

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
 Data File : BP010874.D
 Acq On : 27 Jun 2022 13:08
 Operator : CG/JU
 Sample : N3392-11 10X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXEY6

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-BP062322.M
 Title : SVOA CALIBRATION

Signal : TIC: BP010874.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.205	16	20	23	rBV2	74769	98797	0.80%	0.093%
2	3.240	23	26	39	rVB	1137100	1346716	10.85%	1.264%
3	3.546	74	78	83	rVB	30954	40268	0.32%	0.038%
4	3.616	86	90	95	rBV	550556	722662	5.82%	0.679%
5	3.664	95	98	103	rVB	167113	207541	1.67%	0.195%
6	3.834	123	127	133	rVB	59543	92880	0.75%	0.087%
7	3.916	134	141	148	rVB2	168942	268739	2.17%	0.252%
8	4.075	164	168	174	rVB2	27373	39271	0.32%	0.037%
9	4.193	183	188	195	rVB	201592	251237	2.02%	0.236%
10	4.258	195	199	204	rVV	45625	58123	0.47%	0.055%
11	4.369	213	218	224	rVV	523030	644572	5.19%	0.605%
12	4.434	224	229	235	rVV	239006	303332	2.44%	0.285%
13	4.481	235	237	242	rVB3	11800	13343	0.11%	0.013%
14	4.646	261	265	270	rBV4	7147	12892	0.10%	0.012%
15	4.722	273	278	281	rBV4	11462	15778	0.13%	0.015%
16	4.816	286	294	305	rVB	532975	745962	6.01%	0.700%
17	5.311	374	378	384	rVB3	14427	17866	0.14%	0.017%
18	5.434	395	399	404	rBV	9711	15190	0.12%	0.014%
19	5.675	435	440	446	rBV2	45640	66644	0.54%	0.063%
20	5.734	446	450	454	rVV8	7058	12540	0.10%	0.012%
21	6.887	637	646	654	rBV	488624	746420	6.01%	0.701%
22	7.052	667	674	690	rVB	1981044	2956551	23.82%	2.776%
23	7.246	699	707	717	rBV	1812252	2836895	22.86%	2.664%
24	7.340	719	723	732	rVB9	10781	26056	0.21%	0.024%
25	7.710	779	786	793	rBV	1467974	2245226	18.09%	2.108%
26	7.781	794	798	803	rVV2	21052	33099	0.27%	0.031%
27	8.075	841	848	855	rBV2	15079	29846	0.24%	0.028%
28	8.422	900	907	916	rBV	1386301	2139404	17.24%	2.009%
29	8.534	919	926	934	rVB	80519	138171	1.11%	0.130%
30	8.869	976	983	993	rBV	2067912	3311753	26.68%	3.109%
31	9.375	1064	1069	1071	rBV3	15824	23364	0.19%	0.022%
32	9.434	1076	1079	1088	rVB6	24720	43917	0.35%	0.041%
33	9.593	1097	1106	1114	rBV	1452940	2317710	18.67%	2.176%
34	9.910	1153	1160	1169	rVB	11133	28105	0.23%	0.026%
35	10.040	1178	1182	1189	rVB10	9574	13834	0.11%	0.013%
36	10.128	1189	1197	1205	rBV	2249643	3599160	29.00%	3.379%

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Integrator: RTE

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Filtering: 5

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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-BP062322.M
 Title : SVOA CALIBRATION

37	10.193	1205	1208	1214	rVB7	12032	20702	0.17%	0.019%
38	10.440	1245	1250	1254	rVB5	12704	21132	0.17%	0.020%
39	10.504	1254	1261	1266	rVV	1917066	3091796	24.91%	2.903%
40	10.557	1266	1270	1277	rVB	195187	320485	2.58%	0.301%
41	10.651	1277	1286	1296	rBV	2138955	3625689	29.21%	3.404%
42	11.057	1351	1355	1361	rBV	16030	26553	0.21%	0.025%
43	11.122	1361	1366	1370	rVV	19797	36203	0.29%	0.034%
44	11.163	1370	1373	1377	rVB5	9782	13066	0.11%	0.012%
45	12.098	1522	1532	1539	rBV	49739	90258	0.73%	0.085%
46	12.175	1539	1545	1551	rVB	117826	186008	1.50%	0.175%
47	12.393	1573	1582	1588	rVB	89983	148079	1.19%	0.139%
48	13.198	1713	1719	1724	rVB	16604	23874	0.19%	0.022%
49	13.275	1727	1732	1736	rBV2	24685	39163	0.32%	0.037%
50	13.528	1768	1775	1776	rBV7	12291	19763	0.16%	0.019%
51	13.687	1795	1802	1806	rBV3	23619	45533	0.37%	0.043%
52	13.739	1807	1811	1813	rVV2	12554	15411	0.12%	0.014%
53	13.787	1813	1819	1825	rVV	3983041	5397608	43.49%	5.068%
54	13.828	1825	1826	1832	rVB2	20499	24625	0.20%	0.023%
55	14.057	1854	1865	1882	rBV	4778048	6730194	54.22%	6.319%
56	14.281	1898	1903	1909	rVB3	15899	23531	0.19%	0.022%
57	14.369	1910	1918	1924	rVV	2362917	3530210	28.44%	3.315%
58	14.428	1924	1928	1936	rVB	38694	69466	0.56%	0.065%
59	14.575	1946	1953	1969	rBV	324730	524572	4.23%	0.493%
60	14.798	1989	1991	1996	rVB5	10537	14066	0.11%	0.013%
61	15.151	2044	2051	2056	rBV	83447	129699	1.04%	0.122%
62	15.363	2081	2087	2094	rBV	6060596	8497651	68.46%	7.979%
63	15.422	2094	2097	2101	rVV	40385	57399	0.46%	0.054%
64	15.481	2101	2107	2122	rVV	1245150	1794716	14.46%	1.685%
65	15.604	2124	2128	2136	rVB	74119	111072	0.89%	0.104%
66	16.081	2202	2209	2216	rBV	124385	185683	1.50%	0.174%
67	16.157	2218	2222	2227	rVB4	22769	35033	0.28%	0.033%
68	16.810	2329	2333	2338	rVB8	11861	19035	0.15%	0.018%
69	16.910	2346	2350	2354	rBV5	11801	20521	0.17%	0.019%
70	16.945	2354	2356	2361	rVB6	11032	13807	0.11%	0.013%
71	17.128	2381	2387	2393	rBV	2799067	3965638	31.95%	3.723%
72	17.169	2393	2394	2398	rVV	56136	58141	0.47%	0.055%
73	17.228	2398	2404	2420	rVV	6413296	9189991	74.04%	8.629%
74	17.816	2500	2504	2509	rVB6	31056	39770	0.32%	0.037%
75	19.057	2711	2715	2719	rBV	78533	106543	0.86%	0.100%
76	19.122	2722	2726	2732	rBV	78347	113385	0.91%	0.106%

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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

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77	19.480	2781	2787	2794	rBV	8439161	10800517	87.01%	10.141%
78	21.239	3082	3086	3092	rBV	3616445	4584261	36.93%	4.304%
79	23.457	3456	3463	3474	rVB2	6631857	12412387	100.00%	11.654%
80	23.610	3483	3489	3498	rVB2	2603823	4869067	39.23%	4.572%

Sum of corrected areas: 106506167

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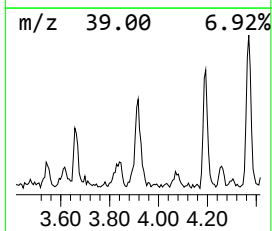
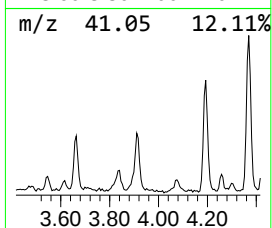
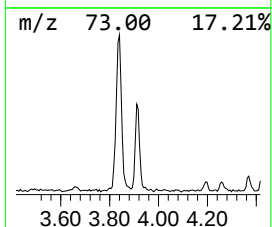
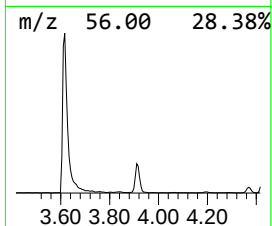
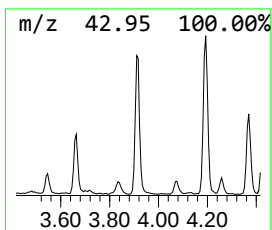
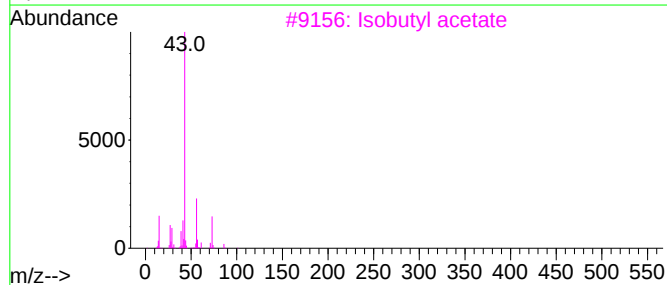
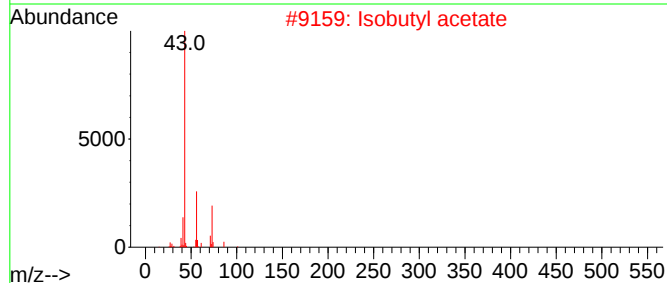
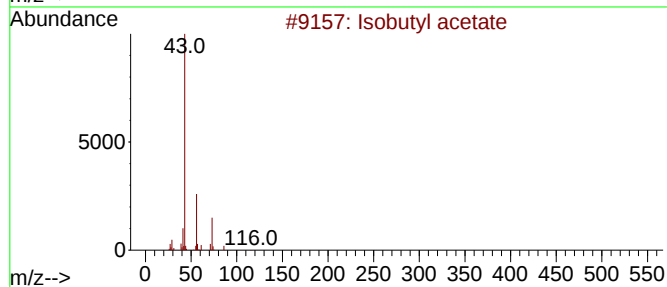
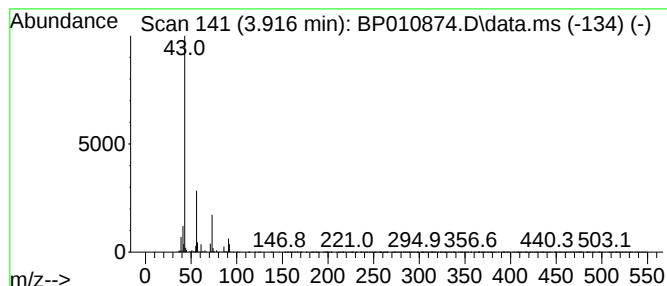
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Isobutyl acetate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.916	2.39 ng/ul	268739	1,4-Dichlorobenzene-d4	7.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isobutyl acetate	116	C6H12O2	000110-19-0	72
2			Isobutyl acetate	116	C6H12O2	000110-19-0	59
3			Isobutyl acetate	116	C6H12O2	000110-19-0	56
4			Isobutyl acetate	116	C6H12O2	000110-19-0	56
5			Isobutyl acetate	116	C6H12O2	000110-19-0	53



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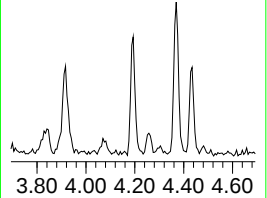
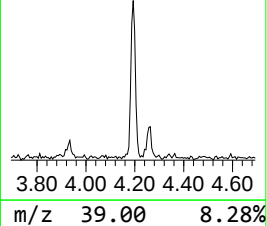
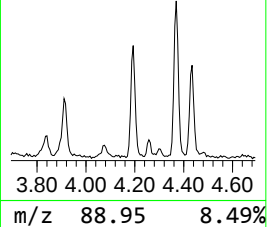
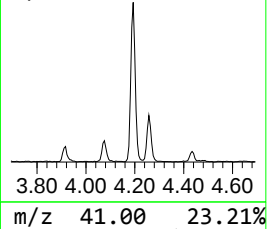
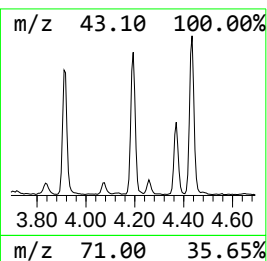
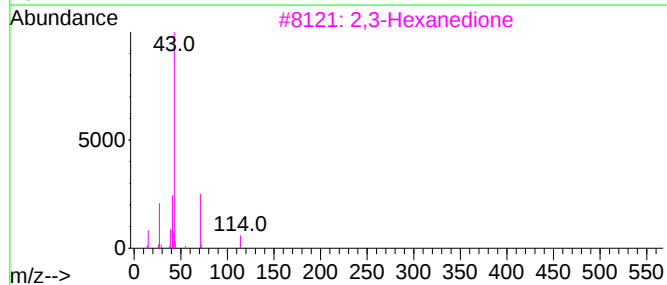
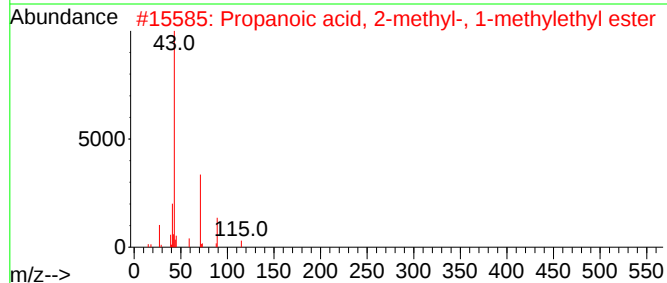
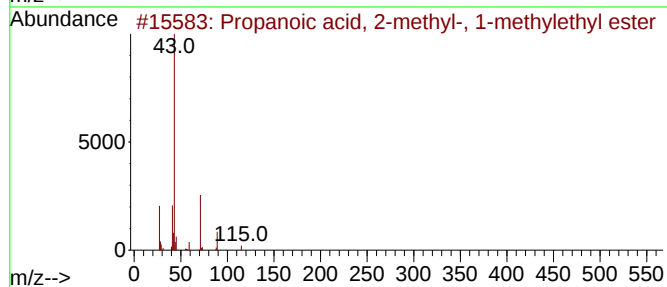
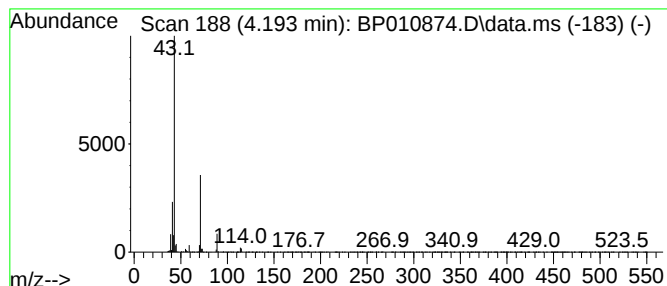
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Propanoic acid, 2-methyl-, ... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.193	2.24 ng/ul	251237	1,4-Dichlorobenzene-d4	7.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	64
2			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	64
3			2,3-Hexanedione	114	C6H10O2	003848-24-6	53
4			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	50
5			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	50



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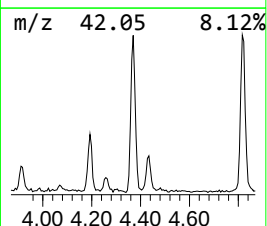
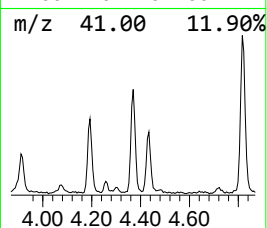
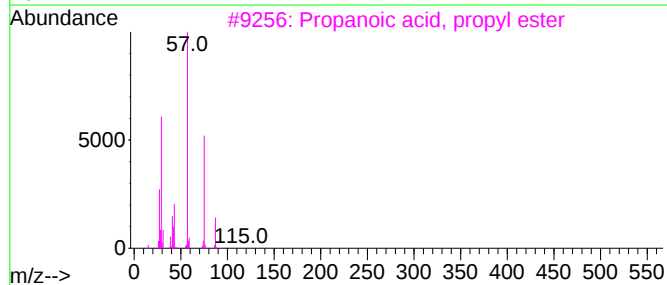
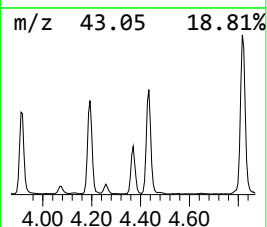
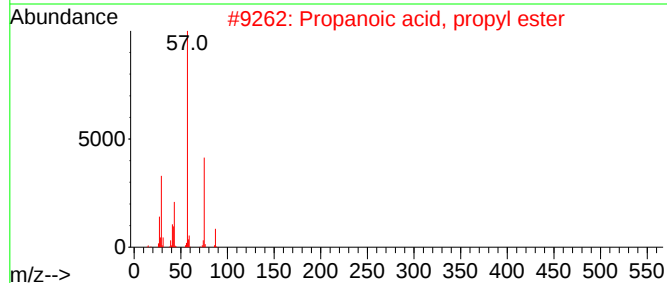
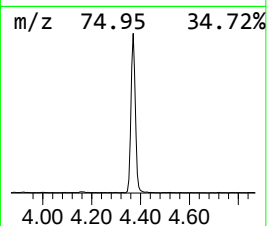
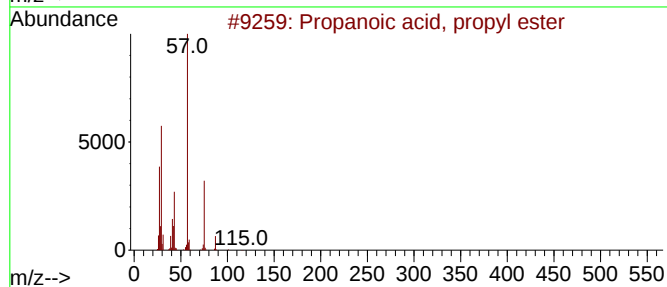
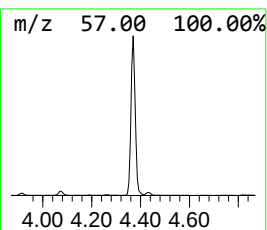
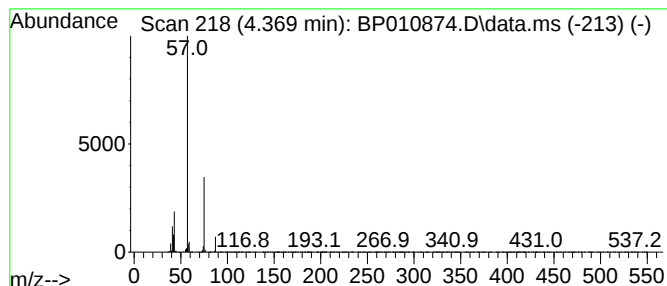
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Propanoic acid, propyl ester Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.369	5.74 ng/ul	644572	1,4-Dichlorobenzene-d4	7.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
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	78



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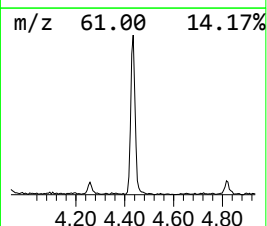
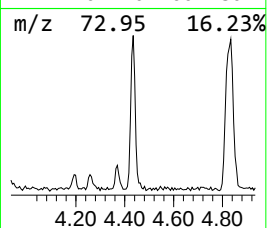
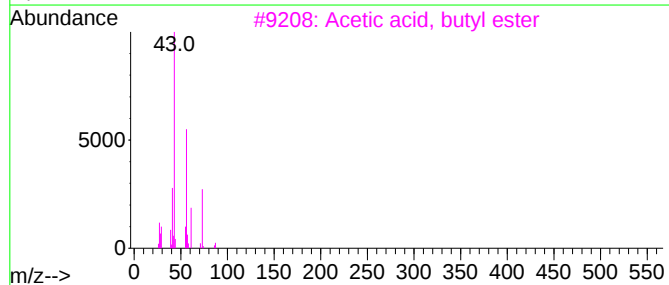
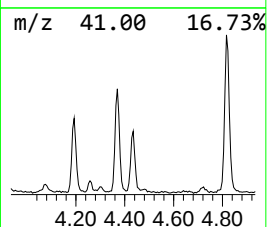
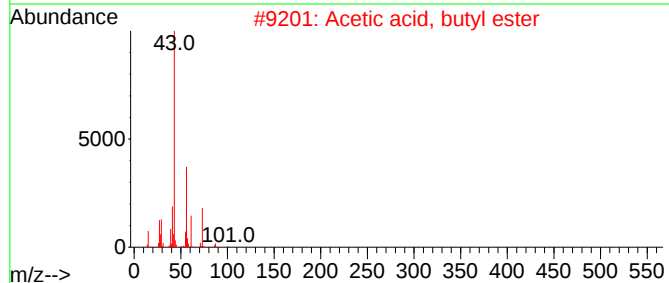
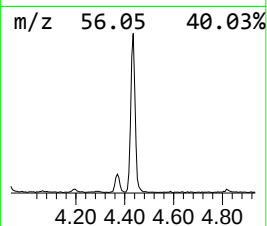
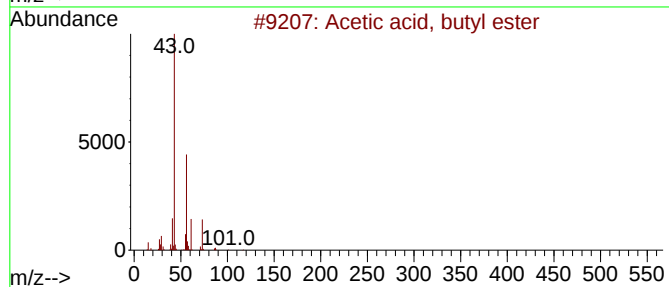
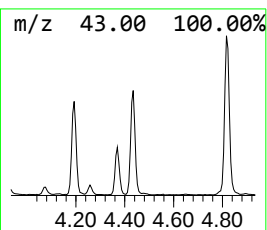
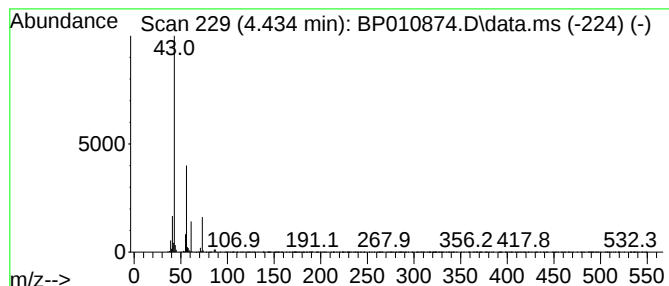
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Acetic acid, butyl ester Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.434	2.70 ng/ul	303332	1,4-Dichlorobenzene-d4	7.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetic acid, butyl ester	116	C6H12O2	000123-86-4	78
2			Acetic acid, butyl ester	116	C6H12O2	000123-86-4	72
3			Acetic acid, butyl ester	116	C6H12O2	000123-86-4	64
4			Acetic acid, butyl ester	116	C6H12O2	000123-86-4	56
5			Di-n-propyl ether	102	C6H14O	000111-43-3	16



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010874.D
Acq On : 27 Jun 2022 13:08
Operator : CG/JU
Sample : N3392-11 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXEY6

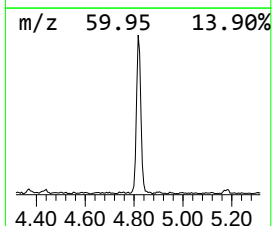
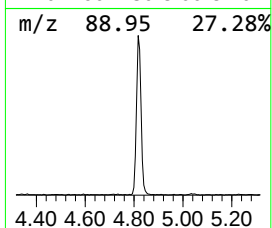
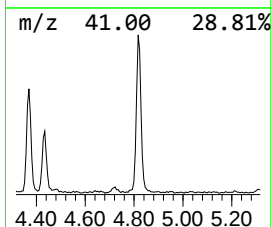
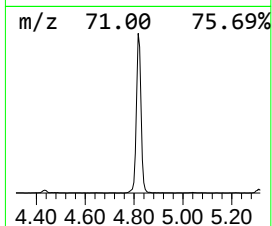
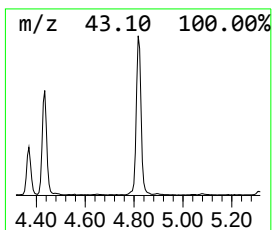
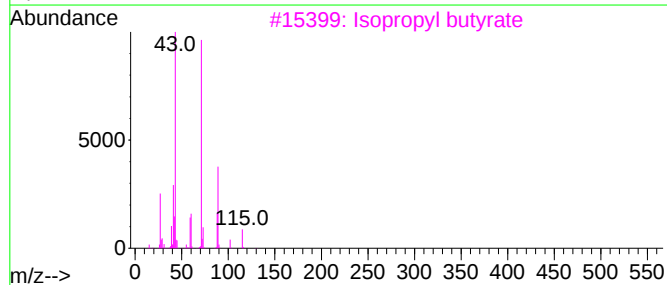
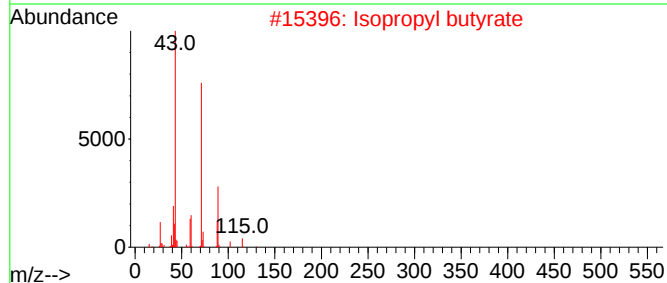
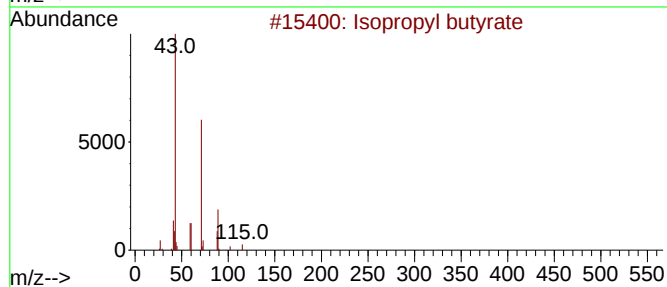
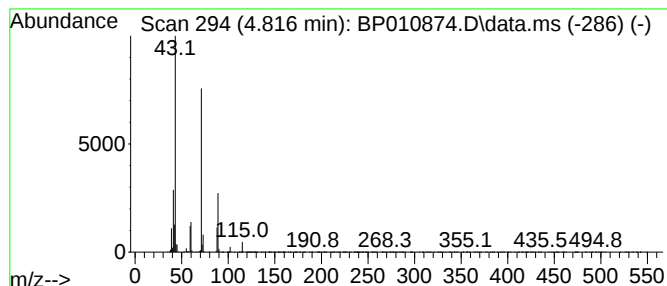
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Isopropyl butyrate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.816	6.64 ng/ul	745962	1,4-Dichlorobenzene-d4	7.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropyl butyrate	130	C7H14O2	000638-11-9	90
2			Isopropyl butyrate	130	C7H14O2	000638-11-9	90
3			Isopropyl butyrate	130	C7H14O2	000638-11-9	90
4			Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	80
5			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010874.D
Acq On : 27 Jun 2022 13:08
Operator : CG/JU
Sample : N3392-11 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXEY6

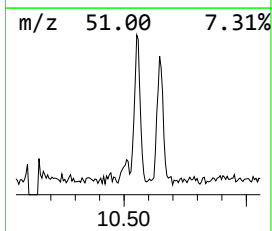
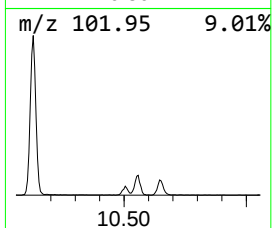
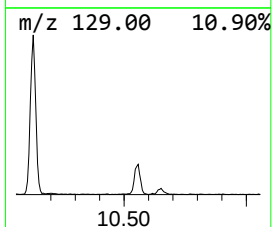
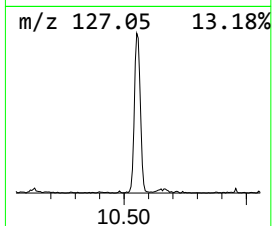
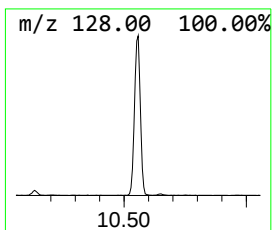
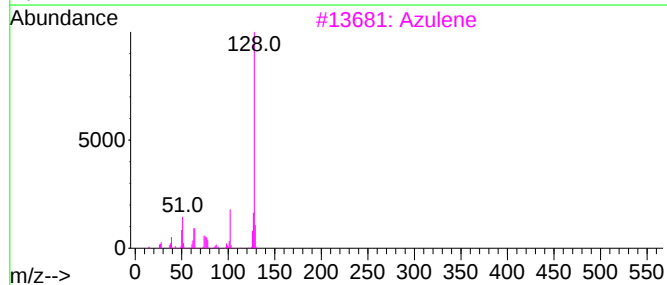
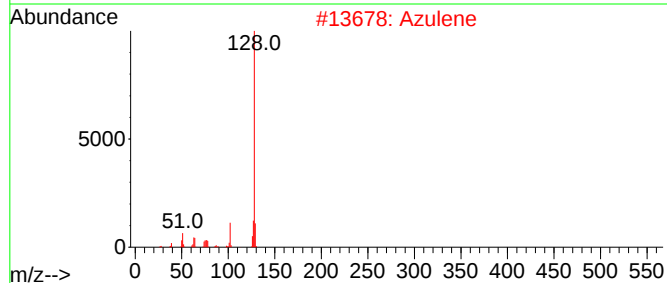
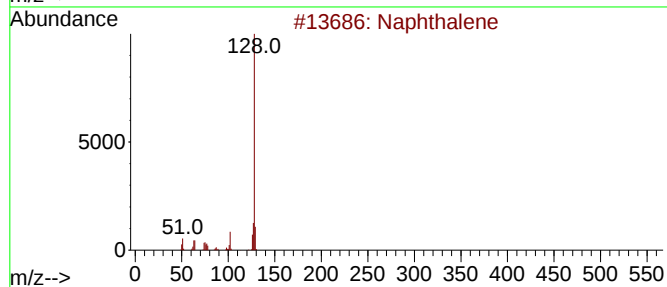
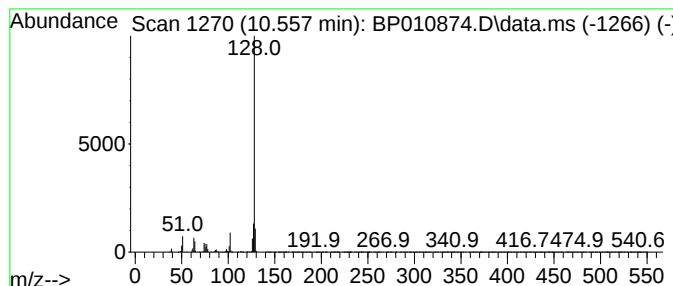
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Naphthalene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.557	2.07 ng/ul	320485	Naphthalene-d8	10.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	94
2			Azulene	128	C10H8	000275-51-4	94
3			Azulene	128	C10H8	000275-51-4	94
4			1H-Indene, 1-methylene-	128	C10H8	002471-84-3	91
5			Naphthalene	128	C10H8	000091-20-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010874.D
Acq On : 27 Jun 2022 13:08
Operator : CG/JU
Sample : N3392-11 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXEY6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Isobutyl acetate	3.916	2.4	ng/ul	268739	1	7.710	2245230	20.0
Propanoic acid,...	4.193	2.2	ng/ul	251237	1	7.710	2245230	20.0
Propanoic acid,...	4.369	5.7	ng/ul	644572	1	7.710	2245230	20.0
Acetic acid, bu...	4.434	2.7	ng/ul	303332	1	7.710	2245230	20.0
Isopropyl butyrate	4.816	6.6	ng/ul	745962	1	7.710	2245230	20.0
Naph thalene	10.557	2.1	ng/ul	320485	2	10.504	3091800	20.0