

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011224\
 Data File : BP019328.D
 Acq On : 12 Jan 2024 14:07
 Operator : MA/JU
 Sample : P1016-06 10X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCMH6

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-BP010824.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP019328.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.223	19	23	25	rBV3	15744	19403	0.33%	0.042%
2	3.258	25	29	40	rVB	162310	210319	3.62%	0.459%
3	3.564	78	81	87	rVB	17954	22581	0.39%	0.049%
4	3.640	91	94	100	rBV	146936	228475	3.93%	0.499%
5	3.864	128	132	137	rVB5	6019	9796	0.17%	0.021%
6	3.946	142	146	153	rBV2	29399	49708	0.85%	0.109%
7	4.064	163	166	170	rBV6	5884	8370	0.14%	0.018%
8	4.105	170	173	178	rVV6	5133	8141	0.14%	0.018%
9	4.222	189	193	200	rVB	30630	43452	0.75%	0.095%
10	4.293	202	205	210	rVB5	9002	12374	0.21%	0.027%
11	4.411	221	225	232	rVB	45794	63883	1.10%	0.139%
12	4.475	232	236	242	rVB	37235	49189	0.85%	0.107%
13	4.864	297	302	307	rBV	69869	97457	1.68%	0.213%
14	4.899	307	308	313	rVB5	8219	9820	0.17%	0.021%
15	5.458	398	403	410	rBV2	25491	40428	0.70%	0.088%
16	5.728	443	449	454	rBV4	7178	12236	0.21%	0.027%
17	5.987	487	493	504	rVB8	4635	13558	0.23%	0.030%
18	6.440	566	570	574	rBV6	4647	6535	0.11%	0.014%
19	6.734	615	620	626	rVV8	3495	9072	0.16%	0.020%
20	6.969	653	660	670	rVB	144409	238043	4.09%	0.520%
21	7.122	678	686	698	rVB	505329	784048	13.48%	1.711%
22	7.316	709	719	728	rBV	465517	768333	13.21%	1.677%
23	7.416	731	736	743	rVB5	7670	14050	0.24%	0.031%
24	7.781	790	798	806	rBV	467113	749711	12.89%	1.637%
25	7.852	806	810	818	rVB5	9195	16730	0.29%	0.037%
26	8.128	851	857	862	rVB2	16310	25383	0.44%	0.055%
27	8.505	913	921	934	rBV	412641	680316	11.70%	1.485%
28	8.616	935	940	947	rVV2	21369	34125	0.59%	0.074%
29	8.705	950	955	960	rBV	12517	20078	0.35%	0.044%
30	8.899	984	988	989	rBV3	6698	6314	0.11%	0.014%
31	8.946	989	996	1006	rVB	643332	1039292	17.87%	2.269%
32	9.175	1029	1035	1045	rBV2	33819	82446	1.42%	0.180%
33	9.293	1047	1055	1060	rVV2	5573	12712	0.22%	0.028%
34	9.340	1060	1063	1067	rVB4	4600	6293	0.11%	0.014%
35	9.505	1086	1091	1094	rBV5	5774	9093	0.16%	0.020%
36	9.663	1111	1118	1133	rBV	534885	886073	15.24%	1.934%

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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-BP010824.MA.M
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37	10.205	1201	1210	1225	rBV	735850	1302053	22.39%	2.842%
38	10.575	1265	1273	1287	rVB	659556	1097995	18.88%	2.397%
39	10.728	1290	1299	1317	rBV	606043	1111411	19.11%	2.426%
40	10.887	1317	1326	1331	rVB9	4203	12491	0.21%	0.027%
41	11.116	1358	1365	1371	rBV	62999	105694	1.82%	0.231%
42	11.181	1371	1376	1385	rVB	75116	134477	2.31%	0.294%
43	11.416	1409	1416	1429	rVB9	13996	38961	0.67%	0.085%
44	11.652	1451	1456	1461	rVB8	4194	6631	0.11%	0.014%
45	11.969	1502	1510	1514	rBV5	10203	24884	0.43%	0.054%
46	11.999	1514	1515	1522	rVB4	7885	11419	0.20%	0.025%
47	12.175	1538	1545	1550	rBV2	12231	23801	0.41%	0.052%
48	12.275	1557	1562	1570	rBV	27104	44717	0.77%	0.098%
49	12.898	1662	1668	1673	rBV8	4377	7664	0.13%	0.017%
50	13.246	1721	1727	1732	rVB10	3804	7313	0.13%	0.016%
51	13.322	1735	1740	1748	rBV2	17266	30861	0.53%	0.067%
52	13.393	1748	1752	1756	rBV6	3759	7760	0.13%	0.017%
53	13.463	1761	1764	1770	rVB7	3242	6191	0.11%	0.014%
54	13.546	1774	1778	1784	rVB8	3676	7196	0.12%	0.016%
55	13.846	1822	1829	1834	rBV	1701772	2316123	39.83%	5.056%
56	13.893	1834	1837	1843	rVB	70886	95055	1.63%	0.207%
57	14.116	1866	1875	1889	rBV	1777393	2547758	43.81%	5.561%
58	14.304	1904	1907	1914	rVB5	6226	11840	0.20%	0.026%
59	14.422	1921	1927	1936	rVB	1133666	1664647	28.63%	3.634%
60	14.563	1946	1951	1960	rVB	3805	9639	0.17%	0.021%
61	14.669	1963	1969	1986	rBV2	127999	279413	4.81%	0.610%
62	14.845	1994	1999	2003	rBV8	4119	8698	0.15%	0.019%
63	14.916	2008	2011	2016	rBV7	3999	6125	0.11%	0.013%
64	15.034	2028	2031	2036	rVB7	4654	6813	0.12%	0.015%
65	15.234	2061	2065	2075	rVB6	10531	27172	0.47%	0.059%
66	15.422	2090	2097	2105	rBV	2582751	3592010	61.77%	7.841%
67	15.557	2114	2120	2134	rBV	978801	1400121	24.08%	3.056%
68	15.663	2134	2138	2142	rVB2	11765	17293	0.30%	0.038%
69	16.210	2227	2231	2235	rVB6	9216	12746	0.22%	0.028%
70	16.857	2336	2341	2349	rVB6	6692	11405	0.20%	0.025%
71	16.963	2351	2359	2365	rBV6	6242	16465	0.28%	0.036%
72	17.181	2389	2396	2403	rBV	1656457	2272398	39.08%	4.960%
73	17.281	2406	2413	2423	rVV	3360229	4578890	78.74%	9.995%
74	17.581	2461	2464	2470	rVB8	8768	13486	0.23%	0.029%
75	17.851	2506	2510	2514	rBV6	5211	6372	0.11%	0.014%
76	18.075	2542	2548	2557	rBV	122797	174404	3.00%	0.381%

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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

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77	18.486	2614	2618	2622	rVB6	18315	21284	0.37%	0.046%
78	18.910	2688	2690	2693	rBV4	5605	6960	0.12%	0.015%
79	19.104	2718	2723	2728	rBV2	48546	72272	1.24%	0.158%
80	19.175	2731	2735	2738	rBV	45663	64415	1.11%	0.141%
81	19.316	2755	2759	2763	rBV2	60488	83347	1.43%	0.182%
82	19.528	2789	2795	2803	rBV	4468053	5815005	100.00%	12.693%
83	21.286	3089	3094	3102	rBV	2087647	2574713	44.28%	5.620%
84	23.492	3461	3469	3481	rVB	2813507	5418135	93.18%	11.827%
85	23.639	3488	3494	3505	rVB	1224585	2417053	41.57%	5.276%

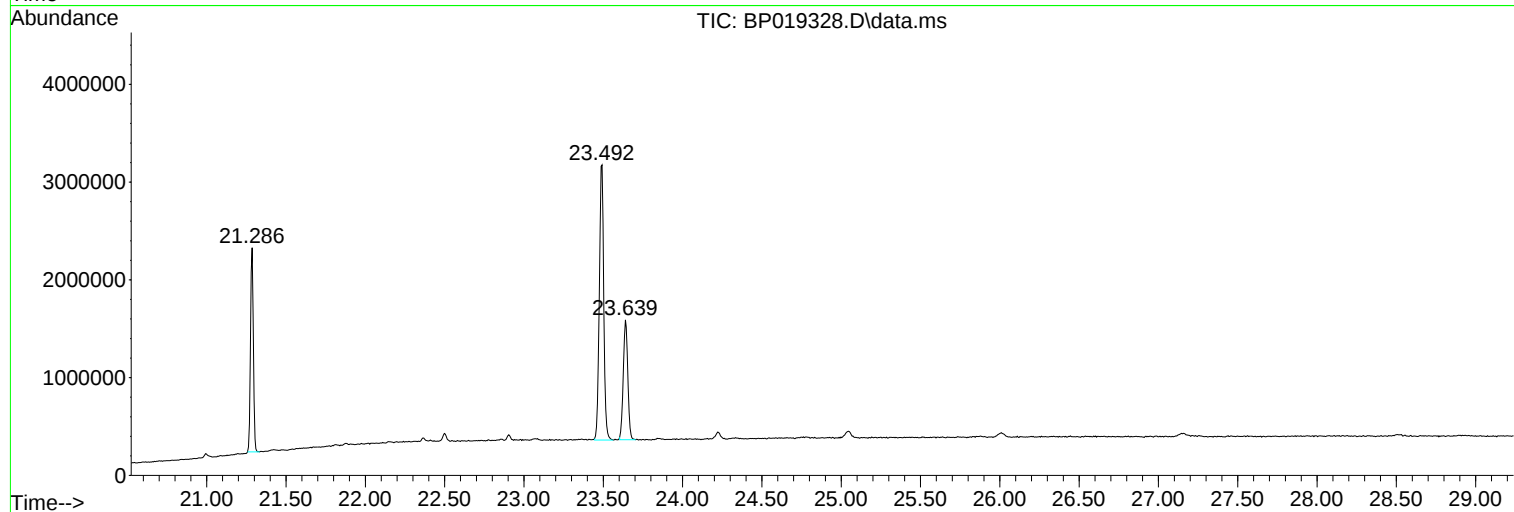
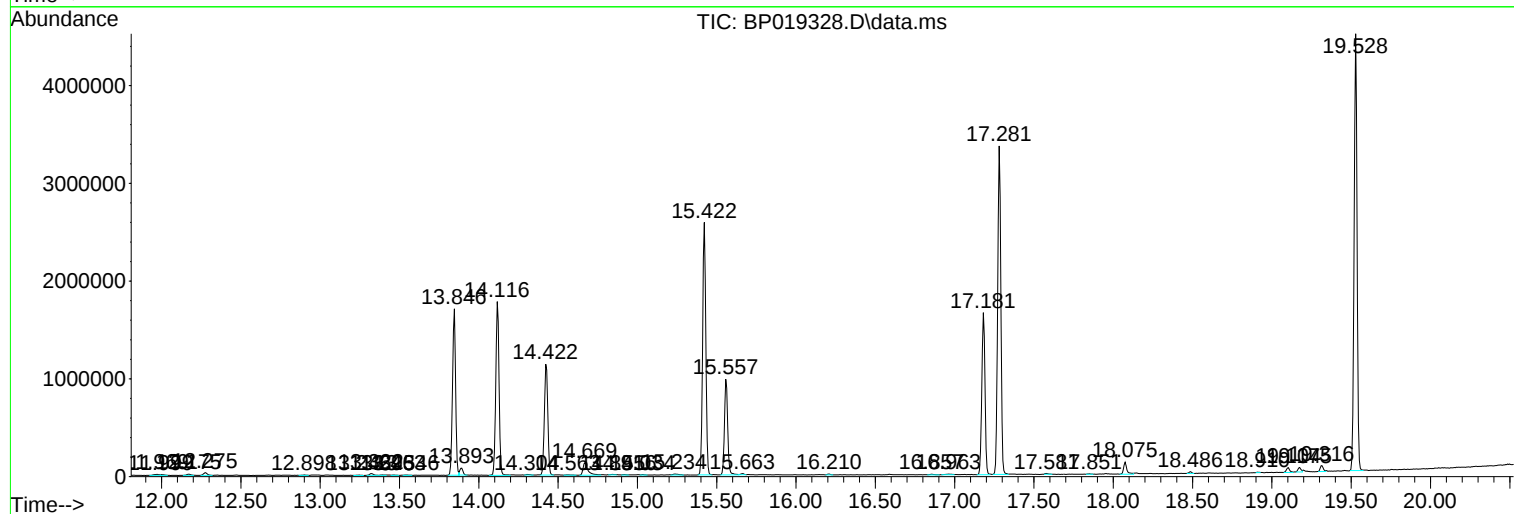
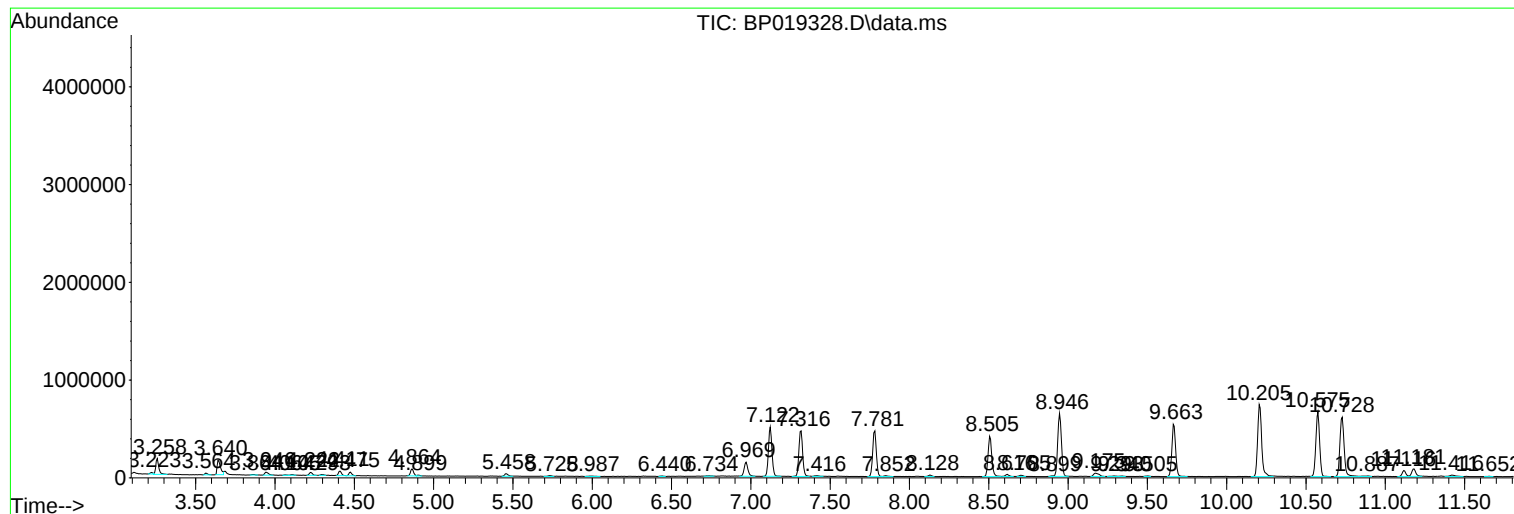
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Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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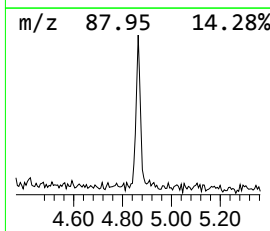
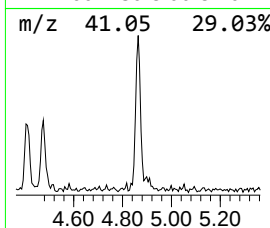
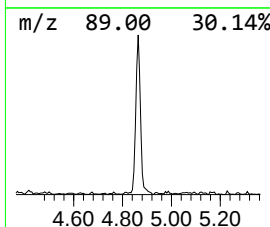
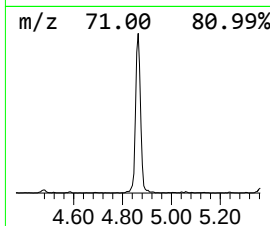
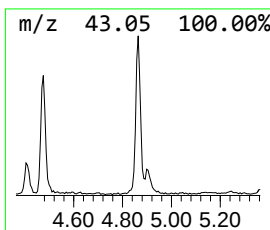
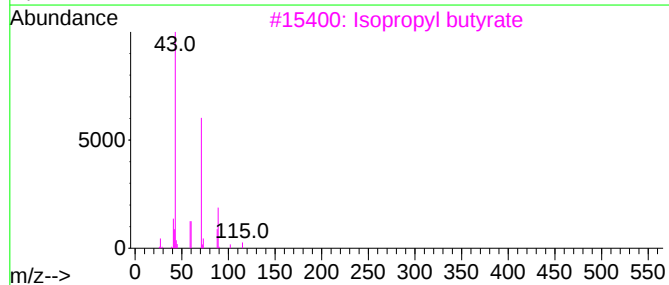
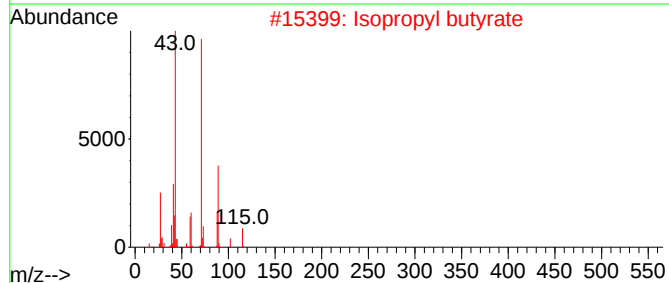
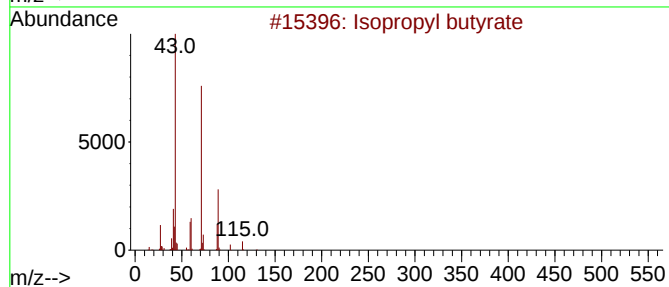
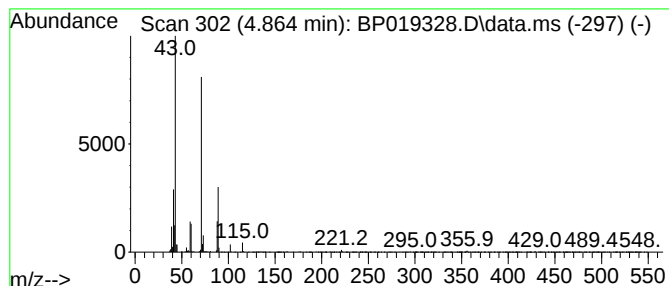
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Isopropyl butyrate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.864	2.60 ng/ul	97457	1,4-Dichlorobenzene-d4	7.781

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-, 1-met...	130	C7H14O2	000617-50-5	72
5			Butanoic acid, pentyl ester	158	C9H18O2	000540-18-1	72



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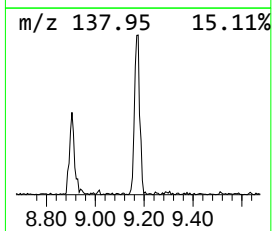
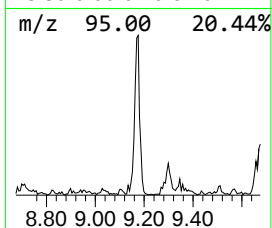
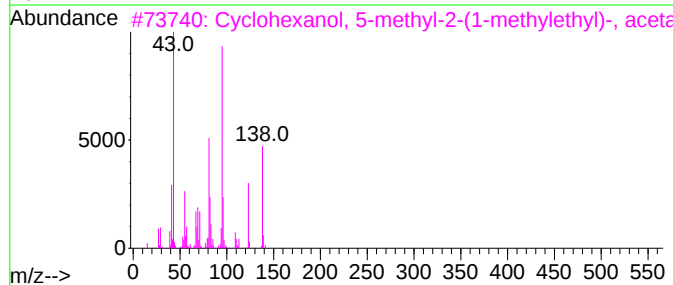
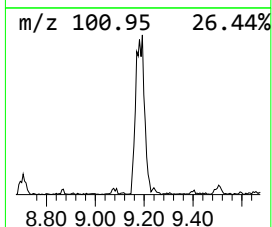
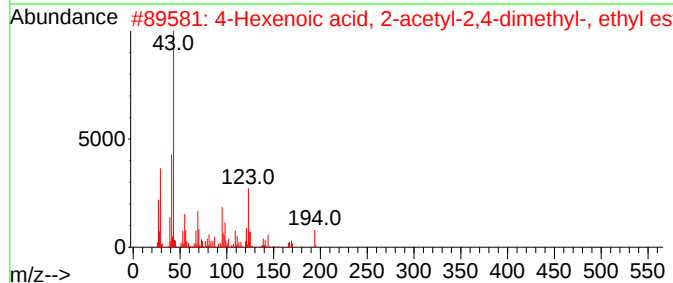
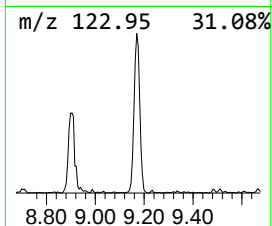
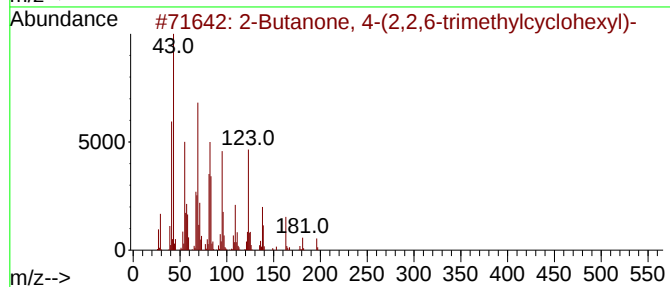
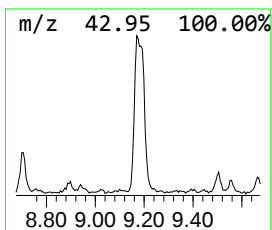
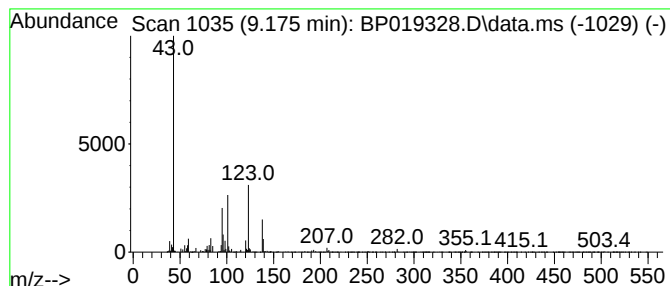
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.175	2.20 ng/ul	82446	1,4-Dichlorobenzene-d4	7.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanone, 4-(2,2,6-trimethylcy...	196	C13H24O	006138-85-8	28		
2	4-Hexenoic acid, 2-acetyl-2,4-di...	212	C12H20O3	1000159-11-3	23		
3	Cyclohexanol, 5-methyl-2-(1-meth...	198	C12H22O2	020777-45-1	12		
4	2-Pyrimidinamine, 4,5-dimethyl-	123	C6H9N3	001193-74-4	11		
5	Methyl 3,4,6-tri-O-acetyl-2-acet...	361	C15H23NO9	002595-39-3	10		



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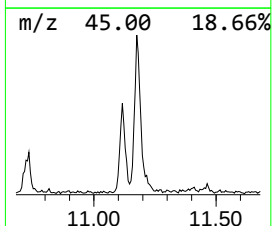
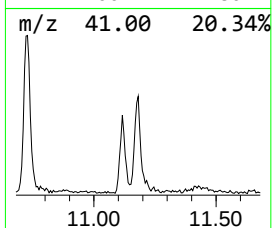
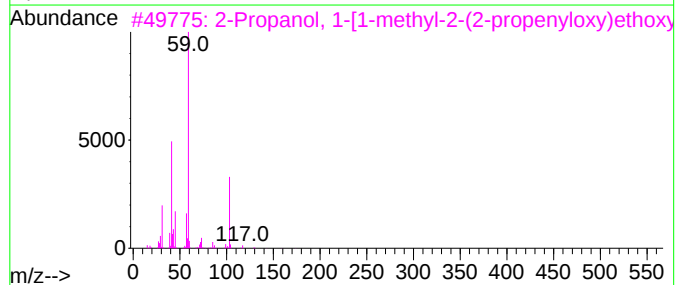
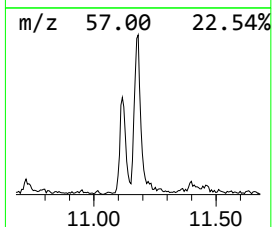
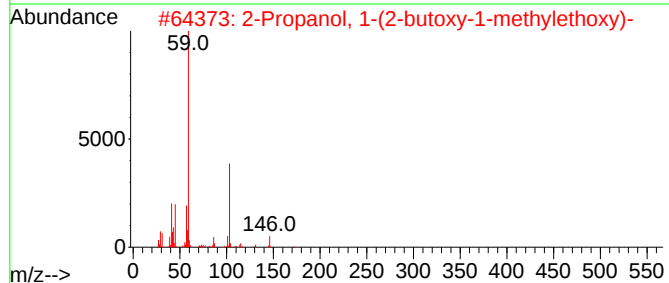
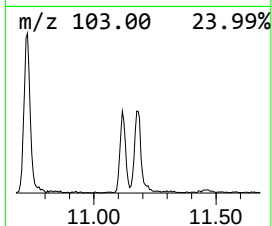
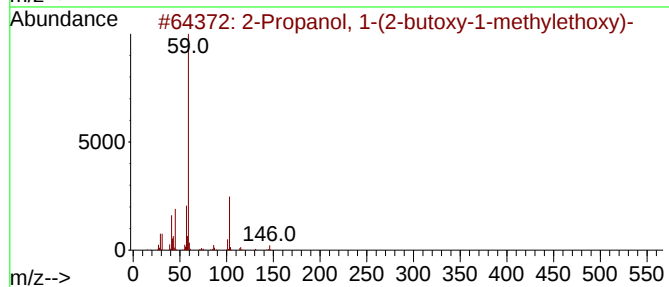
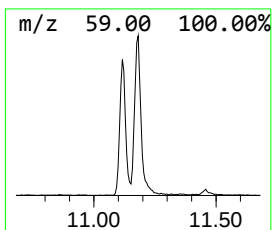
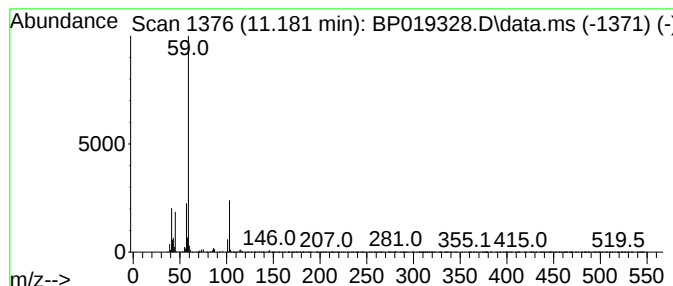
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 2-Propanol, 1-(2-butoxy-1-m... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.181	2.45 ng/ul	134477	Naphthalene-d8	10.575

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	90		
2	2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	74		
3	2-Propanol, 1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	72		
4	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	72		
5	1,4,7-Trimethyl-3,6-dioxaoctane...	192	C9H20O4	1000423-63-2	59		



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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.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
Isopropyl butyrate	4.864	2.6	ng/ul	97457	1	7.781	749711	20.0
unknown-01	9.175	2.2	ng/ul	82446	1	7.781	749711	20.0
2-Propanol, 1-(...	11.181	2.5	ng/ul	134477	2	10.575	1098000	20.0