

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
 Data File : BP014467.D
 Acq On : 01 Apr 2023 04:55
 Operator : CG/JU
 Sample : 02092-15
 Misc :
 ALS Vial : 77 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB982

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-BP032823.M
 Title : SVOA CALIBRATION

Signal : TIC: BP014467.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.387	31	34	35	rVV	51099	50183	1.20%	0.075%
2	3.423	35	40	56	rVB	1541392	2059280	49.28%	3.071%
3	3.670	77	82	87	rBV2	23025	35391	0.85%	0.053%
4	3.834	107	110	122	rBV	128529	233720	5.59%	0.349%
5	4.046	140	146	149	rBV	18764	34116	0.82%	0.051%
6	4.217	170	175	180	rVV	746245	946325	22.65%	1.411%
7	4.264	180	183	188	rVB2	36033	51465	1.23%	0.077%
8	4.599	234	240	248	rBV2	144211	247471	5.92%	0.369%
9	4.946	296	299	309	rBV4	15365	36580	0.88%	0.055%
10	5.075	314	321	327	rBV	19494	30410	0.73%	0.045%
11	5.770	433	439	447	rBV6	12483	36583	0.88%	0.055%
12	6.275	519	525	533	rVV	34863	67486	1.61%	0.101%
13	6.593	575	579	585	rBV2	16775	31971	0.77%	0.048%
14	6.958	637	641	650	rBV	41336	67135	1.61%	0.100%
15	7.111	662	667	671	rVV	25380	45549	1.09%	0.068%
16	7.187	671	680	694	rVB	230845	465297	11.13%	0.694%
17	7.334	699	705	715	rBV	704994	1131407	27.07%	1.688%
18	7.469	723	728	734	rVB3	30268	54666	1.31%	0.082%
19	7.546	734	741	749	rBV	743984	1236631	29.59%	1.844%
20	8.011	813	820	827	rVV	820595	1290710	30.89%	1.925%
21	8.093	827	834	840	rVV4	37976	91088	2.18%	0.136%
22	8.152	840	844	850	rVB4	23636	39527	0.95%	0.059%
23	8.258	855	862	866	rBV8	22758	47372	1.13%	0.071%
24	8.505	898	904	908	rVV	26423	42540	1.02%	0.063%
25	8.734	937	943	952	rVV	635537	1117731	26.75%	1.667%
26	8.805	952	955	968	rVB8	43229	99807	2.39%	0.149%
27	8.958	970	981	985	rBV	99828	214125	5.12%	0.319%
28	9.011	985	990	996	rVV	772174	1264696	30.26%	1.886%
29	9.081	996	1002	1006	rVV3	195291	480359	11.50%	0.716%
30	9.122	1006	1009	1014	rVV	185438	289274	6.92%	0.431%
31	9.187	1014	1020	1028	rVV	918154	1564178	37.43%	2.333%
32	9.269	1029	1034	1039	rVB	97474	166768	3.99%	0.249%
33	9.381	1049	1053	1061	rVB4	92209	164529	3.94%	0.245%
34	9.452	1061	1065	1071	rVB3	21360	29613	0.71%	0.044%
35	9.540	1071	1080	1087	rBV7	65217	172929	4.14%	0.258%
36	9.693	1093	1106	1109	rBV	515013	1209349	28.94%	1.804%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
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37	9.734	1109	1113	1126	rVV2	169885	376220	9.00%	0.561%
38	9.834	1126	1130	1137	rVV9	26898	61352	1.47%	0.092%
39	9.916	1137	1144	1155	rVV	778838	1373895	32.88%	2.049%
40	10.046	1157	1166	1170	rVV4	88477	196762	4.71%	0.293%
41	10.099	1170	1175	1186	rVV5	69000	166157	3.98%	0.248%
42	10.193	1186	1191	1195	rVV4	26826	48451	1.16%	0.072%
43	10.252	1195	1201	1209	rVB10	32185	81372	1.95%	0.121%
44	10.375	1214	1222	1230	rBV5	145970	315174	7.54%	0.470%
45	10.463	1230	1237	1246	rBV	1063702	1965810	47.04%	2.932%
46	10.616	1258	1263	1271	rVV2	117969	223998	5.36%	0.334%
47	10.687	1272	1275	1279	rVB4	20944	27566	0.66%	0.041%
48	10.775	1285	1290	1294	rBV4	82142	156756	3.75%	0.234%
49	10.834	1294	1300	1312	rVB	1221994	2017486	48.28%	3.009%
50	10.981	1317	1325	1335	rBV	574169	1218557	29.16%	1.818%
51	11.099	1340	1345	1356	rBV9	59735	189399	4.53%	0.282%
52	11.275	1369	1375	1376	rBV4	55956	87816	2.10%	0.131%
53	11.310	1377	1381	1386	rVB3	91285	162226	3.88%	0.242%
54	11.416	1394	1399	1402	rBV5	47228	97310	2.33%	0.145%
55	11.469	1405	1408	1412	rVV4	77784	116689	2.79%	0.174%
56	11.522	1412	1417	1423	rVB7	49563	115097	2.75%	0.172%
57	11.675	1435	1443	1444	rBV6	70274	155556	3.72%	0.232%
58	11.822	1464	1468	1475	rVB7	82607	153474	3.67%	0.229%
59	11.887	1475	1479	1482	rBV3	64638	95827	2.29%	0.143%
60	11.963	1489	1492	1496	rVB5	26328	34097	0.82%	0.051%
61	12.040	1502	1505	1515	rVB6	55892	137566	3.29%	0.205%
62	12.181	1526	1529	1535	rVB3	57764	100311	2.40%	0.150%
63	12.287	1543	1547	1551	rBV4	40952	60560	1.45%	0.090%
64	12.346	1551	1557	1562	rBV	184038	373049	8.93%	0.556%
65	12.722	1617	1621	1630	rBV	59975	128163	3.07%	0.191%
66	12.828	1634	1639	1653	rVB3	188424	449163	10.75%	0.670%
67	12.928	1653	1656	1662	rBV5	38549	68875	1.65%	0.103%
68	13.022	1670	1672	1674	rBV3	37725	42507	1.02%	0.063%
69	13.057	1674	1678	1683	rVB6	104878	160271	3.84%	0.239%
70	13.693	1782	1786	1788	rBV3	63266	99199	2.37%	0.148%
71	13.728	1788	1792	1798	rVV4	125814	295257	7.07%	0.440%
72	13.799	1800	1804	1811	rVB3	109321	191794	4.59%	0.286%
73	13.910	1819	1823	1827	rBV5	65840	114652	2.74%	0.171%
74	13.981	1833	1835	1841	rVB6	41921	54097	1.29%	0.081%
75	14.075	1845	1851	1859	rVB	2238297	3169095	75.84%	4.727%
76	14.175	1861	1868	1872	rBV5	88717	221790	5.31%	0.331%

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77	14.251	1878	1881	1885	rVB3	85960	102871	2.46%	0.153%
78	14.363	1894	1900	1907	rBV	2273600	3336871	79.85%	4.977%
79	14.504	1918	1924	1933	rVB	713382	1311413	31.38%	1.956%
80	14.593	1934	1939	1944	rVV	318304	479364	11.47%	0.715%
81	14.663	1946	1951	1961	rVB	1749932	2694364	64.48%	4.019%
82	14.922	1992	1995	2003	rVB3	128302	268144	6.42%	0.400%
83	15.663	2115	2121	2127	rVB	2823035	4158746	99.52%	6.203%
84	15.793	2137	2143	2148	rBV	965600	1461976	34.99%	2.181%
85	16.187	2206	2210	2215	rBV	621677	821785	19.67%	1.226%
86	16.375	2238	2242	2250	rVB5	197285	379614	9.08%	0.566%
87	16.698	2294	2297	2302	rBV	341355	492998	11.80%	0.735%
88	16.828	2316	2319	2323	rBV2	76773	112530	2.69%	0.168%
89	16.940	2335	2338	2340	rBV	108113	140116	3.35%	0.209%
90	17.428	2416	2421	2428	rVB	2020017	2825758	67.62%	4.215%
91	17.528	2433	2438	2448	rVB	2933411	4178802	100.00%	6.233%
92	18.298	2565	2569	2577	rBV2	603571	814209	19.48%	1.214%
93	19.063	2696	2699	2703	rVB	164255	182958	4.38%	0.273%
94	19.522	2774	2777	2785	rVB	243080	363856	8.71%	0.543%
95	19.757	2812	2817	2823	rVB	3155050	4015528	96.09%	5.989%
96	21.116	3046	3048	3057	rBV3	218707	329256	7.88%	0.491%
97	21.192	3058	3061	3066	rVB3	278979	364087	8.71%	0.543%
98	21.516	3112	3116	3122	rVB	1667341	2152632	51.51%	3.211%
99	23.874	3510	3517	3527	rVB	1798371	3561936	85.24%	5.313%
100	24.039	3539	3545	3557	rVB2	1079308	2245789	53.74%	3.350%

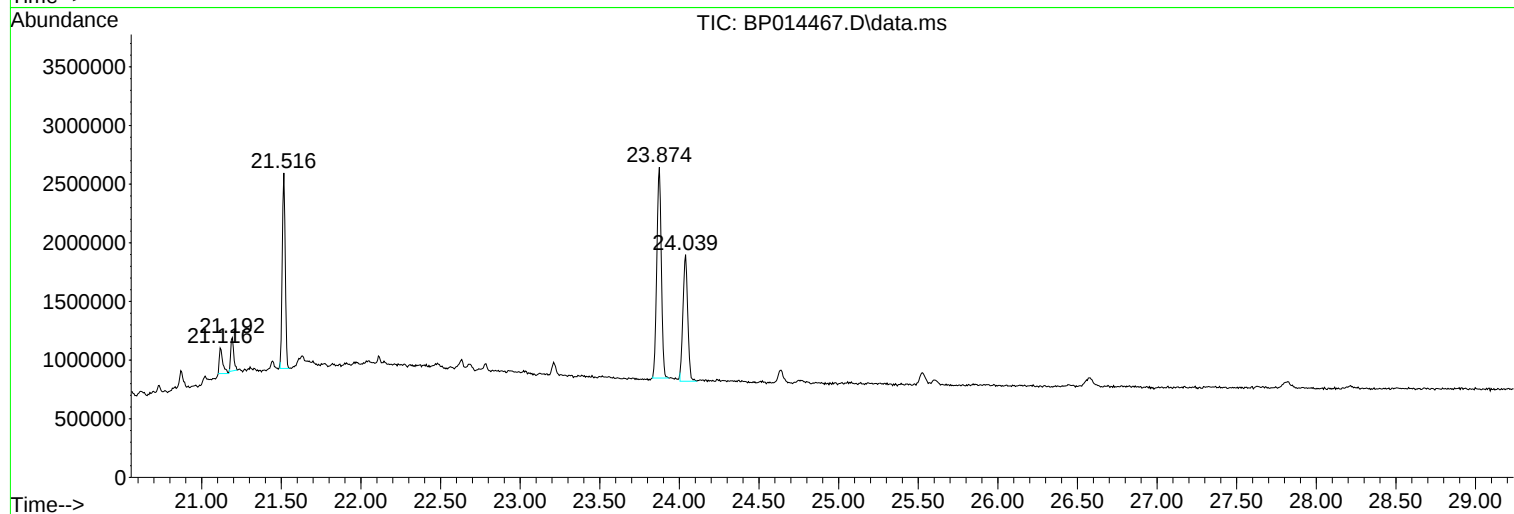
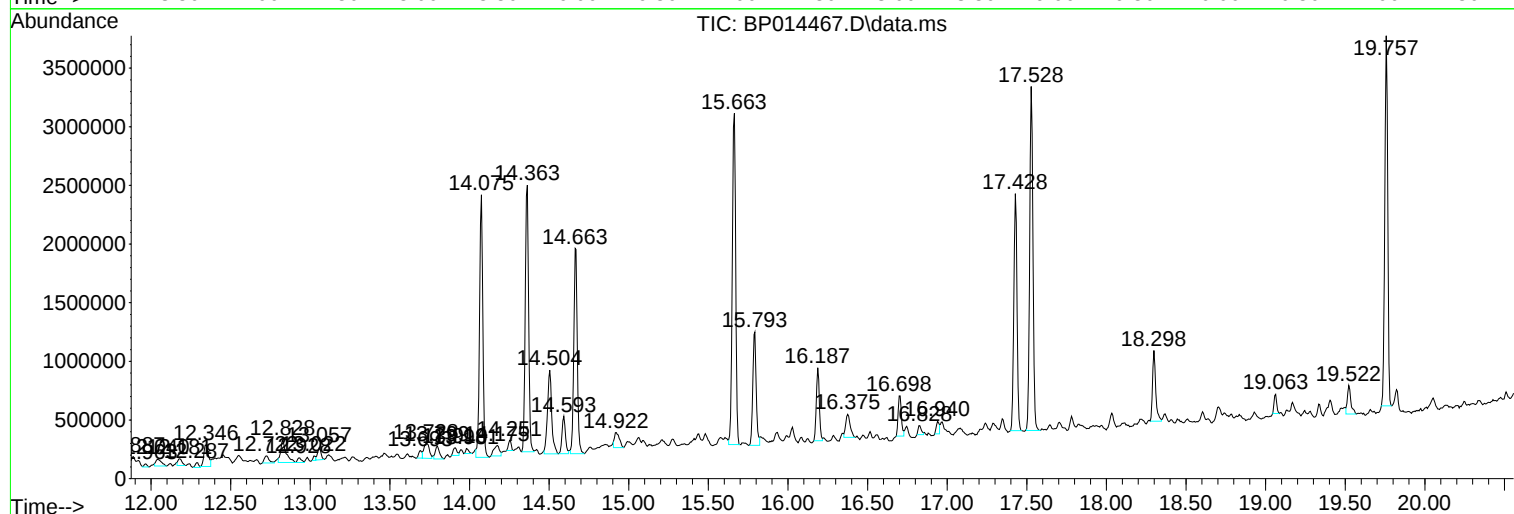
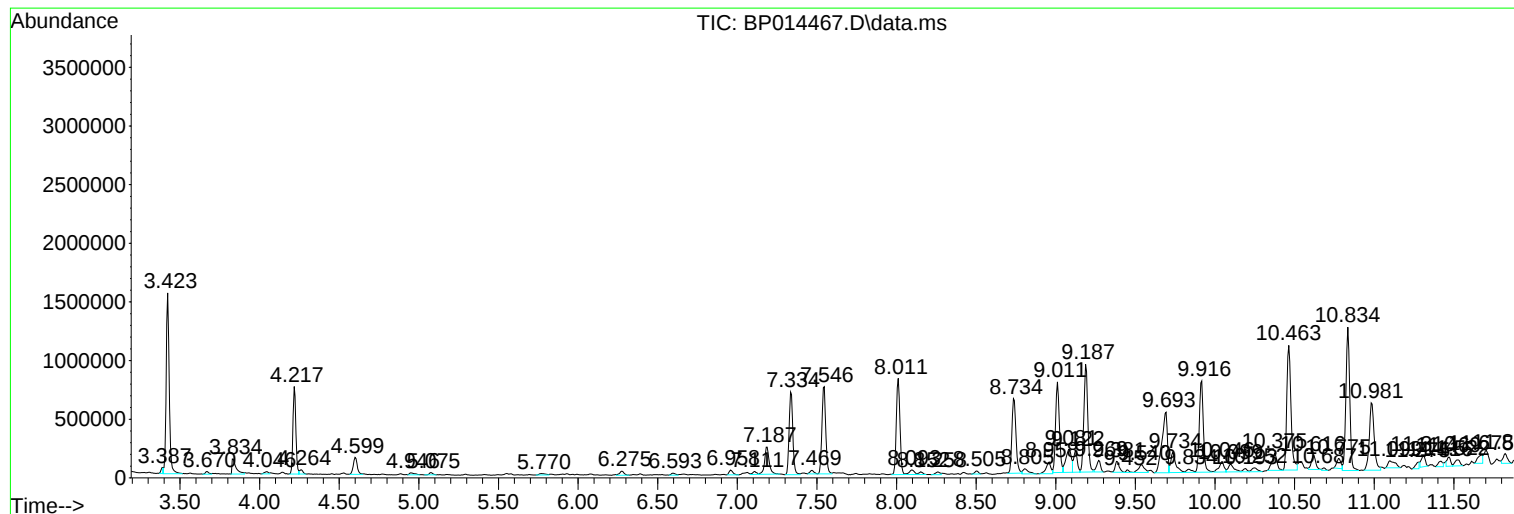
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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



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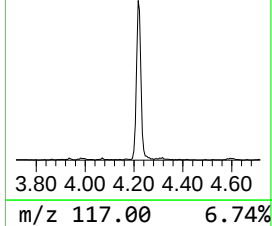
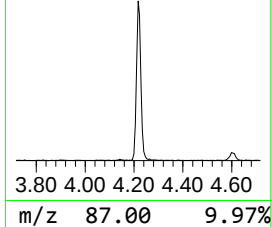
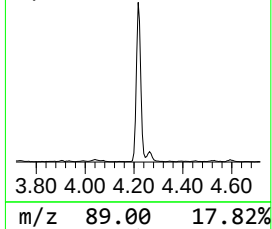
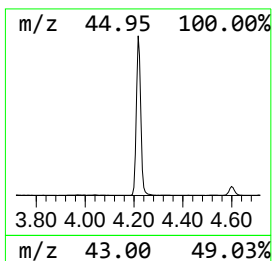
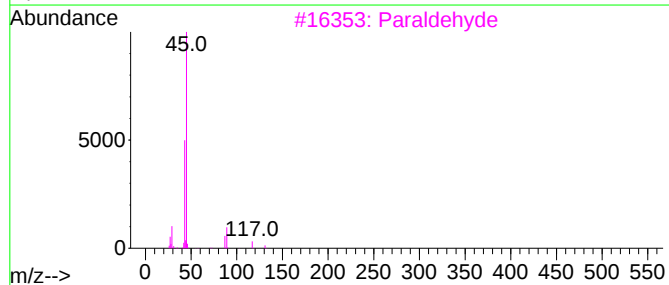
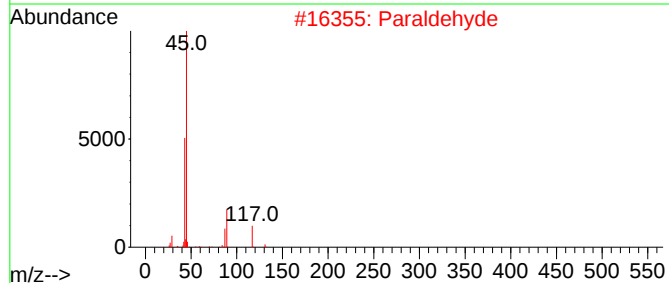
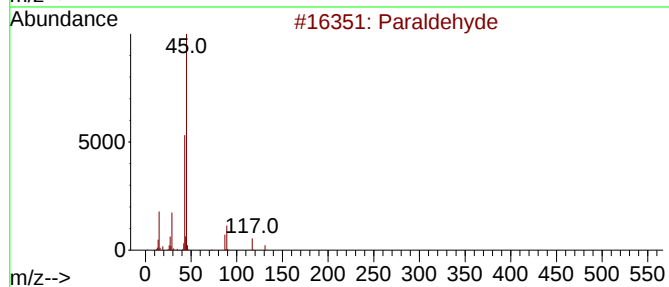
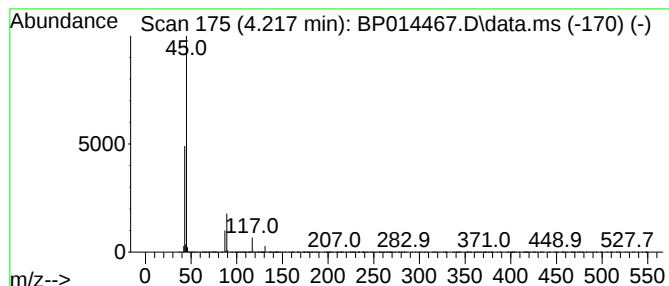
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Paraldehyde Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.217	14.66 ng/ul	946325	1,4-Dichlorobenzene-d4	8.011

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Paraldehyde			132	C6H12O3	000123-63-7	90
2	Paraldehyde			132	C6H12O3	000123-63-7	90
3	Paraldehyde			132	C6H12O3	000123-63-7	83
4	Paraldehyde			132	C6H12O3	000123-63-7	56
5	Propane, 2,2'-[ethylidenebis(oxy...			146	C8H18O2	004285-59-0	56



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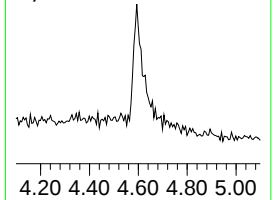
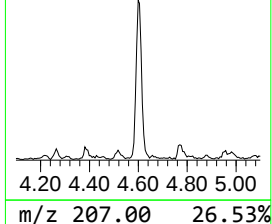
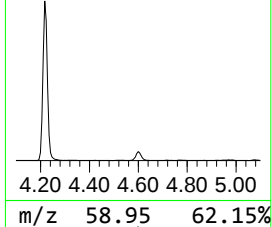
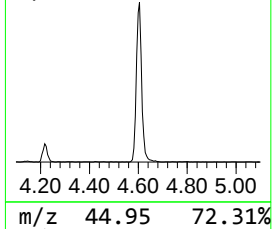
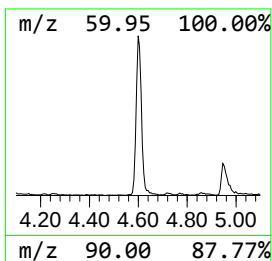
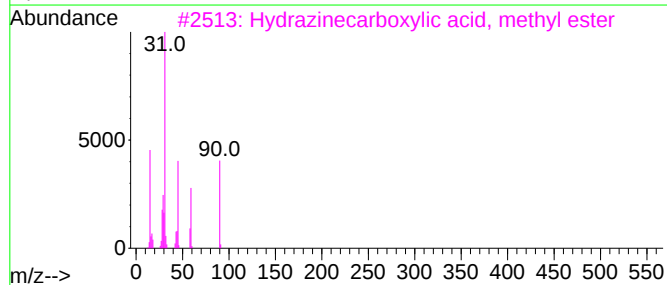
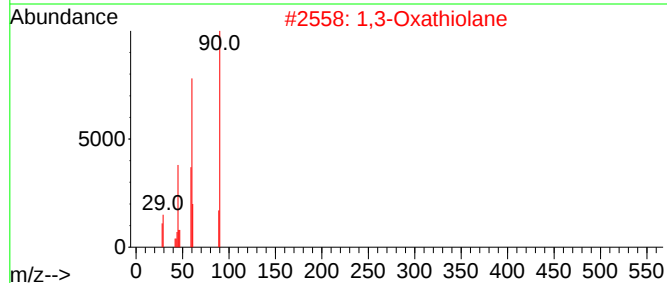
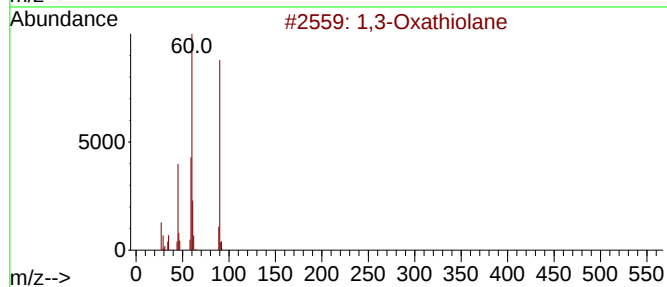
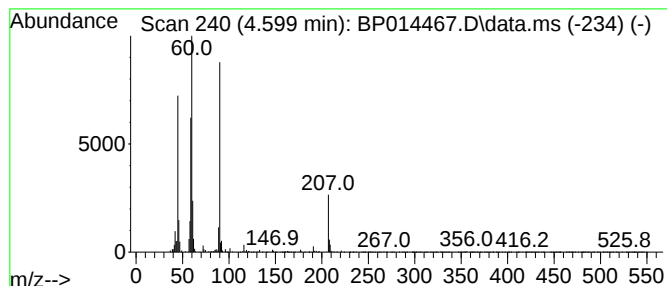
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1,3-Oxathiolane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.599	3.83 ng/ul	247471	1,4-Dichlorobenzene-d4	8.011

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Oxathiolane	90	C3H6OS	002094-97-5	90
2			1,3-Oxathiolane	90	C3H6OS	002094-97-5	59
3			Hydrazinecarboxylic acid, methyl...	90	C2H6N2O2	006294-89-9	53
4			Acrolein, 2-chloro-	90	C3H3ClO	000683-51-2	42
5			Thiourea, methyl-	90	C2H6N2S	000598-52-7	40



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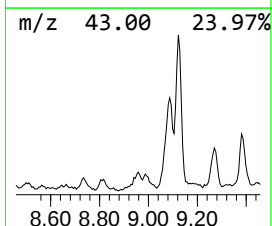
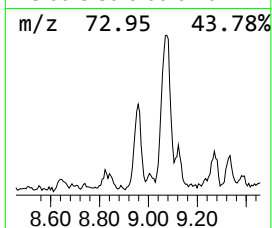
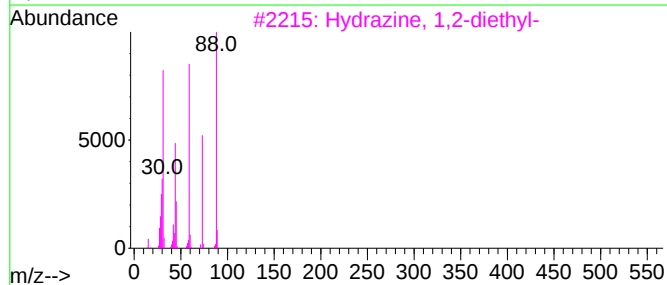
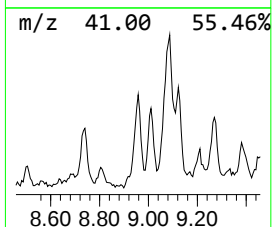
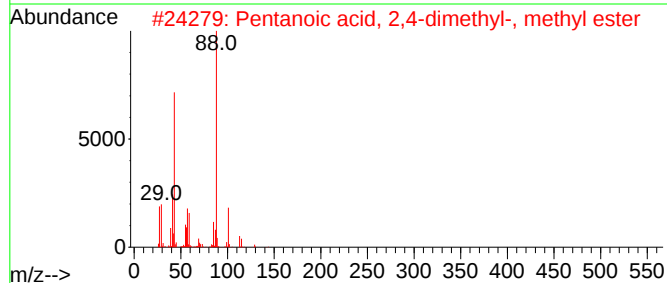
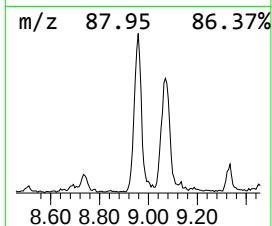
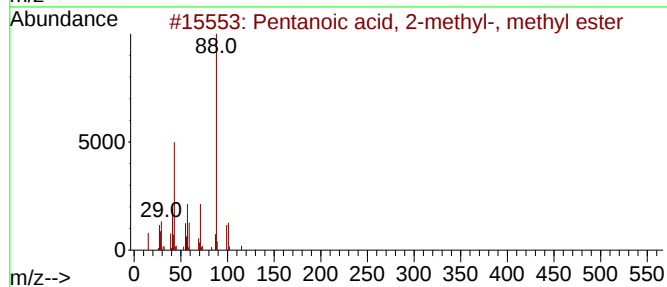
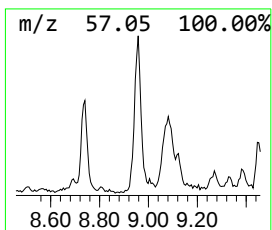
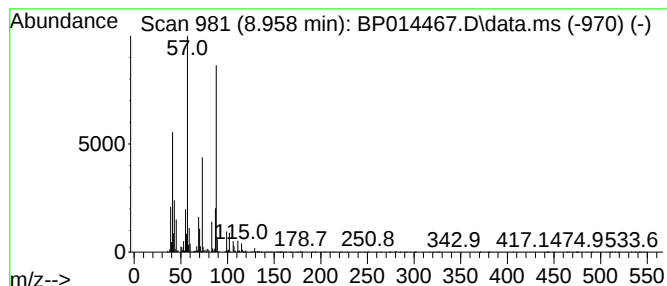
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Pentanoic acid, 2-methyl-, ... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.958	3.32 ng/ul	214125	1,4-Dichlorobenzene-d4	8.011

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentanoic acid, 2-methyl-, methy...	130	C7H14O2	002177-77-7	53
2			Pentanoic acid, 2,4-dimethyl-, m...	144	C8H16O2	071672-33-8	50
3			Hydrazine, 1,2-diethyl-	88	C4H12N2	001615-80-1	46
4			3-Aza-2-oxabicyclo[2.2.2]oct-5-e...	267	C10H12F3NO4	102698-35-1	43
5			Hydrazine, 1,2-diethyl-	88	C4H12N2	001615-80-1	43



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Acq On : 01 Apr 2023 04:55
Operator : CG/JU
Sample : 02092-15
Misc :
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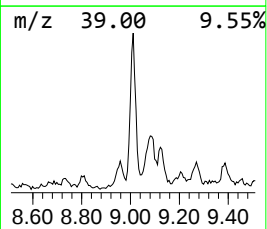
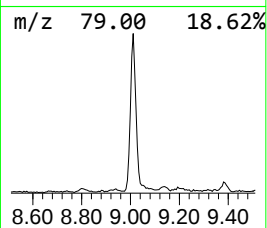
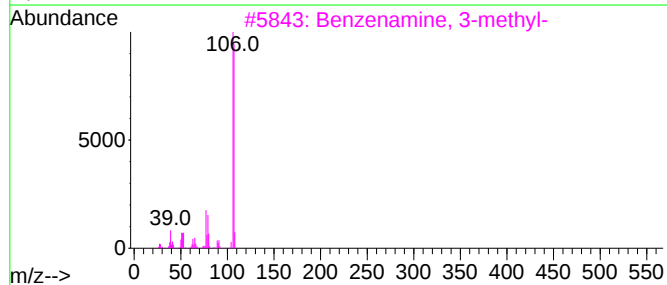
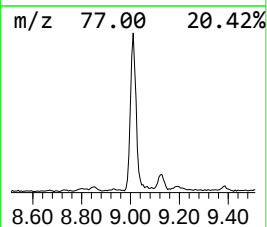
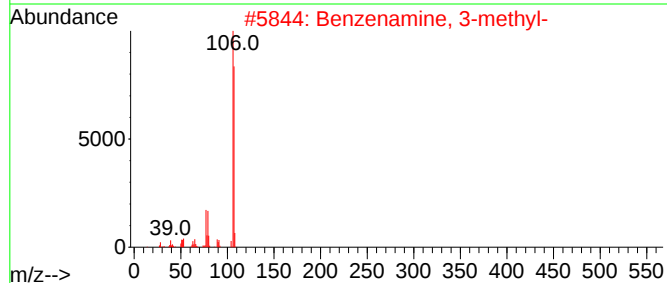
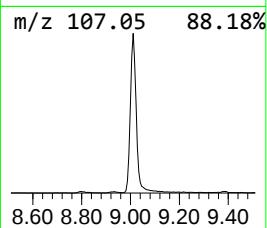
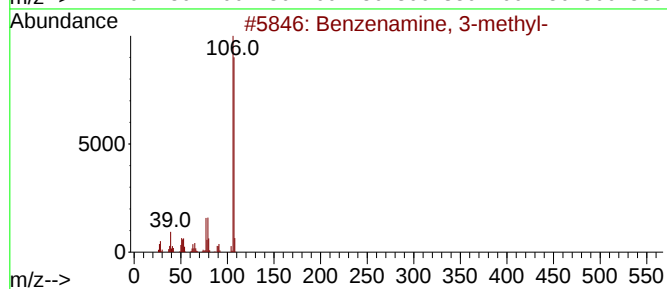
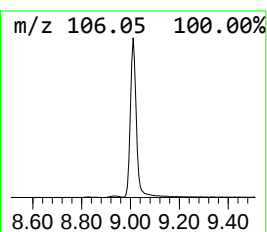
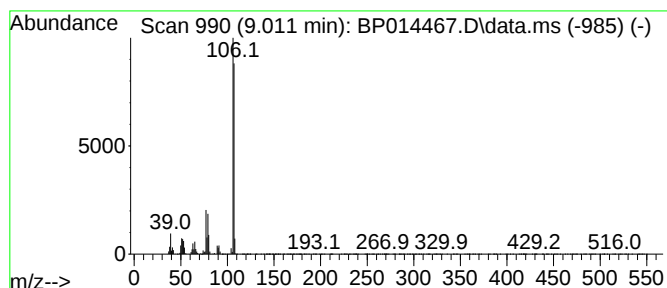
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzenamine, 3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.011	19.60 ng/ul	1264700	1,4-Dichlorobenzene-d4	8.011

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
2			Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
3			Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
4			Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	96
5			p-Aminotoluene	107	C7H9N	000106-49-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014467.D
Acq On : 01 Apr 2023 04:55
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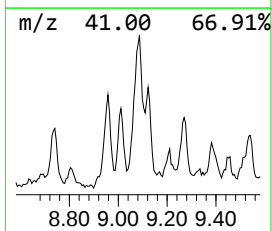
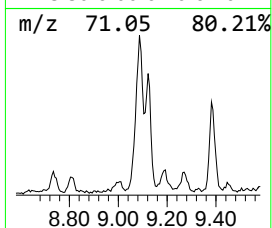
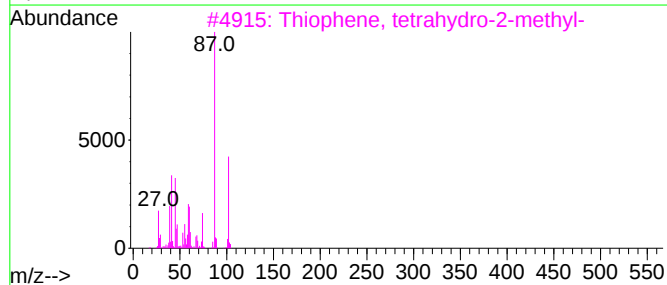
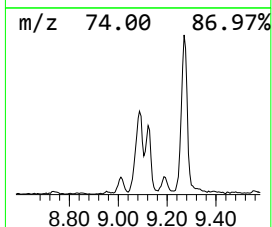
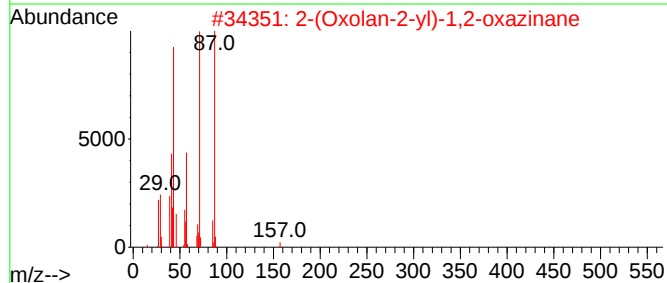
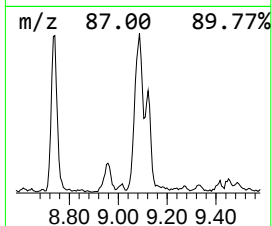
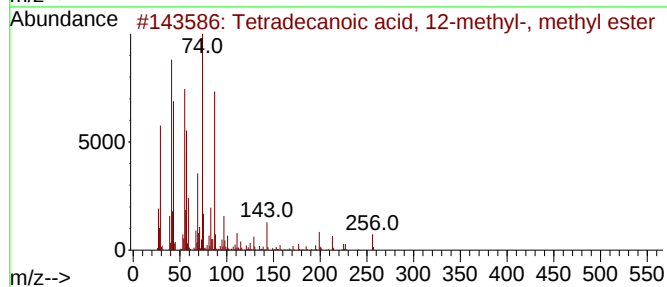
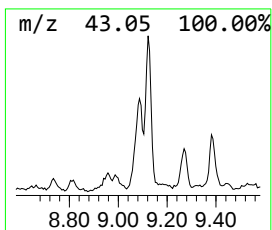
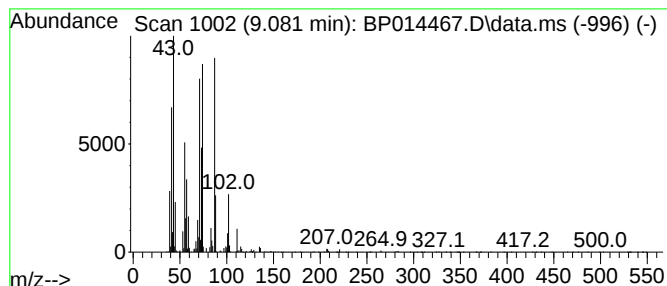
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.081	7.44 ng/ul	480359	1,4-Dichlorobenzene-d4	8.011

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecanoic acid, 12-methyl-, ...	256	C16H32O2	005129-66-8	37
2			2-(Oxolan-2-yl)-1,2-oxazinane	157	C8H15N02	1000445-08-1	32
3			Thiophene, tetrahydro-2-methyl-	102	C5H10S	001795-09-1	27
4			1-Hexene, 4,5-dimethyl-	112	C8H16	016106-59-5	15
5			Ethanedioic acid, bis(3-methylbu...	230	C12H22O4	002051-00-5	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014467.D
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Sample : 02092-15
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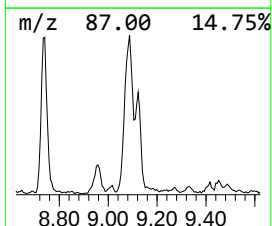
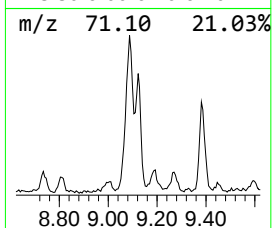
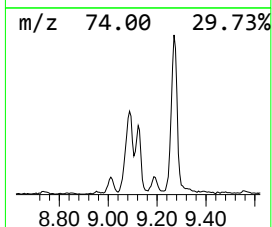
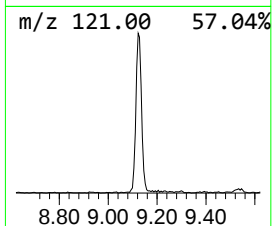
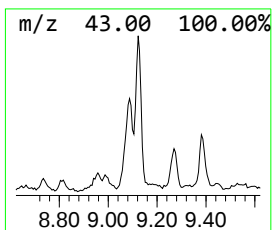
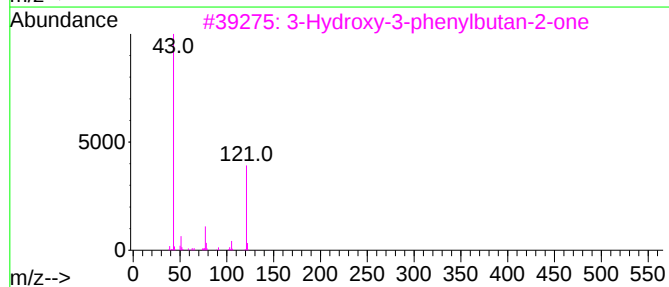
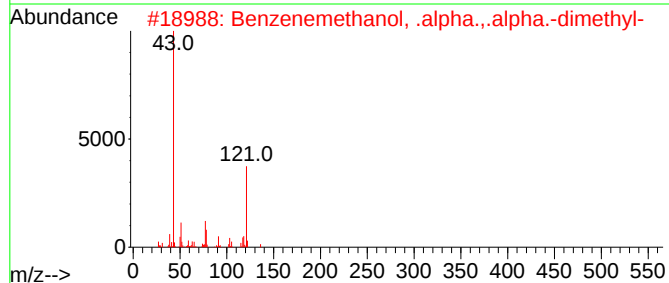
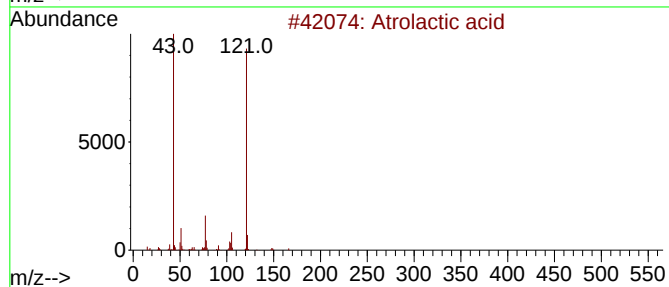
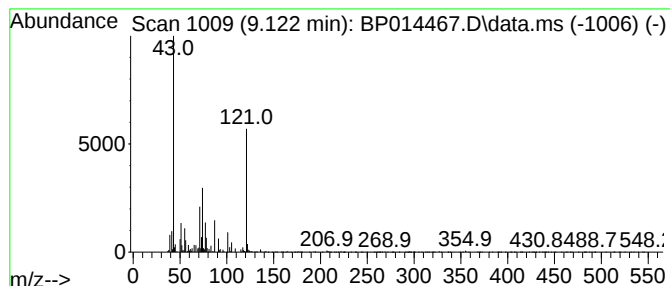
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-02 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.122	4.48 ng/ul	289274	1,4-Dichlorobenzene-d4	8.011

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Atrolactic acid	166	C9H10O3	000515-30-0	38
2			Benzenemethanol, .alpha.,.alpha....	136	C9H12O	000617-94-7	30
3			3-Hydroxy-3-phenylbutan-2-one	164	C10H12O2	003155-01-9	27
4			d1-2-Phenyl-1,2-propanediol	152	C9H12O2	004217-66-7	27
5			Benzenemethanol, .alpha.,.alpha....	136	C9H12O	000617-94-7	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014467.D
Acq On : 01 Apr 2023 04:55
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Sample : 02092-15
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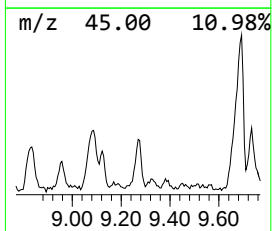
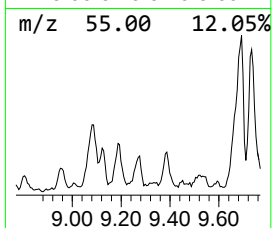
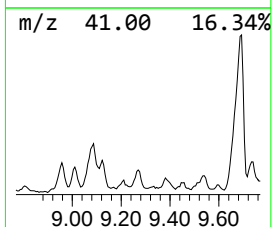
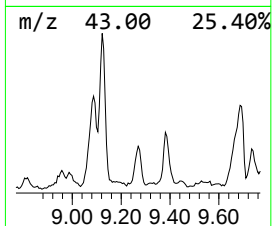
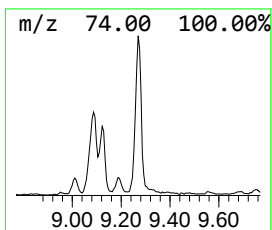
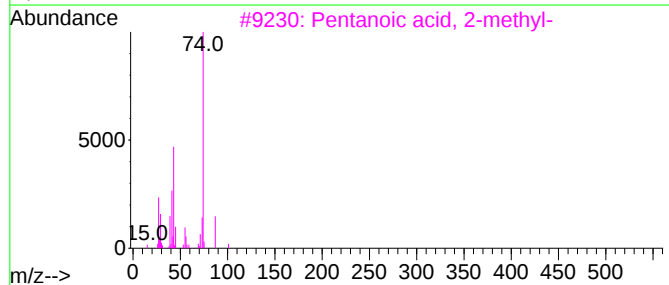
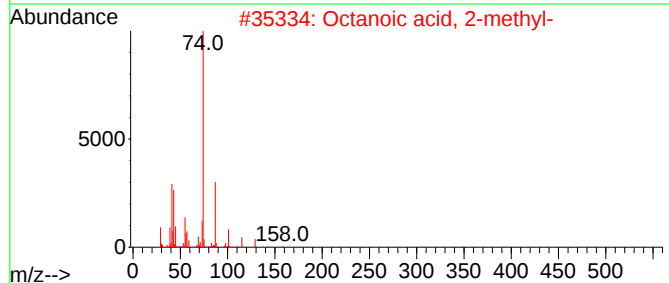
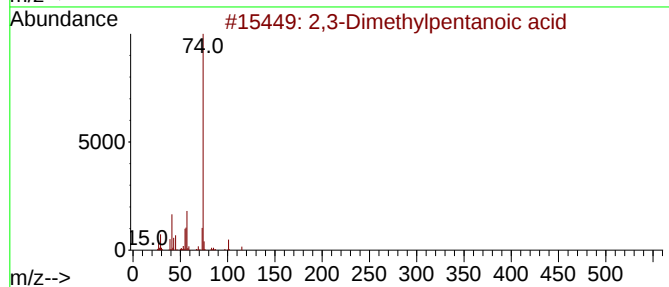
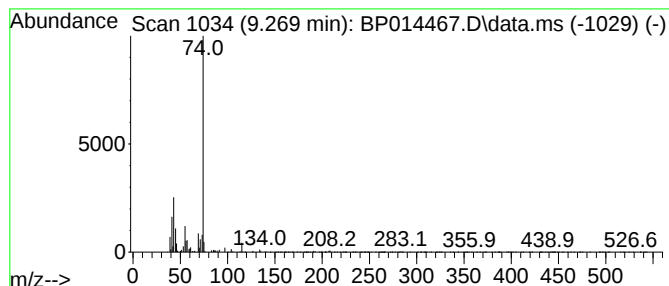
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-03 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.269	2.58 ng/ul	166768	1,4-Dichlorobenzene-d4	8.011

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,3-Dimethylpentanoic acid	130	C7H14O2	082608-03-5	45
2			Octanoic acid, 2-methyl-	158	C9H18O2	003004-93-1	39
3			Pentanoic acid, 2-methyl-	116	C6H12O2	000097-61-0	38
4			Butanoic acid, 2-methyl-	102	C5H10O2	000116-53-0	38
5			1,3,2-Oxathioborolane, 2-ethyl-5...	130	C5H11BOS	1000163-05-5	9



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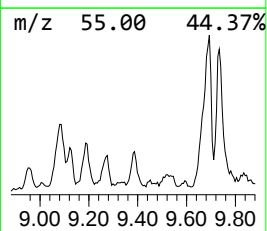
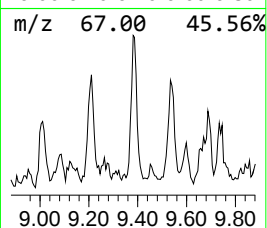
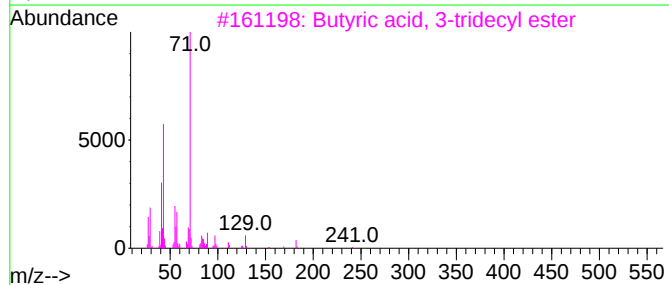
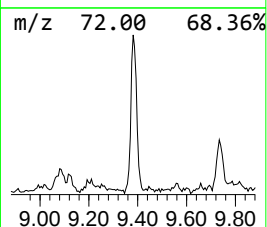
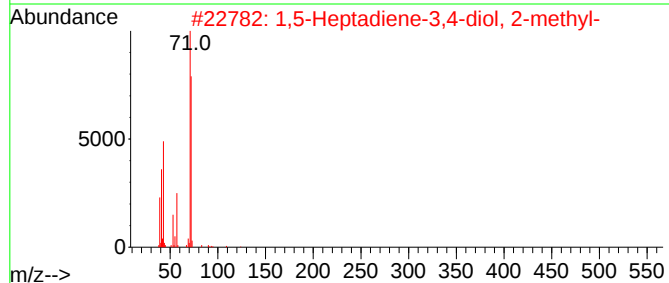
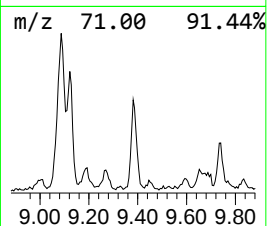
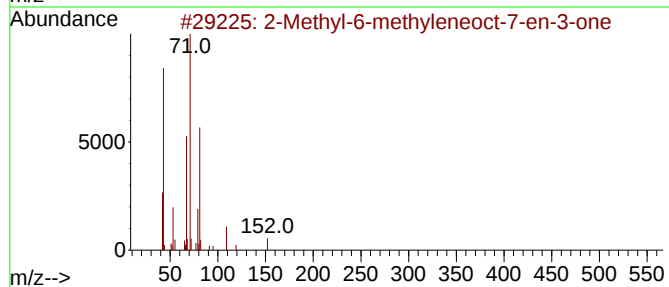
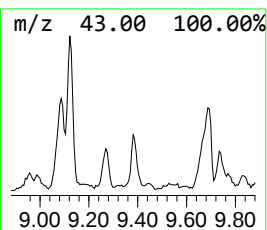
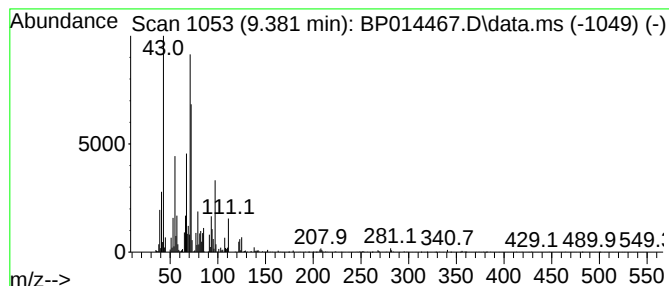
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-04 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.381	2.55 ng/ul	164529	1,4-Dichlorobenzene-d4	8.011

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Methyl-6-methyleneoct-7-en-3-one	152	C10H16O	1000424-30-4	43
2			1,5-Heptadiene-3,4-diol, 2-methyl-	142	C8H14O2	022726-05-2	38
3			Butyric acid, 3-tridecyl ester	270	C17H34O2	1000280-56-8	35
4			5-Nonanol, 2-methylpropionate	214	C13H26O2	1000463-47-6	35
5			Isophytol	296	C20H40O	000505-32-8	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
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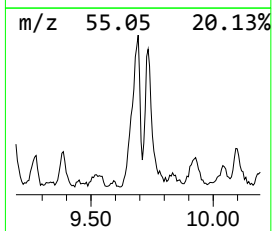
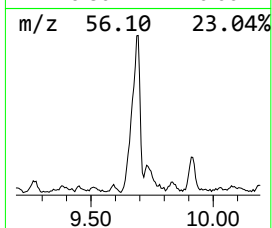
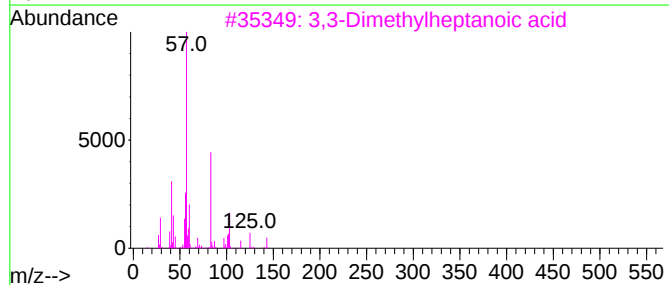
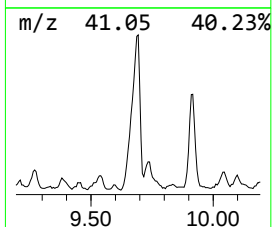
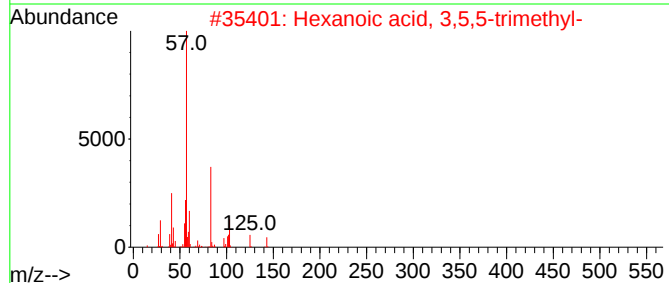
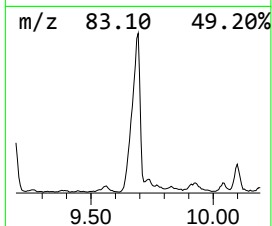
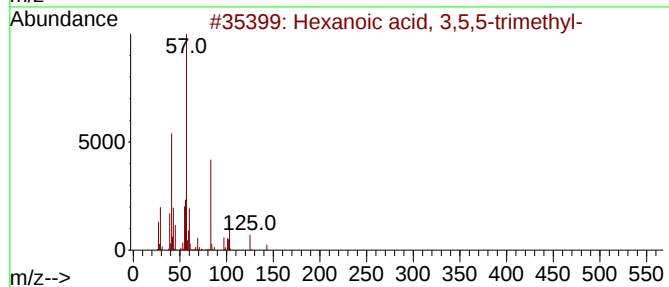
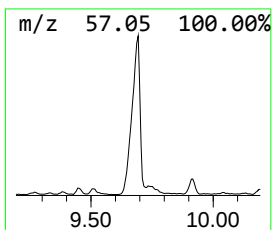
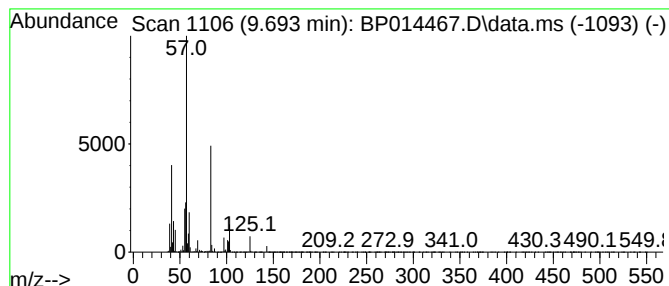
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Hexanoic acid, 3,5,5-trimet... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.693	11.99 ng/ul	1209350	Naphthalene-d8	10.834

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 3,5,5-trimethyl-	158	C9H18O2	003302-10-1	90
2			Hexanoic acid, 3,5,5-trimethyl-	158	C9H18O2	003302-10-1	83
3			3,3-Dimethylheptanoic acid	158	C9H18O2	067061-30-7	83
4			Hexanoic acid, 3,5,5-trimethyl-	158	C9H18O2	003302-10-1	64
5			Hexanoic acid, 3,5,5-trimethyl-	158	C9H18O2	003302-10-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
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ALS Vial : 77 Sample Multiplier: 1

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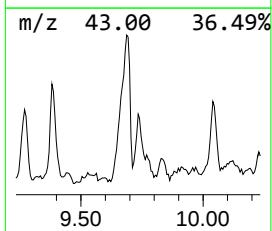
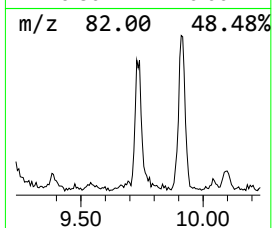
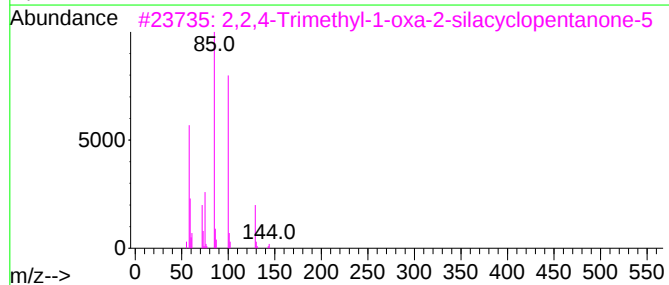
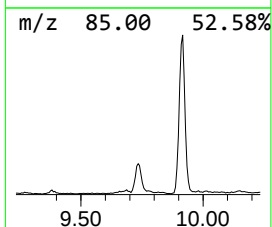
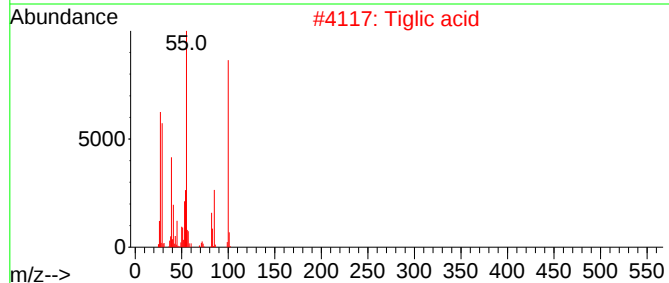
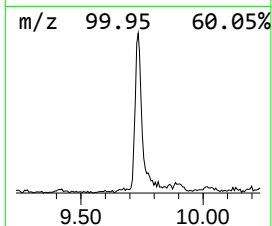
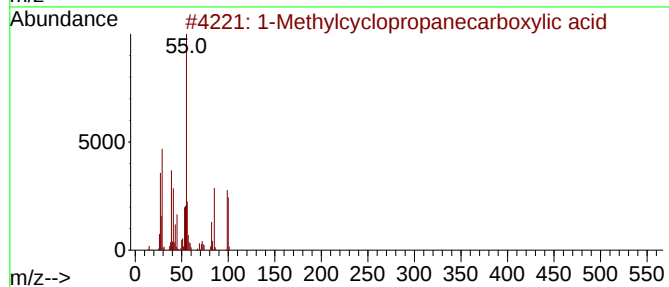
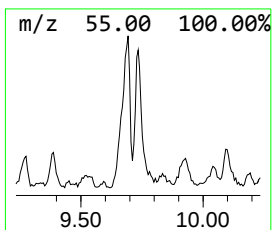
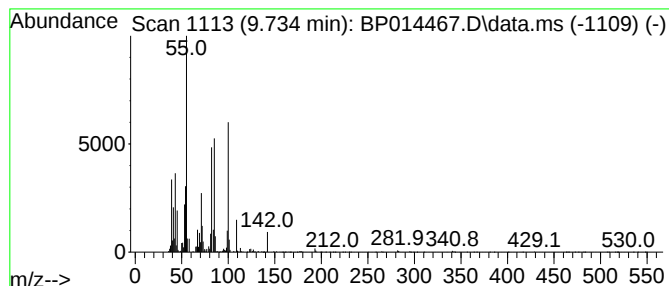
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-05 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.734	3.73 ng/ul	376220	Naphthalene-d8	10.834

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Methylcyclopropanecarboxylic acid	100	C5H8O2	006914-76-7	38
2			Tiglic acid	100	C5H8O2	000080-59-1	35
3			2,2,4-Trimethyl-1-oxa-2-silacycl...	144	C6H12O2Si	020471-80-1	27
4			1-[-]-4-Hydroxy-1-methylproline	145	C6H11NO3	1000251-00-7	27
5			2-Piperazinone	100	C4H8N2O	005625-67-2	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014467.D
Acq On : 01 Apr 2023 04:55
Operator : CG/JU
Sample : 02092-15
Misc :
ALS Vial : 77 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB982

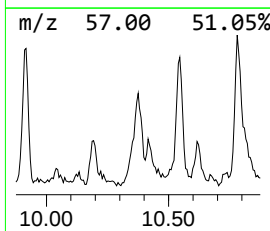
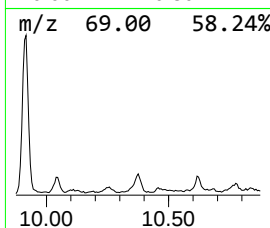
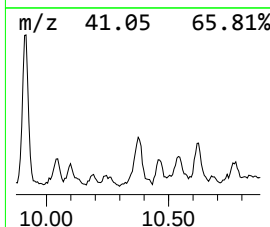
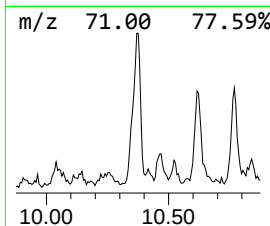
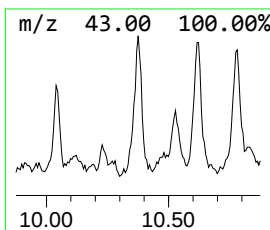
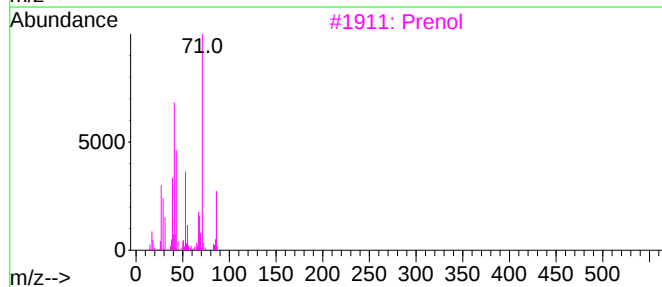
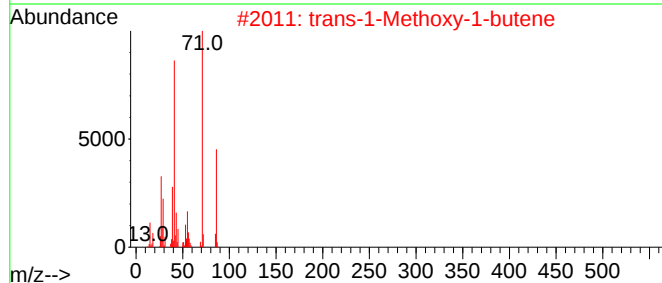
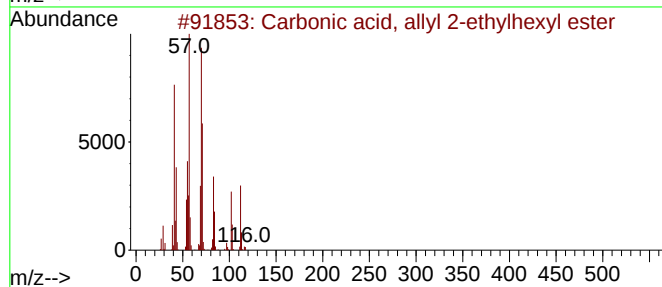
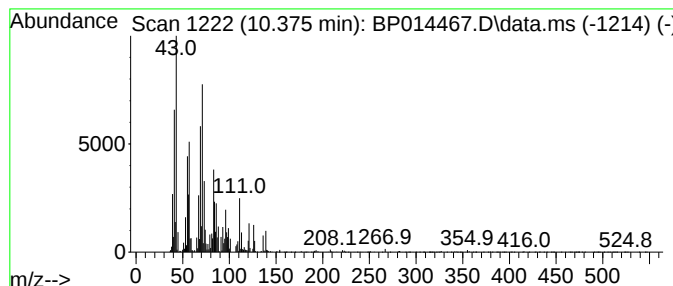
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-06 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.375	3.12 ng/ul	315174	Naphthalene-d8	10.834

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, allyl 2-ethylhexy...	214	C12H22O3	1000357-80-6	43
2			trans-1-Methoxy-1-butene	86	C5H10O	010034-13-6	30
3			Prenol	86	C5H10O	000556-82-1	27
4			Sulfurous acid, butyl nonyl ester	264	C13H28O3S	1000309-17-6	27
5			2-Nonanol, trimethylacetate	228	C14H28O2	1000463-21-6	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014467.D
Acq On : 01 Apr 2023 04:55
Operator : CG/JU
Sample : 02092-15
Misc :
ALS Vial : 77 Sample Multiplier: 1

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ClientSampleId :
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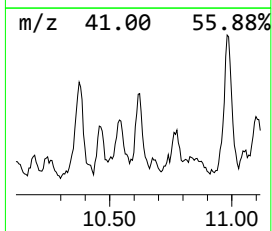
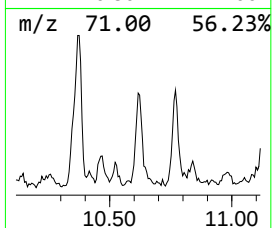
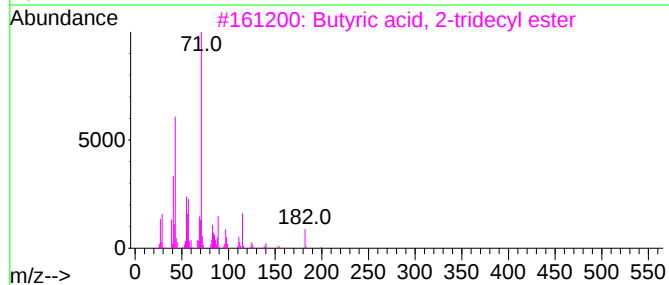
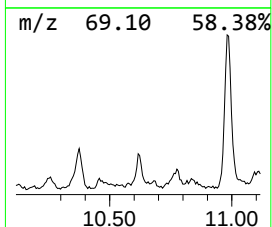
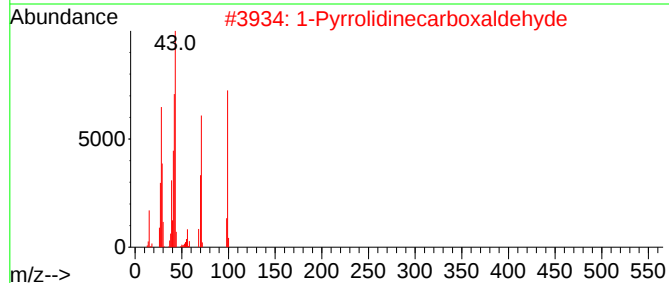
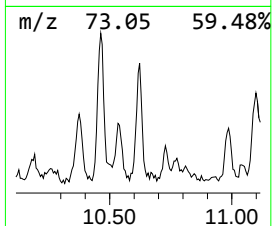
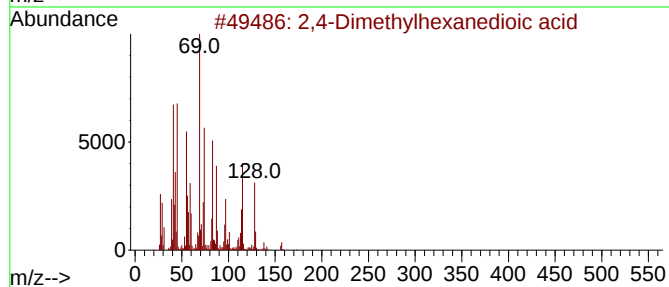
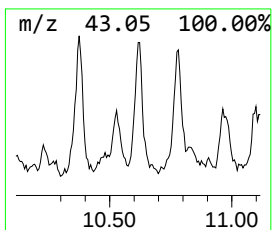
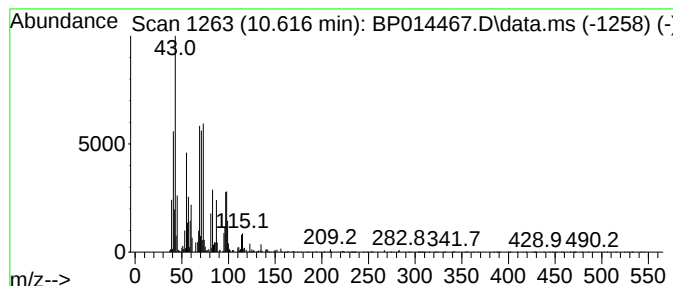
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 unknown-07 Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.616	2.22 ng/ul	223998	Naphthalene-d8	10.834

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,4-Dimethylhexanedioic acid	174	C8H14O4	040933-45-7	22
2			1-Pyrrolidinecarboxaldehyde	99	C5H9NO	003760-54-1	14
3			Butyric acid, 2-tridecyl ester	270	C17H34O2	055193-07-2	14
4			Methyl-4,6-di-O-methylmannoside	222	C9H18O6	1000426-87-7	14
5			Phytol	296	C20H40O	000150-86-7	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014467.D
Acq On : 01 Apr 2023 04:55
Operator : CG/JU
Sample : 02092-15
Misc :
ALS Vial : 77 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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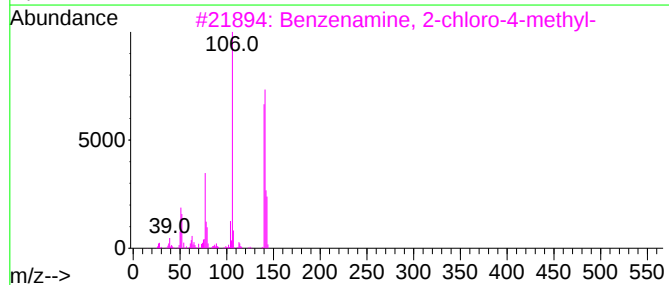
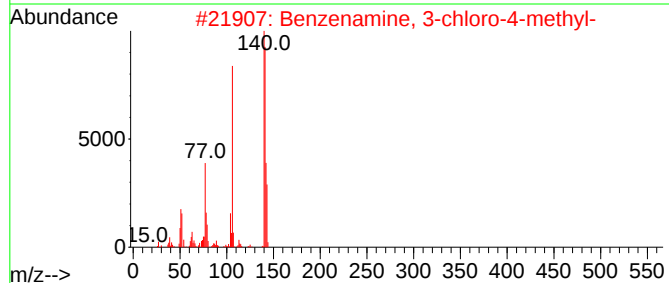
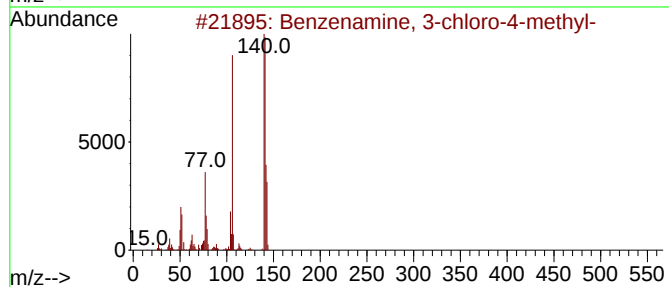
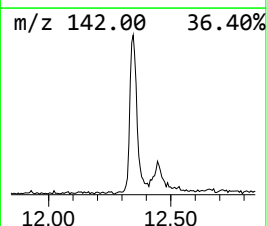
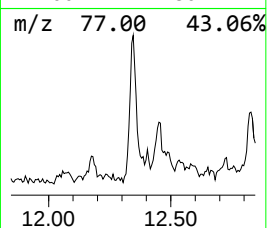
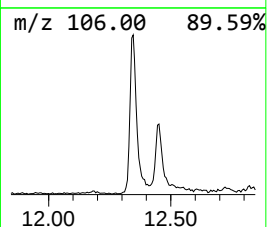
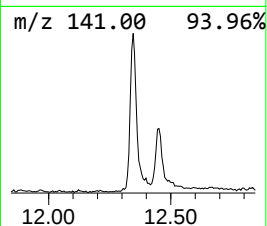
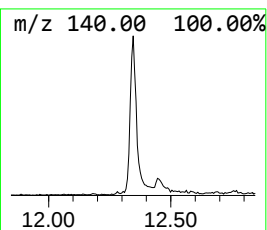
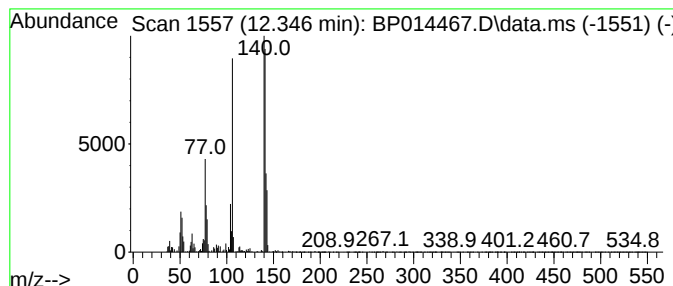
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Benzenamine, 3-chloro-4-met... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.346	3.70 ng/ul	373049	Naphthalene-d8	10.834

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	98
2			Benzenamine, 3-chloro-4-methyl-	141	C7H8ClN	000095-74-9	97
3			Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	93
4			Benzenamine, 3-chloro-2-methyl-	141	C7H8ClN	000087-60-5	89
5			Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014467.D
Acq On : 01 Apr 2023 04:55
Operator : CG/JU
Sample : 02092-15
Misc :
ALS Vial : 77 Sample Multiplier: 1

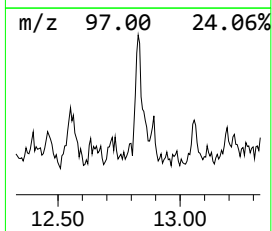
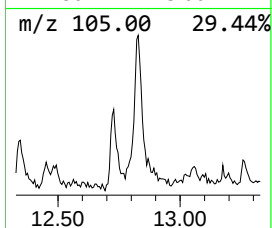
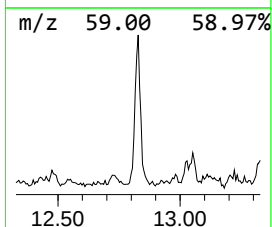
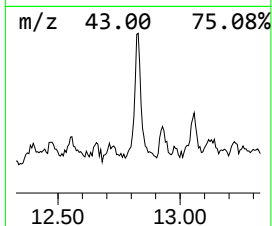
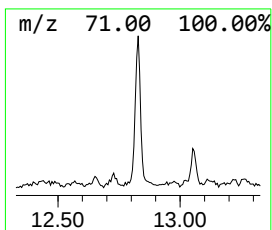
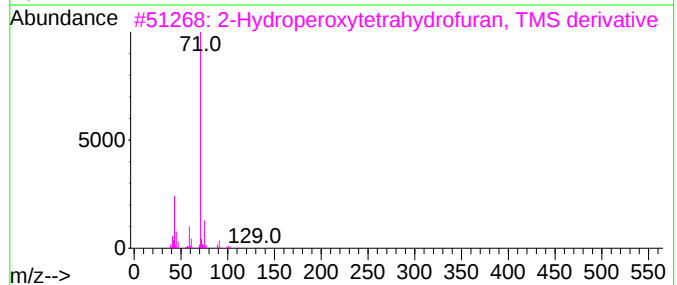
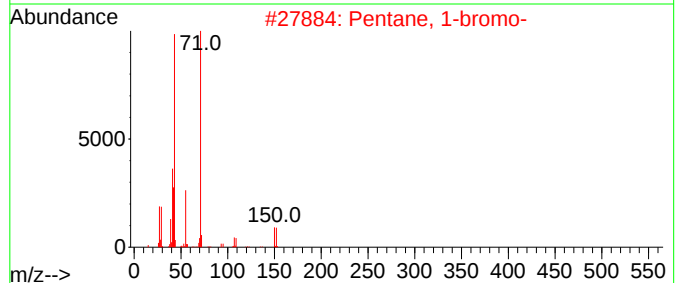
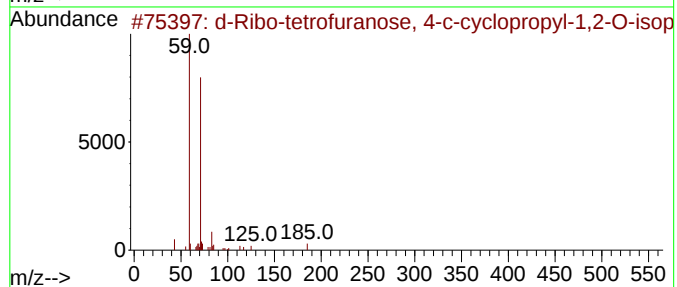
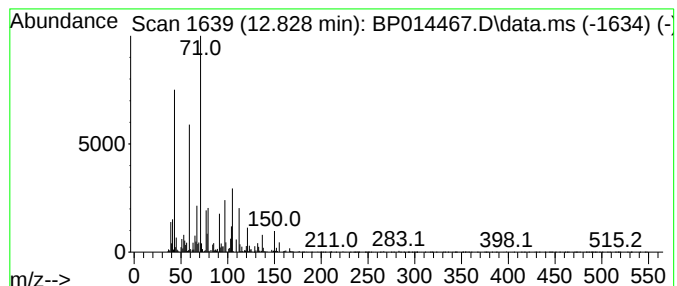
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 unknown-08 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.828	3.33 ng/ul	449163	Acenaphthene-d10	14.669	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	d-Ribo-tetrofuranose, 4-c-cyclop...	200	C10H16O4	030571-50-7	25
2	Pentane, 1-bromo-	150	C5H11Br	000110-53-2	22
3	2-Hydroperoxytetrahydrofuran, TM...	176	C7H16O3Si	1000368-50-8	22
4	3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	16
5	4,4-Dimethyl-1-hexene	112	C8H16	001647-08-1	16



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014467.D
Acq On : 01 Apr 2023 04:55
Operator : CG/JU
Sample : 02092-15
Misc :
ALS Vial : 77 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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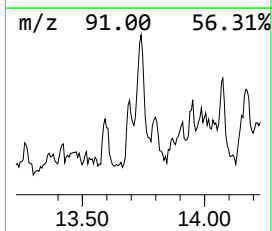
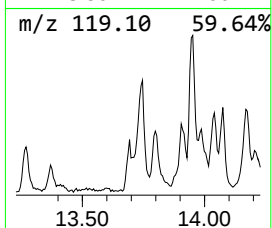
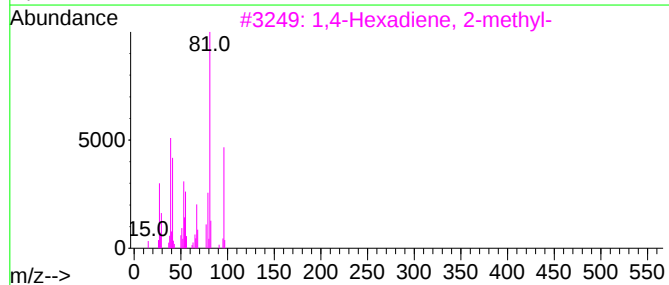
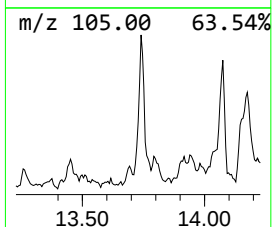
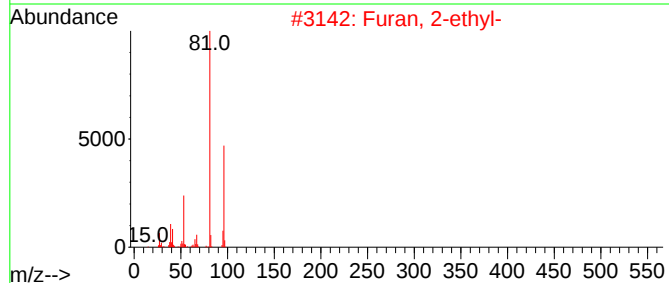
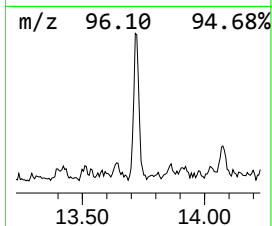
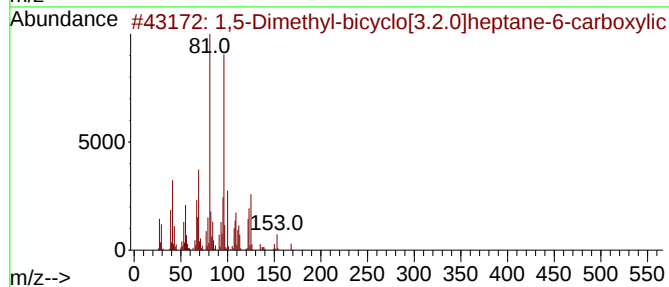
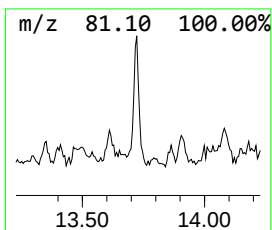
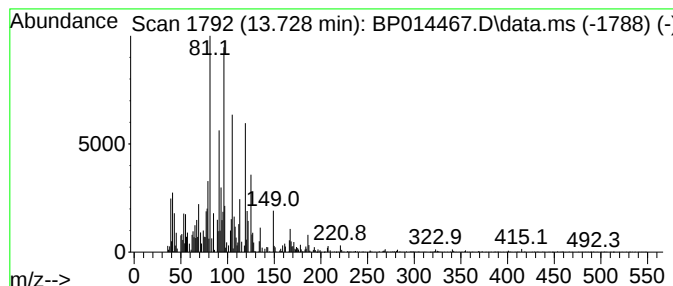
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 1,5-Dimethyl-bicyclo[3.2.0]... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.728	2.19 ng/ul	295257	Acenaphthene-d10	14.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,5-Dimethyl-bicyclo[3.2.0]hepta...	168	C10H16O2	1000190-69-8	53
2			Furan, 2-ethyl-	96	C6H8O	003208-16-0	46
3			1,4-Hexadiene, 2-methyl-	96	C7H12	001119-14-8	46
4			1,3-Pentadiene, 2,3-dimethyl-	96	C7H12	001113-56-0	46
5			2,4-Hexadiene, 3-methyl-	96	C7H12	028823-42-9	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
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Acq On : 01 Apr 2023 04:55
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Misc :
ALS Vial : 77 Sample Multiplier: 1

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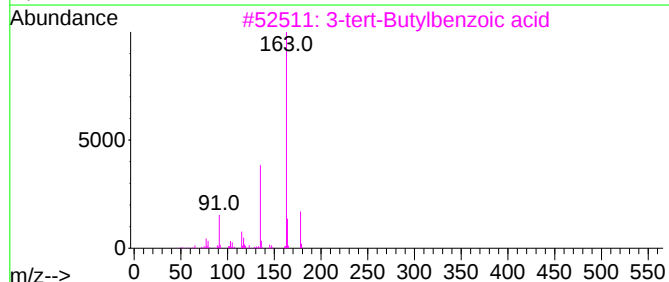
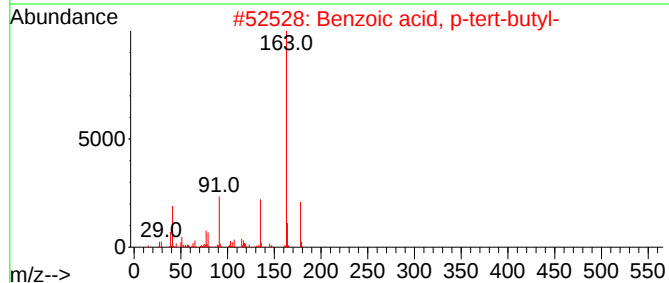
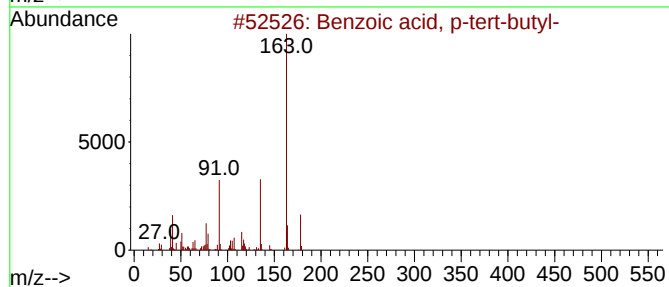
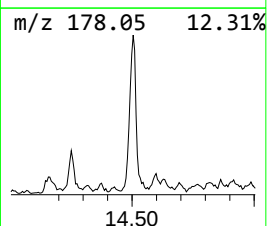
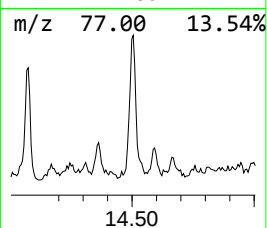
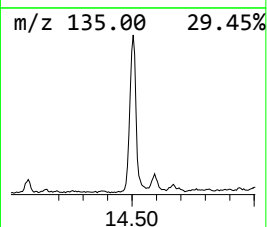
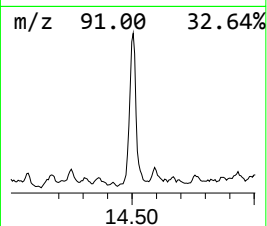
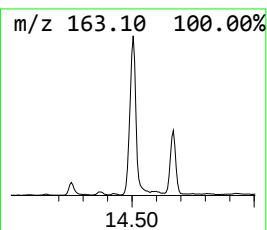
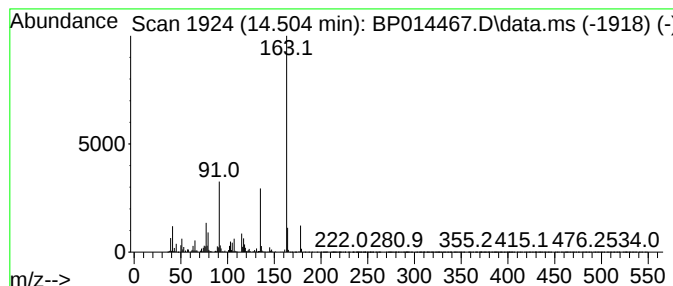
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Benzoic acid, p-tert-butyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.504	9.73 ng/ul	1311410	Acenaphthene-d10	14.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid, p-tert-butyl-	178	C11H14O2	000098-73-7	96
2			Benzoic acid, p-tert-butyl-	178	C11H14O2	000098-73-7	93
3			3-tert-Butylbenzoic acid	178	C11H14O2	007498-54-6	81
4			Benzoic acid, p-tert-butyl-	178	C11H14O2	000098-73-7	76
5			6-tert-Butyl-2,4-dimethylphenol	178	C12H18O	001879-09-0	64



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Data File : BP014467.D
Acq On : 01 Apr 2023 04:55
Operator : CG/JU
Sample : 02092-15
Misc :
ALS Vial : 77 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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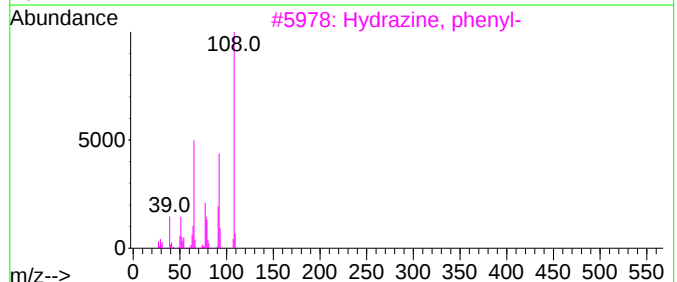
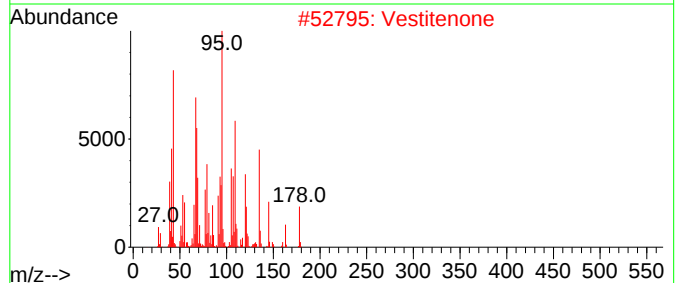
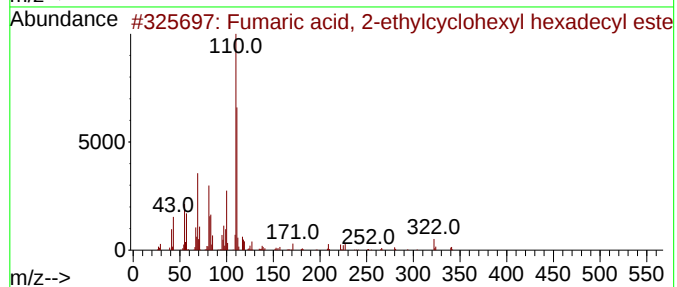
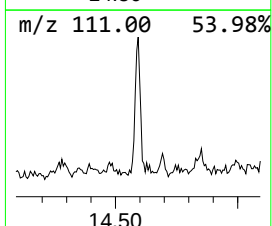
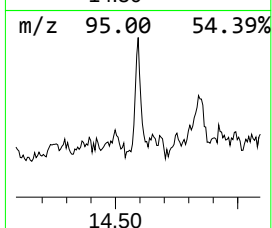
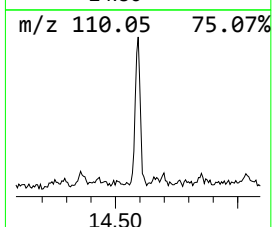
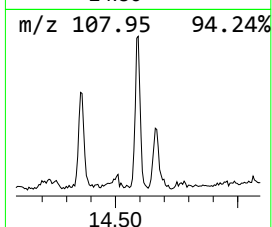
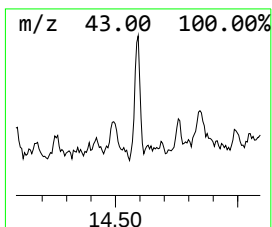
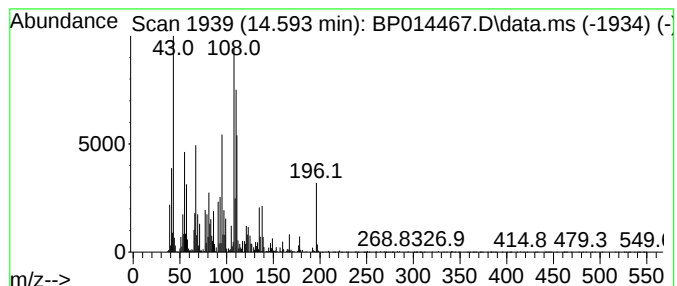
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 unknown-09 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.593	3.56 ng/ul	479364	Acenaphthene-d10	14.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fumaric acid, 2-ethylcyclohexyl ...	450	C28H50O4	1010345-19-6	22
2			Vestitenone	178	C12H18O	069401-36-1	14
3			Hydrazine, phenyl-	108	C6H8N2	000100-63-0	11
4			Hydroquinone	110	C6H6O2	000123-31-9	11
5			Carbamic acid, methyl-, 3-methyl...	165	C9H11NO2	001129-41-5	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014467.D
Acq On : 01 Apr 2023 04:55
Operator : CG/JU
Sample : 02092-15
Misc :
ALS Vial : 77 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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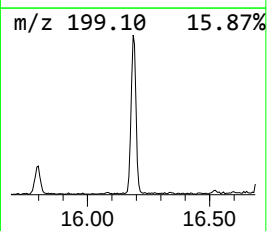
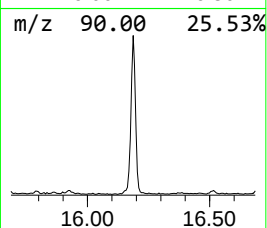
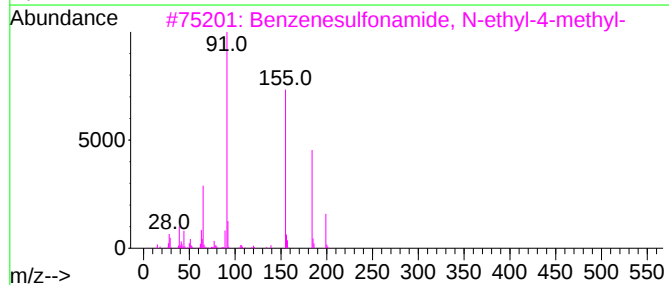
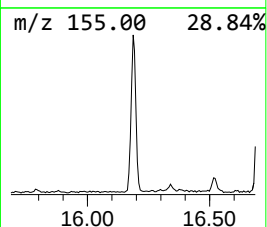
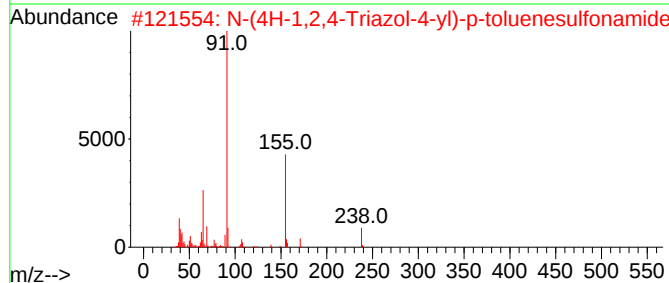
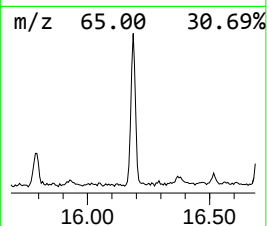
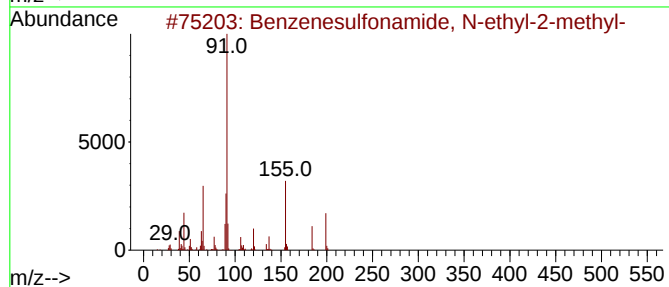
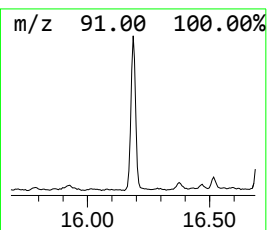
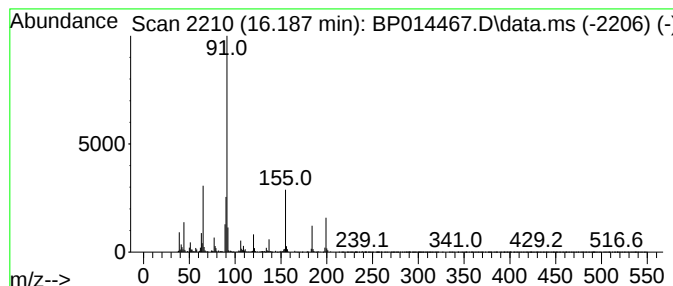
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 Benzenesulfonamide, N-ethyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.187	5.82 ng/ul	821785	Phenanthrene-d10	17.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	98
2			N-(4H-1,2,4-Triazol-4-yl)-p-tolu...	238	C9H10N4O2S	032585-76-5	49
3			Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	49
4			Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	46
5			Benzene, 1-methyl-3-nitro-	137	C7H7NO2	000099-08-1	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014467.D
Acq On : 01 Apr 2023 04:55
Operator : CG/JU
Sample : 02092-15
Misc :
ALS Vial : 77 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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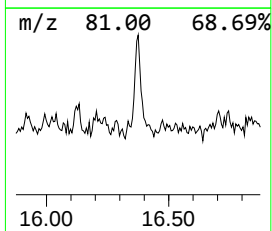
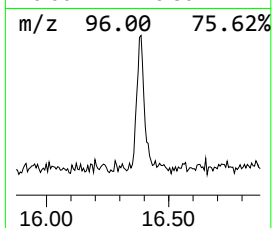
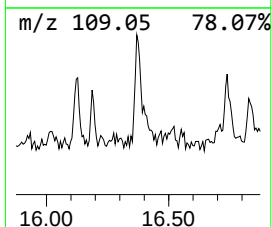
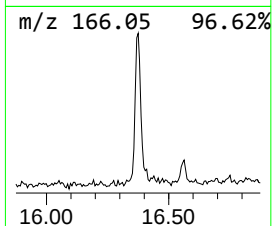
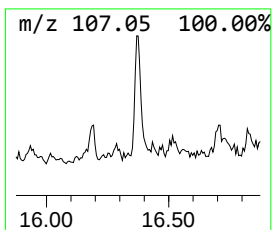
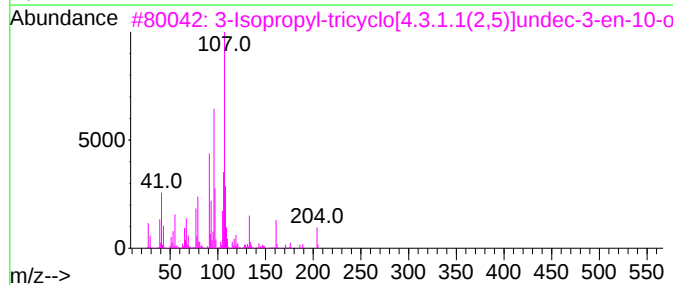
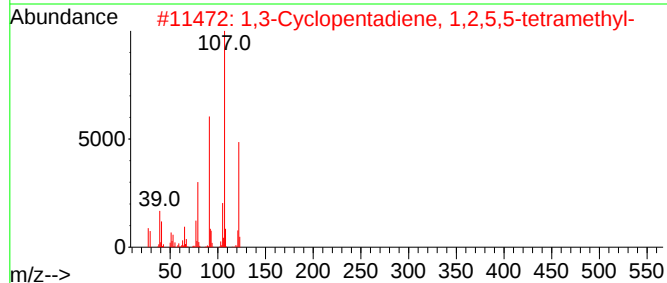
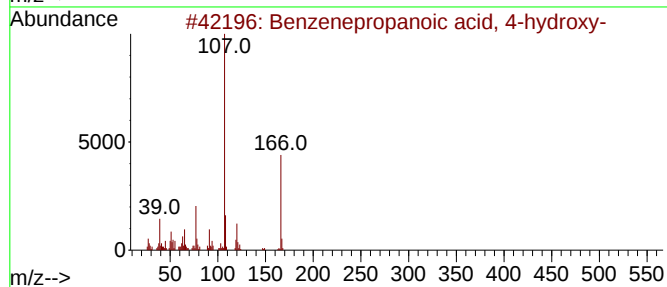
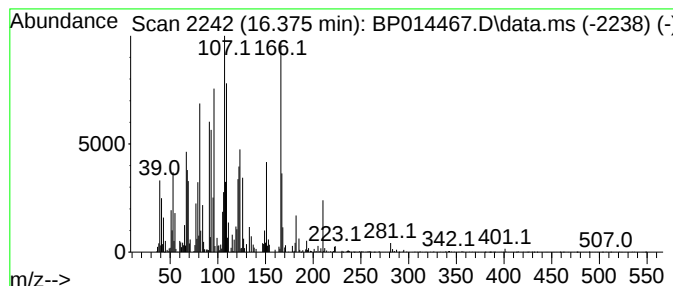
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 unknown-10 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.375	2.69 ng/ul	379614	Phenanthrene-d10	17.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenepropanoic acid, 4-hydroxy-	166	C9H10O3	000501-97-3	22
2			1,3-Cyclopentadiene, 1,2,5,5-tet...	122	C9H14	004249-12-1	15
3			3-Isopropyl-tricyclo[4.3.1.1(2,5...	204	C14H20O	1000191-47-2	14
4			Diacetamate	193	C10H11NO3	002623-33-8	11
5			Phenol, 3-ethyl-	122	C8H10O	000620-17-7	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014467.D
Acq On : 01 Apr 2023 04:55
Operator : CG/JU
Sample : 02092-15
Misc :
ALS Vial : 77 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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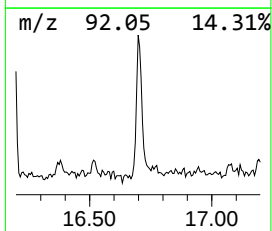
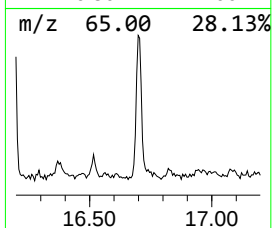
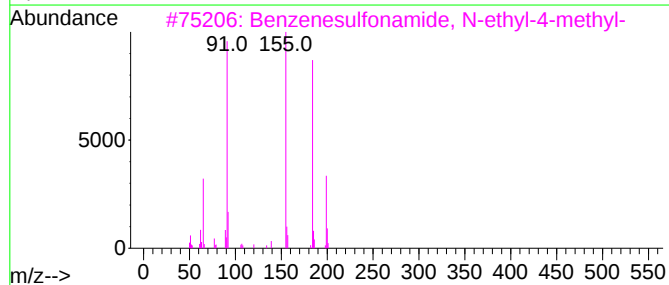
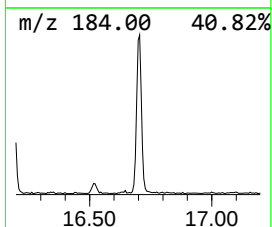
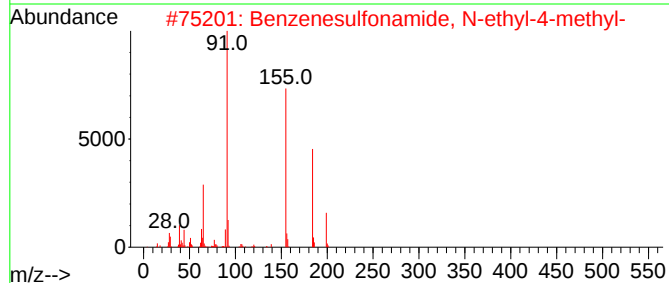
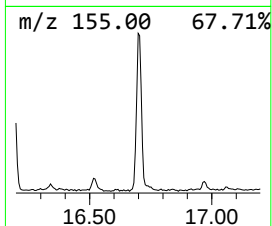
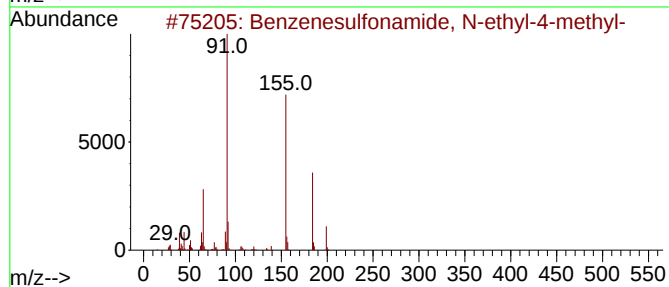
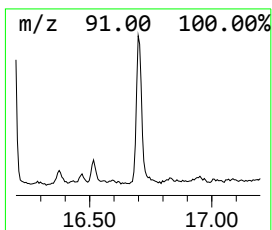
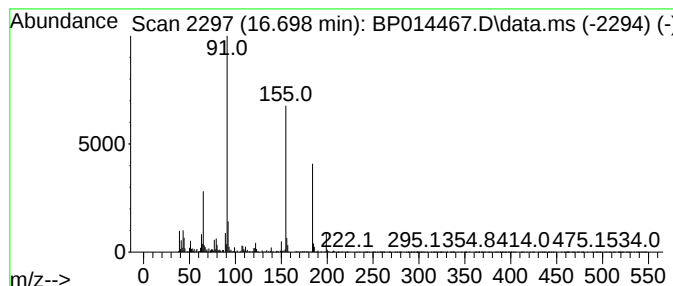
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 Benzenesulfonamide, N-ethyl... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.698	3.49 ng/ul	492998	Phenanthrene-d10	17.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	95
2			Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	94
3			Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	94
4			Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	91
5			4-Methyl-N-(2-oxobutyl)benzenesu...	241	C11H15NO3S	1000192-14-5	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014467.D
Acq On : 01 Apr 2023 04:55
Operator : CG/JU
Sample : 02092-15
Misc :
ALS Vial : 77 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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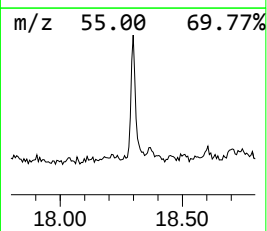
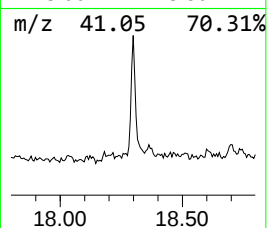
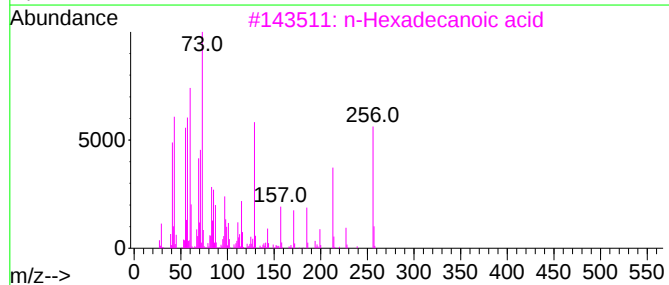
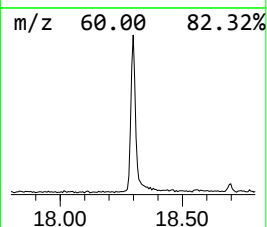
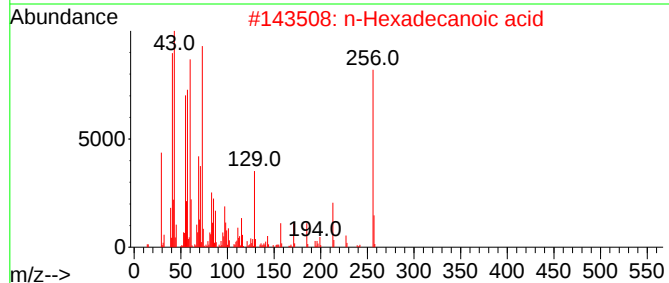
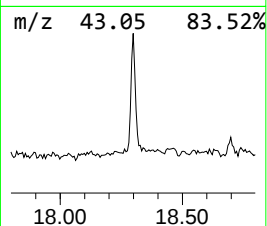
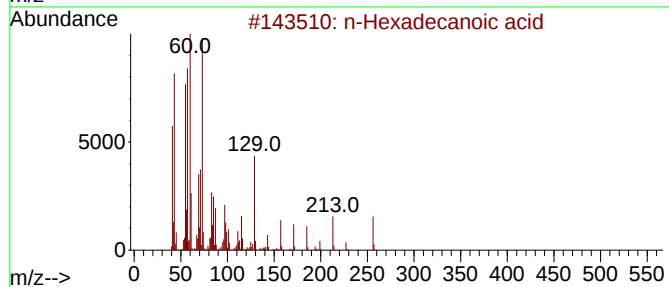
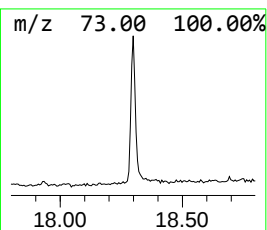
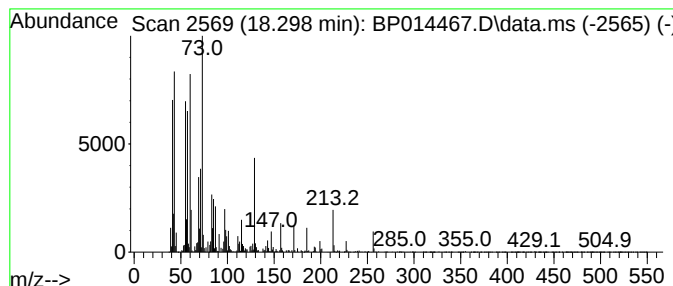
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.298	5.76 ng/ul	814209	Phenanthrene-d10	17.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
5			Tridecanoic acid	214	C13H26O2	000638-53-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014467.D
Acq On : 01 Apr 2023 04:55
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Sample : 02092-15
Misc :
ALS Vial : 77 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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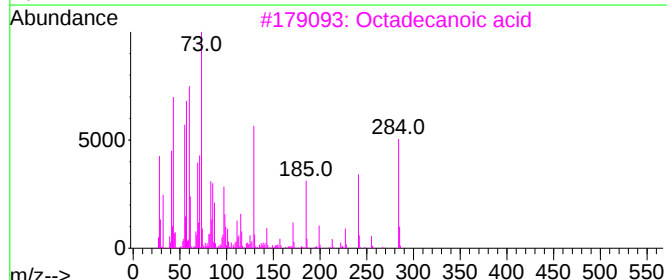
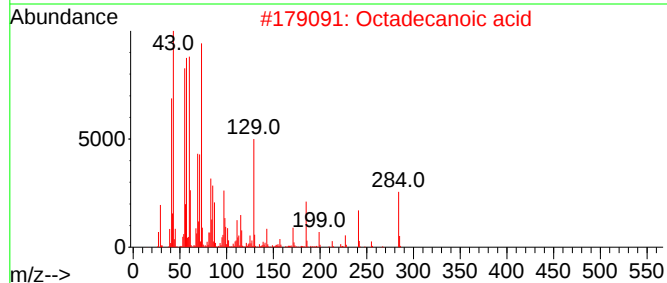
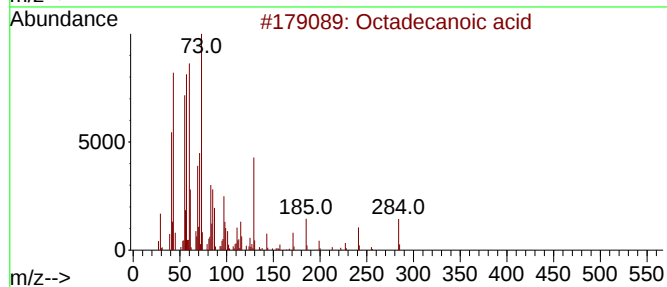
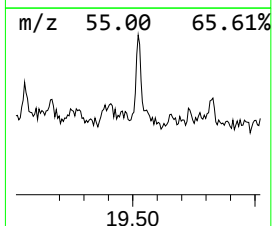
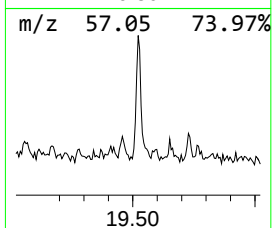
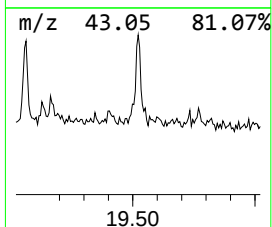
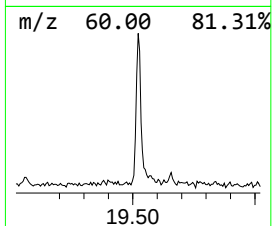
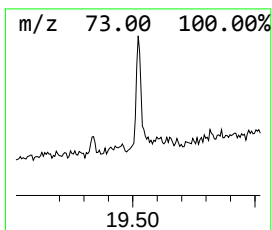
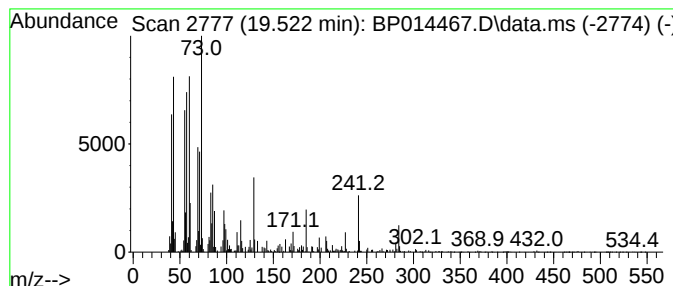
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 Octadecanoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.522	3.38 ng/ul	363856	Chrysene-d12	21.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	98
3			Octadecanoic acid	284	C18H36O2	000057-11-4	96
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014467.D
Acq On : 01 Apr 2023 04:55
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Sample : 02092-15
Misc :
ALS Vial : 77 Sample Multiplier: 1

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ClientSampleId :
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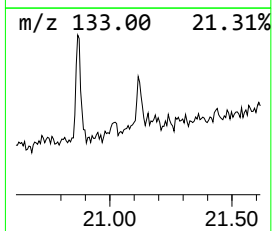
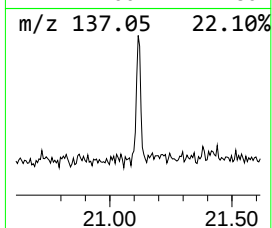
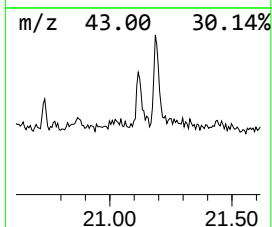
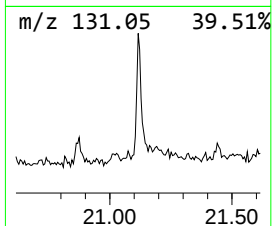
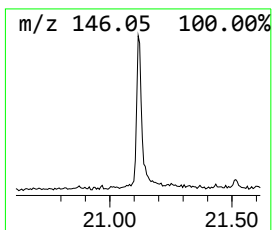
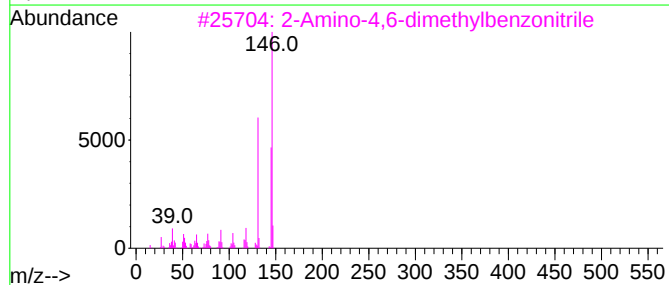
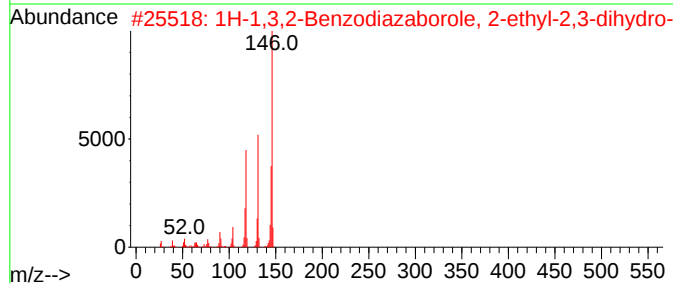
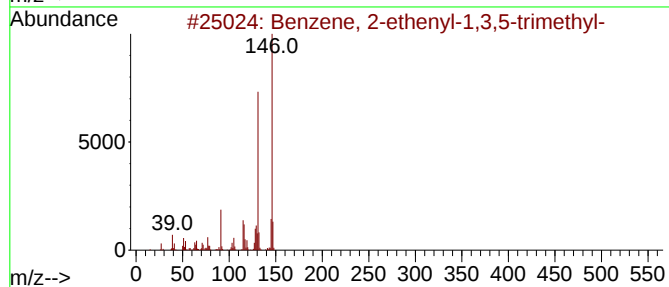
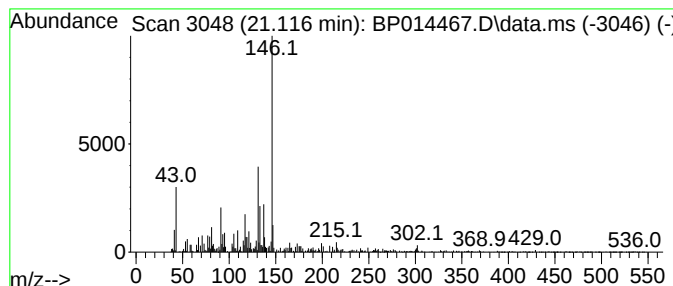
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 Benzene, 2-ethenyl-1,3,5-tr... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.116	3.06 ng/ul	329256	Chrysene-d12	21.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	52
2			1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	52
3			2-Amino-4,6-dimethylbenzonitrile	146	C9H10N2	1010461-05-1	50
4			1H-Indole-3-ethanol, 5-hydroxy-	177	C10H11NO2	000154-02-9	43
5			(4-Methoxy-phenyl)-(3,5,5-trimet...	257	C15H19N3O	1000274-25-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014467.D
Acq On : 01 Apr 2023 04:55
Operator : CG/JU
Sample : 02092-15
Misc :
ALS Vial : 77 Sample Multiplier: 1

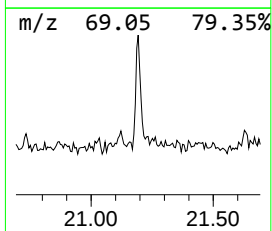
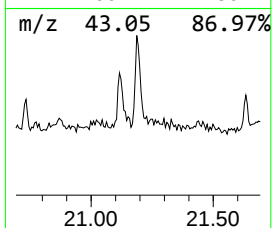
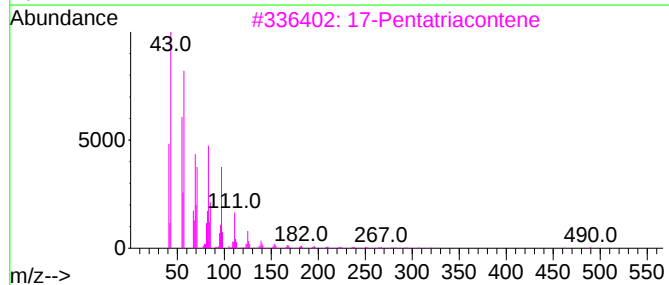
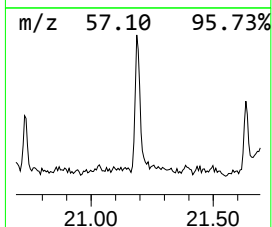
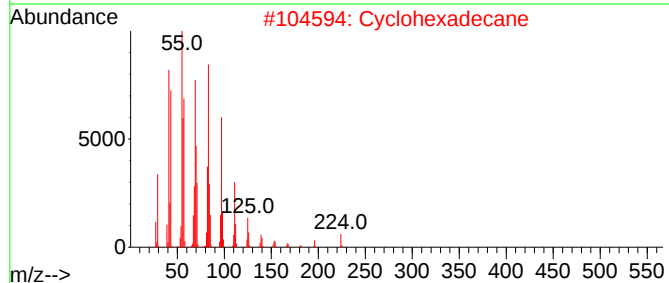
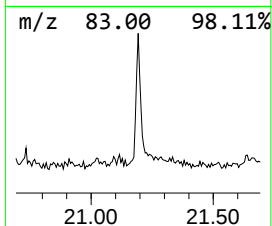
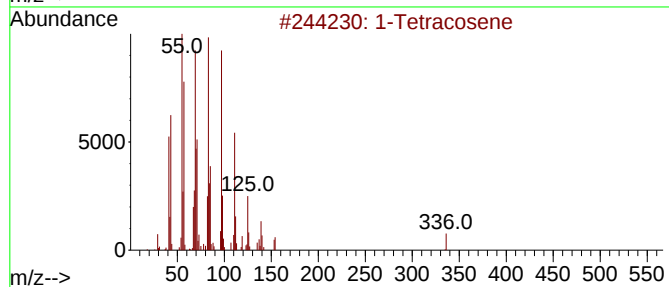
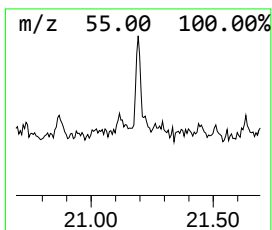
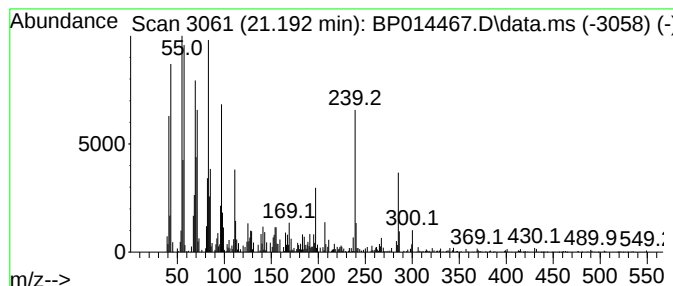
Instrument :
BNA_P
ClientSampleId :
CB982

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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

Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.192	3.38 ng/ul	364087	Chrysene-d12	21.516	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetracosene	336	C24H48	010192-32-2	97
2	Cyclohexadecane	224	C16H32	000295-65-8	97
3	17-Pentatriacontene	491	C35H70	006971-40-0	95
4	Cyclotetradecane	196	C14H28	000295-17-0	91
5	Cyclotetradecane	196	C14H28	000295-17-0	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
 Data File : BP014467.D
 Acq On : 01 Apr 2023 04:55
 Operator : CG/JU
 Sample : 02092-15
 Misc :
 ALS Vial : 77 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB982

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.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
Paraldehyde	4.217	14.7	ng/ul	946325	1	8.011	1290710	20.0
1,3-Oxathiolane	4.599	3.8	ng/ul	247471	1	8.011	1290710	20.0
Pentanoic acid,...	8.958	3.3	ng/ul	214125	1	8.011	1290710	20.0
Benzenamine, 3-...	9.011	19.6	ng/ul	1264700	1	8.011	1290710	20.0
unknown-01	9.081	7.4	ng/ul	480359	1	8.011	1290710	20.0
unknown-02	9.122	4.5	ng/ul	289274	1	8.011	1290710	20.0
unknown-03	9.269	2.6	ng/ul	166768	1	8.011	1290710	20.0
unknown-04	9.381	2.5	ng/ul	164529	1	8.011	1290710	20.0
Hexanoic acid, ...	9.693	12.0	ng/ul	1209350	2	10.834	2017490	20.0
unknown-05	9.734	3.7	ng/ul	376220	2	10.834	2017490	20.0
unknown-06	10.375	3.1	ng/ul	315174	2	10.834	2017490	20.0
unknown-07	10.616	2.2	ng/ul	223998	2	10.834	2017490	20.0
Benzenamine, 3-...	12.346	3.7	ng/ul	373049	2	10.834	2017490	20.0
unknown-08	12.828	3.3	ng/ul	449163	3	14.669	2694360	20.0
1,5-Dimethyl-bi...	13.728	2.2	ng/ul	295257	3	14.669	2694360	20.0
Benzoic acid, p...	14.504	9.7	ng/ul	1311410	3	14.669	2694360	20.0
unknown-09	14.593	3.6	ng/ul	479364	3	14.669	2694360	20.0
Benzenesulfonam...	16.187	5.8	ng/ul	821785	4	17.428	2825760	20.0
unknown-10	16.375	2.7	ng/ul	379614	4	17.428	2825760	20.0
Benzenesulfonam...	16.698	3.5	ng/ul	492998	4	17.428	2825760	20.0
n-Hexadecanoic ...	18.298	5.8	ng/ul	814209	4	17.428	2825760	20.0
Octadecanoic acid	19.522	3.4	ng/ul	363856	5	21.516	2152630	20.0
Benzene, 2-ethe...	21.116	3.1	ng/ul	329256	5	21.516	2152630	20.0
(DEL) Alkane: S...	21.192	3.4	ng/ul	364087	5	21.516	2152630	20.0