

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010845.D
 Acq On : 26 Jun 2022 01:11
 Operator : CG/JU
 Sample : N3392-15 10X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXEZ0

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: BP010845.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.211	17	21	23	rBV2	69110	92990	0.79%	0.081%
2	3.246	23	27	41	rVB	920026	1075892	9.16%	0.938%
3	3.546	75	78	82	rVB	30659	32769	0.28%	0.029%
4	3.617	87	90	95	rVV	636673	793508	6.76%	0.692%
5	3.664	95	98	106	rVB	191717	249360	2.12%	0.217%
6	3.840	122	128	134	rVV	67949	107086	0.91%	0.093%
7	3.917	134	141	152	rVB2	170248	299350	2.55%	0.261%
8	4.011	153	157	164	rVB2	29942	41233	0.35%	0.036%
9	4.075	164	168	175	rBV2	29379	42199	0.36%	0.037%
10	4.193	183	188	196	rBV	253419	334683	2.85%	0.292%
11	4.264	196	200	203	rVV	36986	48759	0.42%	0.042%
12	4.369	214	218	225	rVV	351096	461919	3.93%	0.403%
13	4.434	225	229	235	rVV	173195	226262	1.93%	0.197%
14	4.822	287	295	306	rVB	569990	773558	6.59%	0.674%
15	5.234	359	365	373	rVB2	53486	87655	0.75%	0.076%
16	5.311	373	378	383	rBV2	21087	31935	0.27%	0.028%
17	5.369	383	388	396	rVV	75467	156961	1.34%	0.137%
18	5.681	436	441	444	rBV	51286	68220	0.58%	0.059%
19	5.734	444	450	455	rVB2	50060	85990	0.73%	0.075%
20	6.211	526	531	538	rVB2	18487	29058	0.25%	0.025%
21	6.699	609	614	623	rVB	40560	67193	0.57%	0.059%
22	6.805	627	632	637	rVV	140335	208393	1.77%	0.182%
23	6.893	637	647	652	rVV2	578675	984637	8.38%	0.858%
24	6.940	652	655	663	rVB	104255	150701	1.28%	0.131%
25	7.058	667	675	681	rBV	2141282	3244805	27.63%	2.828%
26	7.105	681	683	691	rVB2	56917	79209	0.67%	0.069%
27	7.252	700	708	716	rBV	1994679	3131329	26.66%	2.729%
28	7.363	720	727	734	rVV	302190	469219	4.00%	0.409%
29	7.716	778	787	796	rBV	1720776	2697996	22.97%	2.352%
30	7.787	796	799	802	rVV2	25903	40904	0.35%	0.036%
31	7.828	802	806	815	rVV	80683	130052	1.11%	0.113%
32	8.087	842	850	857	rVB	119702	187790	1.60%	0.164%
33	8.210	866	871	874	rBV3	20265	32709	0.28%	0.029%
34	8.246	874	877	885	rVB2	44872	75420	0.64%	0.066%
35	8.357	890	896	900	rBV3	28863	46400	0.40%	0.040%
36	8.434	901	909	919	rVV	1551069	2432672	20.71%	2.120%

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

37	8.540	919	927	933	rVB	86410	145729	1.24%	0.127%
38	8.822	964	975	978	rBV	45481	78747	0.67%	0.069%
39	8.875	978	984	1002	rVB	2311924	3696013	31.47%	3.221%
40	9.063	1012	1016	1024	rVB	12993	29559	0.25%	0.026%
41	9.157	1024	1032	1036	rBV	8931	25540	0.22%	0.022%
42	9.304	1052	1057	1060	rBV3	17760	25834	0.22%	0.023%
43	9.410	1066	1075	1076	rVV8	20388	51322	0.44%	0.045%
44	9.434	1076	1079	1086	rVB7	29709	53655	0.46%	0.047%
45	9.599	1098	1107	1121	rBV	1571116	2637518	22.46%	2.299%
46	9.910	1155	1160	1169	rVB2	45423	87398	0.74%	0.076%
47	10.134	1188	1198	1206	rBV	2512983	3992771	34.00%	3.480%
48	10.210	1207	1211	1216	rVB4	13343	26605	0.23%	0.023%
49	10.440	1244	1250	1254	rVV6	17047	40272	0.34%	0.035%
50	10.510	1254	1262	1268	rVV	2292485	3770563	32.10%	3.286%
51	10.557	1268	1270	1277	rVV	82604	124929	1.06%	0.109%
52	10.657	1277	1287	1296	rVV	2276610	3919637	33.37%	3.416%
53	10.734	1296	1300	1307	rVB5	21799	39385	0.34%	0.034%
54	10.981	1336	1342	1349	rVB	38549	65633	0.56%	0.057%
55	11.128	1364	1367	1371	rVB	18543	27501	0.23%	0.024%
56	12.069	1521	1527	1531	rBV	119852	199569	1.70%	0.174%
57	12.104	1531	1533	1537	rVB	48359	59422	0.51%	0.052%
58	12.404	1578	1584	1587	rBV	49596	83688	0.71%	0.073%
59	12.440	1587	1590	1595	rVB5	18102	26018	0.22%	0.023%
60	12.645	1619	1625	1630	rVB6	20183	40164	0.34%	0.035%
61	12.916	1666	1671	1676	rVV5	27668	42903	0.37%	0.037%
62	13.204	1717	1720	1727	rVV6	16301	33173	0.28%	0.029%
63	13.310	1729	1738	1744	rVV5	28446	82846	0.71%	0.072%
64	13.363	1744	1747	1751	rVV3	23657	32573	0.28%	0.028%
65	13.551	1769	1779	1783	rBV7	17482	49094	0.42%	0.043%
66	13.692	1792	1803	1807	rBV4	34449	89927	0.77%	0.078%
67	13.740	1807	1811	1814	rBV2	31070	46776	0.40%	0.041%
68	13.792	1814	1820	1825	rBV	4436941	5836282	49.69%	5.087%
69	13.834	1825	1827	1831	rVB	20847	26246	0.22%	0.023%
70	13.922	1838	1842	1845	rBV	26096	38218	0.33%	0.033%
71	13.957	1845	1848	1854	rVB4	33243	47275	0.40%	0.041%
72	14.063	1854	1866	1881	rBV	5108576	7338863	62.48%	6.397%
73	14.369	1912	1918	1926	rVB	2932128	4271074	36.37%	3.723%
74	14.587	1949	1955	1965	rBV	360609	582148	4.96%	0.507%
75	14.739	1975	1981	1983	rBV5	18131	31697	0.27%	0.028%
76	14.845	1996	1999	2004	rVB5	16297	26064	0.22%	0.023%

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 Title : SVOA CALIBRATION

77	14.992	2020	2024	2029	rBV2	21890	41753	0.36%	0.036%
78	15.039	2029	2032	2037	rVB6	18262	28532	0.24%	0.025%
79	15.104	2037	2043	2047	rBV9	11299	24708	0.21%	0.022%
80	15.157	2047	2052	2057	rBV	94429	148538	1.26%	0.129%
81	15.369	2081	2088	2094	rBV	6400297	8919436	75.94%	7.774%
82	15.422	2094	2097	2101	rVB4	29467	40412	0.34%	0.035%
83	15.486	2101	2108	2116	rBV	1131016	1534221	13.06%	1.337%
84	15.545	2116	2118	2124	rVV6	20693	31137	0.27%	0.027%
85	15.610	2124	2129	2134	rVV2	67599	93307	0.79%	0.081%
86	15.857	2168	2171	2179	rVB5	20286	34057	0.29%	0.030%
87	15.939	2179	2185	2193	rBV5	15954	35830	0.31%	0.031%
88	16.157	2220	2222	2228	rVB4	26854	39807	0.34%	0.035%
89	16.481	2274	2277	2283	rVB4	16280	24071	0.20%	0.021%
90	16.945	2354	2356	2363	rVB5	13764	24302	0.21%	0.021%
91	17.133	2381	2388	2393	rBV	3321761	4558492	38.81%	3.973%
92	17.175	2393	2395	2399	rVV	45019	56056	0.48%	0.049%
93	17.233	2399	2405	2425	rVB2	6709199	9328951	79.43%	8.131%
94	17.816	2501	2504	2507	rBV4	23441	26479	0.23%	0.023%
95	19.057	2711	2715	2721	rVB	85995	114771	0.98%	0.100%
96	19.127	2723	2727	2731	rBV	78611	100575	0.86%	0.088%
97	19.480	2782	2787	2795	rBV	7918872	10580246	90.08%	9.222%
98	21.245	3082	3087	3092	rBV	4132594	5062072	43.10%	4.412%
99	23.457	3456	3463	3473	rVB2	6132729	11745002	100.00%	10.237%
100	23.610	3483	3489	3498	rVB2	2742685	5295655	45.09%	4.616%

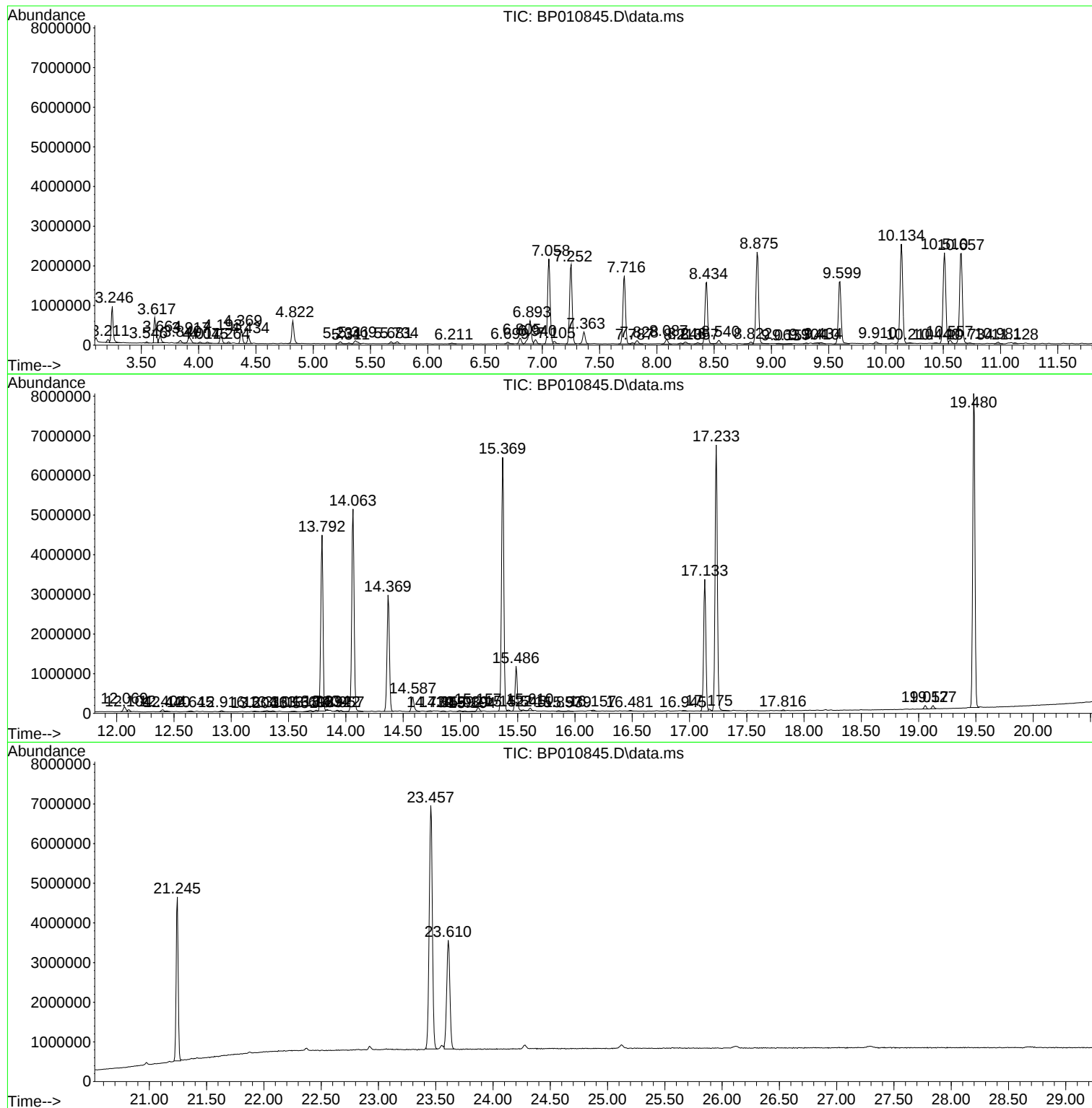
Sum of corrected areas: 114729857

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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



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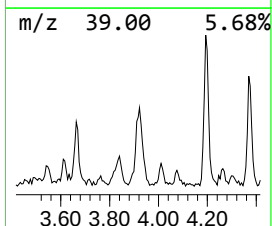
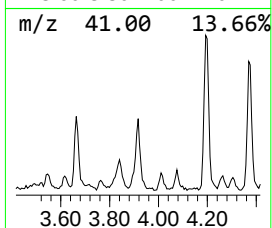
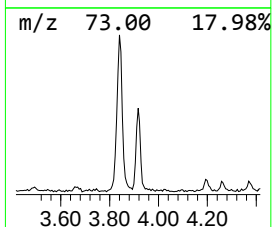
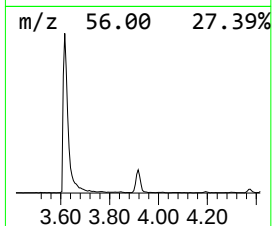
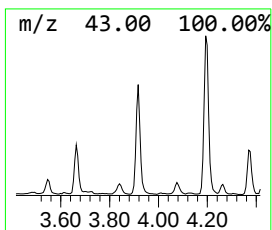
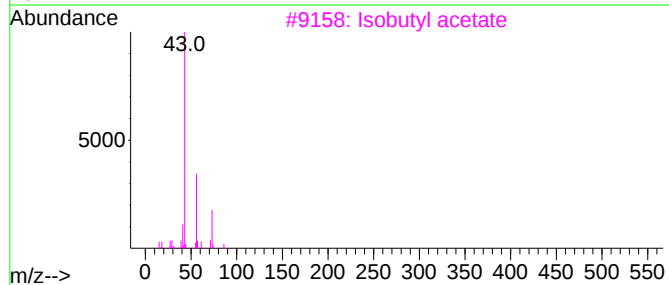
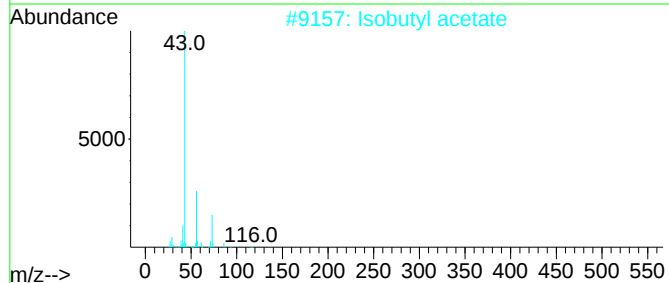
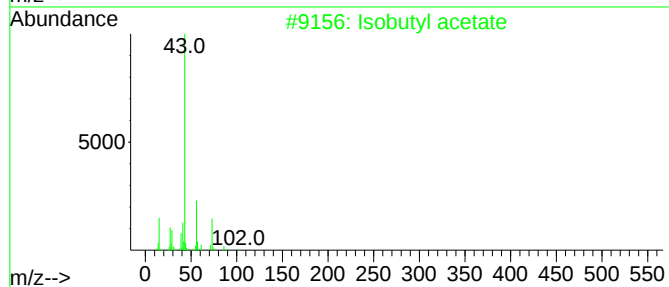
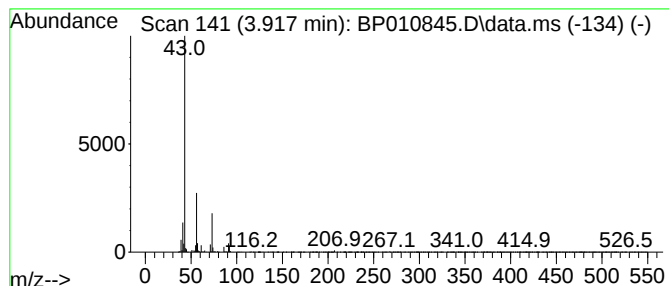
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Isobutyl acetate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.917	2.22 ng/ul	299350	1,4-Dichlorobenzene-d4	7.716

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isobutyl acetate	116	C6H12O2	000110-19-0	78
2			Isobutyl acetate	116	C6H12O2	000110-19-0	72
3			Isobutyl acetate	116	C6H12O2	000110-19-0	72
4			Isobutyl acetate	116	C6H12O2	000110-19-0	64
5			Isobutyl acetate	116	C6H12O2	000110-19-0	59



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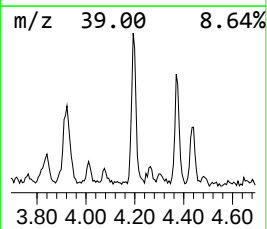
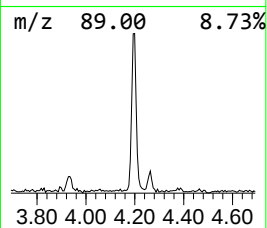
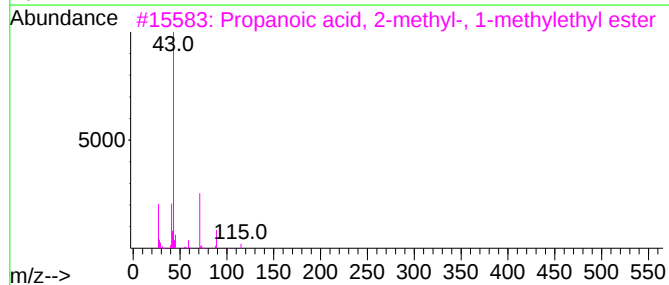
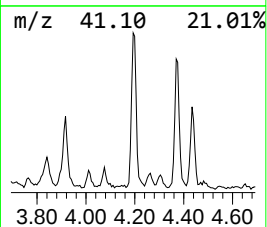
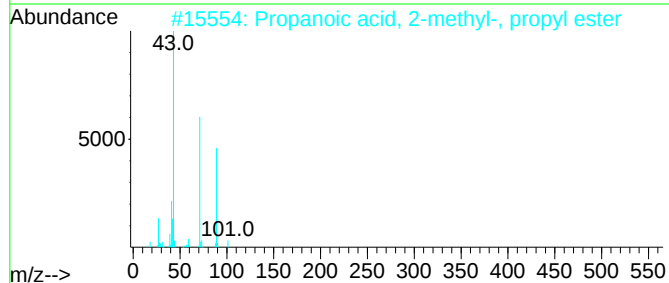
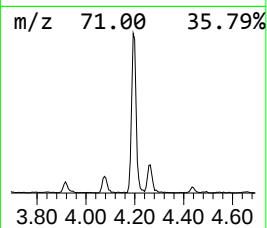
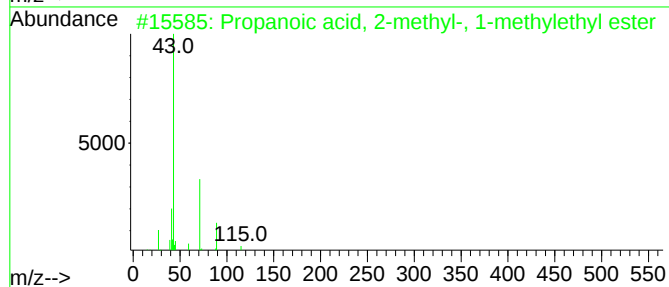
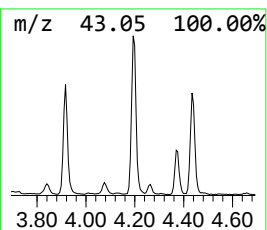
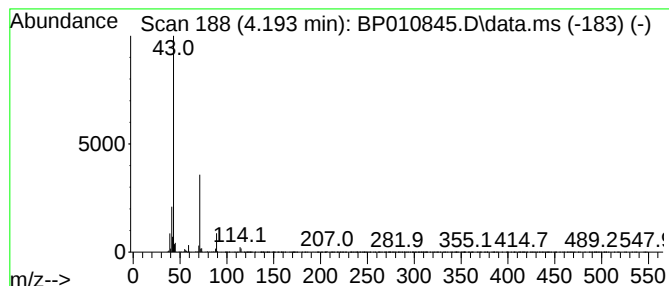
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Propanoic acid, 2-methyl-, ... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.193	2.48 ng/ul	334683	1,4-Dichlorobenzene-d4	7.716

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-, propy...	130	C7H14O2	000644-49-5	72
3			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	64
4			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	64
5			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	59



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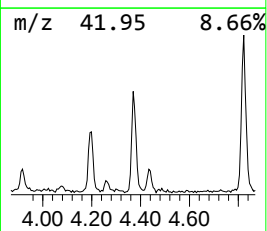
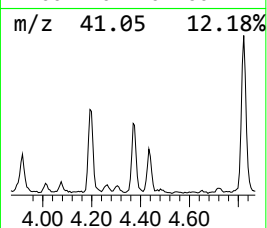
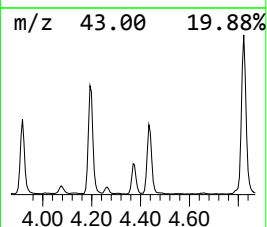
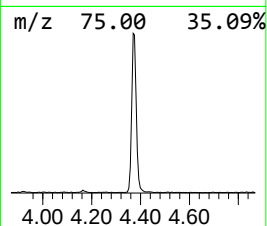
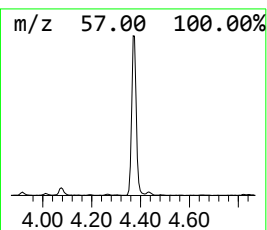
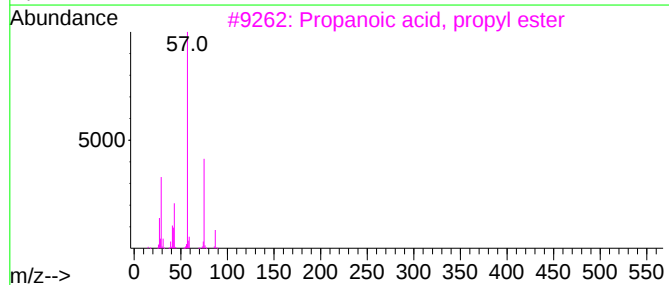
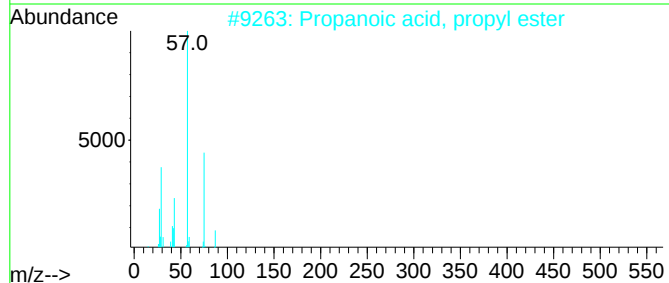
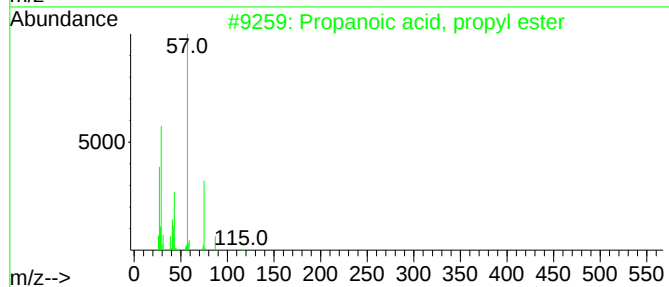
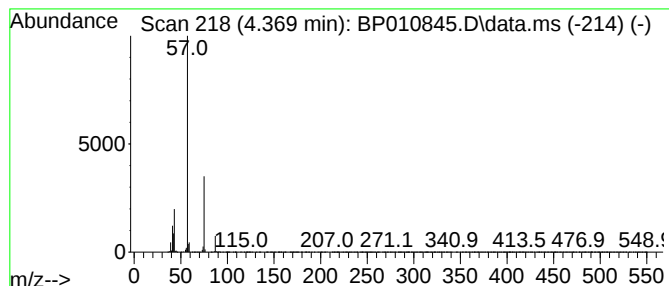
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Propanoic acid, propyl ester Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.369	3.42 ng/ul	461919	1,4-Dichlorobenzene-d4	7.716

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	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
Data File : BP010845.D
Acq On : 26 Jun 2022 01:11
Operator : CG/JU
Sample : N3392-15 10X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXEZ0

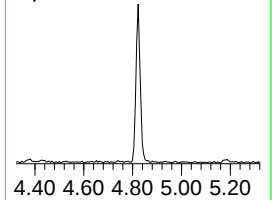
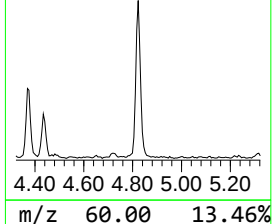
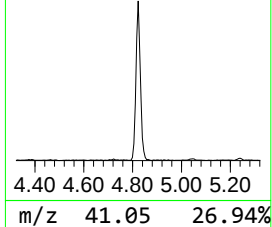
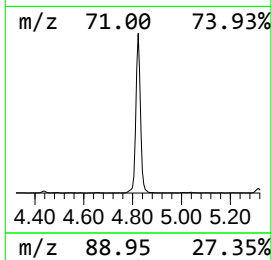
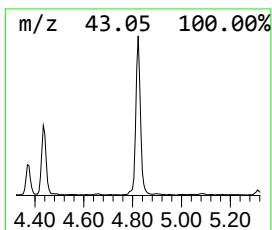
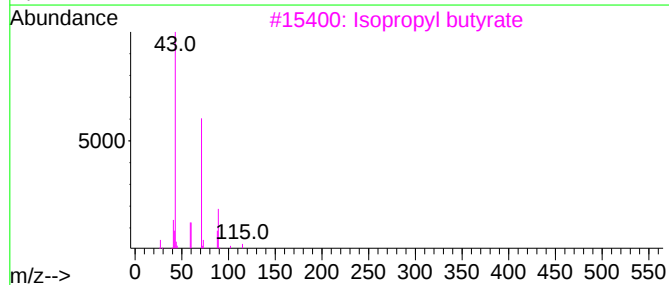
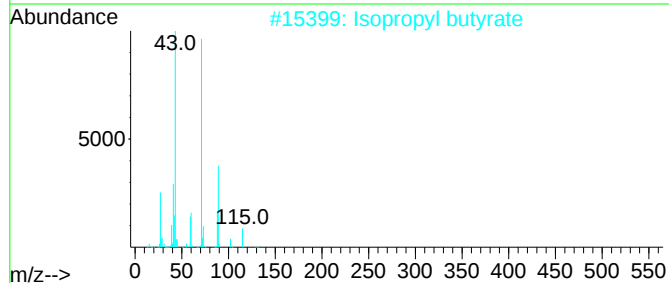
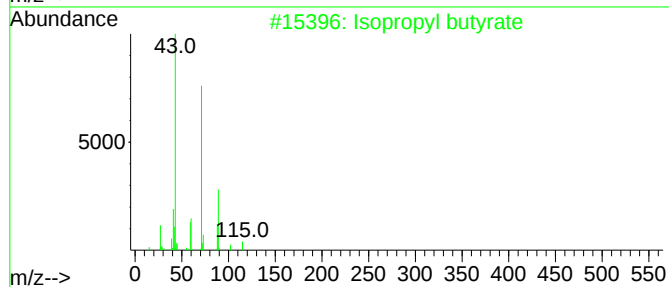
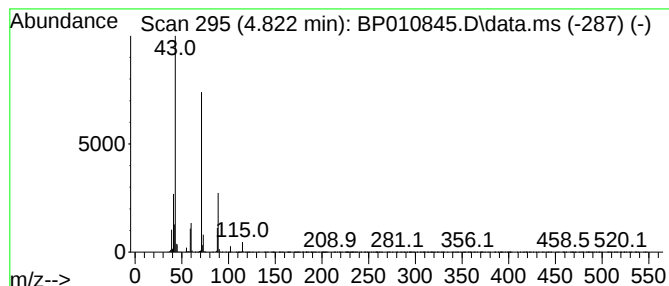
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Isopropyl butyrate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.822	5.73 ng/ul	773558	1,4-Dichlorobenzene-d4	7.716

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	83
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\BP062522\
Data File : BP010845.D
Acq On : 26 Jun 2022 01:11
Operator : CG/JU
Sample : N3392-15 10X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXEZ0

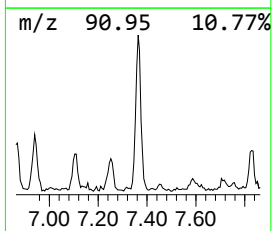
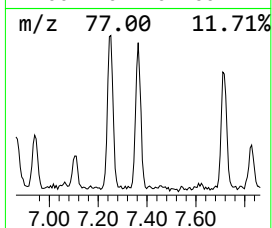
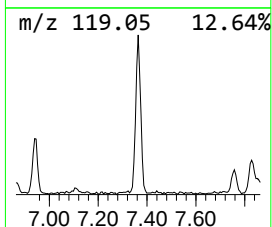
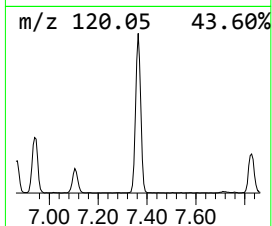
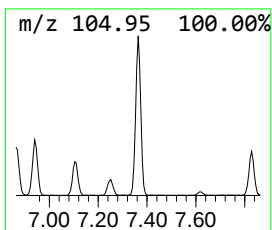
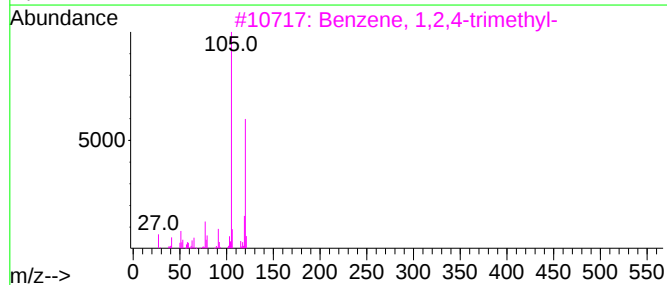
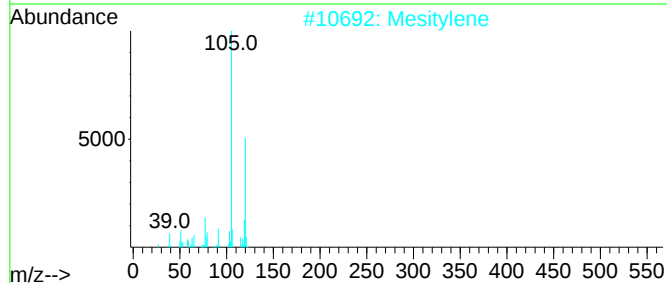
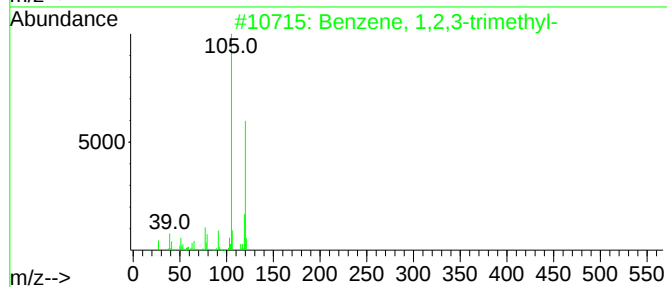
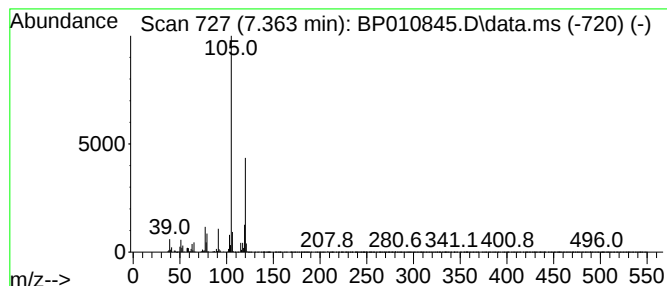
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.363	3.48 ng/ul	469219	1,4-Dichlorobenzene-d4	7.716

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2			Mesitylene	120	C9H12	000108-67-8	97
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
Data File : BP010845.D
Acq On : 26 Jun 2022 01:11
Operator : CG/JU
Sample : N3392-15 10X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXEZ0

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	3.917	2.2	ng/ul	299350	1	7.716	2698000	20.0
Propanoic acid,...	4.193	2.5	ng/ul	334683	1	7.716	2698000	20.0
Propanoic acid,...	4.369	3.4	ng/ul	461919	1	7.716	2698000	20.0
Isopropyl butyrate	4.822	5.7	ng/ul	773558	1	7.716	2698000	20.0
Benzene, 1,2,3-...	7.363	3.5	ng/ul	469219	1	7.716	2698000	20.0