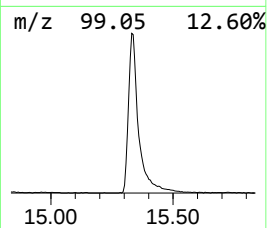
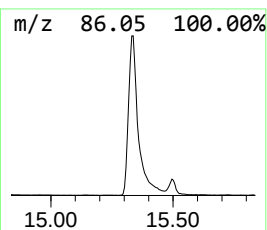
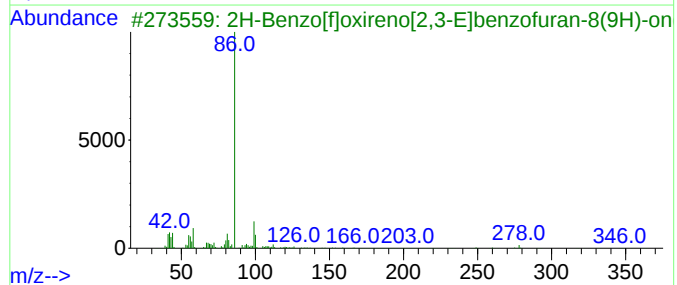
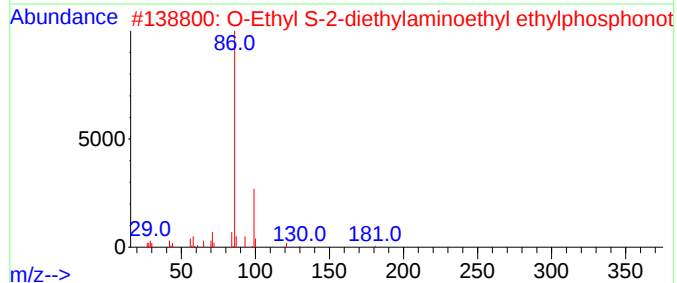
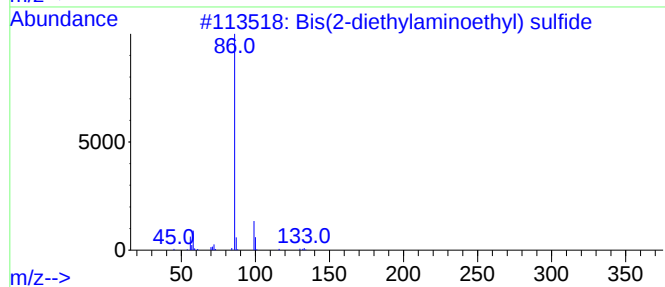
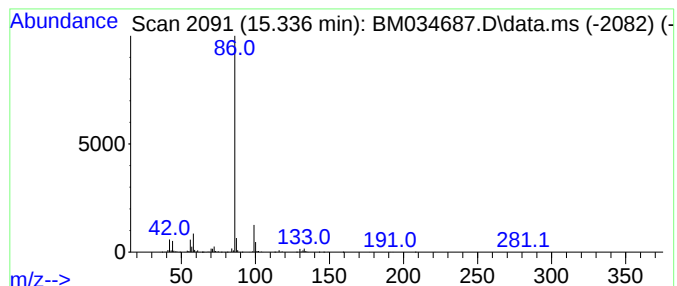
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Bis(2-diethylaminoethyl) su... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.336	39.48 ng/ul	2853400	Acenaphthene-d10	14.507

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bis(2-diethylaminoethyl) sulfide	232	C12H28N2S	006006-58-2	86
2			O-Ethyl S-2-diethylaminoethyl et...	253	C10H24N02PS	021738-25-0	78
3			2H-Benzo[f]oxireno[2,3-E]benzofu...	364	C21H36N2O3	1000327-59-9	78
4			O-Methyl S-2-diethylaminoethyl e...	239	C9H22N02PS	170800-77-8	78
5			O-Methyl S-2-diethylaminoethyl e...	239	C9H22N02PS	170800-77-8	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
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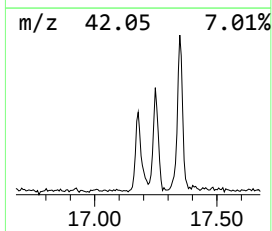
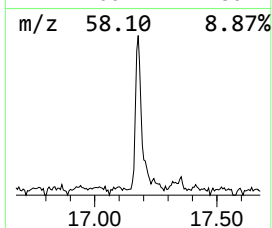
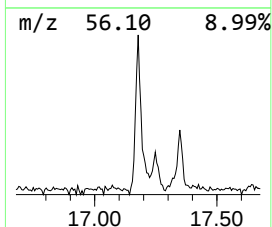
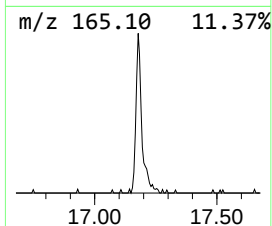
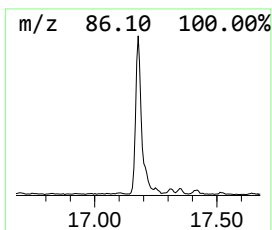
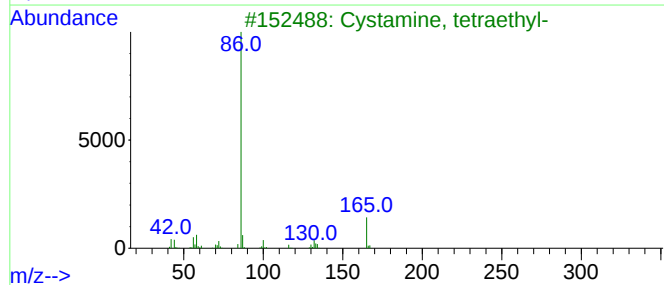
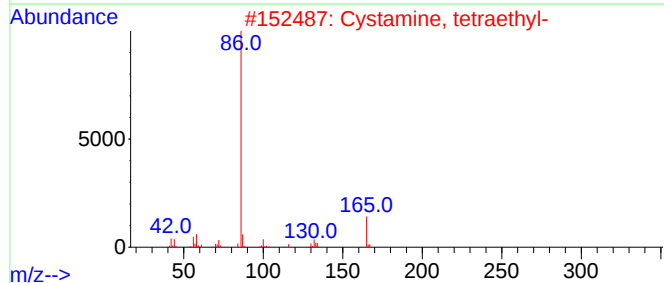
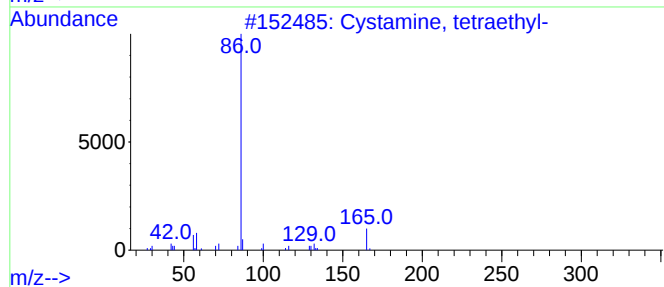
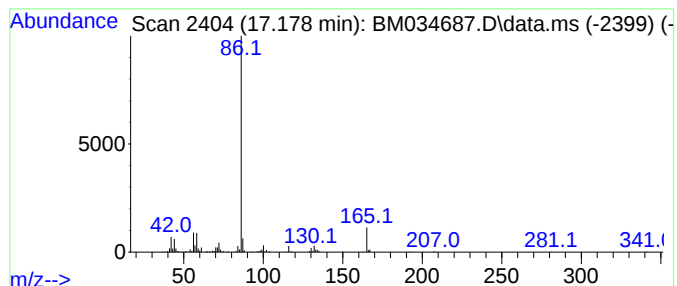
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Cystamine, tetraethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.177	3.11 ng/ul	244516	Phenanthrene-d10	17.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cystamine, tetraethyl-	264	C12H28N2S2	000589-32-2	90
2			Cystamine, tetraethyl-	264	C12H28N2S2	000589-32-2	83
3			Cystamine, tetraethyl-	264	C12H28N2S2	000589-32-2	83
4			1,2-Ethanediamine, N,N-diethyl-	116	C6H16N2	000100-36-7	64
5			1,2-Ethanediamine, N,N-diethyl-	116	C6H16N2	000100-36-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034687.D
Acq On : 15 Apr 2022 03:03
Operator : CG/JU
Sample : N2316-07
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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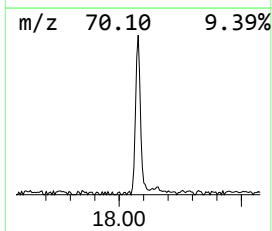
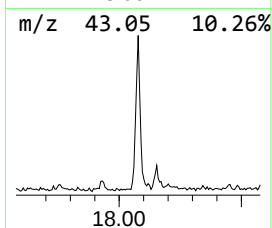
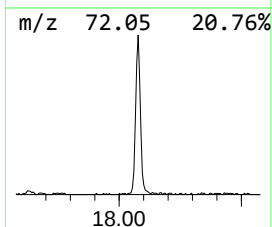
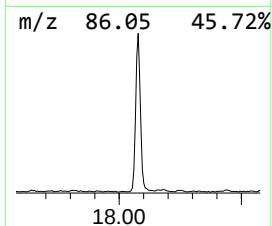
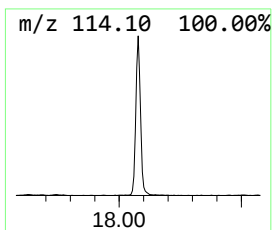
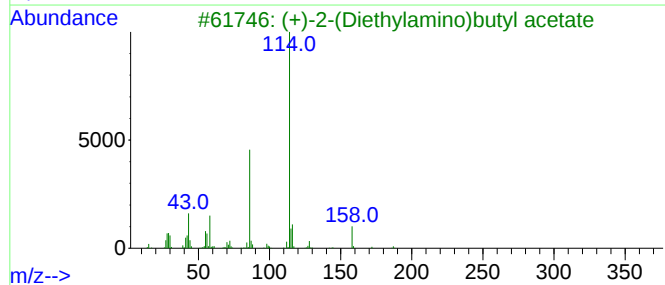
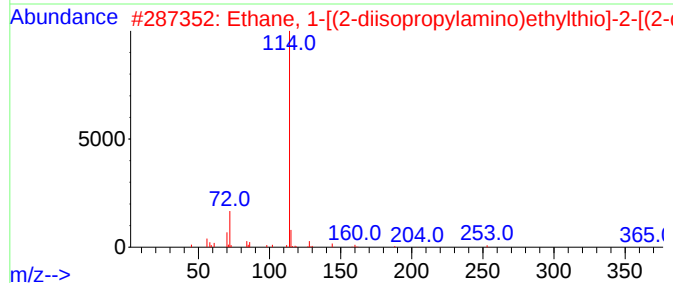
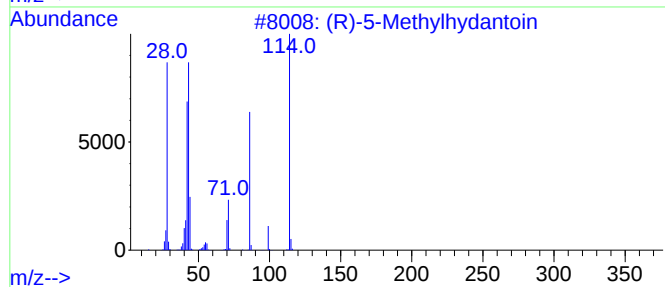
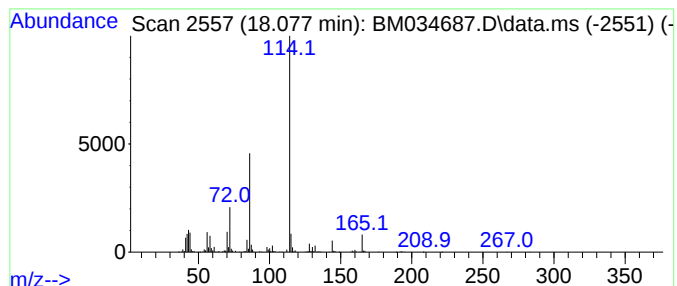
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (R)-5-Methylhydantoin Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.078	6.76 ng/ul	531091	Phenanthrene-d10	17.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(R)-5-Methylhydantoin	114	C4H6N2O2	055147-68-7	53
2			Ethane, 1-[(2-diisopropylamino)e...	380	C18H40N2S3	110501-59-2	53
3			(+)-2-(Diethylamino)butyl acetate	187	C10H21NO2	1000240-85-4	53
4			2-Diisopropylaminoethyl ethyl di...	221	C10H23NS2	310882-85-0	50
5			Tetrapropylammonium iodide	313	C12H28IN	000631-40-3	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034687.D
Acq On : 15 Apr 2022 03:03
Operator : CG/JU
Sample : N2316-07
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
ETNQ6

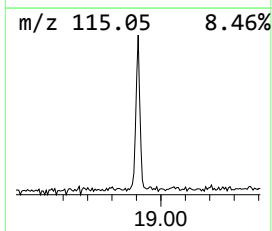
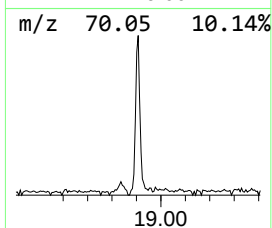
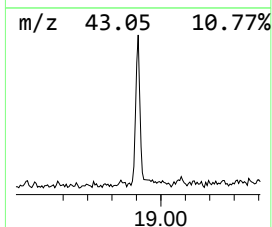
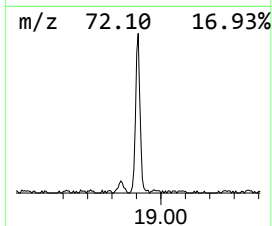
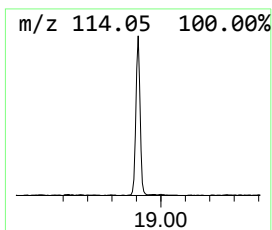
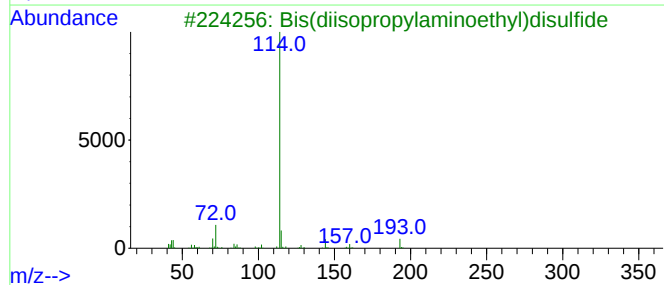
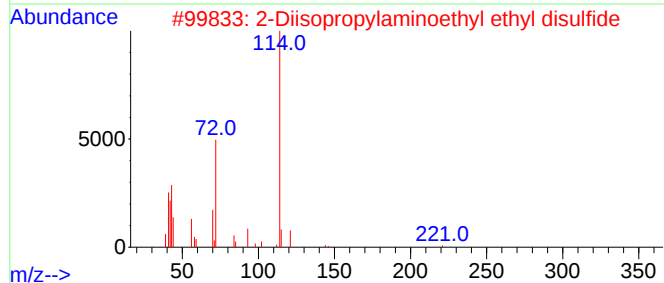
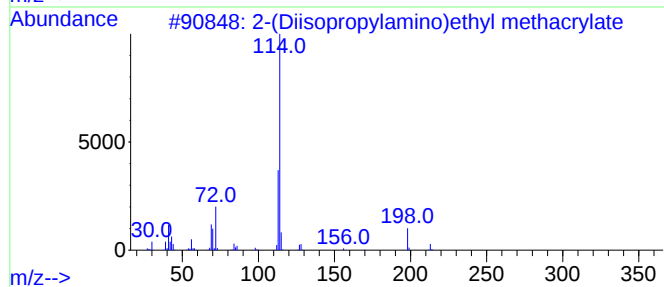
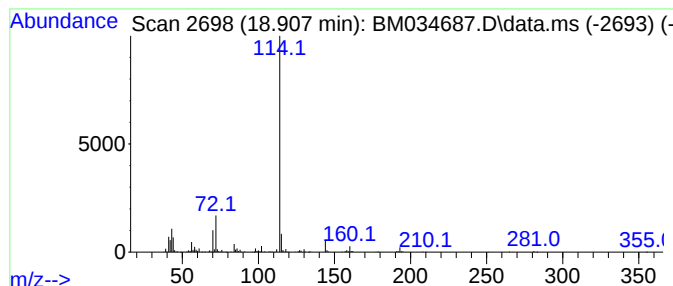
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 2-(Diisopropylamino)ethyl m... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.907	3.18 ng/ul	249854	Phenanthrene-d10	17.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-(Diisopropylamino)ethyl methac...	213	C12H23NO2	016715-83-6	72
2			2-Diisopropylaminoethyl ethyl di...	221	C10H23NS2	310882-85-0	72
3			Bis(diisopropylaminoethyl)disulfide	320	C16H36N2S2	065332-44-7	72
4			Ethane, 1-[(2-diisopropylamino)e...	380	C18H40N2S3	110501-59-2	72
5			Bis(2-diisopropylaminoethyl) sul...	288	C16H36N2S	110501-56-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
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Misc :
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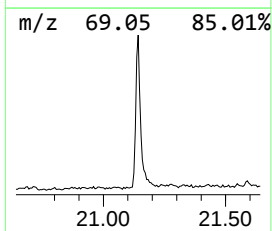
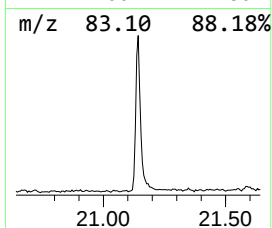
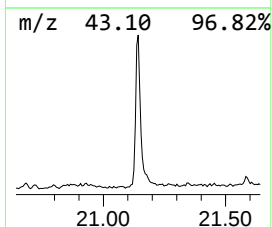
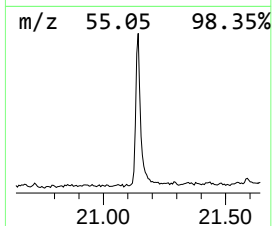
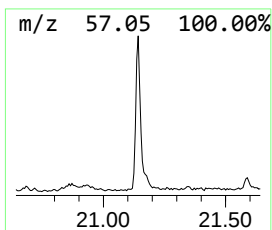
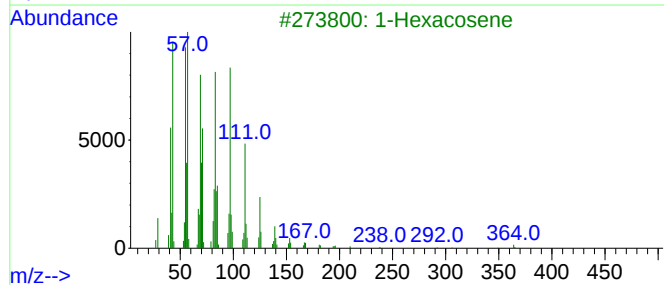
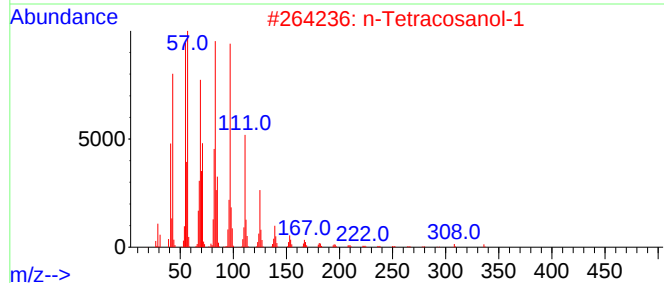
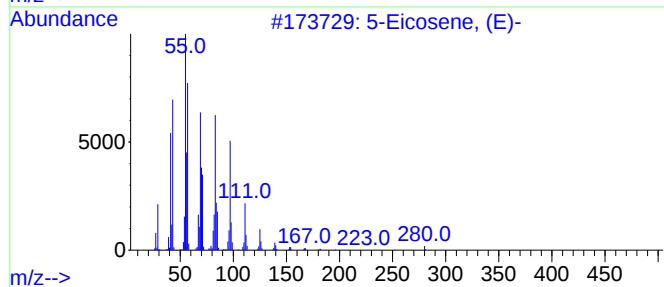
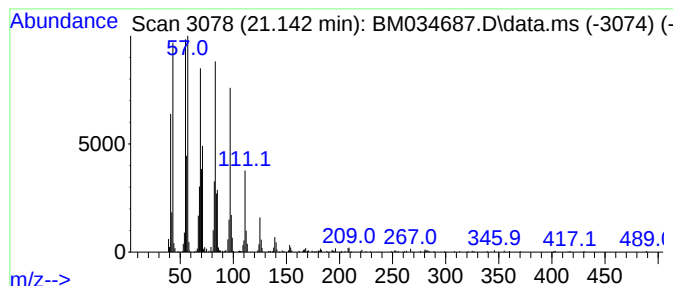
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.142	7.82 ng/ul	610673	Chrysene-d12	21.430

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-		280	C20H40	074685-30-6	99
2	n-Tetracosanol-1		354	C24H50O	000506-51-4	94
3	1-Hexacosene		364	C26H52	018835-33-1	94
4	1-Tetracosene		336	C24H48	010192-32-2	94
5	Pentacos-1-ene		350	C25H50	016980-85-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
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ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
ETNQ6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.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
2-(N,N-Diisopro...	8.319	5.4	ng/ul	213327	1	7.860	796640	20.0
unknown-01	8.484	2.6	ng/ul	102121	1	7.860	796640	20.0
S-Ethyl S-(2-di...	8.878	3.3	ng/ul	132580	1	7.860	796640	20.0
2-Propenoic aci...	11.178	4.3	ng/ul	238314	2	10.672	1120390	20.0
Bis(2-diethylam...	15.336	39.5	ng/ul	2853400	3	14.507	1445480	20.0
Cystamine, tetr...	17.177	3.1	ng/ul	244516	4	17.254	1571120	20.0
(R)-5-Methylhyd...	18.078	6.8	ng/ul	531091	4	17.254	1571120	20.0
2-(Diisopropyla...	18.907	3.2	ng/ul	249854	4	17.254	1571120	20.0
(DEL) Alkane: S...	21.142	7.8	ng/ul	610673	5	21.430	1561260	20.0