LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018101.D
Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH347

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-BP111023.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP018101.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.228	20	24	27	rBV2	17995	22115	0.51%	0.061%
2	3.263	27	30	38	rVB	22436	27265	0.63%	0.076%
3	3.346	39	44	48	rBV	111771	129563	3.00%	0.359%
4	3.669	96	99	113	rBV	82861	147400	3.41%	0.409%
5	3.834	123	127	136	rVB2	41630	60385	1.40%	0.167%
6	4.481	232	237	242	rBV2	22885	34037	0.79%	0.094%
7	5.122	341	346	350	rBV6	9045	15259	0.35%	0.042%
8	5.157	350	352	358	rVV7	6020	8718	0.20%	0.024%
9	5.340	379	383	388	rBV4	10784	15010	0.35%	0.042%
10	5.710	439	446	457	rBV	1561973	2168743	50.14%	6.014%
11	5.916	476	481	488	rVB3	10358	18050	0.42%	0.050%
12	6.034	496	501	508	rBV	75788	105818	2.45%	0.293%
13	6.716	611	617	623	rVB10	4224	7807	0.18%	0.022%
14	7.034	662	671	688	rBV	2911080	4325095	100.00%	11.994%
15	7.192	693	698	710	rVV	339372	524414	12.12%	1.454%
16	7.369	719	728	737	rBV	352525	602992	13.94%	1.672%
17	7.440	737	740	746	rVB3	15559	24555	0.57%	0.068%
18	7.781	791	798	802	rBV7	8936	16375	0.38%	0.045%
19	7.845	802	809	813	rBV	217676	381361	8.82%	1.058%
20	7.981	829	832	839	rVB4	9322	15488	0.36%	0.043%
21	8.198	861	869	874	rBV10	8066	16790	0.39%	0.047%
22	8.386	893	901	911	rVB	53023	84124	1.95%	0.233%
23	8.528	917	925	928	rBV2	45316	75015	1.73%	0.208%
24	8.575	928	933	941	rVV2	237176	425029	9.83%	1.179%
25	8.639	941	944	956	rVB3	27731	68300	1.58%	0.189%
26	8.763	958	965	969	rBV3	20116	29242	0.68%	0.081%
27	8.804	969	972	977	rVB3	8803	11638	0.27%	0.032%
28	8.857	977	981	987	rBV3	17923	32326	0.75%	0.090%
29	8.916	987	991	995	rBV2	18725	27775	0.64%	0.077%
30	8.963	995	999	1005	rBV	19200	28600	0.66%	0.079%
31	9.051	1007	1014	1020	rBV	473902	796438	18.41%	2.209%
32	9.116	1021	1025	1031	rVB	105753	149017	3.45%	0.413%
33	9.192	1031	1038	1041	rBV3	34364	71534	1.65%	0.198%
34	9.257	1046	1049	1055	rVB4	41732	66995	1.55%	0.186%
35	9.351	1055	1065	1074	rVB4	53730	132768	3.07%	0.368%
36	9.445	1074	1081	1093	rBV6	69840	224286	5.19%	0.622%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018101.D
Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH347

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-BP111023.MA.M
Title : SVOA CALIBRATION

37	9.539	1093	1097	1101	rBV2	28928	45849	1.06%	0.127%
38	9.581	1101	1104	1105	rVV3	8339	7167	0.17%	0.020%
39	9.622	1105	1111	1118	rVB6	59244	132754	3.07%	0.368%
40	9.692	1118	1123	1127	rBV4	15534	26037	0.60%	0.072%
41	9.763	1128	1135	1147	rBV	404986	706752	16.34%	1.960%
42	9.904	1155	1159	1161	rBV2	6942	9489	0.22%	0.026%
43	9.945	1161	1166	1169	rVV3	24466	38188	0.88%	0.106%
44	9.986	1169	1173	1178	rVV2	43223	91146	2.11%	0.253%
45	10.045	1178	1183	1194	rVB6	53831	150338	3.48%	0.417%
46	10.204	1202	1210	1215	rBV3	94163	231693	5.36%	0.643%
47	10.298	1219	1226	1238	rVV2	500169	1233556	28.52%	3.421%
48	10.381	1238	1240	1245	rVV2	60723	122846	2.84%	0.341%
49	10.433	1245	1249	1257	rVV4	61181	149875	3.47%	0.416%
50	10.663	1279	1288	1299	rVB	307830	576566	13.33%	1.599%
51	10.845	1311	1319	1340	rVB	429030	1026282	23.73%	2.846%
52	10.992	1340	1344	1351	rBV5	27936	61195	1.41%	0.170%
53	11.092	1357	1361	1365	rBV4	26147	53084	1.23%	0.147%
54	11.339	1398	1403	1407	rBV5	31626	58058	1.34%	0.161%
55	11.404	1407	1414	1417	rBV3	35956	71250	1.65%	0.198%
56	11.445	1417	1421	1422	rBV	34900	40680	0.94%	0.113%
57	11.557	1434	1440	1445	rBV6	43313	111550	2.58%	0.309%
58	11.692	1458	1463	1469	rVB2	142578	239762	5.54%	0.665%
59	12.169	1539	1544	1556	rVB2	43069	99303	2.30%	0.275%
60	12.480	1590	1597	1602	rBV5	14751	33640	0.78%	0.093%
61	12.863	1658	1662	1666	rVV7	5963	10766	0.25%	0.030%
62	12.916	1668	1671	1675	rVB5	4865	7702	0.18%	0.021%
63	13.104	1699	1703	1709	rBV3	16928	29580	0.68%	0.082%
64	13.357	1738	1746	1753	rBV	120263	189614	4.38%	0.526%
65	13.480	1763	1767	1771	rVB6	6294	7763	0.18%	0.022%
66	13.827	1823	1826	1830	rBV5	15533	20604	0.48%	0.057%
67	13.904	1835	1839	1840	rBV2	15983	17882	0.41%	0.050%
68	13.939	1840	1845	1851	rVV	1153576	1543603	35.69%	4.281%
69	14.127	1873	1877	1883	rVB5	18546	31930	0.74%	0.089%
70	14.216	1885	1892	1904	rBV	1164108	1742613	40.29%	4.833%
71	14.333	1906	1912	1919	rVV6	12895	32678	0.76%	0.091%
72	14.521	1936	1944	1951	rBV	426608	650171	15.03%	1.803%
73	14.604	1954	1958	1967	rVB8	12344	27502	0.64%	0.076%
74	14.827	1991	1996	2005	rBV4	21279	61858	1.43%	0.172%
75	14.910	2008	2010	2014	rVB5	8032	9188	0.21%	0.025%
76	15.269	2067	2071	2075	rBV3	10504	16870	0.39%	0.047%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018101.D
Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BH347

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-BP111023.MA.M
Title : SVOA CALIBRATION

77	15.439	2097	2100	2104	rBV5	11129	18027	0.42%	0.050%
78	15.521	2107	2114	2122	rVV	1615775	2241468	51.82%	6.216%
79	15.674	2132	2140	2151	rBV	559276	876127	20.26%	2.430%
80	15.757	2151	2154	2159	rVB6	15814	24680	0.57%	0.068%
81	15.845	2166	2169	2172	rVB5	10510	11377	0.26%	0.032%
82	16.233	2230	2235	2243	rVB5	15237	31212	0.72%	0.087%
83	16.327	2243	2251	2254	rBV4	24960	62972	1.46%	0.175%
84	16.363	2254	2257	2259	rVV4	14838	19072	0.44%	0.053%
85	16.386	2259	2261	2268	rVB2	17570	29651	0.69%	0.082%
86	16.457	2268	2273	2277	rBV4	22321	36758	0.85%	0.102%
87	16.598	2294	2297	2301	rVB5	9379	11605	0.27%	0.032%
88	16.727	2315	2319	2323	rBV2	16566	20160	0.47%	0.056%
89	16.798	2328	2331	2336	rVB7	8379	13189	0.30%	0.037%
90	17.298	2410	2416	2423	rBV	575929	820275	18.97%	2.275%
91	17.398	2427	2433	2445	rBV	1794457	2596864	60.04%	7.202%
92	18.145	2556	2560	2567	rBV	62341	94130	2.18%	0.261%
93	18.351	2593	2595	2598	rBV4	7828	8692	0.20%	0.024%
94	19.221	2739	2743	2746	rBV3	24412	33180	0.77%	0.092%
95	19.292	2753	2755	2760	rVB	26631	35354	0.82%	0.098%
96	19.392	2768	2772	2779	rBV5	43291	66733	1.54%	0.185%
97	19.651	2811	2816	2824	rBV	2392533	3157497	73.00%	8.756%
98	21.427	3113	3118	3124	rBV	696051	955370	22.09%	2.649%
99	23.762	3507	3515	3525	rVB2	1653530	3314576	76.64%	9.192%
100	23.927	3537	3543	3555	rVB	455977	972357	22.48%	2.697%

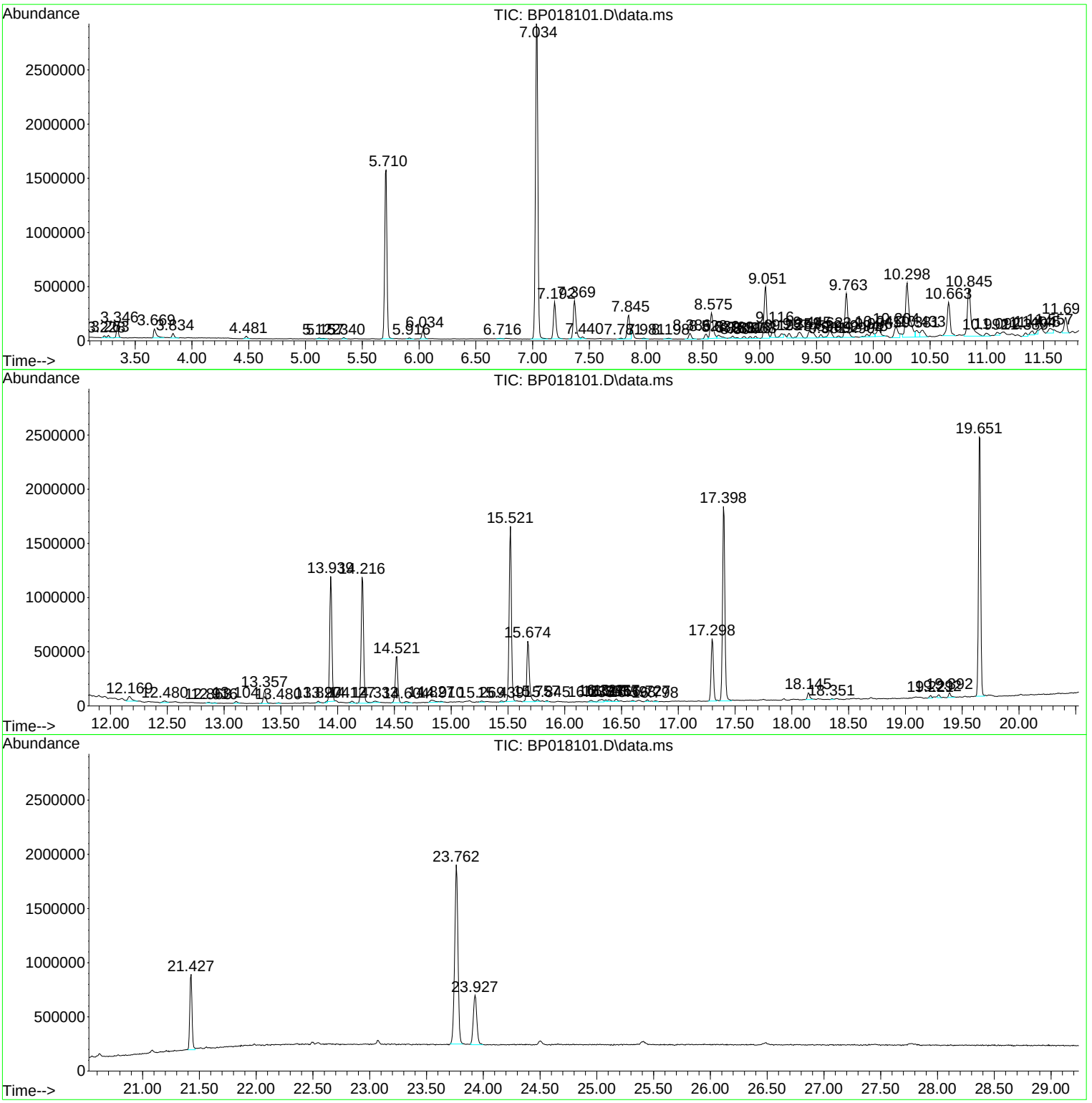
Sum of corrected areas: 36059037

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018101.D
Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH347

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018101.D
Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH347

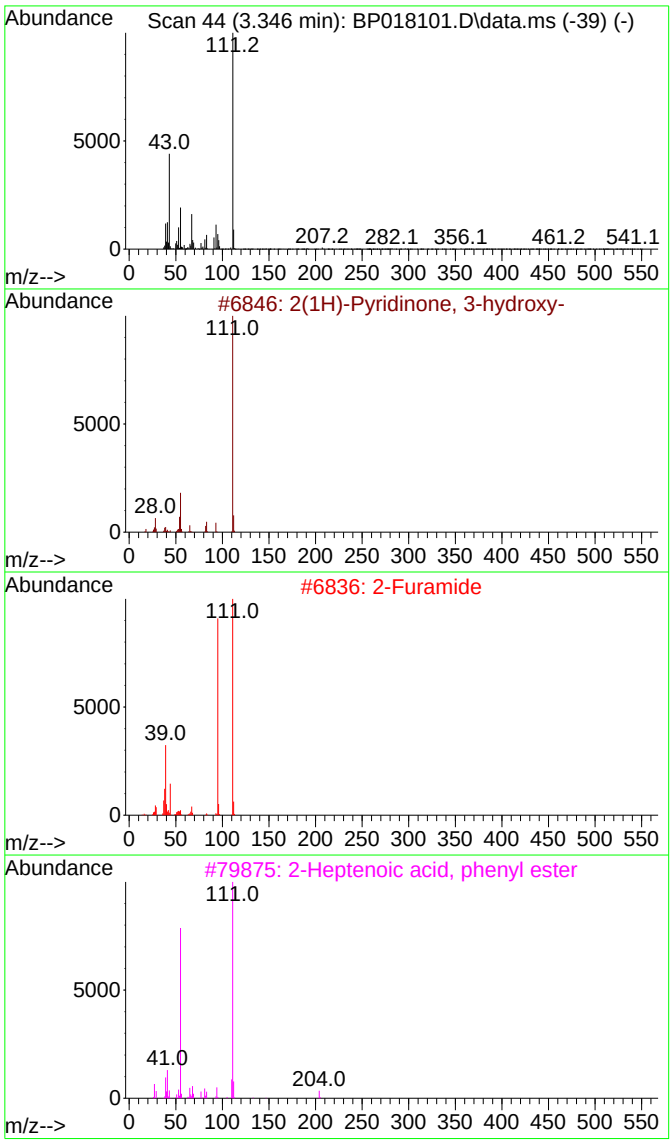
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

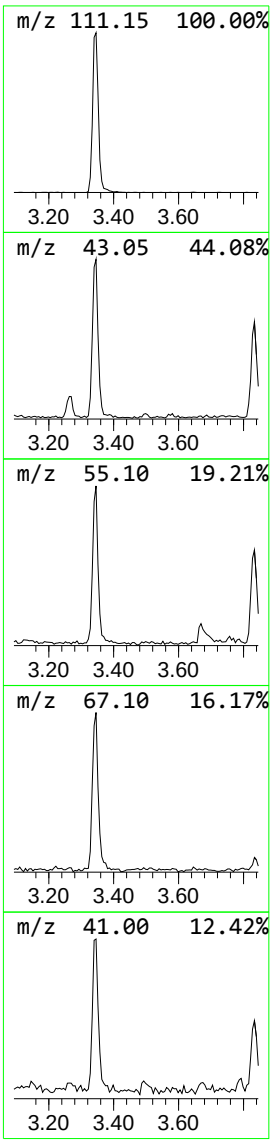
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.346	6.79 ng/ul	129563	1,4-Dichlorobenzene-d4	7.845

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(1H)-Pyridinone, 3-hydroxy-	111	C5H5NO2	016867-04-2	45
2	2-Furamide	111	C5H5NO2	000609-38-1	42
3	2-Heptenoic acid, phenyl ester	204	C13H16O2	1000406-90-4	40
4	1-Cyclohexanol, 2-[1-(phenylsulf...	252	C13H16O3S	1000196-78-3	39
5	2-Thiophenecarboxylic acid, 2-et...	238	C13H18O2S	1010278-88-5	39





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018101.D
Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH347

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

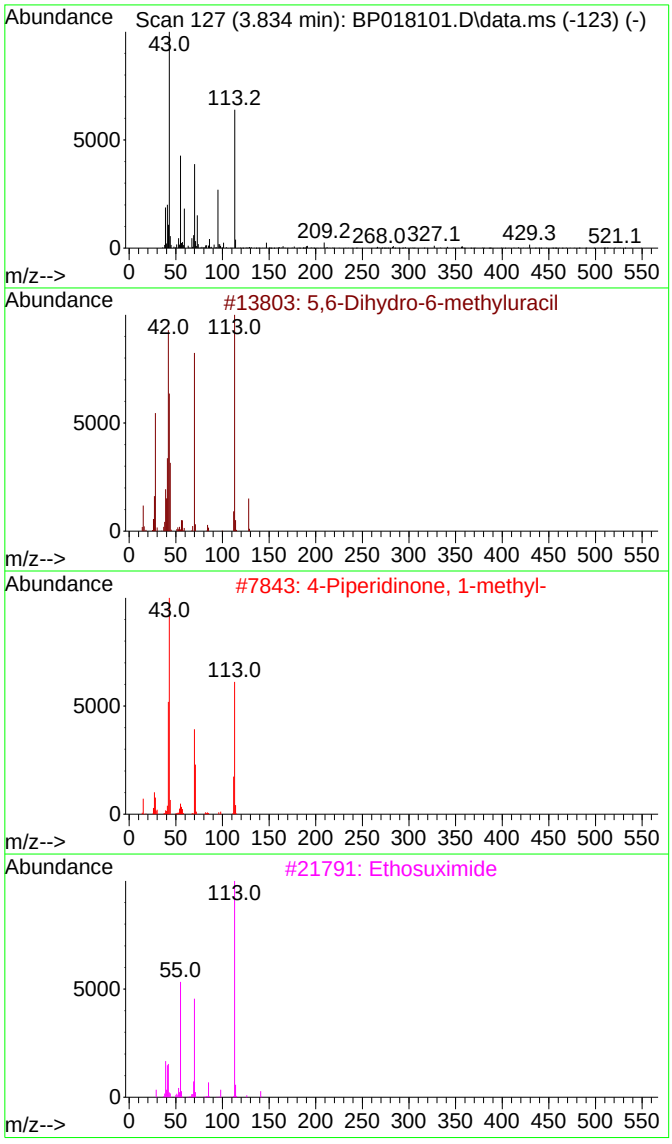
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

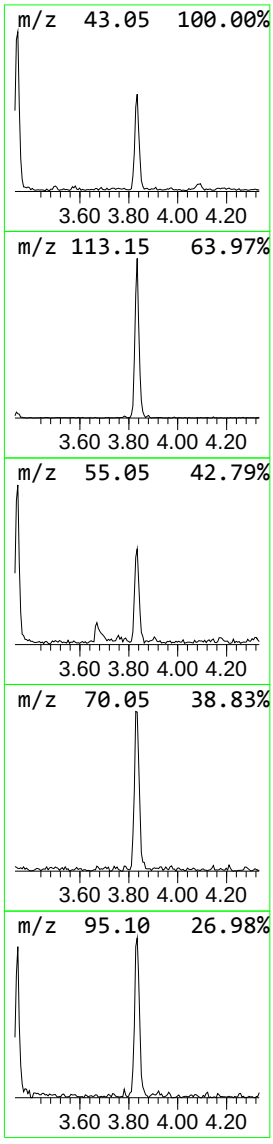
****

Peak Number 2 unknown-02 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.834	3.17 ng/ul	60385	1,4-Dichlorobenzene-d4	7.845

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,6-Dihydro-6-methyluracil	128	C5H8N2O2	002434-49-3	47
2	4-Piperidinone, 1-methyl-	113	C6H11NO	001445-73-4	47
3	Ethosuximide	141	C7H11NO2	000077-67-8	47
4	1-[2-(3-Chloro-phenoxy)-ethyl]-4...	254	C13H19ClN2O	337312-00-2	47
5	Furan, tetrahydro-2,2,5,5-tetram...	128	C8H16O	015045-43-9	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018101.D
Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH347

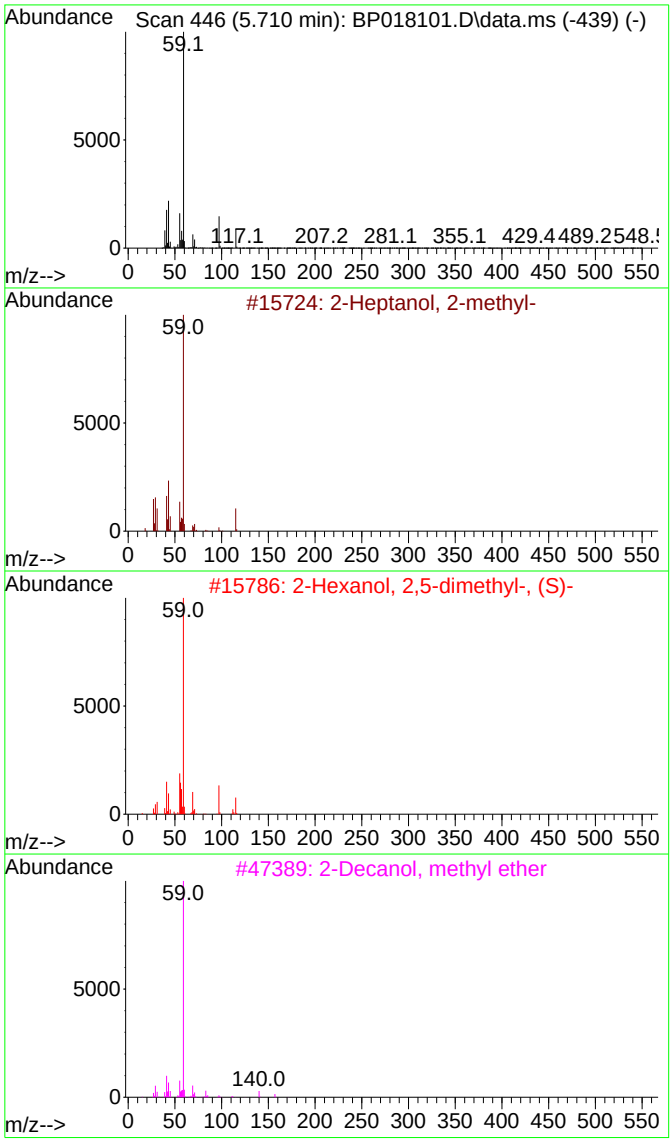
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

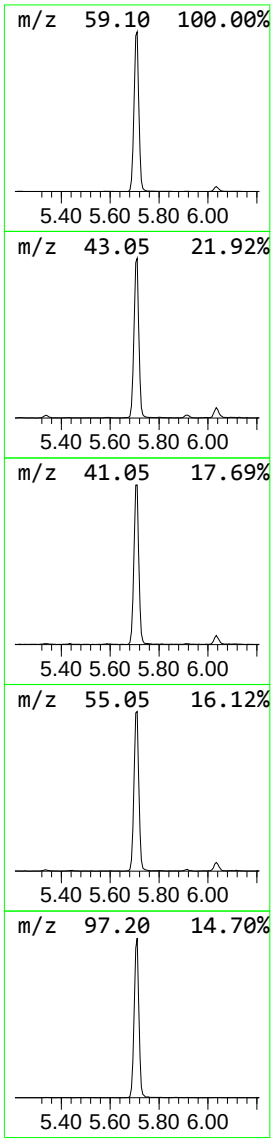
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 2-Heptanol, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.710	113.74 ng/ul	2168740	1,4-Dichlorobenzene-d4	7.845

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanol, 2-methyl-	130	C8H18O	000625-25-2	64
2	2-Hexanol, 2,5-dimethyl-, (S)-	130	C8H18O	003730-60-7	64
3	2-Decanol, methyl ether	172	C11H24O	1000333-83-9	50
4	2-Hexanol, methyl ether	116	C7H16O	1000333-92-0	45
5	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	45





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018101.D
Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH347

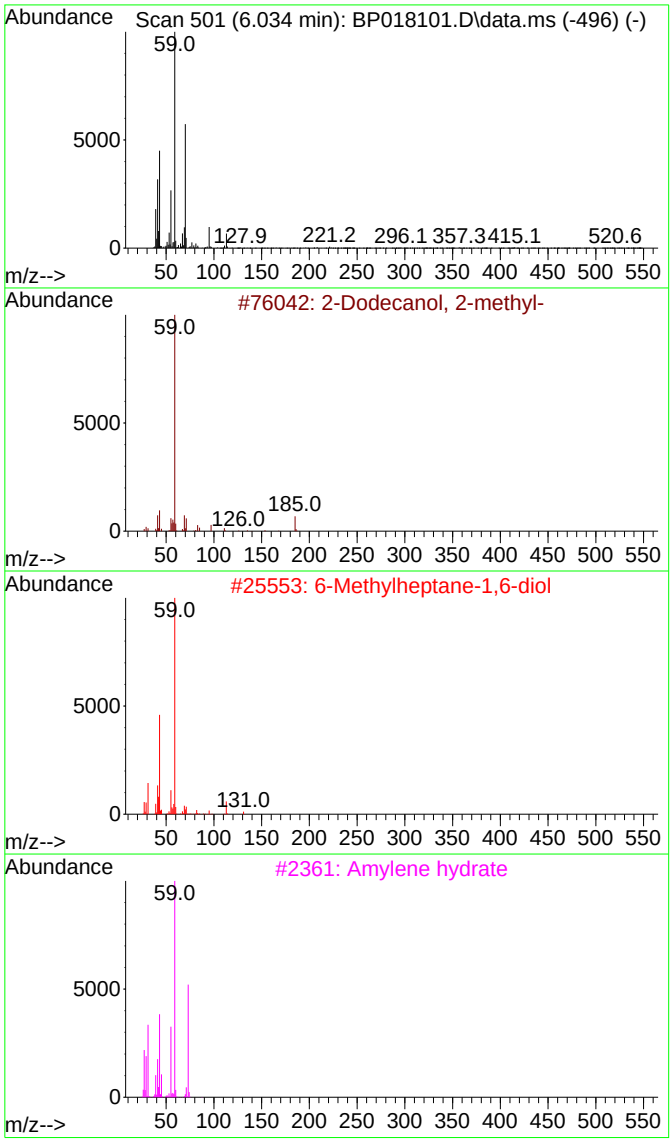
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

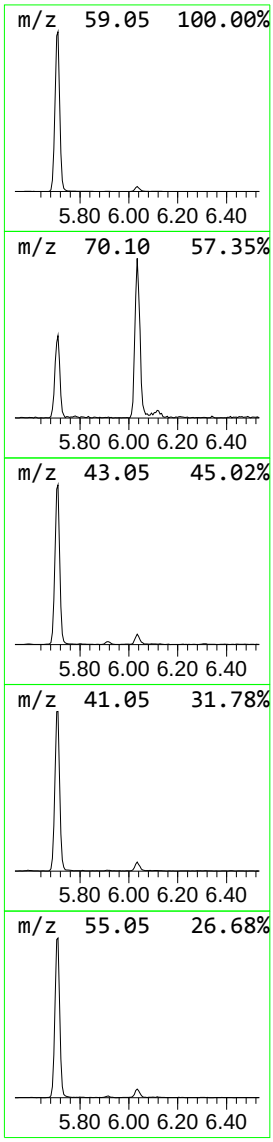
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 unknown-03 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.034	5.55 ng/ul	105818	1,4-Dichlorobenzene-d4	7.845

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Dodecanol, 2-methyl-	200	C13H28O	001653-37-8	47
2	6-Methylheptane-1,6-diol	146	C8H18O2	005392-57-4	38
3	Amylene hydrate	88	C5H12O	000075-85-4	37
4	2-Hexanol, 2,3-dimethyl-	130	C8H18O	019550-03-9	28
5	2-Hexanol, 2,3-dimethyl-	130	C8H18O	019550-03-9	28





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018101.D
Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH347

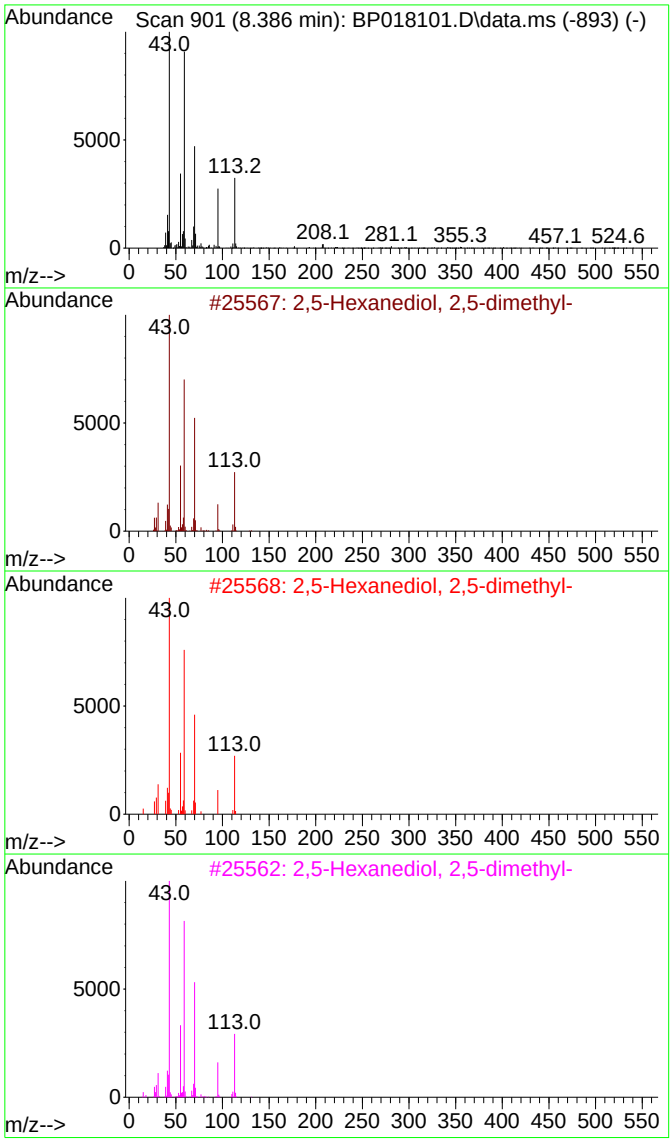
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

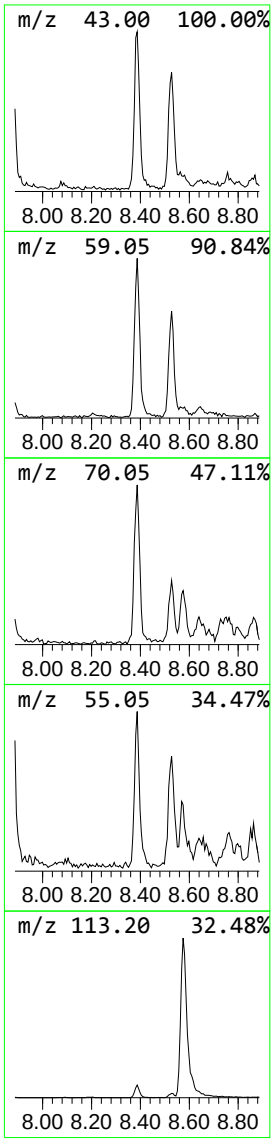
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 2,5-Hexanediol, 2,5-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.386	4.41 ng/ul	84124	1,4-Dichlorobenzene-d4	7.845

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,5-Hexanediol, 2,5-dimethyl-	146	C8H18O2	000110-03-2	78
2	2,5-Hexanediol, 2,5-dimethyl-	146	C8H18O2	000110-03-2	78
3	2,5-Hexanediol, 2,5-dimethyl-	146	C8H18O2	000110-03-2	78
4	2,5-Hexanediol, 2,5-dimethyl-	146	C8H18O2	000110-03-2	50
5	2,5-Hexanediol, 2,5-dimethyl-	146	C8H18O2	000110-03-2	47





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Data File : BP018101.D
Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH347

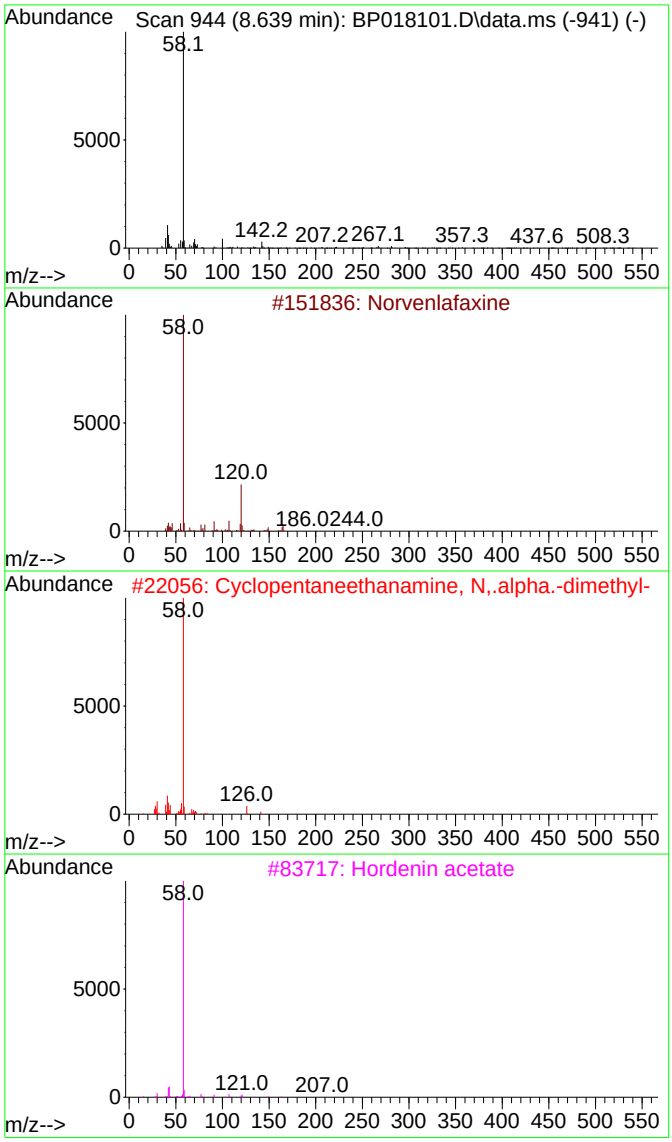
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

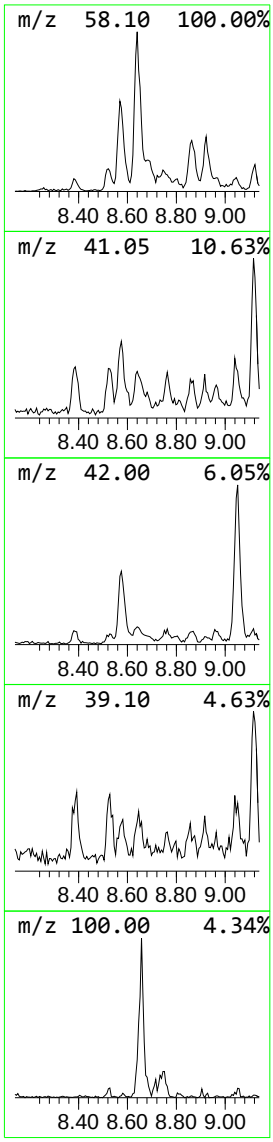
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Norvenlafaxine Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.639	3.58 ng/ul	68300	1,4-Dichlorobenzene-d4	7.845

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Norvenlafaxine	263	C16H25NO2	130198-38-8	64
2	Cyclopentaneethanamine, N,.alpha...	141	C9H19N	000102-45-4	50
3	Hordenin acetate	207	C12H17NO2	095469-40-2	50
4	.alpha.-Pinene, 10-(dimethylamin...	193	C13H23N	015063-97-5	50
5	2,3 Methylenedioxymethcathinone	207	C11H13NO3	1010386-29-7	50





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Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH347

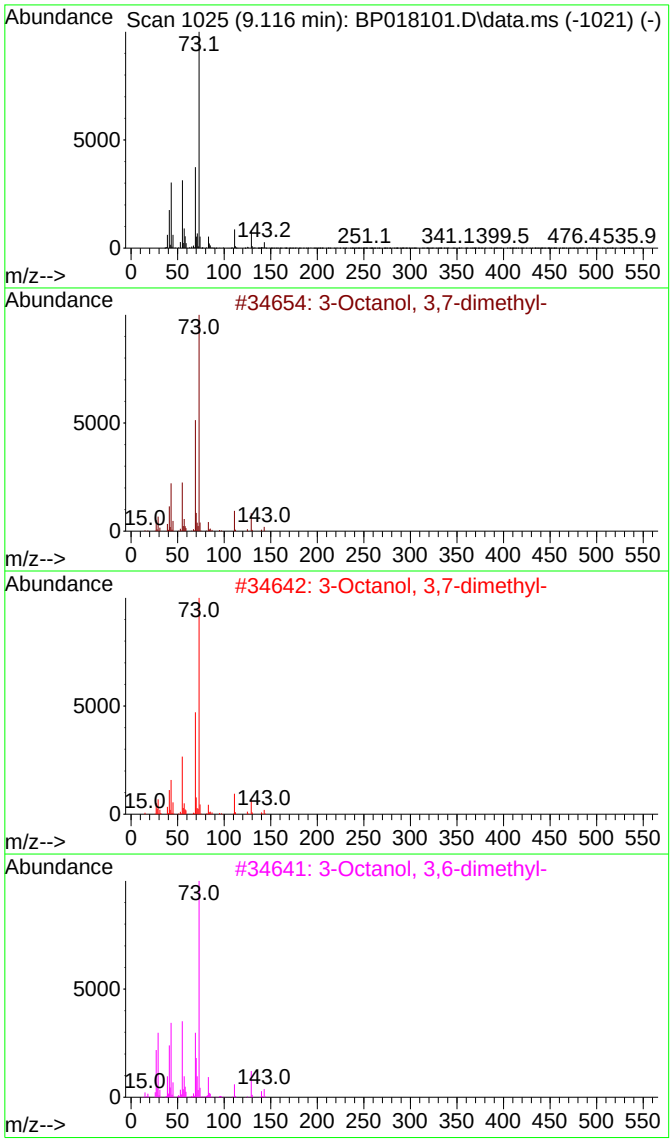
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

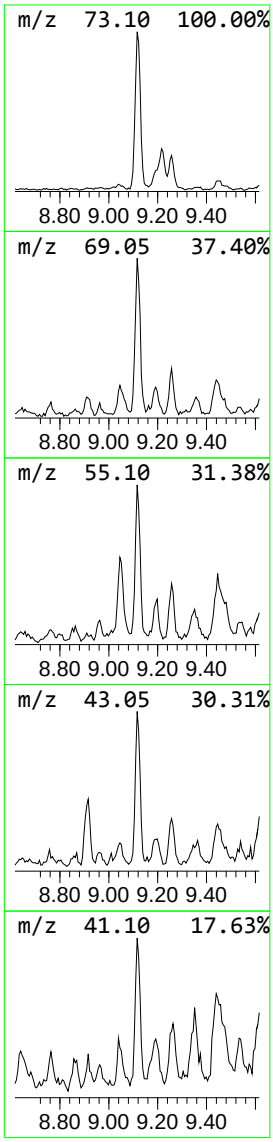
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 3-Octanol, 3,7-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.116	7.82 ng/ul	149017	1,4-Dichlorobenzene-d4	7.845

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	64
2	3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	50
3	3-Octanol, 3,6-dimethyl-	158	C10H22O	000151-19-9	50
4	3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	43
5	3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	40





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Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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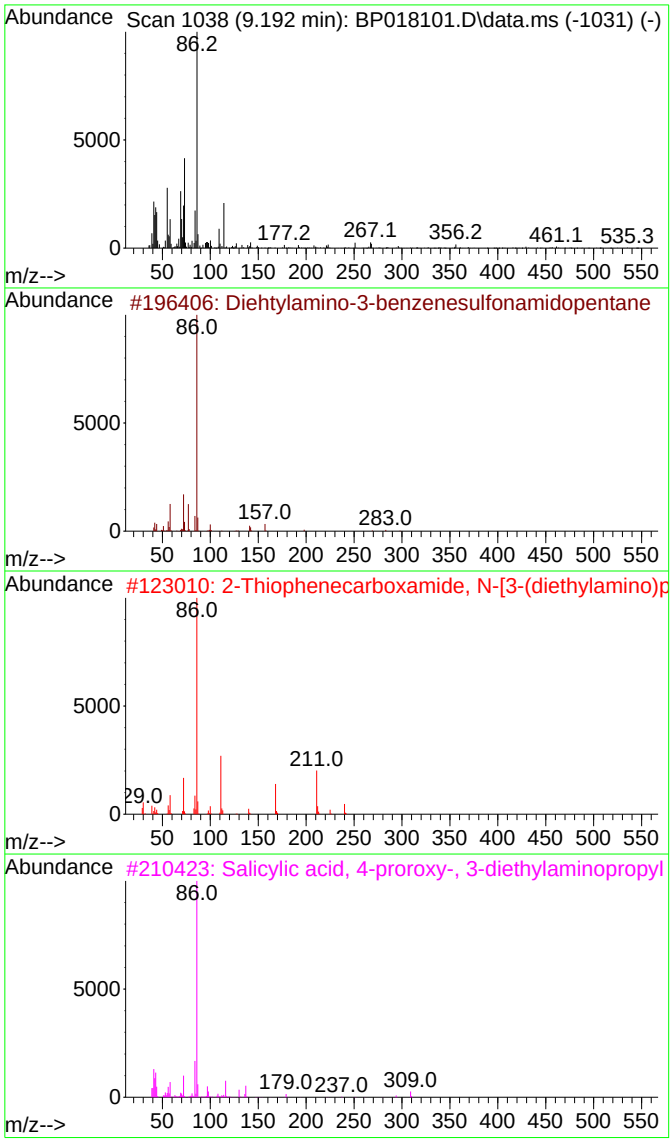
Quant Title : SVOA CALIBRATION

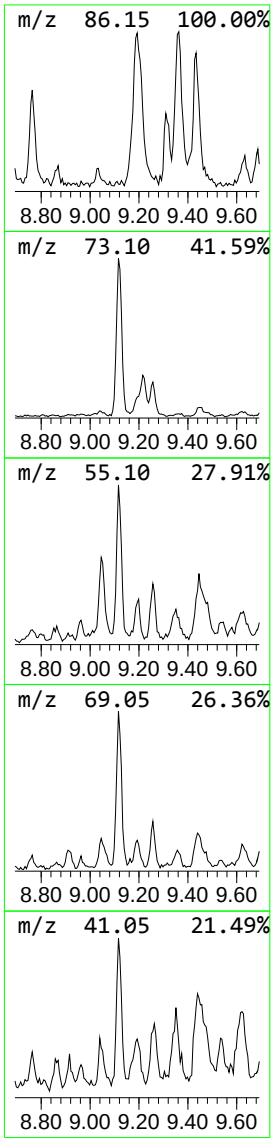
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Diehtylamino-3-benzenesulfo... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.192	3.75 ng/ul	71534	1,4-Dichlorobenzene-d4	7.845

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diehtylamino-3-benzenesulfonamid...	298	C15H26N2O2S	025132-10-9	58
2	2-Thiophenecarboxamide, N-[3-(di...	240	C12H20N2OS	1000319-44-0	53
3	Salicylic acid, 4-proroxy-, 3-di...	309	C17H27NO4	055828-24-5	52
4	2-Aminooctane, N-acetyl-	171	C10H21NO	023602-00-8	50
5	o-(iso-Pentyl) S-(2-diethylamino...	295	C13H30NO2PS	1000273-62-3	50





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Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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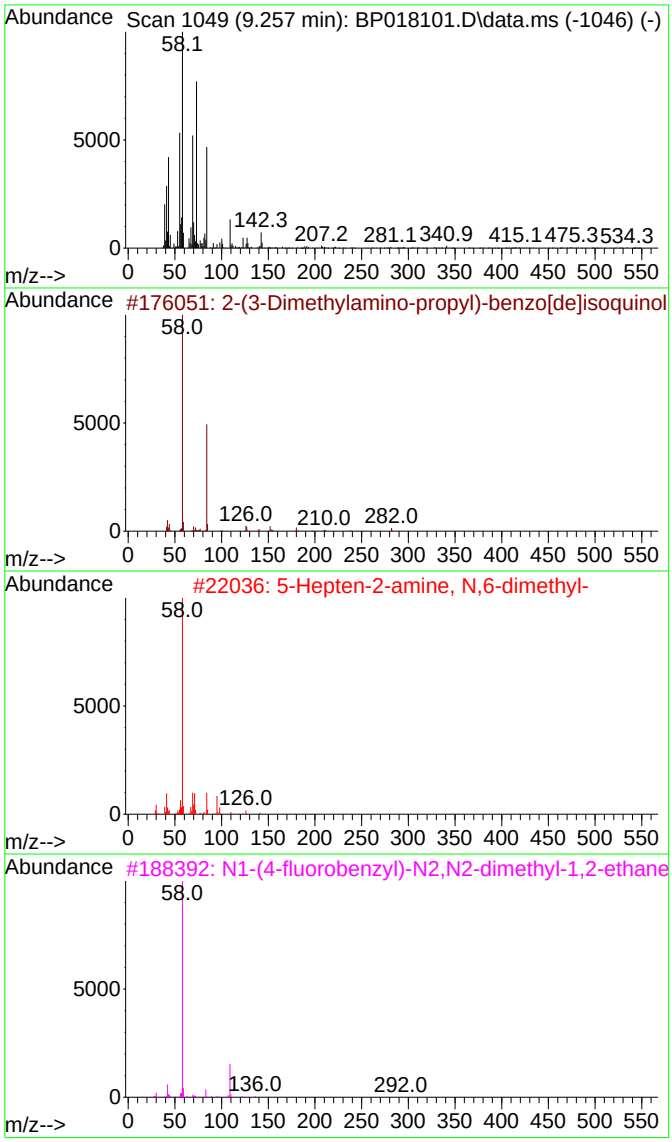
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

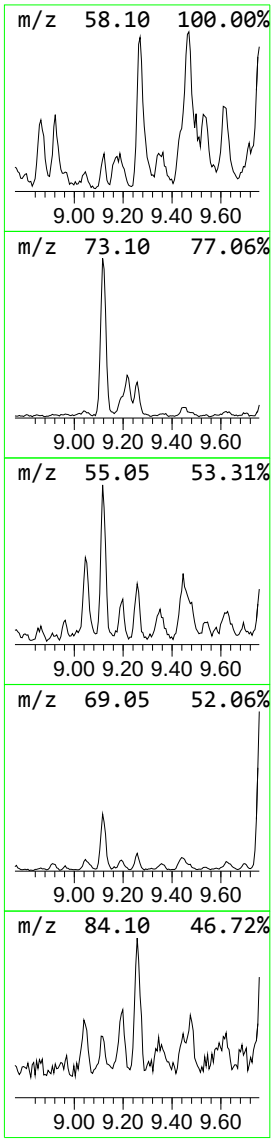
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 2-(3-Dimethylamino-propyl)-... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.257	2.32 ng/ul	66995	Naphthalene-d8	10.663

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(3-Dimethylamino-propyl)-benzo...	282	C17H18N2O2	1000318-08-6	53
2	5-Hepten-2-amine, N,6-dimethyl-	141	C9H19N	000503-01-5	43
3	N1-(4-fluorobenzyl)-N2,N2-dimeth...	292	C13H16F4N2O	1000511-02-0	43
4	4-Piperidineamin, N-(2-(dimethyl...	281	C12H22F3N3O	1000470-40-0	43
5	Methanamine, 1-(dicyclohexylphos...	255	C15H30NP	066830-54-4	38





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Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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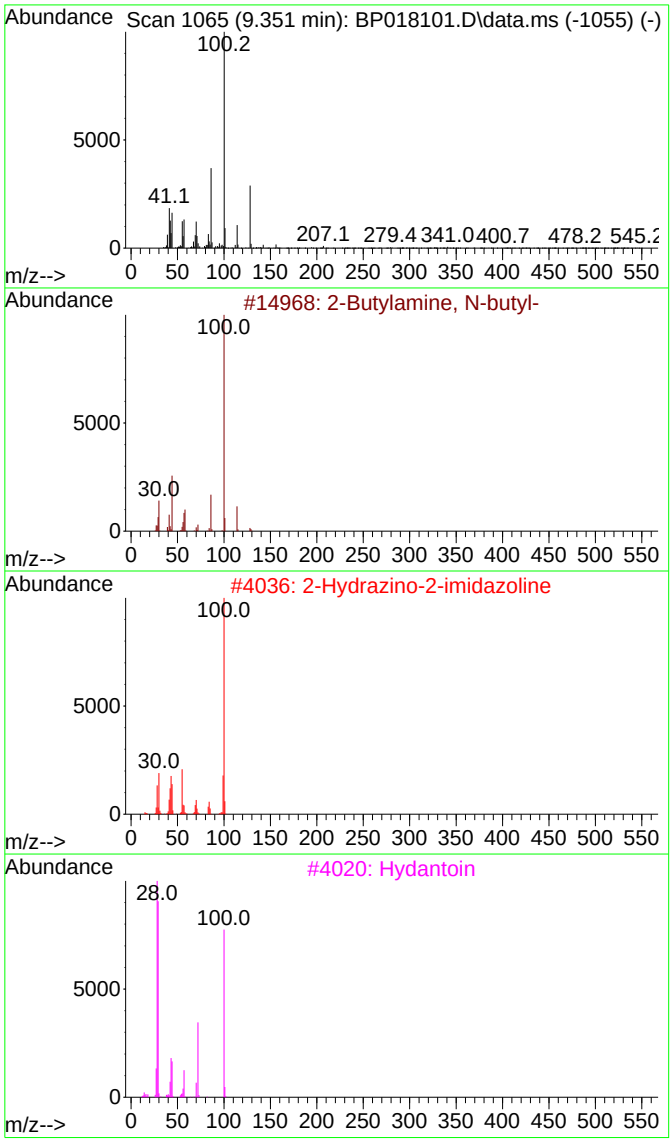
Quant Title : SVOA CALIBRATION

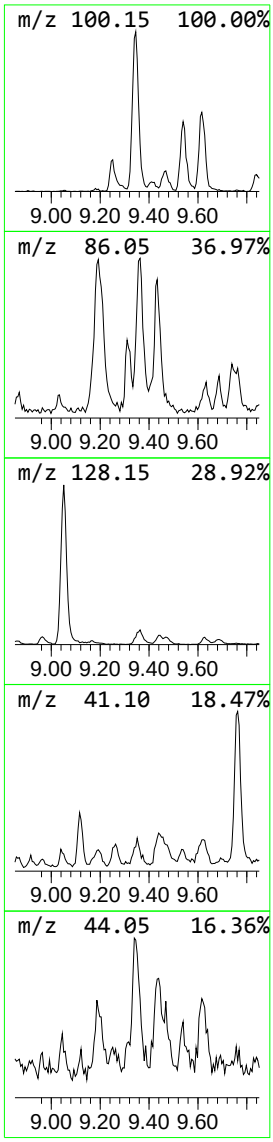
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 2-Butylamine, N-butyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.351	4.61 ng/ul	132768	Naphthalene-d8	10.663

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butylamine, N-butyl-	129	C8H19N	1000463-86-0	64
2	2-Hydrazino-2-imidazoline	100	C3H8N4	1000239-54-7	52
3	Hydantoin	100	C3H4N2O2	000461-72-3	52
4	4-Bromo-5-[3-dibutylaminopropyla...	358	C15H27BrN4O	1000256-15-1	50
5	4-Chloro-N-butylcathinone	239	C13H18ClNO	1000466-85-1	47





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Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH347

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Quant Title : SVOA CALIBRATION

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

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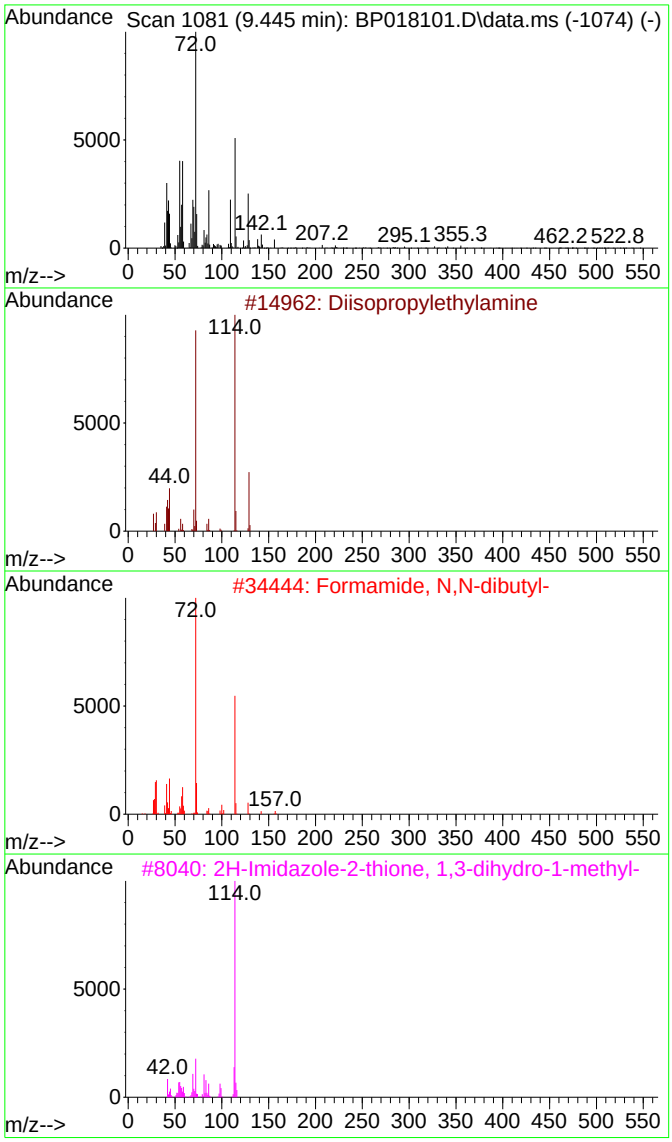
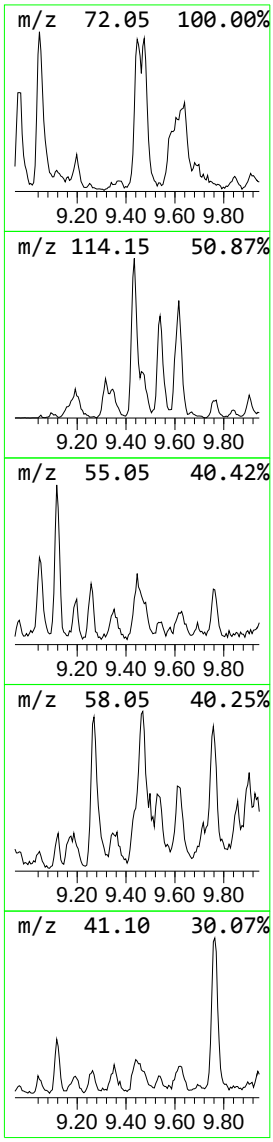
Peak Number 11 unknown-04 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.445	7.78 ng/ul	224286	Naphthalene-d8	10.663

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diisopropylethylamine	129	C8H19N	007087-68-5	43
2	Formamide, N,N-dibutyl-	157	C9H19NO	000761-65-9	37
3	2H-Imidazole-2-thione, 1,3-dihyd...	114	C4H6N2S	000060-56-0	37
4	Fenfluramine, N-acetyl-	273	C14H18F3NO	1000448-49-5	32
5	N-Ethylamphetamine acetate	205	C13H19NO	072687-64-0	25





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Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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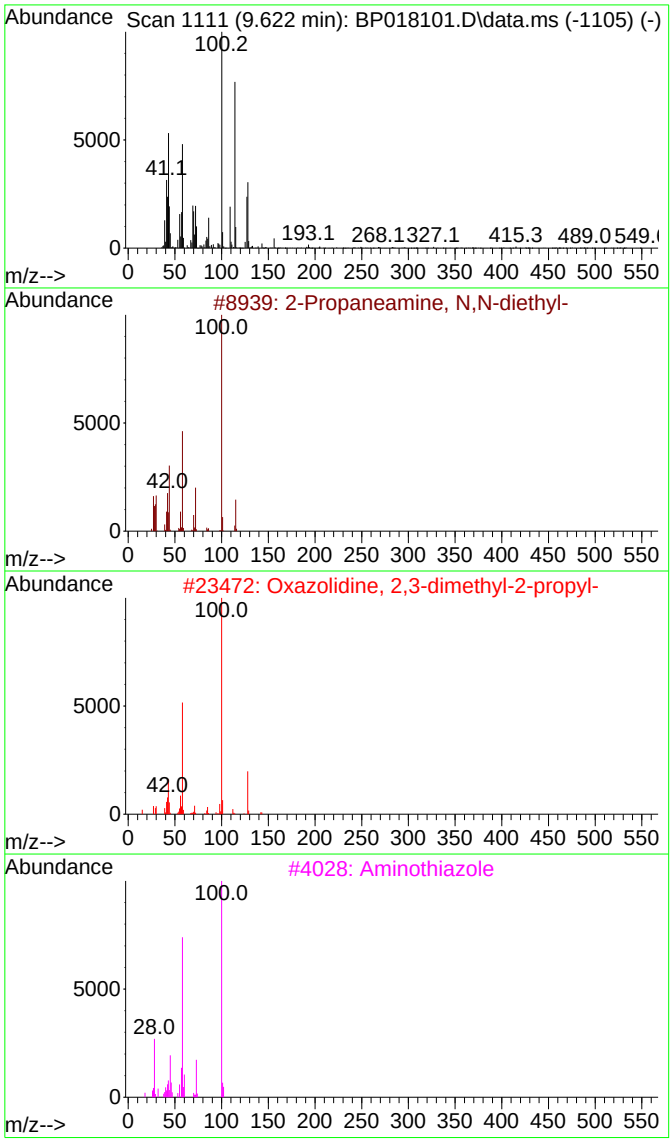
Quant Title : SVOA CALIBRATION

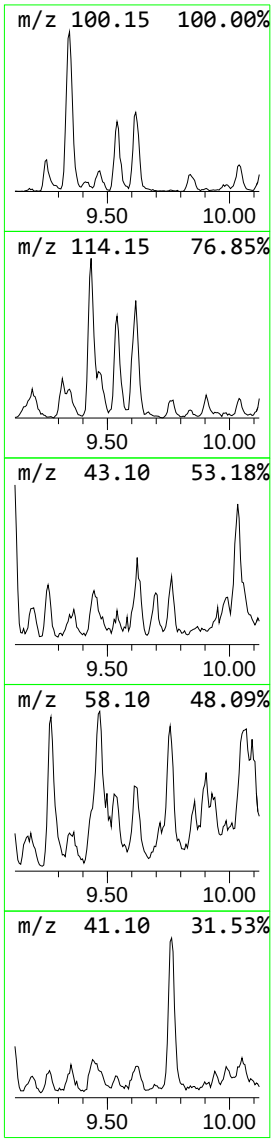
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 unknown-05 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.622	4.60 ng/ul	132754	Naphthalene-d8	10.663

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propaneamine, N,N-diethyl-	115	C7H17N	006006-15-1	35
2	Oxazolidine, 2,3-dimethyl-2-propyl-	143	C8H17NO	161500-41-0	35
3	Aminothiazole	100	C3H4N2S	000096-50-4	30
4	4-Heptanone	114	C7H14O	000123-19-3	27
5	Isobutyl S-2-dipropylaminoethyl ...	323	C15H34NO2PS	1010510-46-2	27





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Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Title : SVOA CALIBRATION

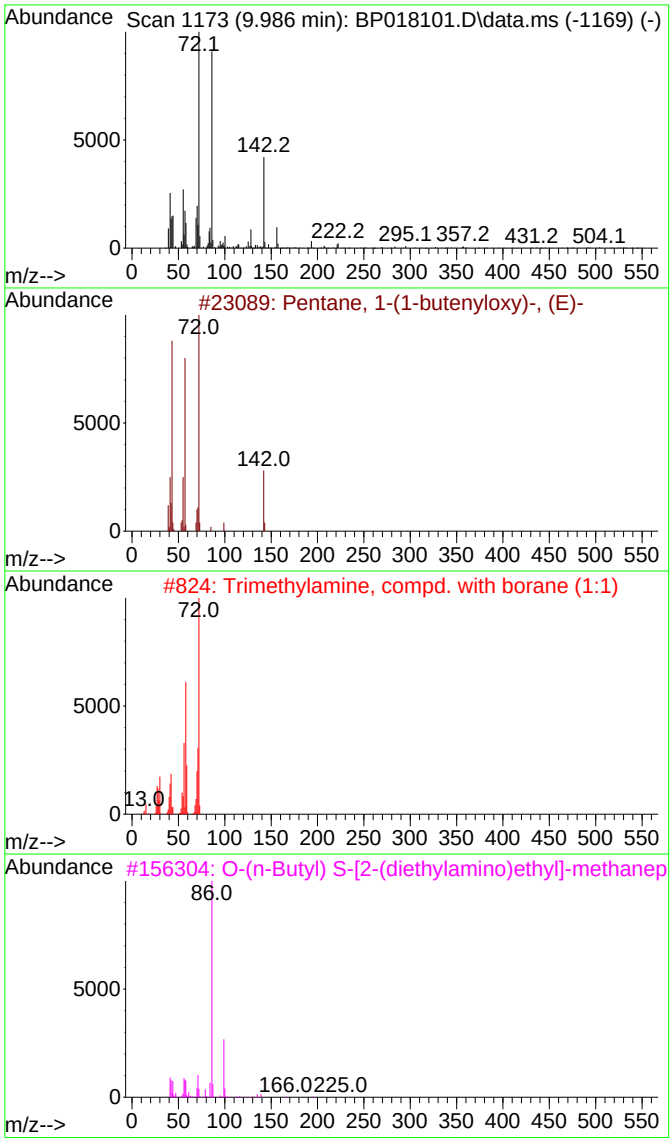
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

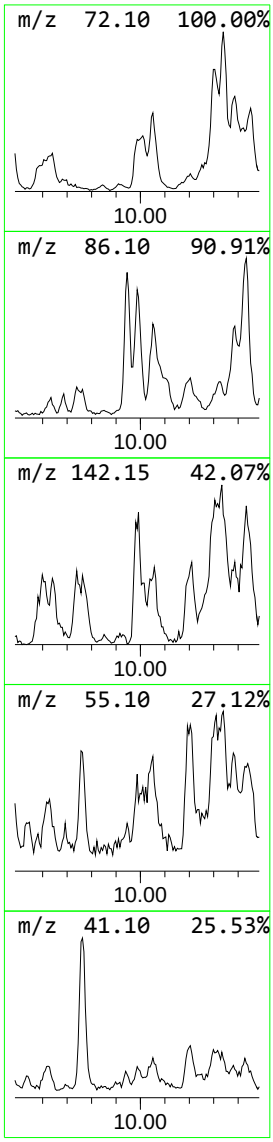
****

Peak Number 13 unknown-06 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.986	3.16 ng/ul	91146	Naphthalene-d8	10.663

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 1-(1-butenyloxy)-, (E)-	142	C9H18O	054004-25-0	38
2	Trimethylamine, compd. with bora...	73	C3H12BN	000075-22-9	27
3	O-(n-Butyl) S-[2-(diethylamino)e...	267	C11H26NO2PS	468712-10-9	27
4	Urea, N,N-dimethyl-N',N'-dibutyl-	200	C11H24N2O	1000439-30-5	27
5	9H-1,3,4,9-Tetraazafluorene-2-th...	301	C15H19N5S	1000304-53-7	22





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Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

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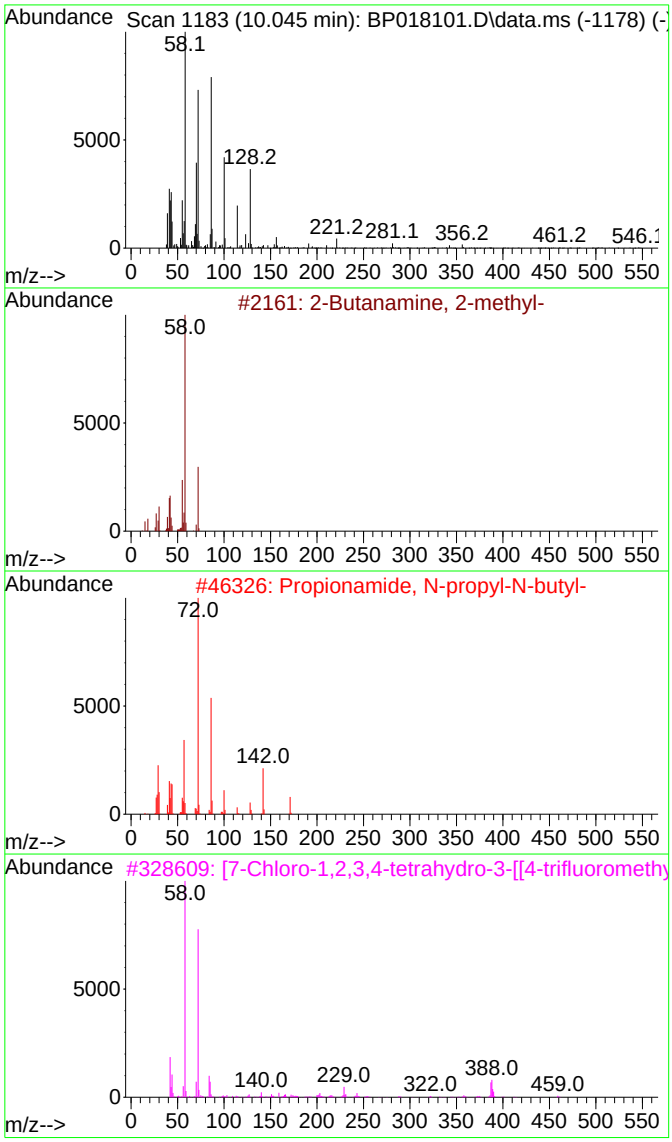
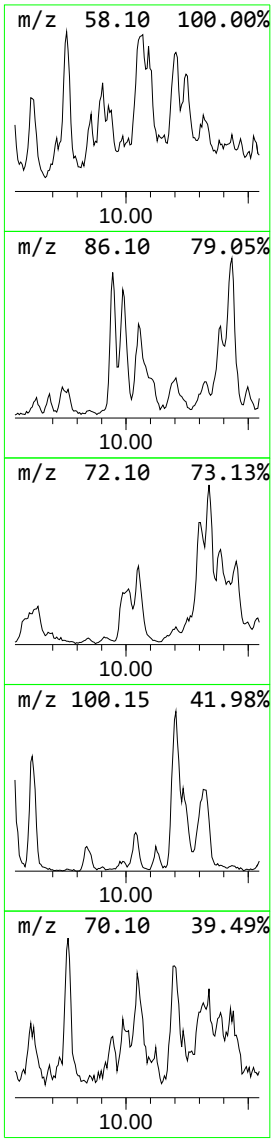
Peak Number 14 unknown-07 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.045	5.21 ng/ul	150338	Naphthalene-d8	10.663

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanamine, 2-methyl-	87	C5H13N	000594-39-8	35
2	Propionamide, N-propyl-N-butyl-	171	C10H21NO	1000438-44-6	35
3	[7-Chloro-1,2,3,4-tetrahydro-3-[...	459	C25H25ClF3N3	144155-87-3	27
4	Undec-10-ynoylamide, N-(2-butyl)...	265	C17H31NO	1000491-72-1	14
5	Methanetricarboxaldehyde	100	C4H4O3	018655-47-5	14





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ALS Vial : 81 Sample Multiplier: 1

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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

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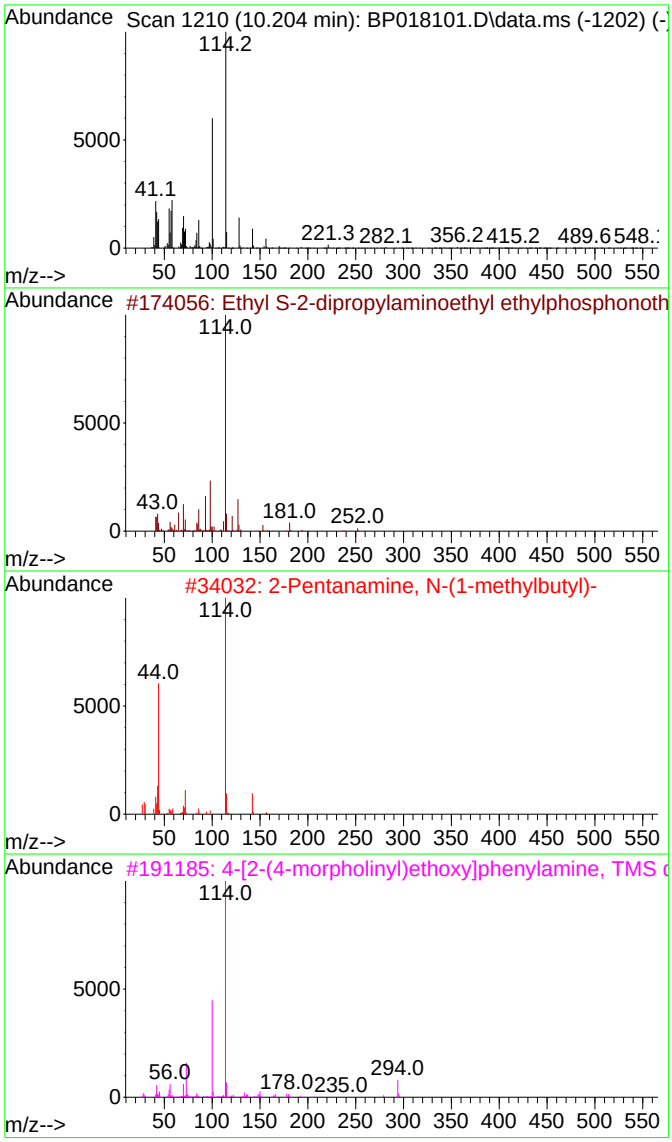
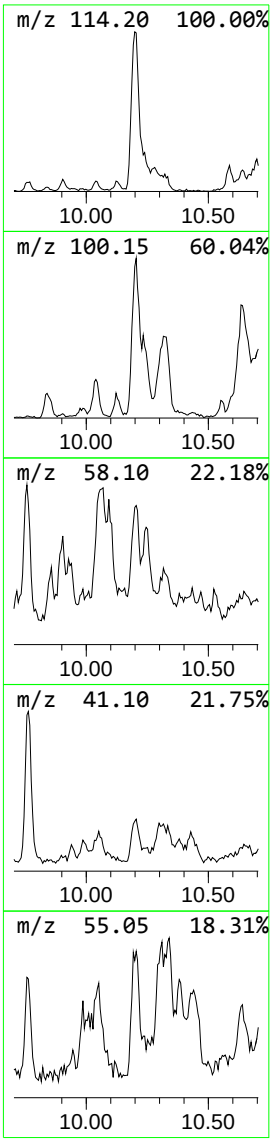
Peak Number 15 unknown-08 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.204	8.04 ng/ul	231693	Naphthalene-d8	10.663

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethyl S-2-dipropylaminoethyl eth...	281	C12H28NO2PS	1000510-44-7	43
2	2-Pentanamine, N-(1-methylbutyl)-	157	C10H23N	040221-44-1	43
3	4-[2-(4-morpholinyl)ethoxy]pheny...	294	C15H26N2O2Si	1000479-02-9	40
4	5-Fluoro-4(3H)-pyrimidinone	114	C4H3FN2O	000671-35-2	38
5	Piperazine, 1,4-dimethyl-	114	C6H14N2	000106-58-1	38





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ClientSampleId :
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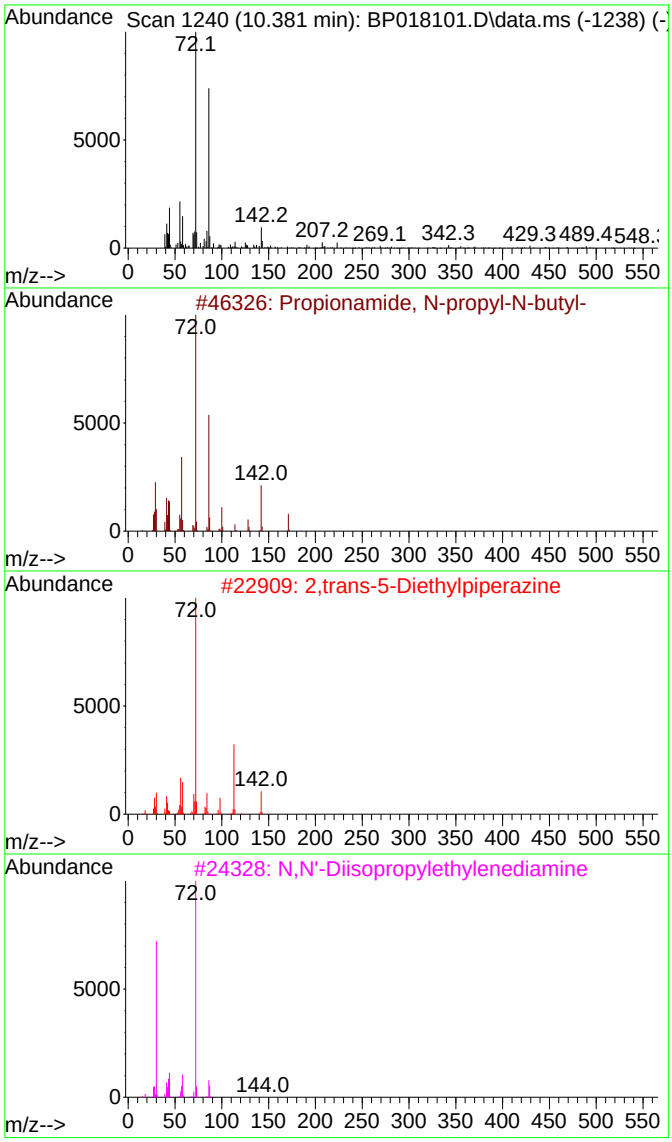
Quant Title : SVOA CALIBRATION

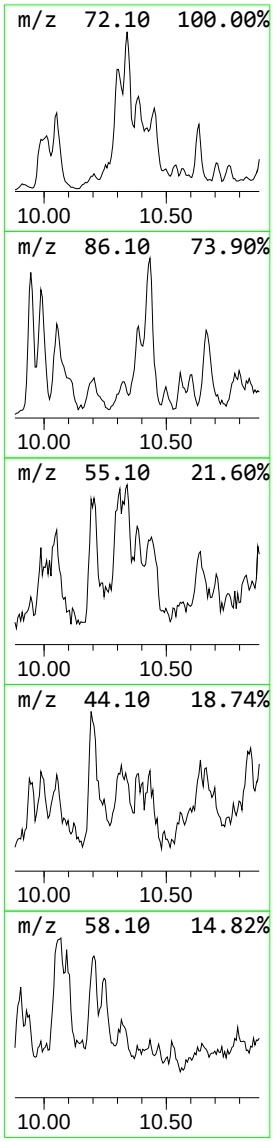
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Propionamide, N-propyl-N-bu... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.381	4.26 ng/ul	122846	Naphthalene-d8	10.663

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propionamide, N-propyl-N-butyl-	171	C10H21NO	1000438-44-6	53
2	2,trans-5-Diethylpiperazine	142	C8H18N2	006189-24-8	38
3	N,N'-Diisopropylethylenediamine	144	C8H20N2	004013-94-9	35
4	O-(n-Butyl) S-[2-(diethylamino)e...	267	C11H26NO2PS	468712-10-9	35
5	L-Valine, N-(3-cyclopentylpropio...	311	C18H33NO3	1000346-65-6	35





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018101.D
Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH347

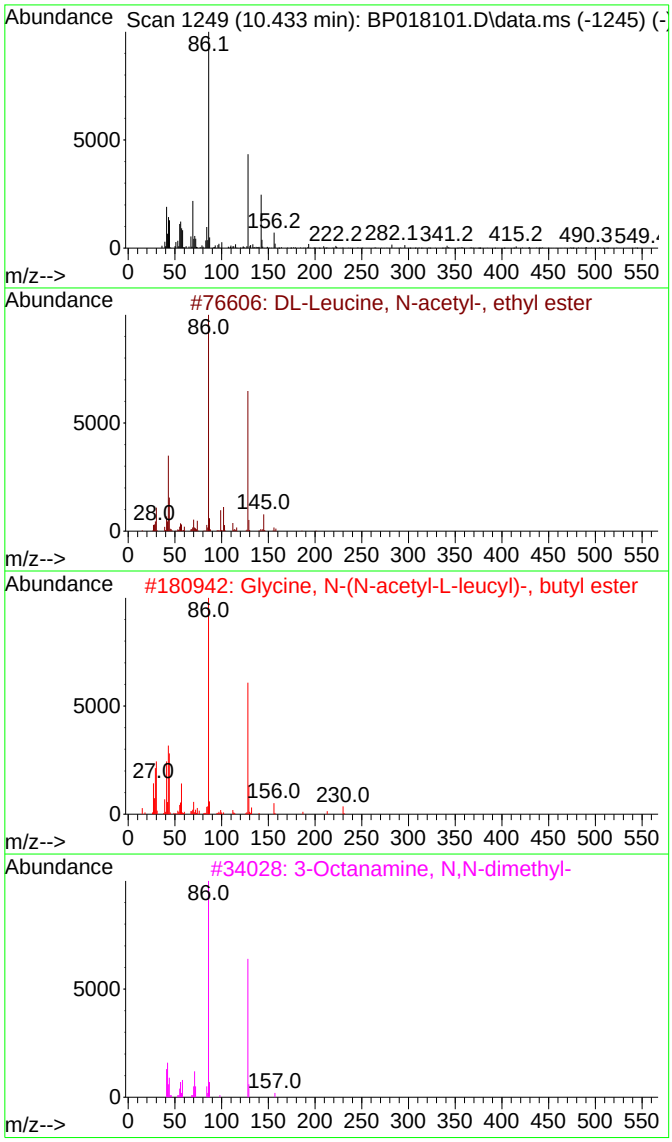
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

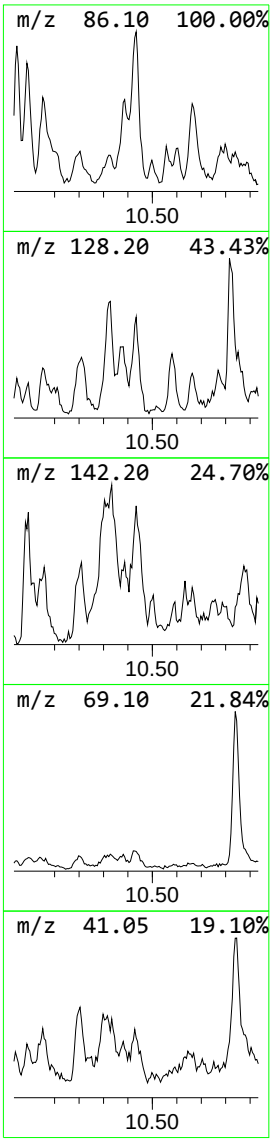
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 DL-Leucine, N-acetyl-, ethy... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.434	5.20 ng/ul	149875	Naphthalene-d8	10.663

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	DL-Leucine, N-acetyl-, ethyl ester	201	C10H19NO3	004071-36-7	59
2	Glycine, N-(N-acetyl-L-leucyl)-,...	286	C14H26N2O4	055712-44-2	59
3	3-Octanamine, N,N-dimethyl-	157	C10H23N	024539-82-0	58
4	2-(4-Tert-butylphenoxy)ethanamin...	277	C16H23NO3	1000502-14-3	53
5	Methyl S-2-diethylaminoethyl met...	225	C8H20NO2PS	203385-78-8	43





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Data File : BP018101.D
Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH347

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

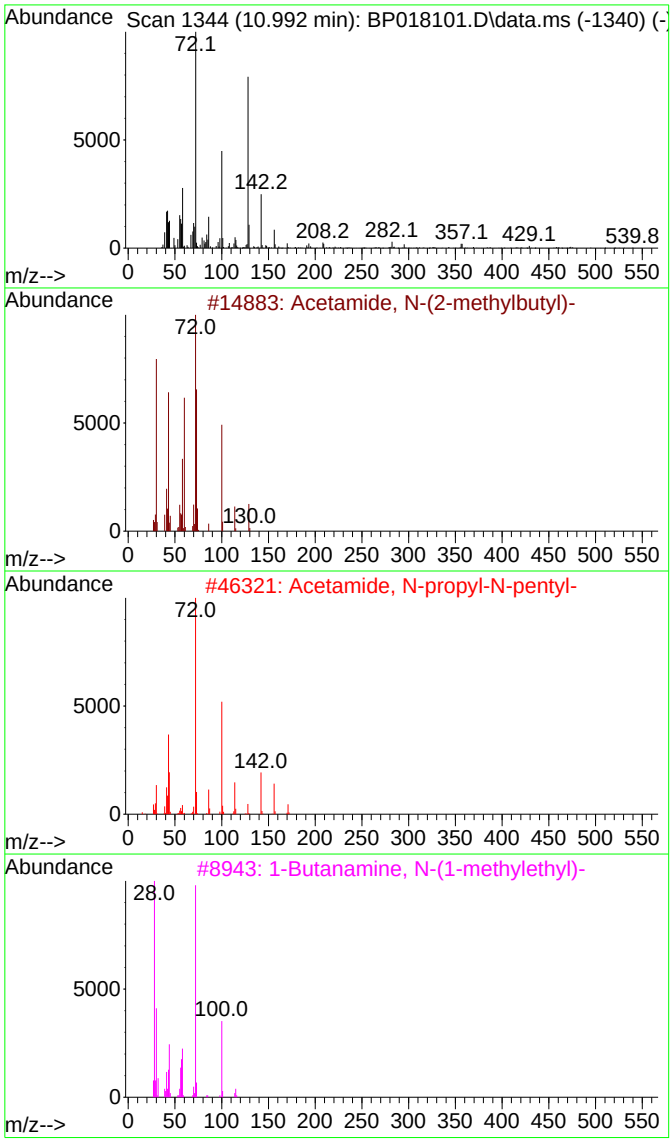
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

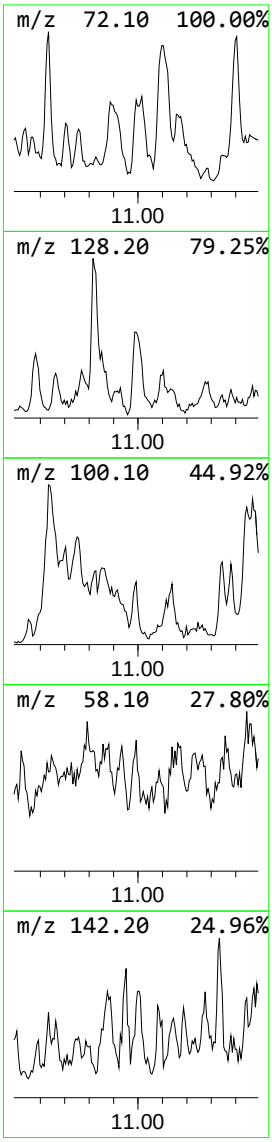
****

Peak Number 18 unknown-09 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.992	2.12 ng/ul	61195	Naphthalene-d8	10.663

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(2-methylbutyl)-	129	C7H15NO	054824-90-7	46
2	Acetamide, N-propyl-N-pentyl-	171	C10H21NO	1000438-05-1	43
3	1-Butanamine, N-(1-methylethyl)-	115	C7H17N	039099-23-5	35
4	1,3-Dimethyl-3,4,5,6-tetrahydro-...	128	C6H12N2O	007226-23-5	30
5	2,4(1H,3H)-Pyrimidinedione, dihy...	128	C5H8N2O2	001672-04-4	27





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Data File : BP018101.D
Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH347

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

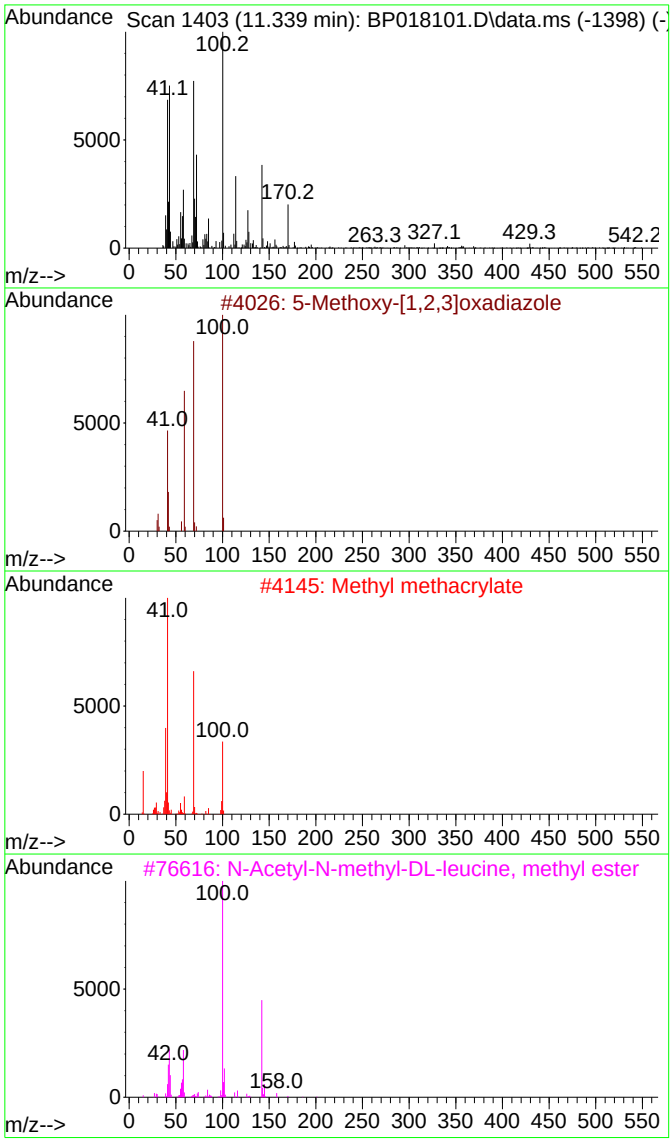
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

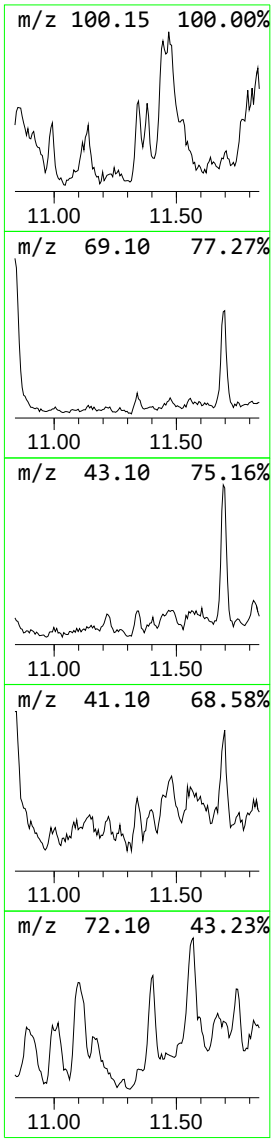
****

Peak Number 19 unknown-10 Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.339	2.01 ng/ul	58058	Naphthalene-d8	10.663

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Methoxy-[1,2,3]oxadiazole	100	C3H4N2O2	1000322-54-7	43
2	Methyl methacrylate	100	C5H8O2	000080-62-6	43
3	N-Acetyl-N-methyl-DL-leucine, me...	201	C10H19NO3	1000453-80-5	30
4	2-Acetamidothiazole	142	C5H6N2OS	002719-23-5	27
5	2(5H)-Thiophenone	100	C4H4OS	003354-32-3	27





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018101.D
Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH347

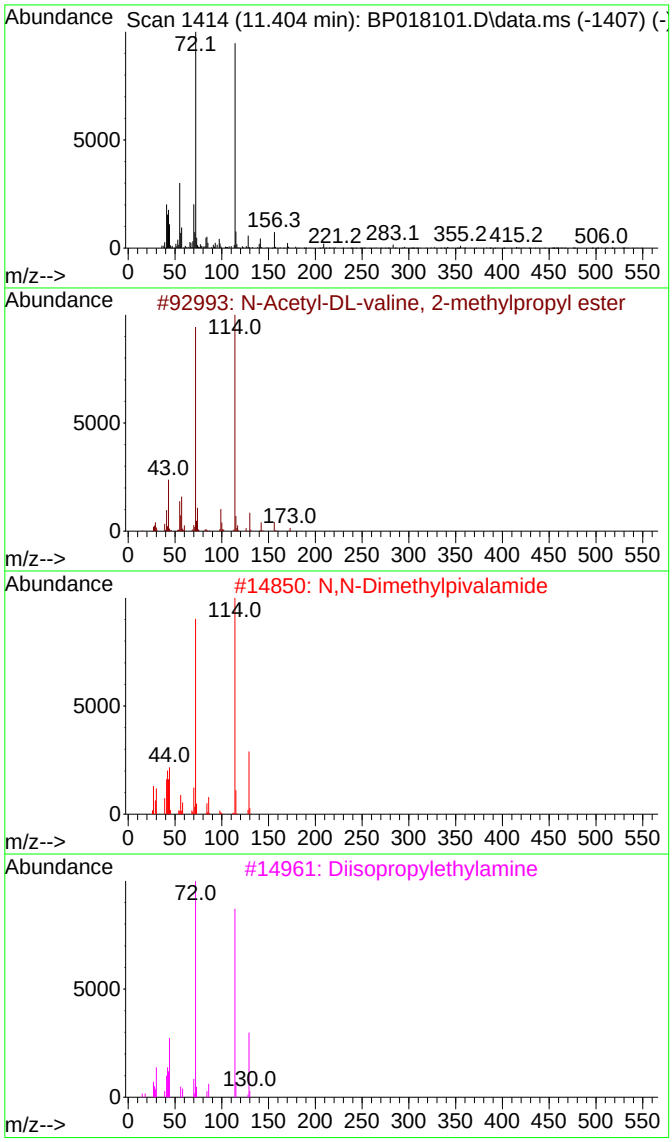
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

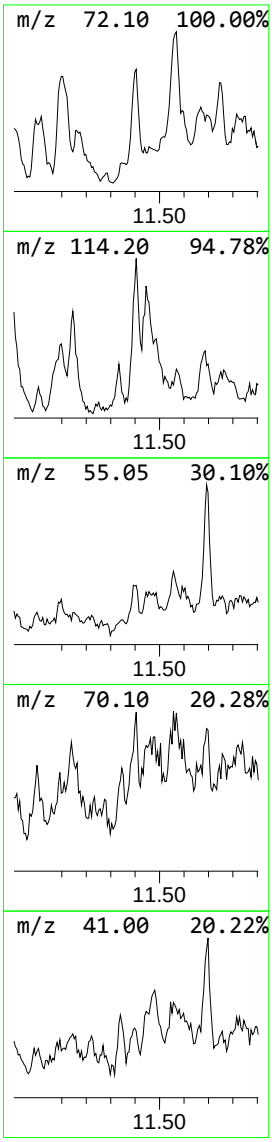
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 N-Acetyl-DL-valine, 2-methy... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.404	2.47 ng/ul	71250	Naphthalene-d8	10.663

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	N-Acetyl-DL-valine, 2-methylprop...	215	C11H21NO3	1000453-78-9	72
2	N,N-Dimethylpivalamide	129	C7H15NO	024331-71-3	64
3	Diisopropylethylamine	129	C8H19N	007087-68-5	64
4	Diisopropylethylamine	129	C8H19N	007087-68-5	64
5	DL-Valine, N-acetyl-, methyl ester	173	C8H15NO3	052152-47-3	45





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018101.D
Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH347

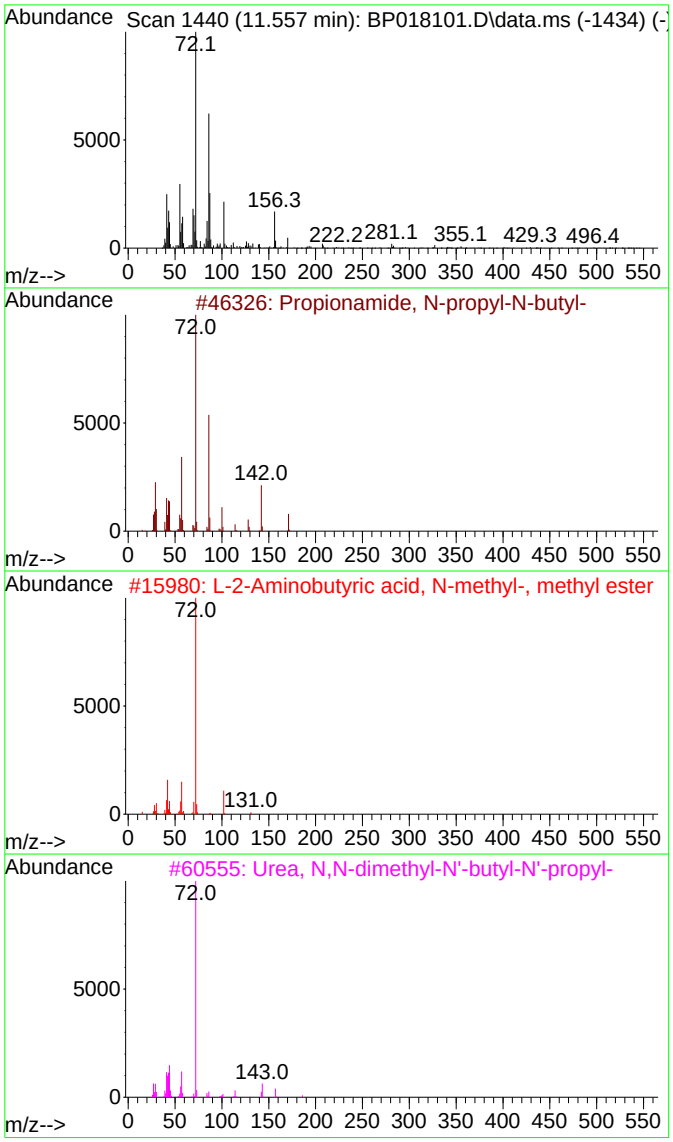
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

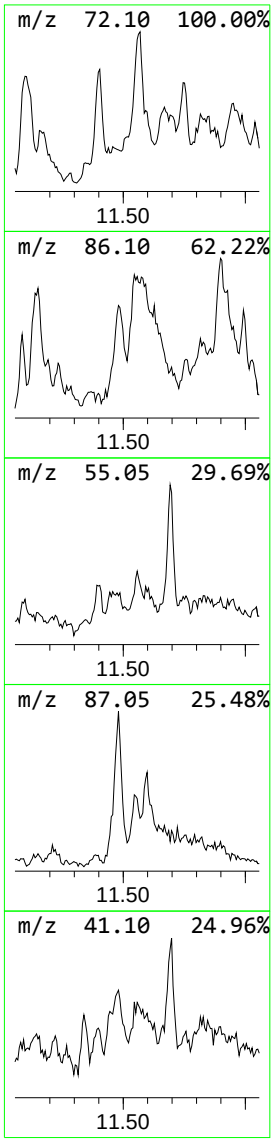
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 unknown-11 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.557	3.87 ng/ul	111550	Naphthalene-d8	10.663

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propionamide, N-propyl-N-butyl-	171	C10H21NO	1000438-44-6	50
2	L-2-Aminobutyric acid, N-methyl-...	131	C6H13NO2	1000332-98-7	38
3	Urea, N,N-dimethyl-N'-butyl-N'-p...	186	C10H22N2O	1000439-30-3	35
4	Acetaldehyde, methylhydrazone	72	C3H8N2	017167-73-6	27
5	1-Butanamine, N-(1-methylethyl)-	115	C7H17N	039099-23-5	27





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Data File : BP018101.D
Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH347

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

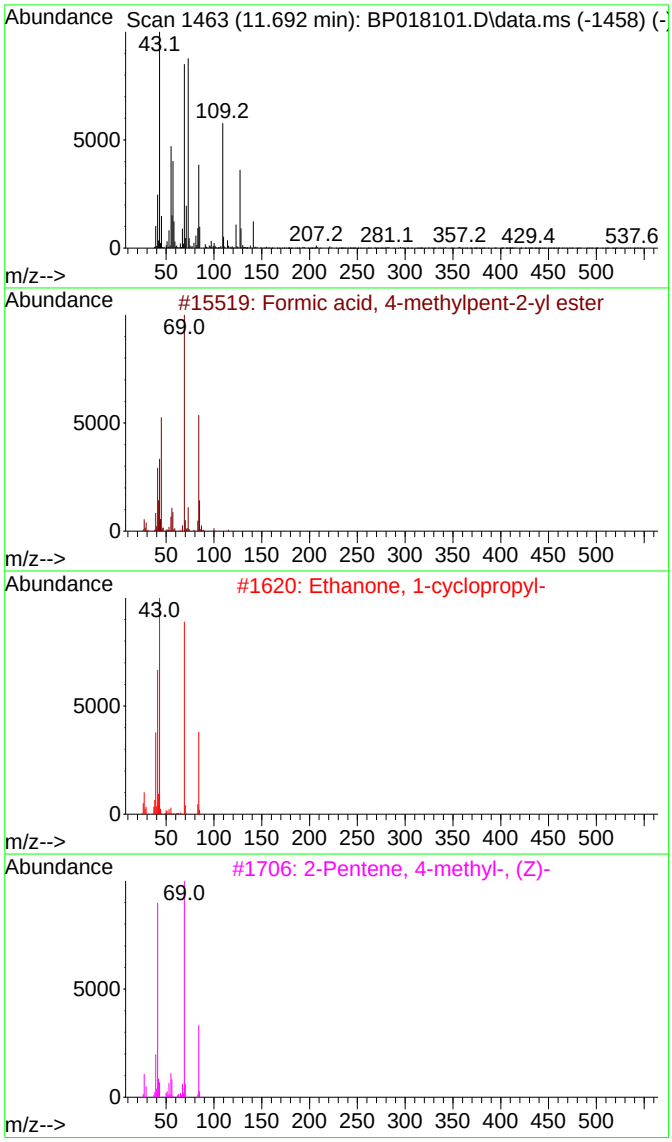
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

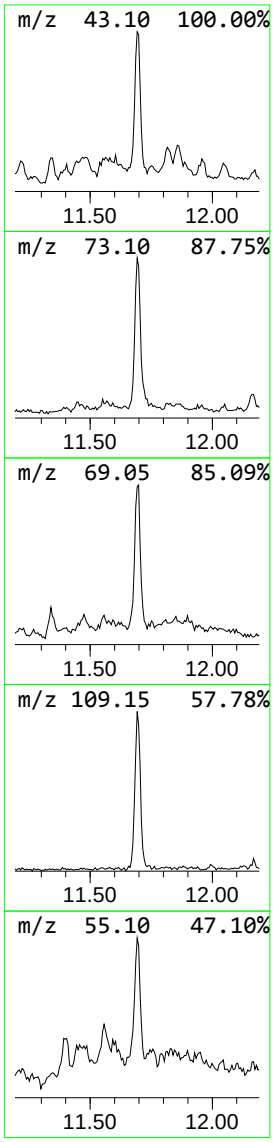
****

Peak Number 22 unknown-12 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.692	8.32 ng/ul	239762	Naphthalene-d8	10.663

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Formic acid, 4-methylpent-2-yl e...	130	C7H14O2	1000367-94-0	43
2	Ethanone, 1-cyclopropyl-	84	C5H8O	000765-43-5	38
3	2-Pentene, 4-methyl-, (Z)-	84	C6H12	000691-38-3	38
4	Ethanone, 1-cyclopropyl-	84	C5H8O	000765-43-5	38
5	3,3-Diethylglutaric acid	188	C9H16O4	004160-95-6	37





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Data File : BP018101.D
Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH347

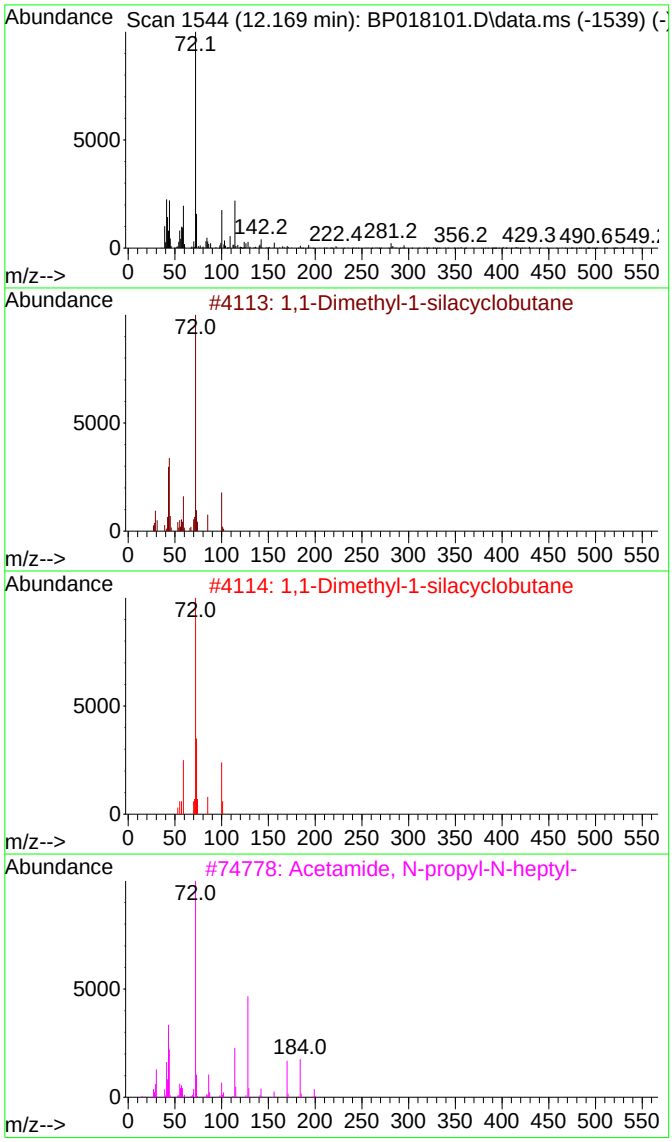
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

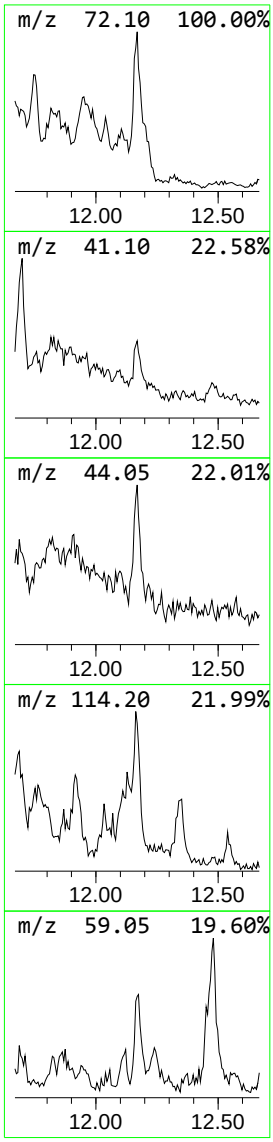
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 1,1-Dimethyl-1-silacyclobutane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.169	3.44 ng/ul	99303	Naphthalene-d8	10.663

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1-Dimethyl-1-silacyclobutane	100	C5H12Si	002295-12-7	59
2	1,1-Dimethyl-1-silacyclobutane	100	C5H12Si	002295-12-7	58
3	Acetamide, N-propyl-N-heptyl-	199	C12H25NO	1000438-05-5	50
4	3,4-Dimethylethcathinone	205	C13H19NO	1000386-29-1	47
5	2-[Ethyl(methyl)amino]ethyl cycl...	319	C19H29NO3	751416-59-8	43





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Data File : BP018101.D
Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH347

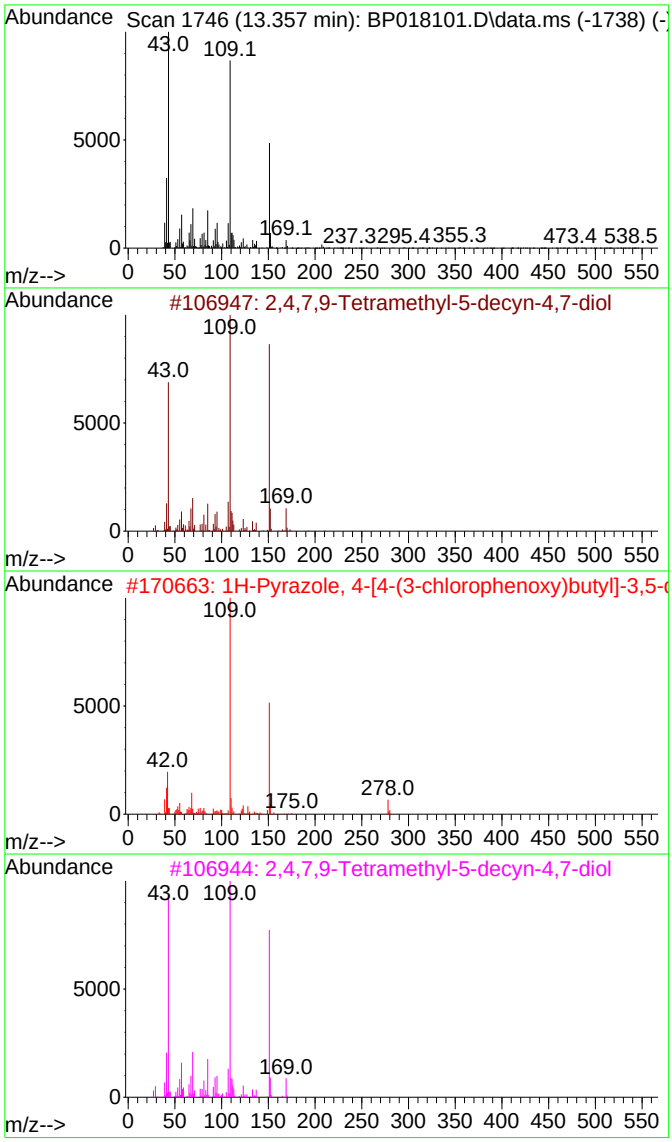
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

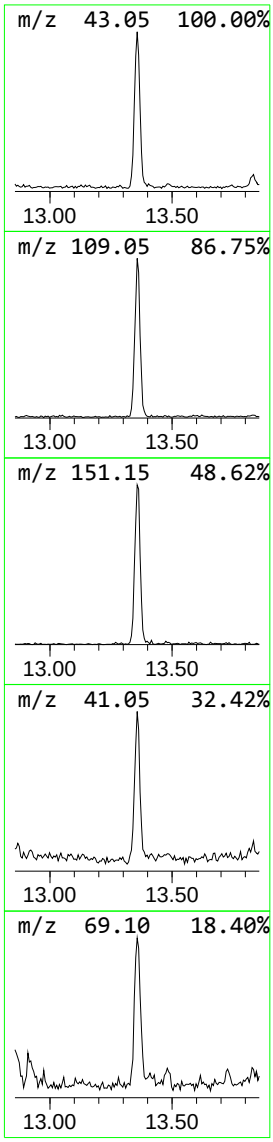
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.357	5.83 ng/ul	189614	Acenaphthene-d10	14.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	90
2	1H-Pyrazole, 4-[4-(3-chloropheno...	278	C15H19ClN2O	1010302-33-9	59
3	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	58
4	N-(4-tert-Butoxyphenyl)acetamide	207	C12H17NO2	002109-73-1	53
5	Butanamide, N-acetyl-N-(4-hydrox...	221	C12H15NO3	1000107-08-7	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018101.D
Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH347

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

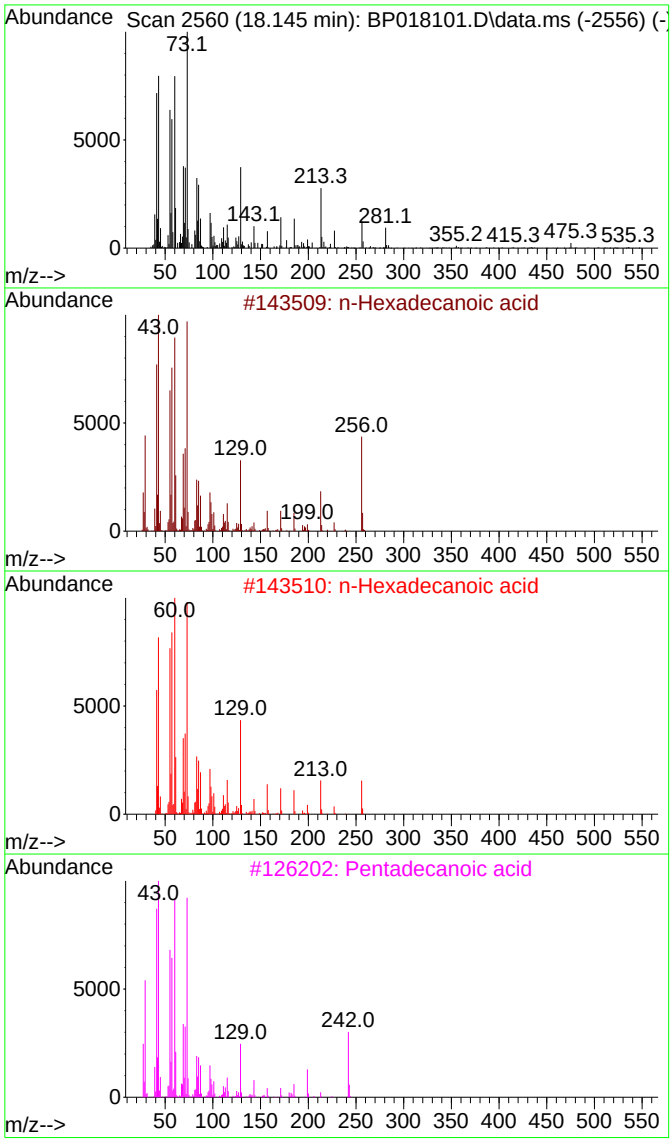
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

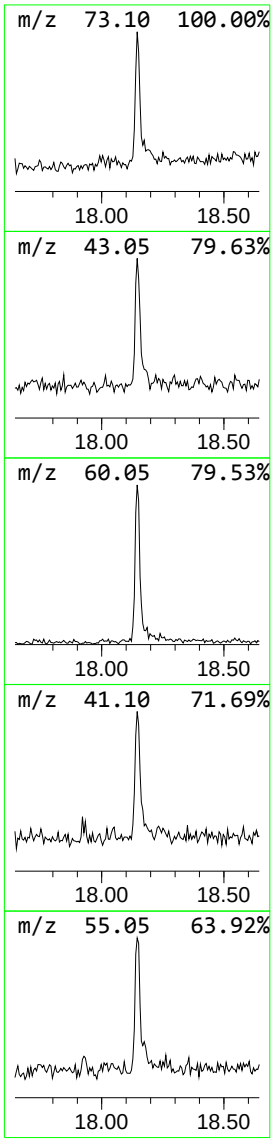
****

Peak Number 25 n-Hexadecanoic acid Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.145	2.30 ng/ul	94130	Phenanthrene-d10	17.298

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	86
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	76
5	8-Methylnonanoic acid	172	C10H20O2	005963-14-4	70





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018101.D
Acq On : 12 Nov 2023 12:41
Operator : MA/JU
Sample : 05158-10
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH347

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
unknown-01	3.346	6.8	ng/ul	129563	1	7.845	381361	20.0
unknown-02	3.834	3.2	ng/ul	60385	1	7.845	381361	20.0
2-Heptanol, 2-m...	5.710	113.7	ng/ul	2168740	1	7.845	381361	20.0
unknown-03	6.034	5.5	ng/ul	105818	1	7.845	381361	20.0
2,5-Hexanediol,...	8.386	4.4	ng/ul	84124	1	7.845	381361	20.0
Norvenlafaxine	8.639	3.6	ng/ul	68300	1	7.845	381361	20.0
3-Octanol, 3,7-...	9.116	7.8	ng/ul	149017	1	7.845	381361	20.0
Diehtylamino-3-...	9.192	3.8	ng/ul	71534	1	7.845	381361	20.0
2-(3-Dimethylam...	9.257	2.3	ng/ul	66995	2	10.663	576566	20.0
2-Butylamine, N...	9.351	4.6	ng/ul	132768	2	10.663	576566	20.0
unknown-04	9.445	7.8	ng/ul	224286	2	10.663	576566	20.0
unknown-05	9.622	4.6	ng/ul	132754	2	10.663	576566	20.0
unknown-06	9.986	3.2	ng/ul	91146	2	10.663	576566	20.0
unknown-07	10.045	5.2	ng/ul	150338	2	10.663	576566	20.0
unknown-08	10.204	8.0	ng/ul	231693	2	10.663	576566	20.0
Propionamide, N...	10.381	4.3	ng/ul	122846	2	10.663	576566	20.0
DL-Leucine, N-a...	10.434	5.2	ng/ul	149875	2	10.663	576566	20.0
unknown-09	10.992	2.1	ng/ul	61195	2	10.663	576566	20.0
unknown-10	11.339	2.0	ng/ul	58058	2	10.663	576566	20.0
N-Acetyl-DL-val...	11.404	2.5	ng/ul	71250	2	10.663	576566	20.0
unknown-11	11.557	3.9	ng/ul	111550	2	10.663	576566	20.0
unknown-12	11.692	8.3	ng/ul	239762	2	10.663	576566	20.0
1,1-Dimethyl-1-...	12.169	3.4	ng/ul	99303	2	10.663	576566	20.0
2,4,7,9-Tetrame...	13.357	5.8	ng/ul	189614	3	14.522	650171	20.0
n-Hexadecanoic ...	18.145	2.3	ng/ul	94130	4	17.298	820275	20.0