

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
 Data File : BP010873.D
 Acq On : 27 Jun 2022 12:34
 Operator : CG/JU
 Sample : N3392-07 10X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXEY2

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: BP010873.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.204	16	20	23	rBV2	58102	79700	0.89%	0.098%
2	3.240	23	26	37	rVB	987442	1099152	12.29%	1.351%
3	3.540	74	77	82	rVB	27830	34898	0.39%	0.043%
4	3.616	87	90	95	rVV	271341	366762	4.10%	0.451%
5	3.657	95	97	112	rVB	149441	211440	2.36%	0.260%
6	3.834	123	127	133	rVB	53793	82549	0.92%	0.101%
7	3.910	134	140	149	rVB2	143374	230648	2.58%	0.283%
8	4.069	164	167	177	rVB3	22065	32856	0.37%	0.040%
9	4.193	181	188	196	rBV	178916	229364	2.56%	0.282%
10	4.257	196	199	203	rBV	36156	41818	0.47%	0.051%
11	4.369	213	218	224	rBV	435242	529462	5.92%	0.651%
12	4.434	224	229	234	rVV	190039	236231	2.64%	0.290%
13	4.481	234	237	241	rVB3	7875	9579	0.11%	0.012%
14	4.646	260	265	269	rBV5	6000	9650	0.11%	0.012%
15	4.722	274	278	281	rBV5	9168	12924	0.14%	0.016%
16	4.816	287	294	303	rBV	456718	615604	6.88%	0.756%
17	5.310	374	378	382	rVB4	13541	18193	0.20%	0.022%
18	5.434	394	399	403	rBV	9901	14939	0.17%	0.018%
19	5.675	435	440	445	rBV	37508	50349	0.56%	0.062%
20	5.781	454	458	462	rBV7	6863	11720	0.13%	0.014%
21	6.392	557	562	569	rBV2	21917	36560	0.41%	0.045%
22	6.716	613	617	622	rVB8	5476	9480	0.11%	0.012%
23	6.887	636	646	654	rBV	378842	582722	6.51%	0.716%
24	7.051	667	674	682	rBV	1562252	2374077	26.54%	2.917%
25	7.128	684	687	693	rVB	14535	21367	0.24%	0.026%
26	7.245	700	707	719	rVB	1417400	2209519	24.70%	2.715%
27	7.345	719	724	728	rBV8	7679	17021	0.19%	0.021%
28	7.710	779	786	794	rBV	1318913	2014942	22.53%	2.476%
29	7.787	794	799	803	rBV2	20244	32731	0.37%	0.040%
30	8.422	899	907	918	rBV	1055597	1647628	18.42%	2.025%
31	8.534	918	926	932	rVB	69393	116947	1.31%	0.144%
32	8.810	969	973	976	rBV6	5324	9010	0.10%	0.011%
33	8.869	976	983	992	rVB	1634602	2631123	29.42%	3.233%
34	8.992	1001	1004	1011	rVB6	8520	11909	0.13%	0.015%
35	9.298	1051	1056	1062	rVB4	13422	22739	0.25%	0.028%
36	9.381	1064	1070	1071	rBV4	13977	21256	0.24%	0.026%

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Integrator: RTE

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
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37	9.434	1076	1079	1087	rVB6	19685	37586	0.42%	0.046%
38	9.592	1099	1106	1117	rBV	1084818	1777243	19.87%	2.184%
39	9.833	1143	1147	1153	rVB8	6067	9464	0.11%	0.012%
40	10.039	1175	1182	1187	rBV8	7493	18848	0.21%	0.023%
41	10.128	1190	1197	1205	rVV	1738860	2782704	31.11%	3.419%
42	10.192	1205	1208	1215	rVB9	12603	24639	0.28%	0.030%
43	10.433	1243	1249	1253	rVB6	10426	17637	0.20%	0.022%
44	10.504	1254	1261	1268	rBV	1668502	2788977	31.18%	3.427%
45	10.557	1268	1270	1274	rVB	17703	20336	0.23%	0.025%
46	10.651	1277	1286	1297	rBV	1613713	2709970	30.30%	3.330%
47	11.063	1349	1356	1360	rBV	11960	21408	0.24%	0.026%
48	11.122	1361	1366	1370	rVV2	14130	26296	0.29%	0.032%
49	11.163	1370	1373	1378	rVB5	8140	12068	0.13%	0.015%
50	11.310	1393	1398	1403	rBV10	6967	12919	0.14%	0.016%
51	12.098	1523	1532	1538	rBV2	37520	63848	0.71%	0.078%
52	12.175	1540	1545	1550	rVB3	11988	17340	0.19%	0.021%
53	12.392	1577	1582	1587	rBV5	8626	13860	0.15%	0.017%
54	13.086	1696	1700	1704	rBV6	5325	9724	0.11%	0.012%
55	13.198	1714	1719	1724	rVB7	10753	17625	0.20%	0.022%
56	13.274	1726	1732	1739	rBV4	18202	36801	0.41%	0.045%
57	13.692	1798	1803	1807	rBV7	5666	9803	0.11%	0.012%
58	13.786	1813	1819	1825	rBV	2939074	3952707	44.19%	4.857%
59	14.057	1857	1865	1882	rBV	3515601	5046870	56.42%	6.202%
60	14.369	1908	1918	1924	rBV	2117980	3102579	34.69%	3.812%
61	14.427	1925	1928	1938	rVB3	22555	42459	0.47%	0.052%
62	14.580	1947	1954	1963	rBV	209360	344113	3.85%	0.423%
63	14.692	1968	1973	1976	rVV5	8640	16888	0.19%	0.021%
64	14.727	1976	1979	1982	rVV5	8511	12076	0.14%	0.015%
65	14.769	1982	1986	1988	rVV	13943	20819	0.23%	0.026%
66	14.798	1989	1991	1996	rVB6	9249	11098	0.12%	0.014%
67	15.151	2046	2051	2057	rBV	50425	79009	0.88%	0.097%
68	15.363	2081	2087	2094	rBV	4327608	6234702	69.70%	7.661%
69	15.416	2094	2096	2101	rVV4	23045	32123	0.36%	0.039%
70	15.486	2101	2108	2120	rVV	806907	1168929	13.07%	1.436%
71	15.604	2124	2128	2134	rVB2	53071	80289	0.90%	0.099%
72	16.157	2218	2222	2232	rVB6	16378	31068	0.35%	0.038%
73	16.798	2326	2331	2338	rBV9	12628	27719	0.31%	0.034%
74	16.910	2346	2350	2354	rBV6	8481	12173	0.14%	0.015%
75	17.133	2380	2388	2393	rBV	2358824	3367031	37.64%	4.137%
76	17.233	2398	2405	2422	rVV	4640474	6705708	74.97%	8.240%

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77	17.815	2501	2504	2511	rVB9	15737	25477	0.28%	0.031%
78	19.057	2711	2715	2722	rBV3	58671	90415	1.01%	0.111%
79	19.121	2723	2726	2730	rVV	52475	67045	0.75%	0.082%
80	19.480	2781	2787	2794	rBV	6186834	7764498	86.80%	9.541%
81	19.533	2794	2796	2800	rVB	38657	50045	0.56%	0.061%
82	21.245	3082	3087	3092	rBV	3101262	3859094	43.14%	4.742%
83	23.456	3456	3463	3473	rVB2	4837982	8944800	100.00%	10.991%
84	23.609	3483	3489	3498	rVB2	2128565	4185174	46.79%	5.143%

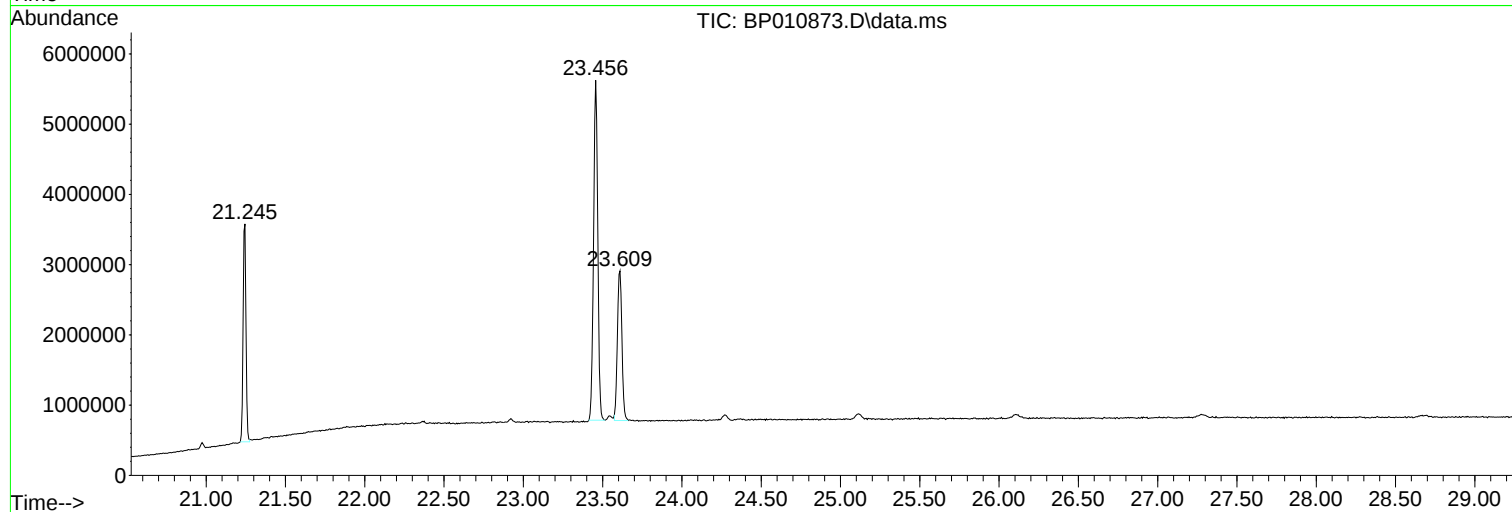
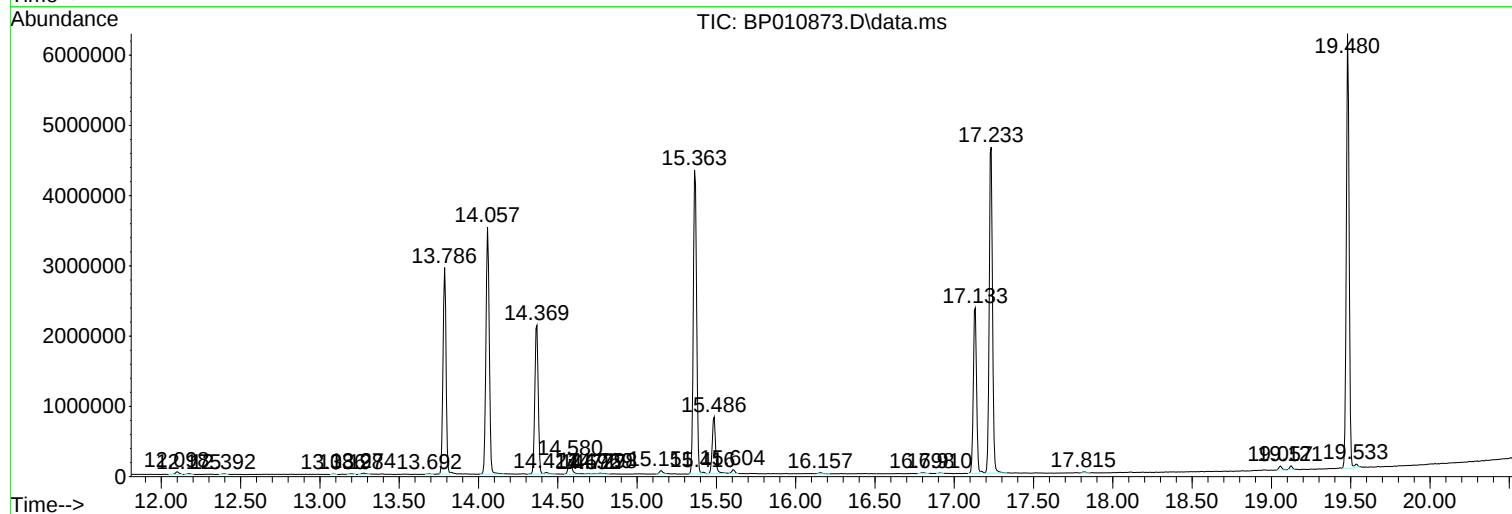
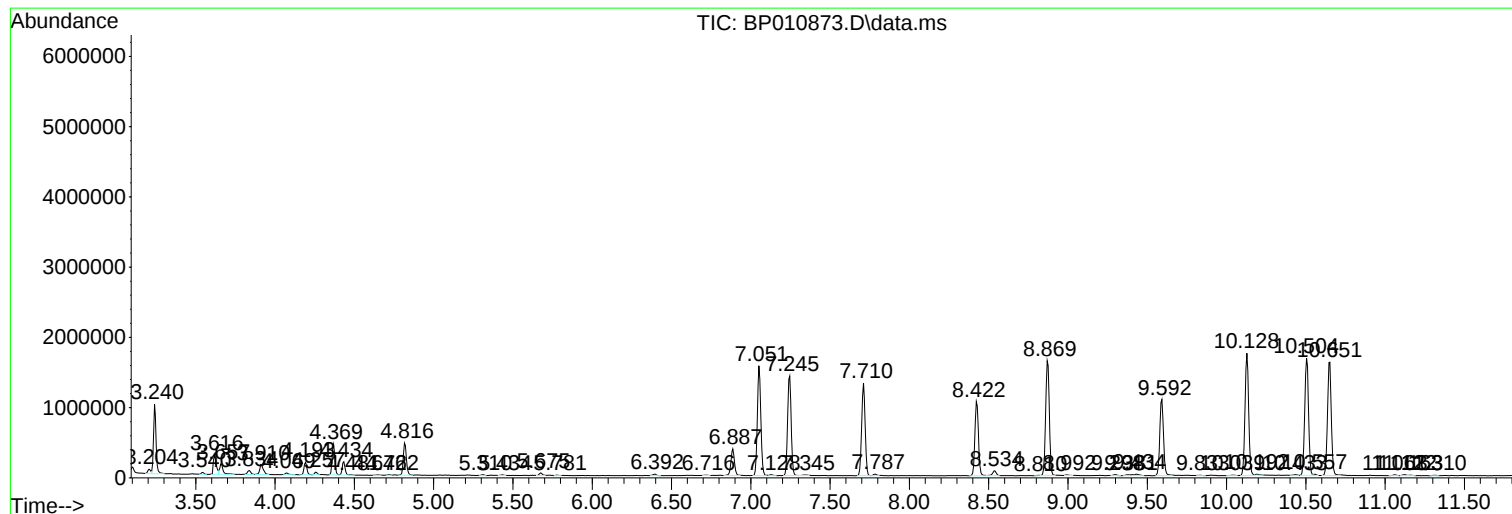
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Instrument :
BNA_P
ClientSampleId :
EXEY2

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



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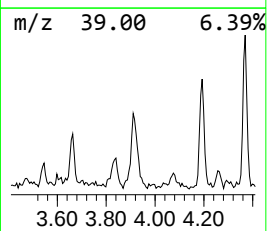
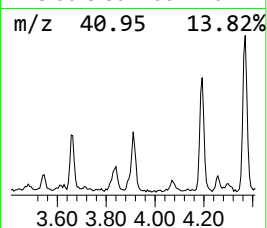
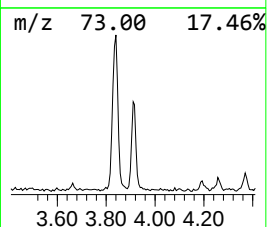
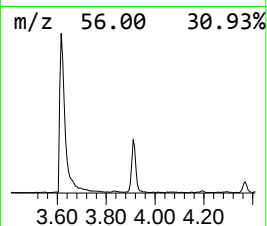
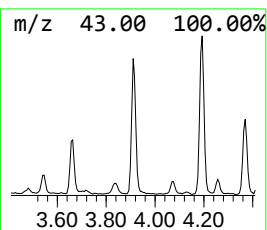
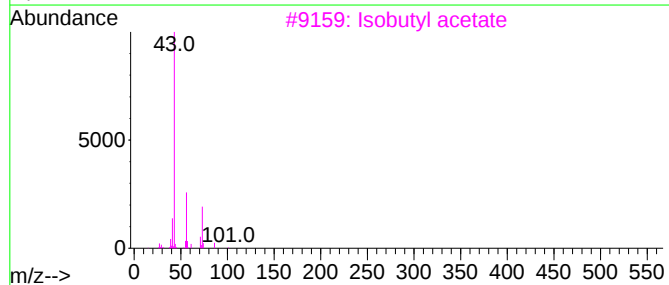
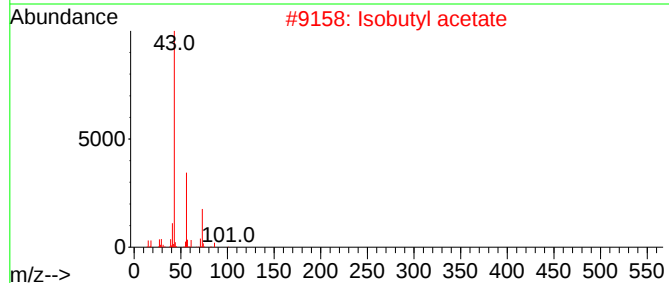
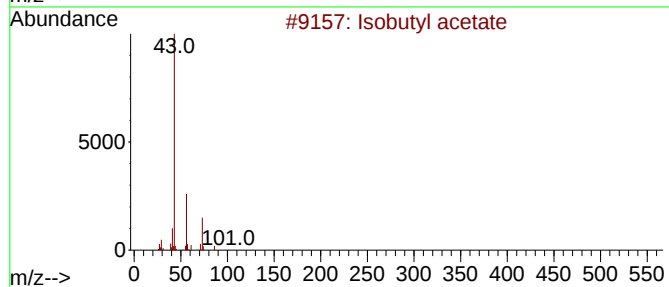
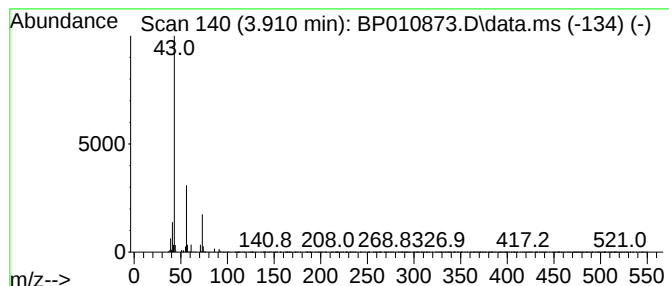
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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.
3.910	2.29 ng/ul	230648	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	56
3			Isobutyl acetate	116	C6H12O2	000110-19-0	45
4			Isobutyl acetate	116	C6H12O2	000110-19-0	43
5			Pentanoic acid, 4-oxo-	116	C5H8O3	000123-76-2	39



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BNA_P
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EXEY2

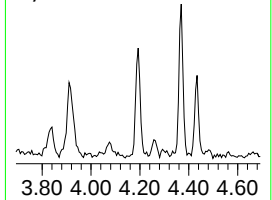
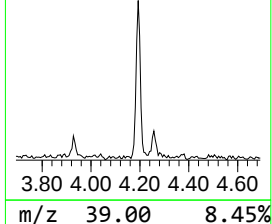
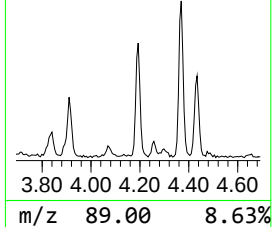
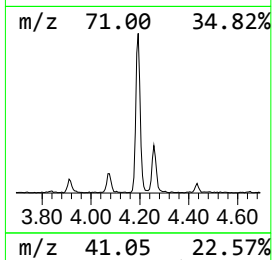
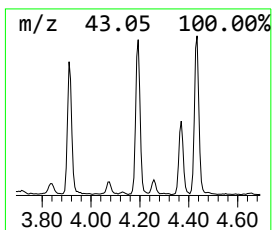
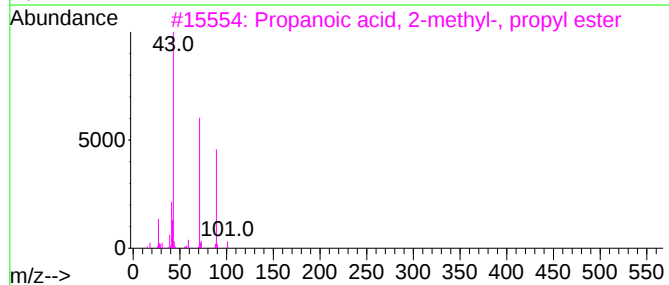
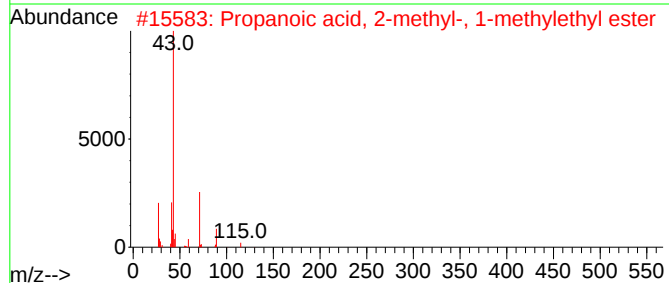
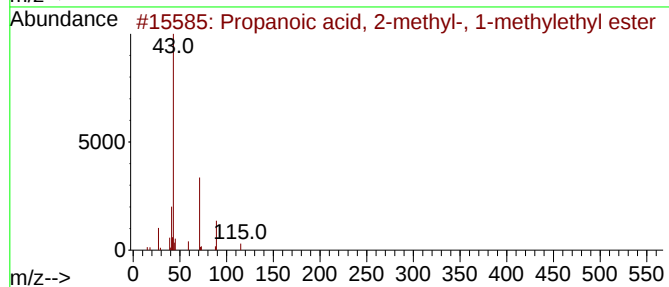
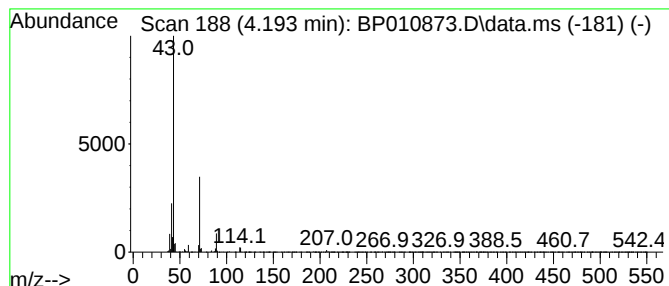
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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	2.28 ng/ul	229364	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	83
2			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	72
3			Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	72
4			Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	56
5			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	50



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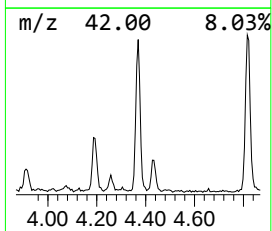
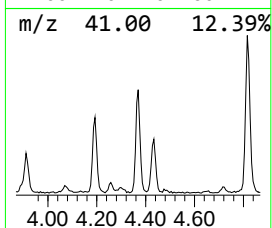
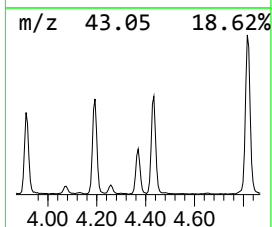
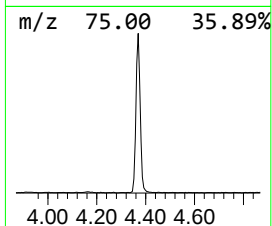
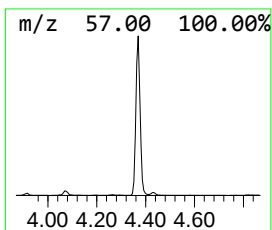
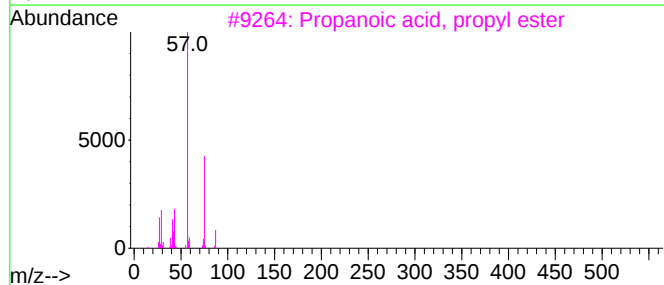
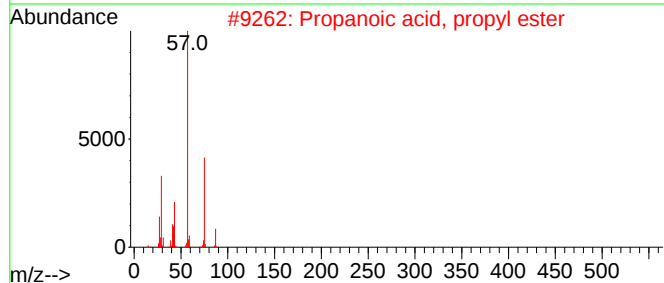
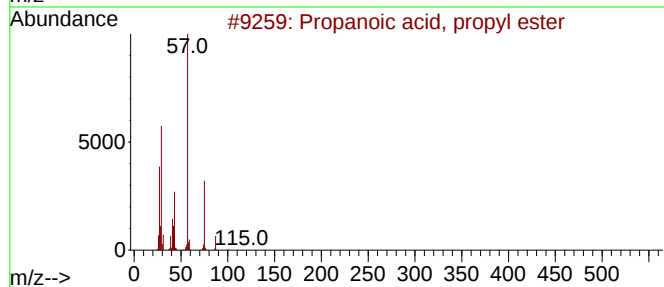
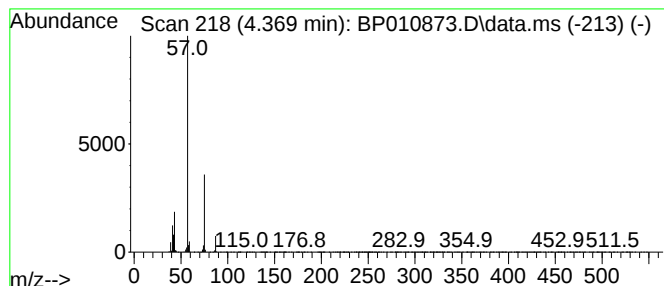
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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	5.26 ng/ul	529462	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	83
3			Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
4			Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	72
5			Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	72



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EXEY2

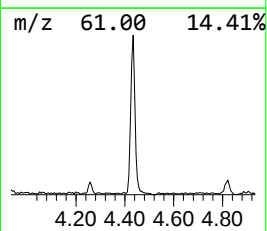
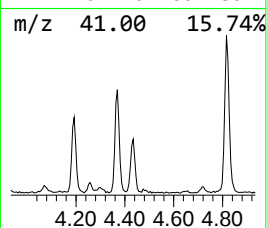
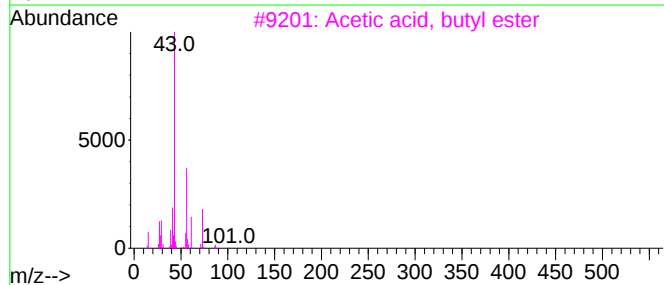
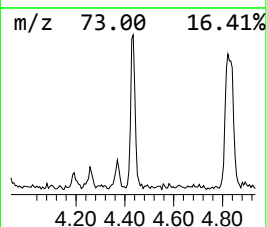
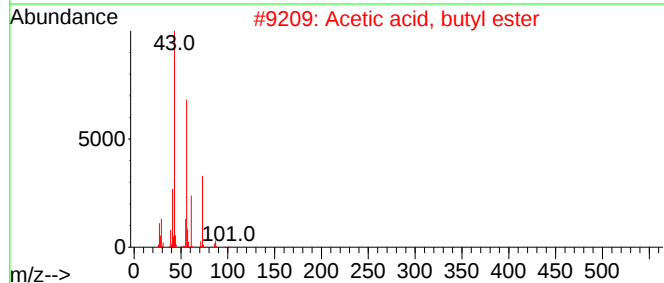
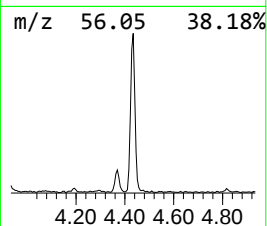
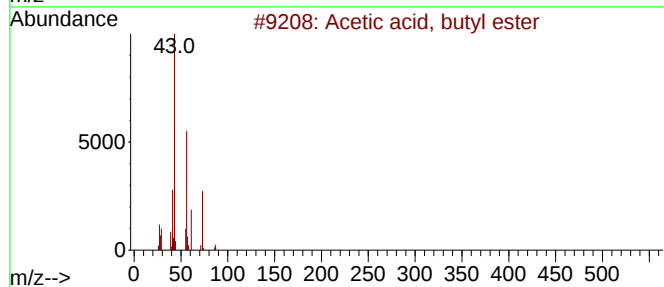
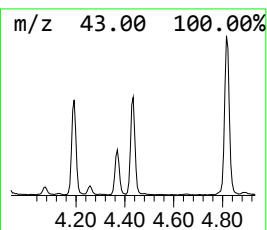
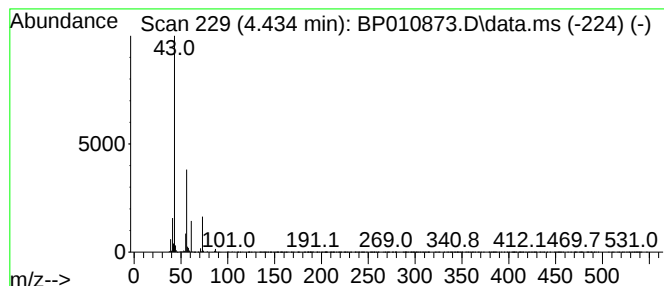
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Quant Title : SVOA CALIBRATION

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	2.34 ng/ul	236231	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	64
2			Acetic acid, butyl ester	116	C6H12O2	000123-86-4	50
3			Acetic acid, butyl ester	116	C6H12O2	000123-86-4	50
4			Acetic acid, butyl ester	116	C6H12O2	000123-86-4	32
5			Isobutyl acetate	116	C6H12O2	000110-19-0	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010873.D
Acq On : 27 Jun 2022 12:34
Operator : CG/JU
Sample : N3392-07 10X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXEY2

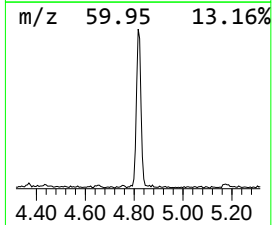
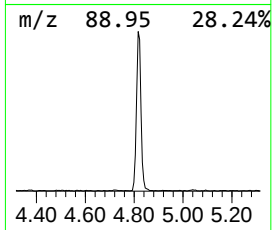
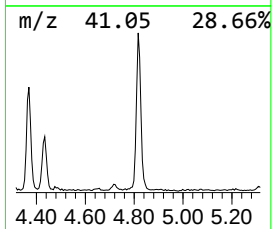
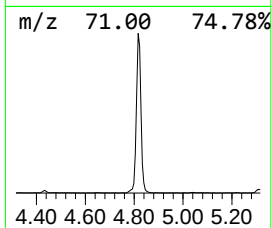
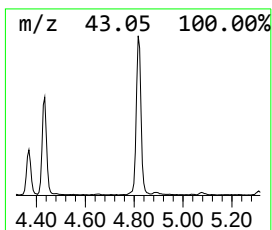
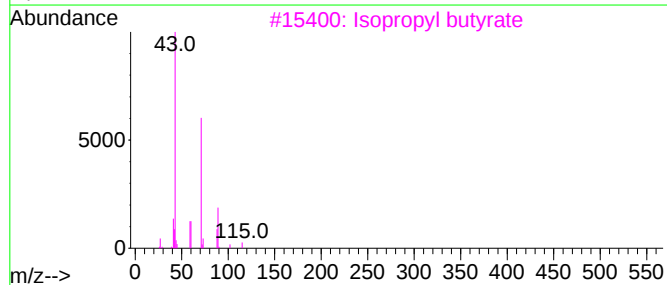
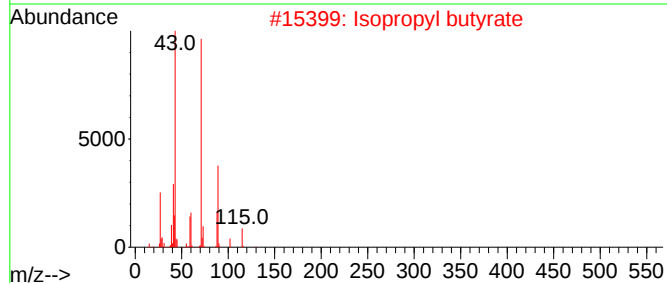
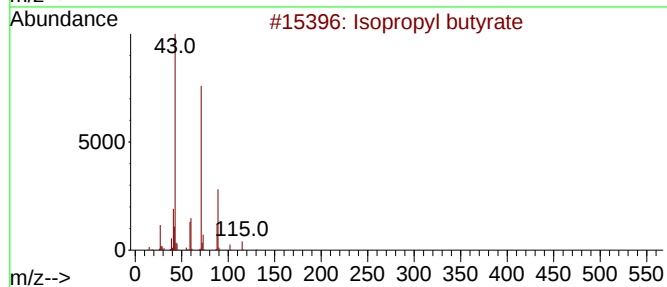
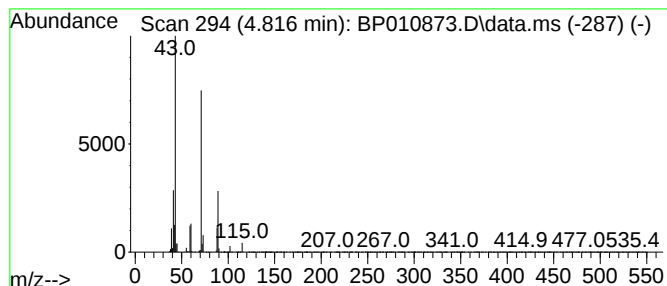
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Isopropyl butyrate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.816	6.11 ng/ul	615604	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	83
3			Isopropyl butyrate	130	C7H14O2	000638-11-9	83
4			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	78
5			Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010873.D
Acq On : 27 Jun 2022 12:34
Operator : CG/JU
Sample : N3392-07 10X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXEY2

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	3.910	2.3	ng/ul	230648	1	7.710	2014940	20.0
Propanoic acid,...	4.193	2.3	ng/ul	229364	1	7.710	2014940	20.0
Propanoic acid,...	4.369	5.3	ng/ul	529462	1	7.710	2014940	20.0
Acetic acid, bu...	4.434	2.3	ng/ul	236231	1	7.710	2014940	20.0
Isopropyl butyrate	4.816	6.1	ng/ul	615604	1	7.710	2014940	20.0