

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121823\
 Data File : BP018939.D
 Acq On : 19 Dec 2023 10:01
 Operator : MA/JU
 Sample : 05682-01 10X
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCLD6

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-BP120123.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP018939.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.152	8	11	19	rVB	41152	55764	0.69%	0.083%
2	3.258	26	29	32	rBV2	29837	41236	0.51%	0.061%
3	3.299	32	36	49	rVB	222734	298185	3.69%	0.444%
4	3.605	85	88	92	rVB2	21484	27002	0.33%	0.040%
5	3.681	97	101	106	rBV	374926	522909	6.48%	0.779%
6	3.728	106	109	115	rVB	74461	99390	1.23%	0.148%
7	3.905	136	139	145	rVB2	19356	28960	0.36%	0.043%
8	3.987	149	153	161	rVB2	33258	57213	0.71%	0.085%
9	4.105	168	173	178	rBV	56816	82301	1.02%	0.123%
10	4.152	178	181	188	rVB2	14799	23144	0.29%	0.034%
11	4.270	197	201	208	rVB	105003	135684	1.68%	0.202%
12	4.340	209	213	217	rVB3	21467	27471	0.34%	0.041%
13	4.452	228	232	238	rVB	98528	124832	1.55%	0.186%
14	4.523	239	244	252	rBV2	31491	65706	0.81%	0.098%
15	4.911	305	310	320	rVB	199571	287234	3.56%	0.428%
16	5.499	404	410	415	rBV	19372	31755	0.39%	0.047%
17	5.775	454	457	462	rVB4	10997	15813	0.20%	0.024%
18	6.481	572	577	585	rVB5	9654	19031	0.24%	0.028%
19	6.846	635	639	645	rBV7	7419	11262	0.14%	0.017%
20	7.011	656	667	675	rVB	263856	453843	5.62%	0.676%
21	7.169	687	694	702	rVB	902125	1415178	17.53%	2.107%
22	7.364	717	727	737	rBV	943283	1483790	18.38%	2.210%
23	7.446	737	741	745	rVB6	10284	16429	0.20%	0.024%
24	7.587	760	765	769	rBV7	7028	14946	0.19%	0.022%
25	7.828	798	806	813	rBV	726129	1118594	13.86%	1.666%
26	7.899	813	818	824	rVB7	8871	16839	0.21%	0.025%
27	8.087	845	850	856	rVV9	9222	19406	0.24%	0.029%
28	8.175	860	865	872	rVV3	10787	17926	0.22%	0.027%
29	8.546	921	928	939	rBV	695732	1147527	14.22%	1.709%
30	8.658	942	947	951	rVV	27451	40944	0.51%	0.061%
31	8.752	961	963	968	rVB6	7574	10374	0.13%	0.015%
32	8.946	990	996	998	rBV2	22666	36315	0.45%	0.054%
33	8.993	998	1004	1012	rVB	1056718	1693998	20.99%	2.523%
34	9.122	1021	1026	1030	rBV5	8001	15825	0.20%	0.024%
35	9.228	1035	1044	1052	rVV	97586	173327	2.15%	0.258%
36	9.552	1093	1099	1103	rVV8	8501	18215	0.23%	0.027%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
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37	9.710	1116	1126	1138	rBV	801091	1369028	16.96%	2.039%
38	9.893	1151	1157	1161	rBV9	5539	11934	0.15%	0.018%
39	10.252	1209	1218	1225	rBV	1231302	2060415	25.53%	3.068%
40	10.552	1265	1269	1272	rBV6	5256	8420	0.10%	0.013%
41	10.622	1272	1281	1288	rVV	860238	1402950	17.38%	2.089%
42	10.769	1298	1306	1326	rBV	1026472	1790192	22.18%	2.666%
43	11.034	1349	1351	1357	rVB6	5034	8803	0.11%	0.013%
44	11.163	1368	1373	1378	rBV	131022	200096	2.48%	0.298%
45	11.222	1378	1383	1390	rVV	167427	271673	3.37%	0.405%
46	11.275	1390	1392	1400	rVB9	8155	14418	0.18%	0.021%
47	11.428	1414	1418	1422	rVV5	12390	24099	0.30%	0.036%
48	11.475	1422	1426	1434	rVB5	14361	36077	0.45%	0.054%
49	11.981	1507	1512	1516	rBV2	12723	21473	0.27%	0.032%
50	12.034	1516	1521	1526	rVB	13550	23506	0.29%	0.035%
51	12.152	1538	1541	1545	rVV4	9757	15686	0.19%	0.023%
52	12.210	1547	1551	1556	rVB2	23308	36226	0.45%	0.054%
53	12.304	1562	1567	1573	rBV	39516	67785	0.84%	0.101%
54	12.828	1651	1656	1659	rBV7	6433	11032	0.14%	0.016%
55	12.922	1669	1672	1678	rVB3	13633	19777	0.25%	0.029%
56	12.993	1678	1684	1692	rBV4	17632	38614	0.48%	0.058%
57	13.199	1715	1719	1726	rBV6	6155	9262	0.11%	0.014%
58	13.363	1740	1747	1753	rBV	33385	59310	0.73%	0.088%
59	13.428	1753	1758	1766	rVB	8156	21477	0.27%	0.032%
60	13.651	1792	1796	1800	rVB7	6148	9651	0.12%	0.014%
61	13.881	1828	1835	1844	rBV	2184893	2967699	36.77%	4.419%
62	14.157	1872	1882	1891	rBV	2624261	3877266	48.04%	5.774%
63	14.351	1909	1915	1916	rBV5	7370	12440	0.15%	0.019%
64	14.463	1923	1934	1941	rBV	1204614	1848003	22.90%	2.752%
65	14.522	1941	1944	1951	rVB3	15279	28049	0.35%	0.042%
66	14.687	1964	1972	1984	rBV	174325	333431	4.13%	0.497%
67	14.828	1992	1996	1998	rBV5	7887	12326	0.15%	0.018%
68	14.863	1999	2002	2011	rVB2	21578	47340	0.59%	0.070%
69	15.028	2028	2030	2034	rVB5	7483	9978	0.12%	0.015%
70	15.069	2034	2037	2040	rBV5	6021	8243	0.10%	0.012%
71	15.134	2044	2048	2053	rVB8	8529	11691	0.14%	0.017%
72	15.210	2056	2061	2064	rVV	38201	54737	0.68%	0.082%
73	15.246	2064	2067	2077	rVB	50687	80284	0.99%	0.120%
74	15.457	2094	2103	2109	rBV	3323404	4858261	60.19%	7.235%
75	15.510	2109	2112	2118	rVV3	43598	74480	0.92%	0.111%
76	15.581	2119	2124	2138	rVV	527194	762015	9.44%	1.135%

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77	15.693	2139	2143	2149	rVB5	17831	30285	0.38%	0.045%
78	16.151	2217	2221	2223	rBV5	7302	10706	0.13%	0.016%
79	16.234	2231	2235	2241	rBV2	23859	39041	0.48%	0.058%
80	16.757	2320	2324	2328	rVB5	10854	16317	0.20%	0.024%
81	17.210	2395	2401	2406	rBV	1541918	2254040	27.93%	3.357%
82	17.257	2406	2409	2413	rVV	142177	214138	2.65%	0.319%
83	17.310	2413	2418	2433	rVB	3837891	5677684	70.35%	8.455%
84	17.481	2443	2447	2450	rBV6	18287	32200	0.40%	0.048%
85	17.881	2513	2515	2519	rVB5	20027	22678	0.28%	0.034%
86	18.104	2548	2553	2562	rBV2	100805	177129	2.19%	0.264%
87	19.128	2723	2727	2733	rBV2	62880	89526	1.11%	0.133%
88	19.198	2735	2739	2741	rBV	62392	87841	1.09%	0.131%
89	19.228	2741	2744	2749	rVB	99364	138016	1.71%	0.206%
90	19.339	2759	2763	2769	rBV2	118704	181306	2.25%	0.270%
91	19.551	2793	2799	2810	rBV	5386559	7298953	90.43%	10.869%
92	21.304	3092	3097	3110	rVB	2325108	3221643	39.92%	4.798%
93	21.845	3186	3189	3195	rVB	420416	533813	6.61%	0.795%
94	22.445	3287	3291	3298	rVB	646850	966884	11.98%	1.440%
95	23.122	3401	3406	3414	rVB	723355	1318997	16.34%	1.964%
96	23.522	3467	3474	3489	rVB2	3898807	8071147	100.00%	12.019%
97	23.674	3494	3500	3508	rVB	1547993	2947373	36.52%	4.389%
98	23.904	3534	3539	3549	rVB2	726771	1655649	20.51%	2.466%

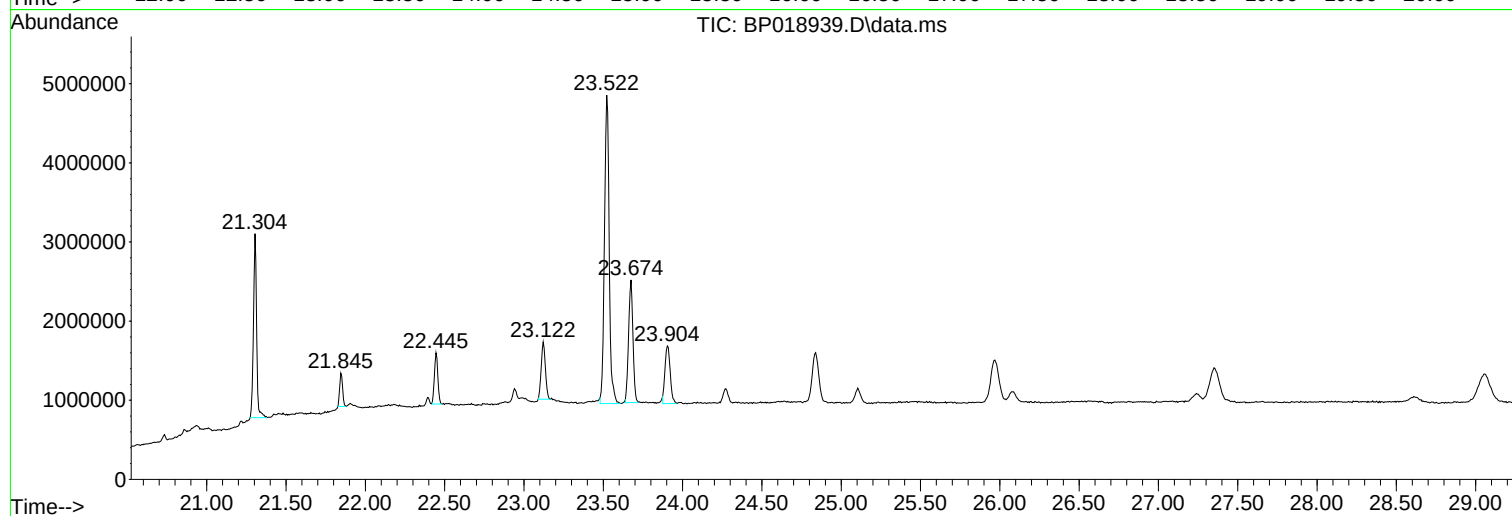
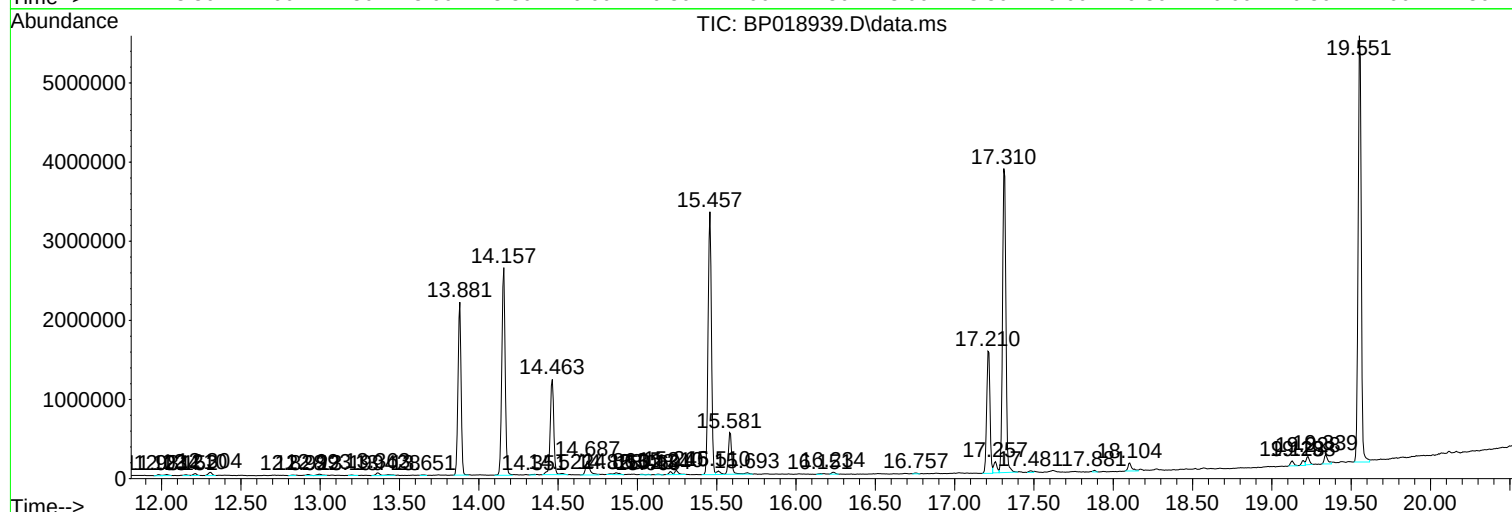
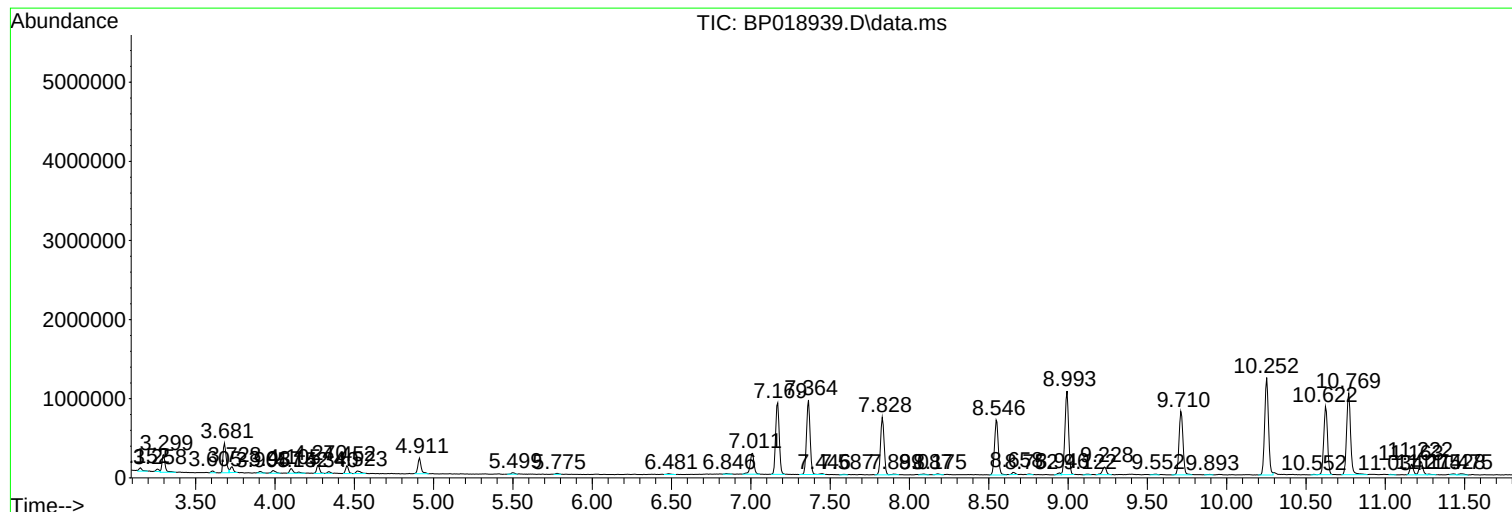
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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



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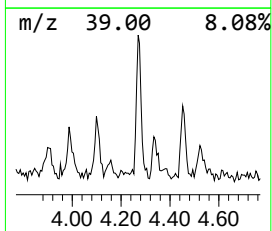
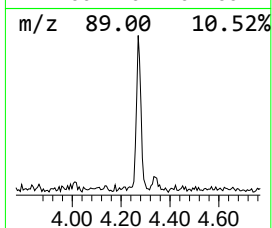
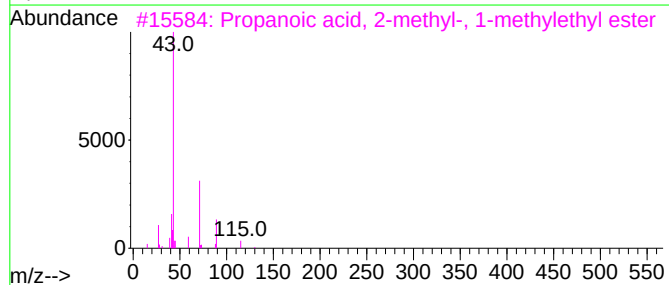
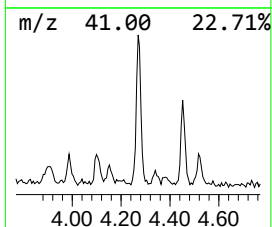
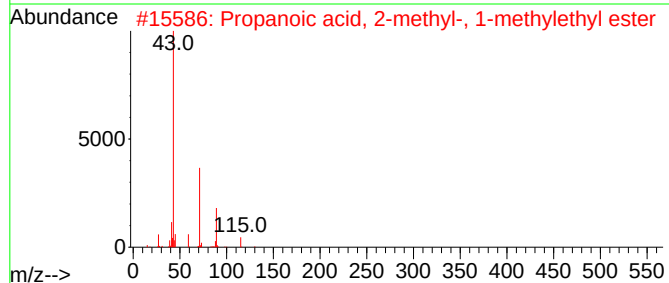
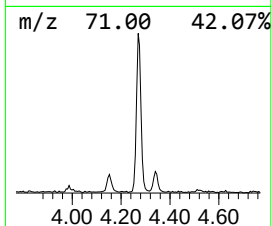
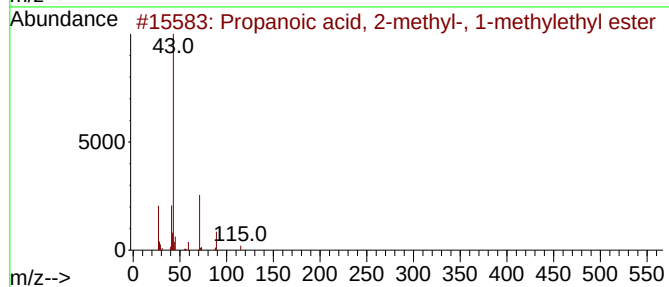
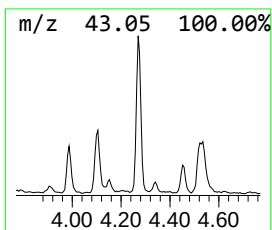
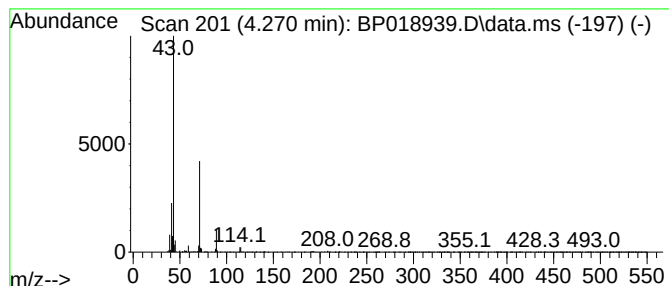
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propanoic acid, 2-methyl-, ... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.270	2.43 ng/ul	135684	1,4-Dichlorobenzene-d4	7.828

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	56
2			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	56
3			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	56
4			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	50
5			3-Hexanone, 2-methyl-	114	C7H14O	007379-12-6	38



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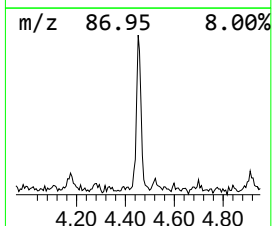
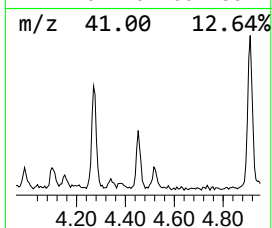
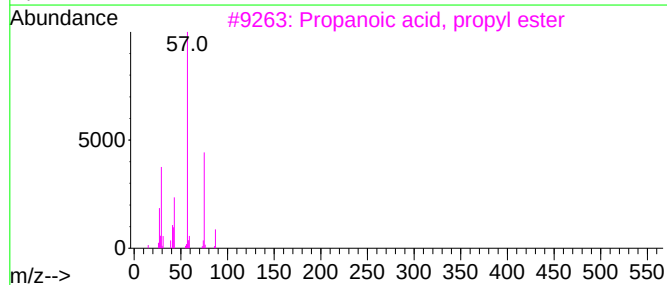
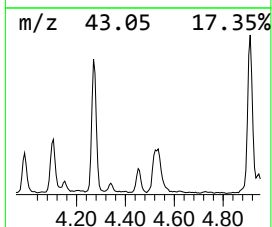
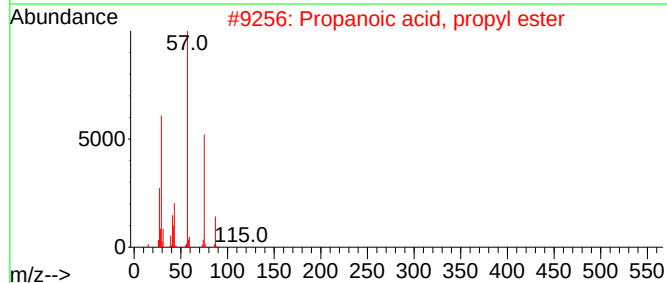
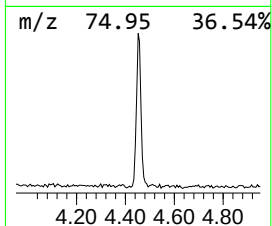
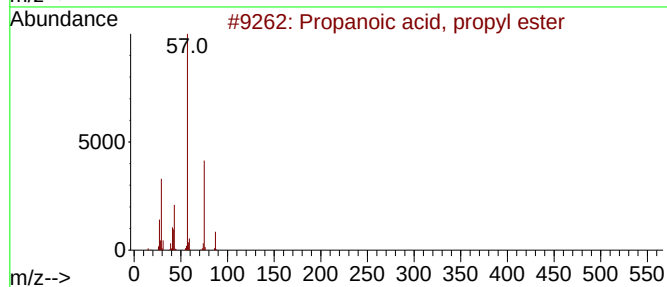
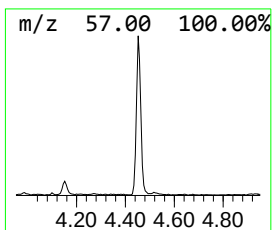
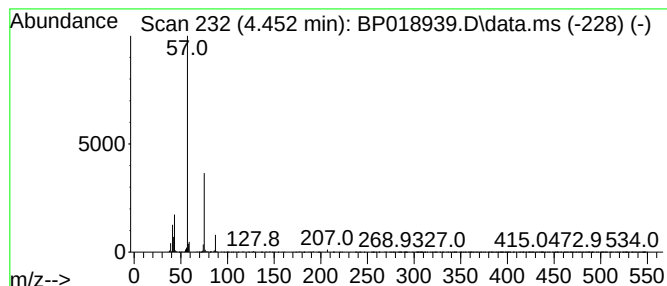
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Propanoic acid, propyl ester Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.452	2.23 ng/ul	124832	1,4-Dichlorobenzene-d4	7.828

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	78
2			Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	78
3			Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	78
4			Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	78
5			Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	56



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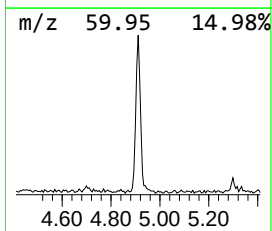
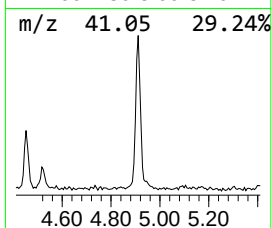
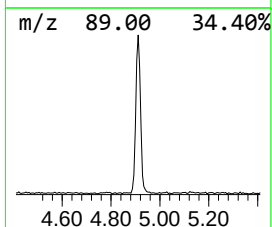
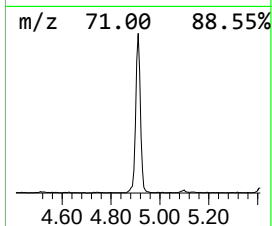
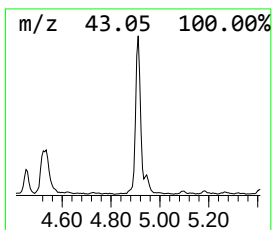
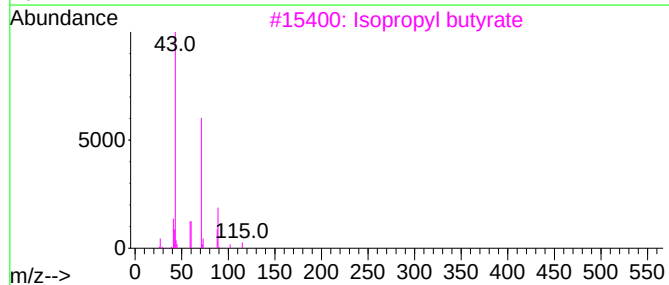
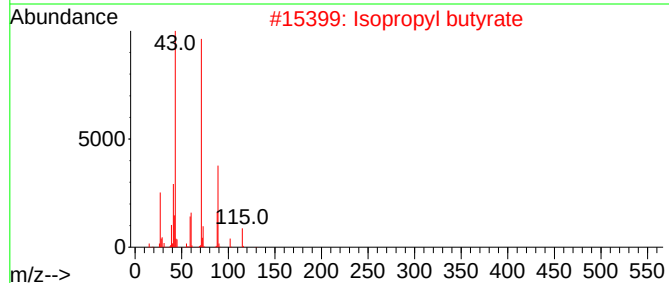
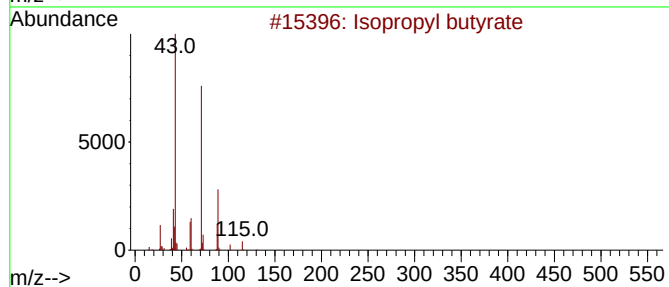
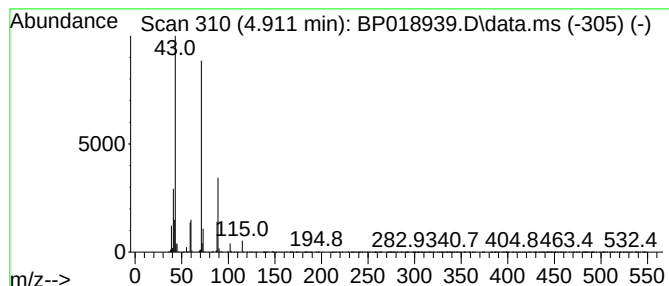
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Isopropyl butyrate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.911	5.14 ng/ul	287234	1,4-Dichlorobenzene-d4	7.828

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropyl butyrate	130	C7H14O2	000638-11-9	94
2			Isopropyl butyrate	130	C7H14O2	000638-11-9	90
3			Isopropyl butyrate	130	C7H14O2	000638-11-9	72
4			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	72
5			Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121823\
Data File : BP018939.D
Acq On : 19 Dec 2023 10:01
Operator : MA/JU
Sample : 05682-01 10X
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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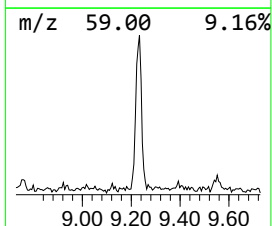
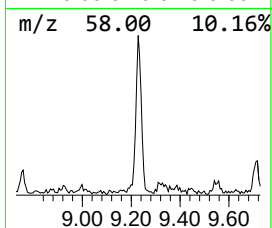
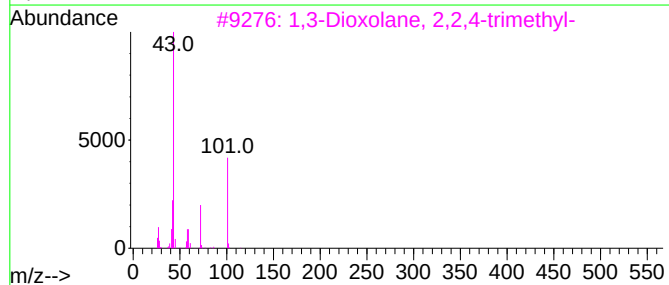
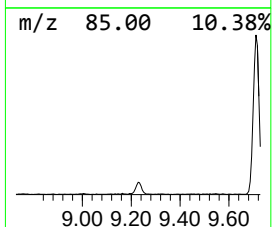
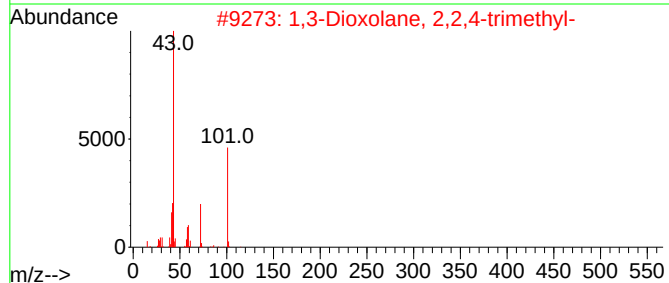
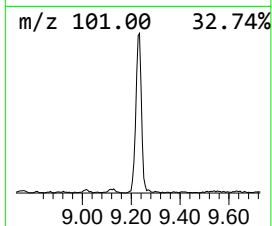
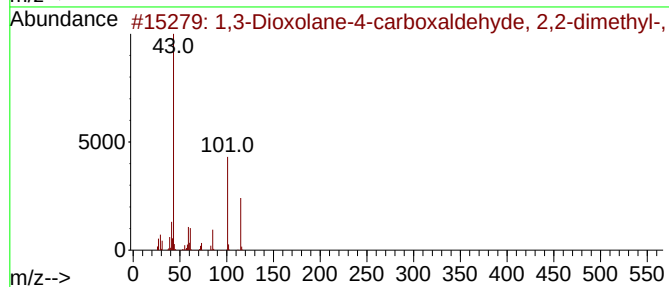
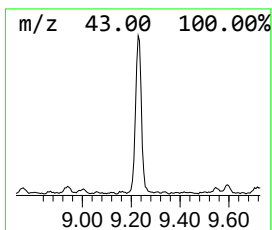
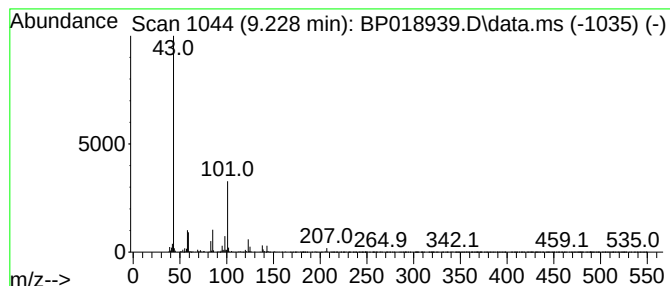
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.228	2.47 ng/ul	173327	Naphthalene-d8	10.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Dioxolane-4-carboxaldehyde, ...	130	C6H10O3	015186-48-8	40
2			1,3-Dioxolane, 2,2,4-trimethyl-	116	C6H12O2	001193-11-9	33
3			1,3-Dioxolane, 2,2,4-trimethyl-	116	C6H12O2	001193-11-9	33
4			Thiocyanic acid, propyl ester	101	C4H7NS	004251-16-5	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121823\
Data File : BP018939.D
Acq On : 19 Dec 2023 10:01
Operator : MA/JU
Sample : 05682-01 10X
Misc :
ALS Vial : 43 Sample Multiplier: 1

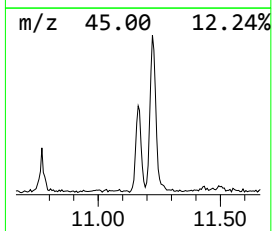
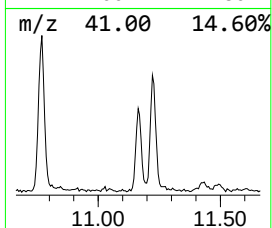
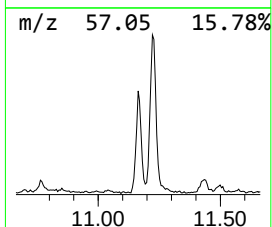
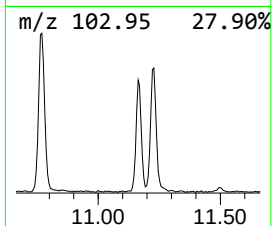
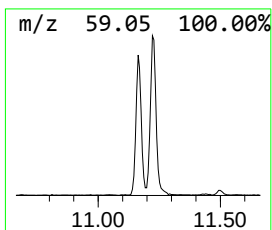
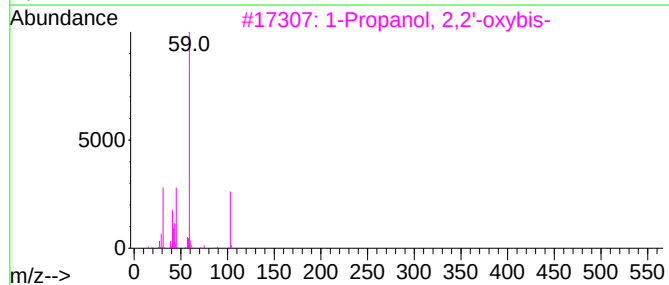
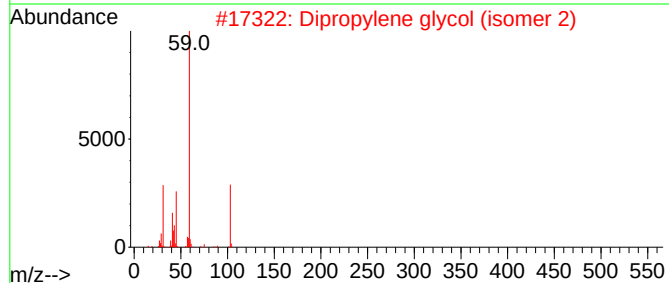
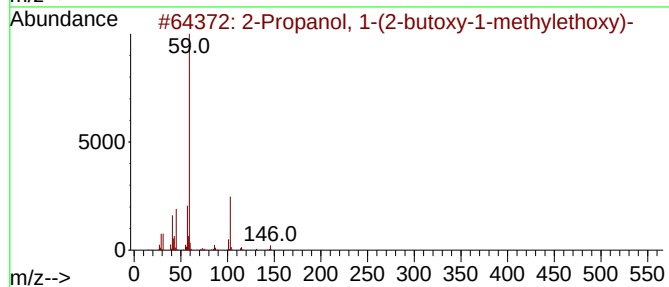
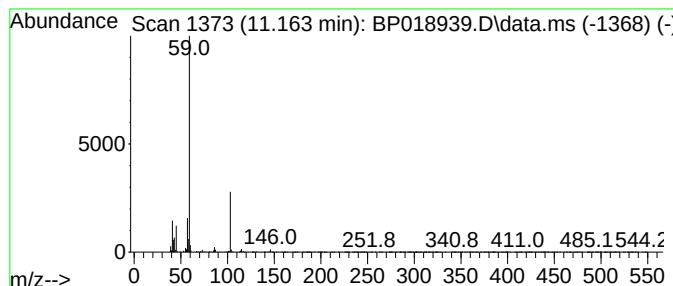
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 2-Propanol, 1-(2-butoxy-1-m... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.163	2.85 ng/ul	200096	Naphthalene-d8	10.622	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	83
2	Dipropylene glycol (isomer 2)	134	C6H14O3	1000506-25-4	72
3	1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	72
4	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	72
5	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121823\
Data File : BP018939.D
Acq On : 19 Dec 2023 10:01
Operator : MA/JU
Sample : 05682-01 10X
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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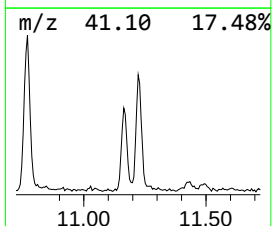
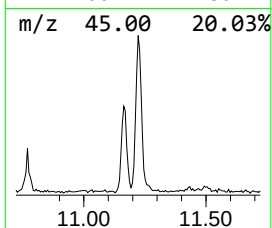
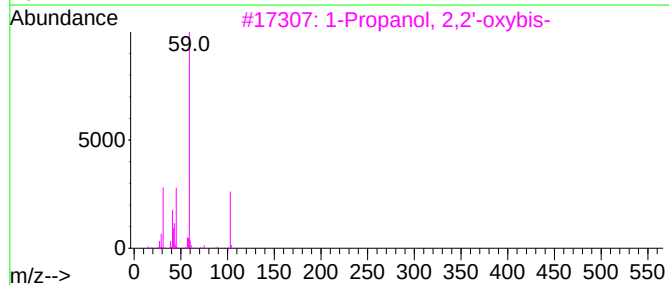
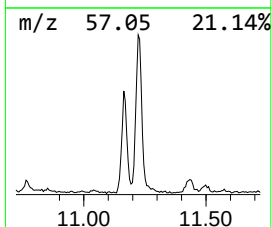
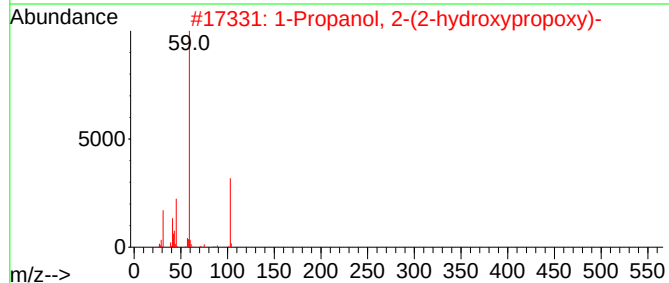
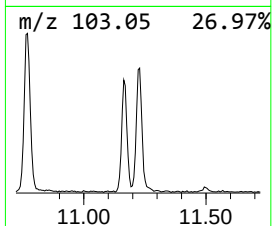
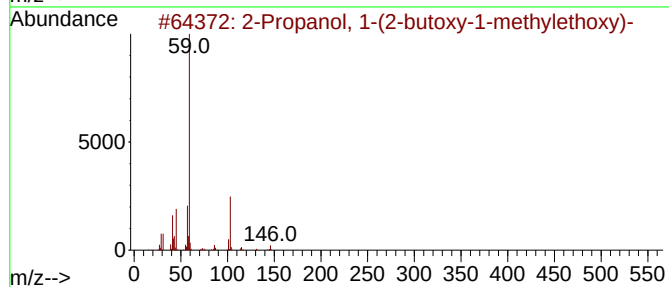
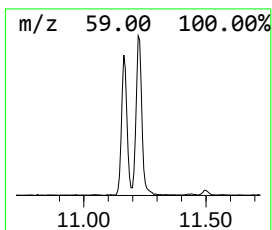
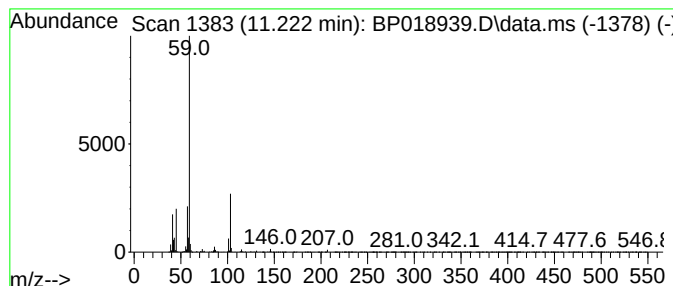
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1-Propanol, 2-(2-hydroxypro... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.222	3.87 ng/ul	271673	Naphthalene-d8	10.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	90		
2	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	78		
3	1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	64		
4	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	64		
5	1,4,7-Trimethyl-3,6-dioxaoctane-...	192	C9H20O4	1000423-63-2	64		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121823\
Data File : BP018939.D
Acq On : 19 Dec 2023 10:01
Operator : MA/JU
Sample : 05682-01 10X
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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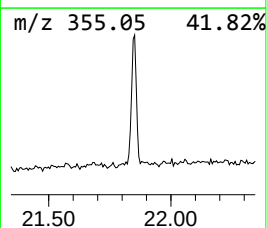
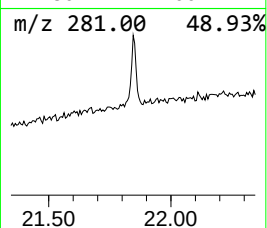
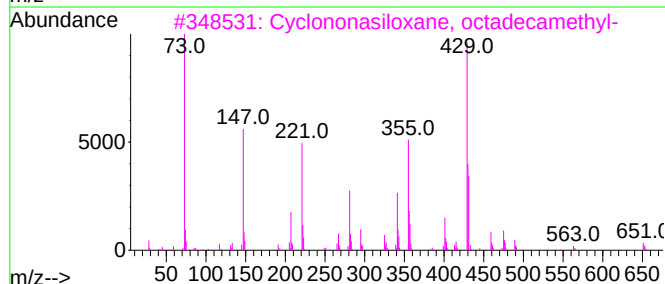
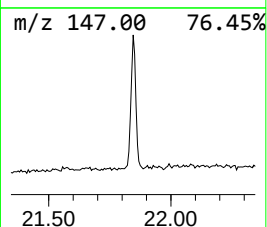
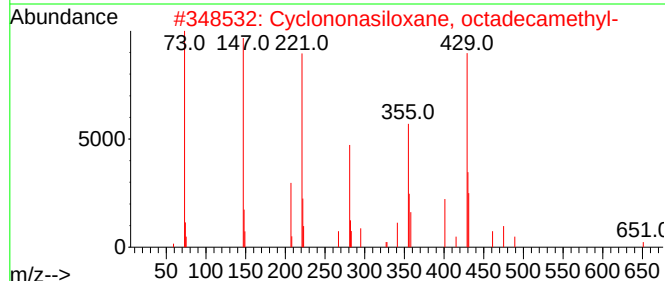
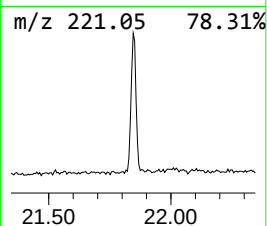
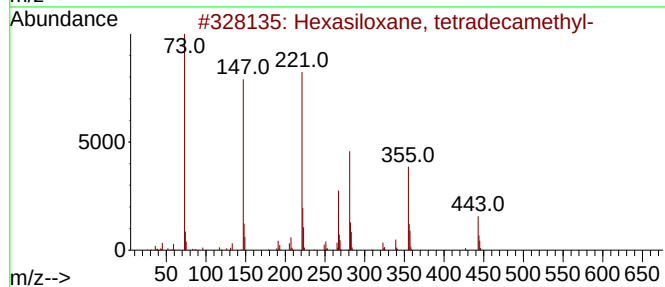
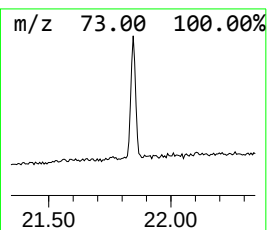
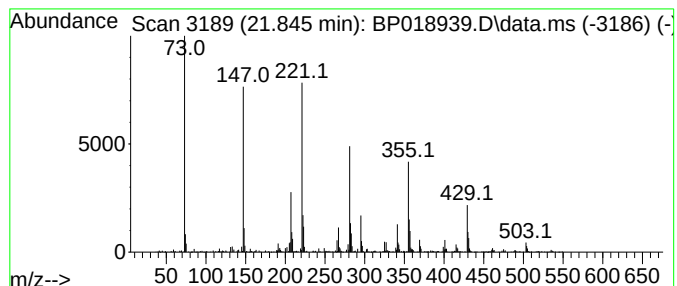
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Hexasiloxane, tetradecamethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.845	3.31 ng/ul	533813	Chrysene-d12	21.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	80
2			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	50
3			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	27
4			N-Benzyl-N-ethyl-p-isopropylbenz...	281	C19H23NO	015089-22-2	27
5			Pentasiloxane, dodecamethyl-	384	C12H36O4Si5	000141-63-9	16



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121823\
Data File : BP018939.D
Acq On : 19 Dec 2023 10:01
Operator : MA/JU
Sample : 05682-01 10X
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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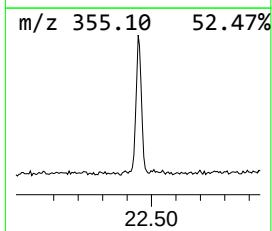
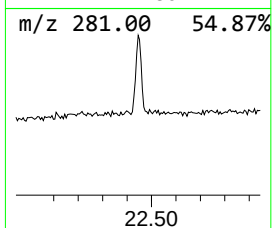
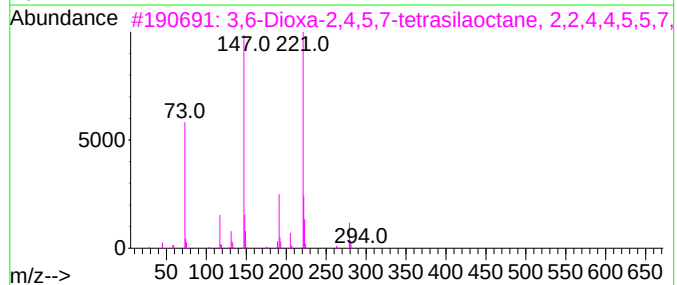
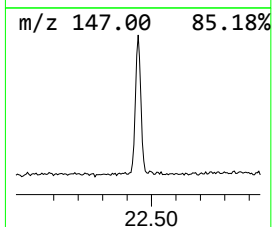
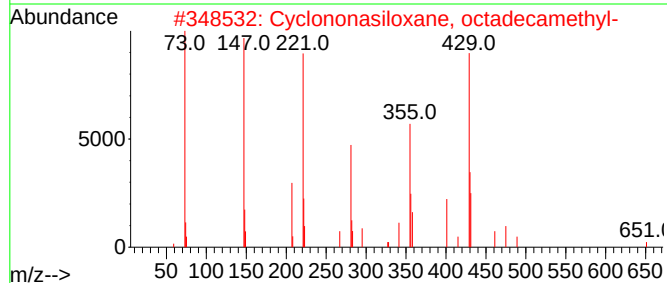
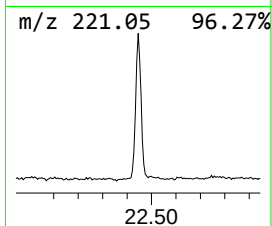
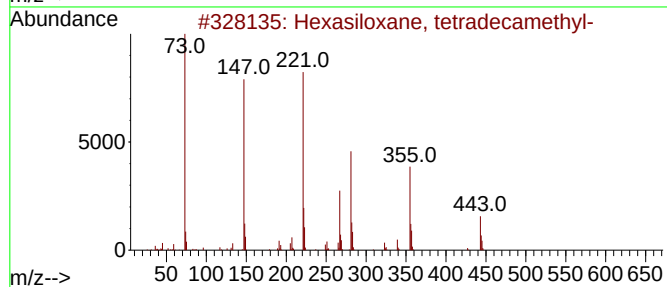
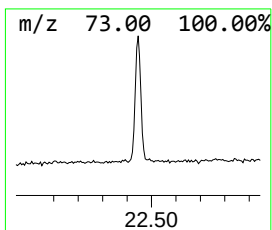
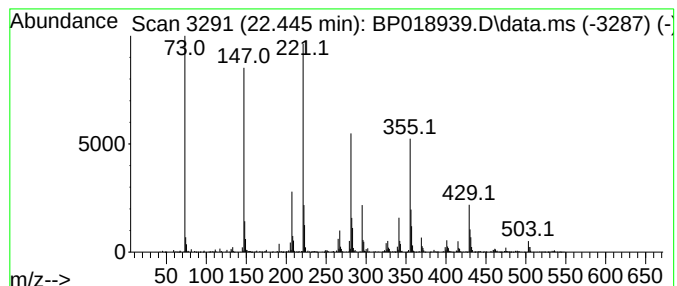
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Cyclononasiloxane, octadeca... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.445	6.00 ng/ul	966884	Chrysene-d12	21.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	74
2			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	72
3			3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	46
4			2,5-Dihydroxybenzoic acid, 3TMS ...	370	C16H30O4Si3	003618-20-0	44
5			Cyclodecasiloxane, eicosamethyl-	740	C20H60O10Si10	018772-36-6	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121823\
Data File : BP018939.D
Acq On : 19 Dec 2023 10:01
Operator : MA/JU
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Misc :
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Instrument :
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ClientSampleId :
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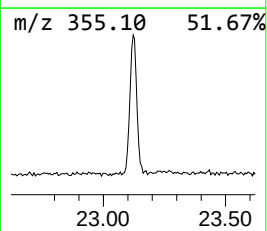
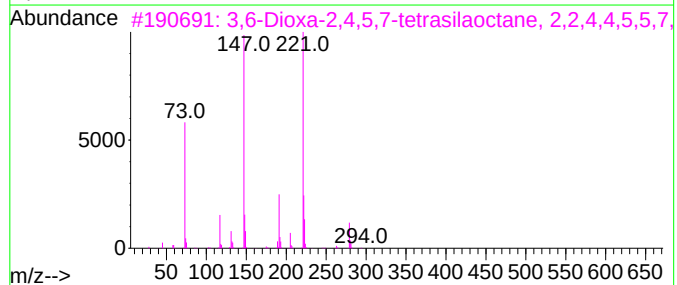
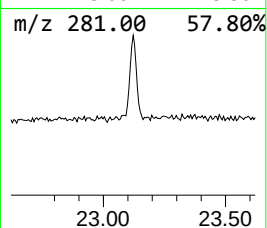
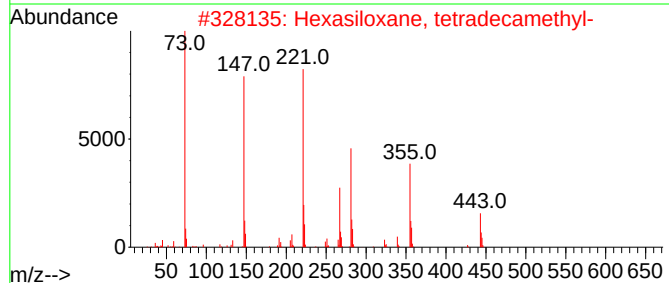
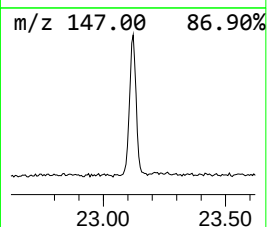
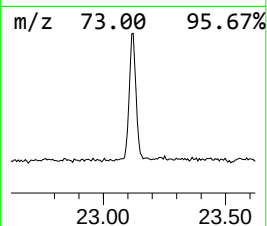
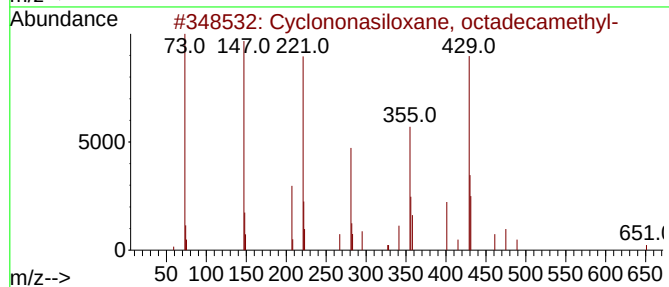
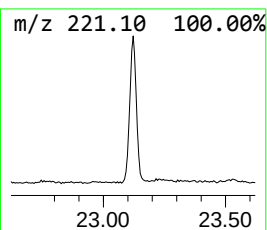
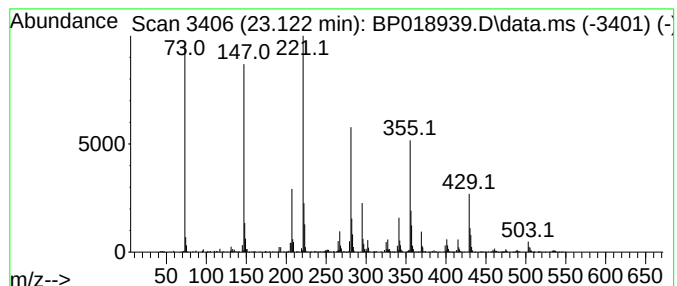
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-02 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.122	8.95 ng/ul	1319000	Perylene-d12	23.674

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	90
2			Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	74
3			3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	37
4			Heptasiloxane, hexadecamethyl-	532	C16H48O6Si7	000541-01-5	25
5			Thioglycolic acid, S-(tert-butyl...	320	C14H32O2SSi2	082112-29-6	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121823\
Data File : BP018939.D
Acq On : 19 Dec 2023 10:01
Operator : MA/JU
Sample : 05682-01 10X
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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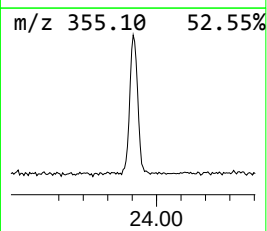
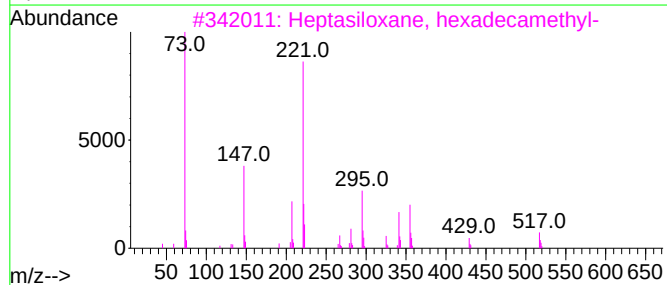
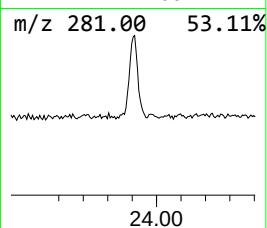
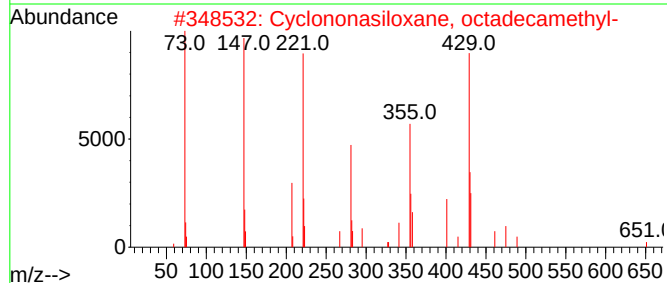
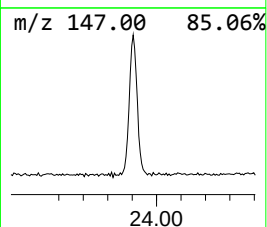
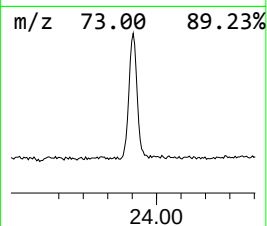
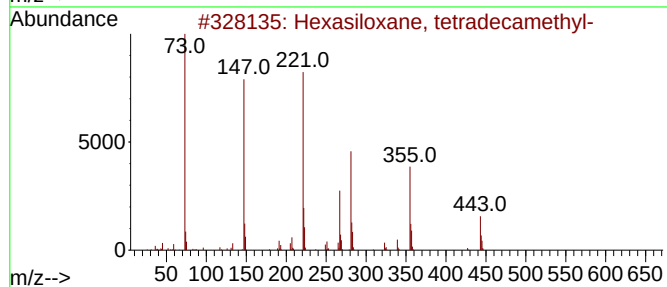
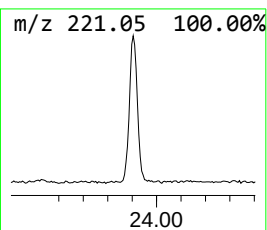
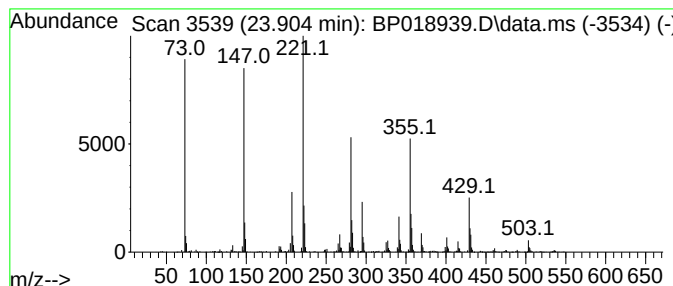
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-03 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.904	11.23 ng/ul	1655650	Perylene-d12	23.674

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	81
2			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	52
3			Heptasiloxane, hexadecamethyl-	532	C16H48O6Si7	000541-01-5	43
4			4-Hydroxybenzyl alcohol, 2TBDS ...	352	C19H36O2Si2	1000364-43-9	25
5			3-Oxobutan-2-yl trimethylsilyl p...	308	C15H20O5Si	1000373-49-5	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121823\
Data File : BP018939.D
Acq On : 19 Dec 2023 10:01
Operator : MA/JU
Sample : 05682-01 10X
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCLD6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propanoic acid,...	4.270	2.4	ng/ul	135684	1	7.828	1118590	20.0
Propanoic acid,...	4.452	2.2	ng/ul	124832	1	7.828	1118590	20.0
Isopropyl butyrate	4.911	5.1	ng/ul	287234	1	7.828	1118590	20.0
unknown-01	9.228	2.5	ng/ul	173327	2	10.622	1402950	20.0
2-Propanol, 1-(...	11.163	2.9	ng/ul	200096	2	10.622	1402950	20.0
1-Propanol, 2-(...	11.222	3.9	ng/ul	271673	2	10.622	1402950	20.0
Hexasiloxane, t...	21.845	3.3	ng/ul	533813	5	21.304	3221640	20.0
Cyclononasiloxa...	22.445	6.0	ng/ul	966884	5	21.304	3221640	20.0
unknown-02	23.122	8.9	ng/ul	1319000	6	23.674	2947370	20.0
unknown-03	23.904	11.2	ng/ul	1655650	6	23.674	2947370	20.0