

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM110519\
 Data File : BM023601.D
 Acq On : 05 Nov 2019 11:13
 Operator : JU
 Sample : K5536-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BEK32

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.176	28	32	46	rVB	3890650	4890516	47.08%	4.434%
2	3.529	89	92	98	rVB2	150641	192779	1.86%	0.175%
3	3.587	99	102	109	rVB	483678	559579	5.39%	0.507%
4	3.840	142	145	151	rVB2	166582	272919	2.63%	0.247%
5	4.111	187	191	199	rVV	982426	1294242	12.46%	1.174%
6	4.182	200	203	212	rVB3	144103	215935	2.08%	0.196%
7	4.287	217	221	229	rBV	1490252	1887877	18.17%	1.712%
8	4.734	292	297	303	rVB	1877151	2405658	23.16%	2.181%
9	5.587	438	442	447	rVB	271354	369874	3.56%	0.335%
10	6.781	641	645	661	rVB	332858	668919	6.44%	0.607%
11	6.946	667	673	689	rVB	1092377	1851832	17.83%	1.679%
12	7.134	700	705	714	rVB	1206522	1961315	18.88%	1.778%
13	7.599	778	784	792	rVB	1397532	2202161	21.20%	1.997%
14	8.311	899	905	916	rBV	915926	1495710	14.40%	1.356%
15	8.746	973	979	991	rVB	1447881	2476425	23.84%	2.245%
16	9.205	1055	1057	1059	rBV2	13919	15177	0.15%	0.014%
17	9.469	1095	1102	1110	rBV	1145887	1886699	18.16%	1.711%
18	10.005	1186	1193	1207	rBV	1697344	2860514	27.54%	2.594%
19	10.375	1249	1256	1265	rBV	1985973	3273339	31.51%	2.968%
20	10.516	1273	1280	1294	rBV	1467482	2558761	24.63%	2.320%
21	11.034	1365	1368	1375	rVB5	13078	24296	0.23%	0.022%
22	11.193	1390	1395	1400	rVV3	16671	28757	0.28%	0.026%
23	11.905	1512	1516	1520	rBV4	8028	12721	0.12%	0.012%
24	11.975	1524	1528	1534	rVB4	31093	52180	0.50%	0.047%
25	12.057	1538	1542	1544	rBV4	9337	14652	0.14%	0.013%
26	12.252	1569	1575	1578	rBV4	8847	18125	0.17%	0.016%
27	12.863	1676	1679	1683	rBV5	12673	14543	0.14%	0.013%
28	13.099	1715	1719	1722	rBV4	11999	18134	0.17%	0.016%
29	13.169	1726	1731	1733	rBV6	6790	10994	0.11%	0.010%
30	13.669	1809	1816	1821	rVV	3286423	4667766	44.93%	4.232%
31	13.710	1821	1823	1830	rVB	233014	319917	3.08%	0.290%
32	13.940	1853	1862	1876	rBV	3971617	5872595	56.53%	5.325%
33	14.175	1897	1902	1906	rBV6	7400	11133	0.11%	0.010%
34	14.251	1908	1915	1920	rBV2	2983071	4478700	43.11%	4.061%

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35	14.293	1920	1922	1929	rVB	116964	157968	1.52%	0.143%
36	14.469	1946	1952	1969	rBV4	150144	404672	3.90%	0.367%
37	14.687	1983	1989	1994	rVV6	56412	93113	0.90%	0.084%
38	14.740	1994	1998	2001	rVB5	10161	16664	0.16%	0.015%
39	14.881	2018	2022	2029	rVB7	8871	16224	0.16%	0.015%
40	15.040	2042	2049	2062	rBV2	101757	216496	2.08%	0.196%
41	15.128	2062	2064	2070	rVB3	15309	20523	0.20%	0.019%
42	15.246	2077	2084	2097	rBV	5085703	7305434	70.32%	6.624%
43	15.375	2100	2106	2120	rVV	1223292	1770259	17.04%	1.605%
44	15.487	2120	2125	2130	rVB	61342	87815	0.85%	0.080%
45	16.034	2214	2218	2224	rVB7	24443	40172	0.39%	0.036%
46	16.157	2234	2239	2245	rBV9	6157	13818	0.13%	0.013%
47	16.687	2323	2329	2333	rVB6	12005	19688	0.19%	0.018%
48	16.787	2341	2346	2352	rVB3	23633	42347	0.41%	0.038%
49	16.998	2375	2382	2391	rBV	3667373	5128001	49.36%	4.650%
50	17.098	2392	2399	2407	rBV	5722881	8377207	80.64%	7.596%
51	17.304	2430	2434	2439	rVB5	10604	20449	0.20%	0.019%
52	17.381	2442	2447	2449	rBV5	7263	10676	0.10%	0.010%
53	17.528	2469	2472	2475	rBV4	14666	16212	0.16%	0.015%
54	17.692	2493	2500	2502	rBV6	8597	17484	0.17%	0.016%
55	17.916	2533	2538	2546	rBV2	188823	296488	2.85%	0.269%
56	18.222	2587	2590	2592	rBV3	19619	18754	0.18%	0.017%
57	18.863	2695	2699	2702	rBV4	33510	45910	0.44%	0.042%
58	19.034	2724	2728	2730	rBV	60606	85462	0.82%	0.077%
59	19.204	2753	2757	2763	rBV4	95729	147045	1.42%	0.133%
60	19.286	2767	2771	2775	rVV2	59113	89902	0.87%	0.082%
61	19.398	2784	2790	2796	rBV	7559919	9833068	94.65%	8.916%
62	19.439	2796	2797	2800	rVV2	85884	83922	0.81%	0.076%
63	19.475	2801	2803	2809	rVB3	58144	66653	0.64%	0.060%
64	19.980	2886	2889	2893	rVB	141813	145887	1.40%	0.132%
65	20.475	2970	2973	2977	rBV	246643	272980	2.63%	0.248%
66	20.939	3048	3052	3059	rBV	730243	1089585	10.49%	0.988%
67	21.198	3091	3096	3104	rBV	4747614	5881844	56.62%	5.333%
68	21.380	3123	3127	3131	rBV	594525	677998	6.53%	0.615%
69	21.810	3196	3200	3209	rBV	562332	938130	9.03%	0.851%
70	22.263	3274	3277	3285	rVB	506487	652756	6.28%	0.592%
71	22.757	3358	3361	3367	rVB	422767	569393	5.48%	0.516%

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72	23.251	3438	3445	3451	rBV	5739615	10388372	100.00%	9.419%
73	23.386	3462	3468	3478	rVB	3409522	6342527	61.05%	5.751%

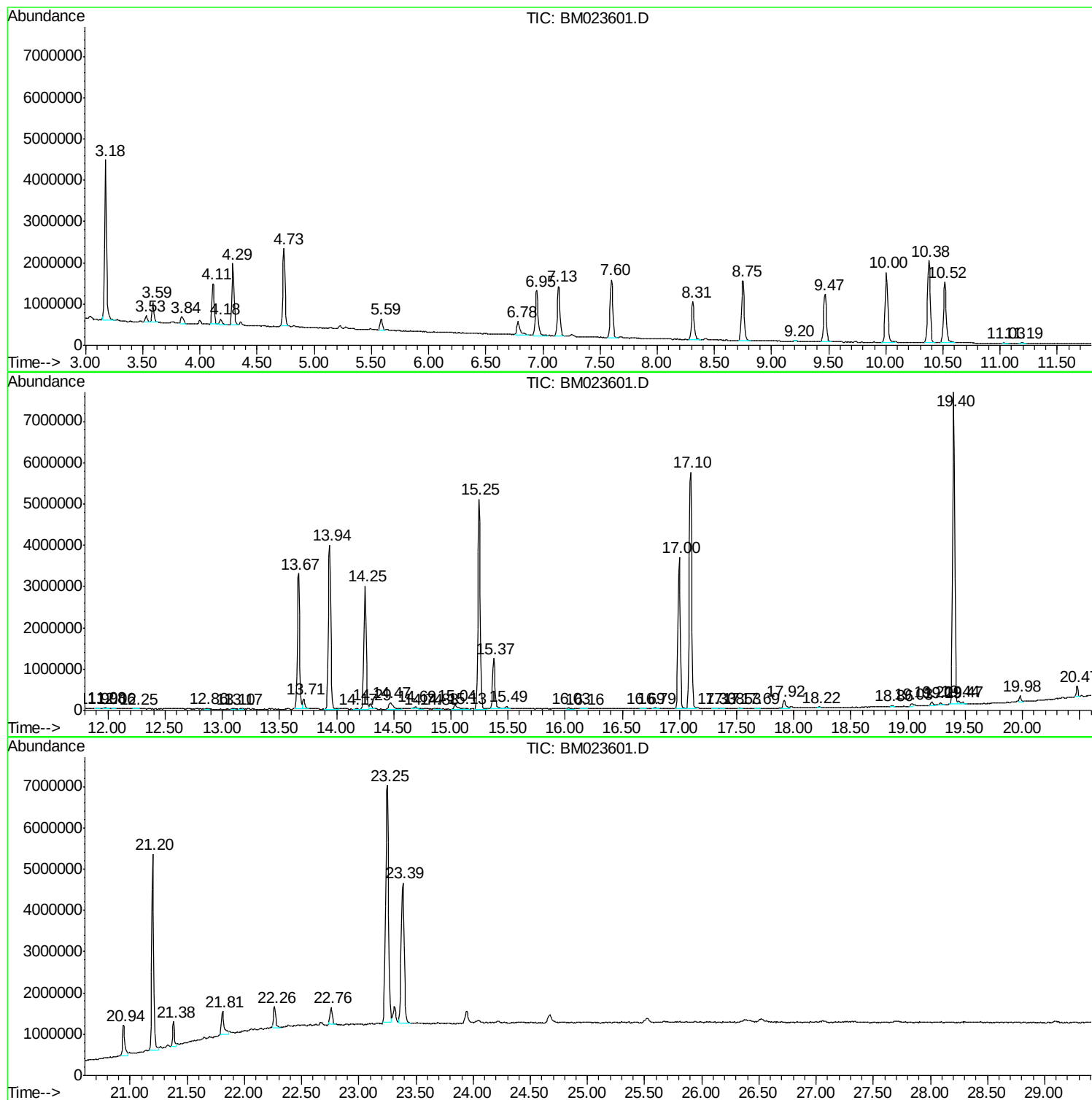
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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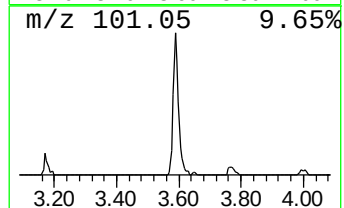
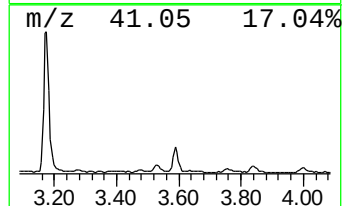
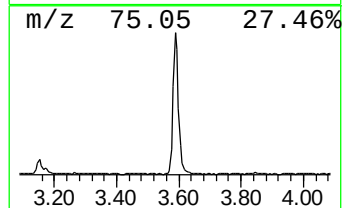
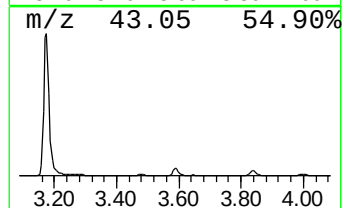
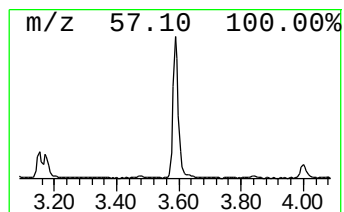
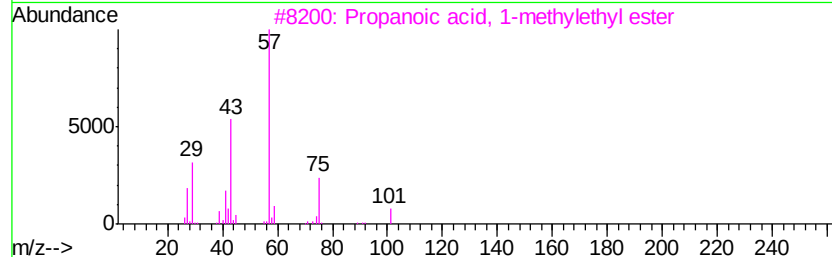
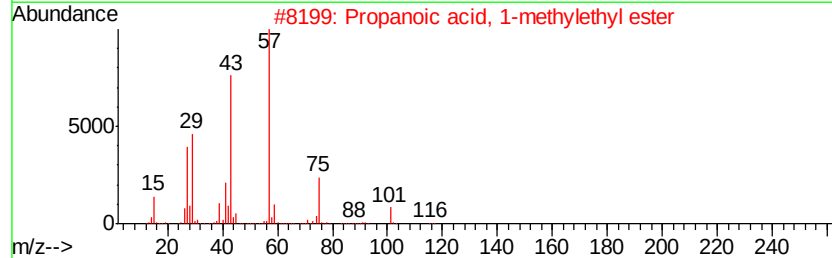
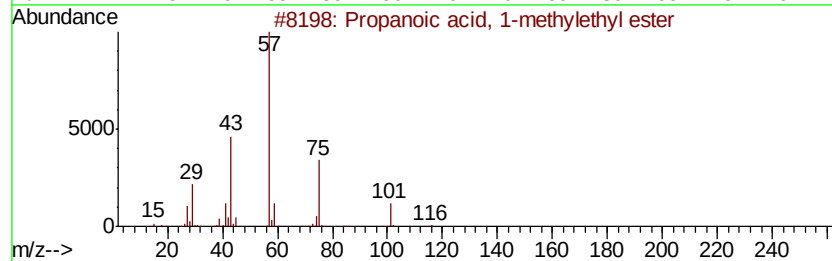
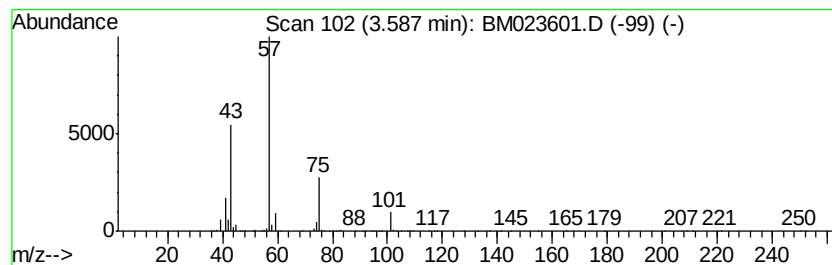
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Propanoic acid, 1-methyleth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.59	5.08 ng/ul	559579	1,4-Dichlorobenzene-d4	7.60
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Propanoic acid, 1-methylethyl ester	116 C6H12O2	000637-78-5	83
2	Propanoic acid, 1-methylethyl ester	116 C6H12O2	000637-78-5	83
3	Propanoic acid, 1-methylethyl ester	116 C6H12O2	000637-78-5	53
4	Propanoic acid, 1-methylethyl ester	116 C6H12O2	000637-78-5	38
5	(+)-4-Amino-4,5-dihydro-2(3H)-f...	101 C4H7NO2	016504-58-8	9



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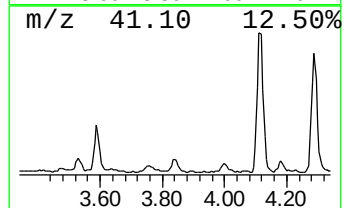
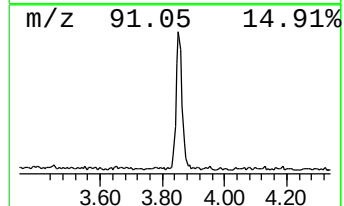
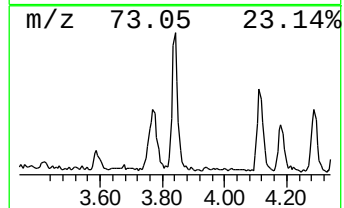
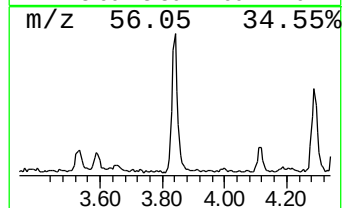
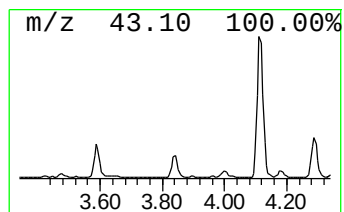
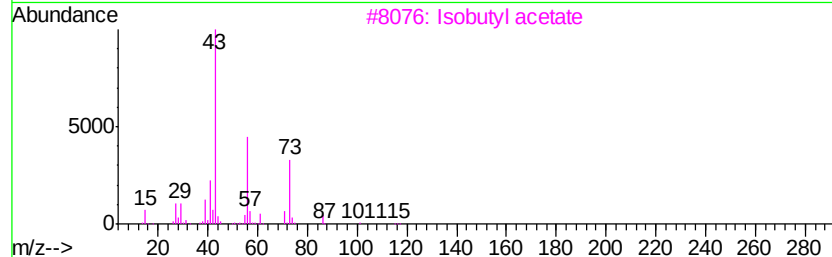
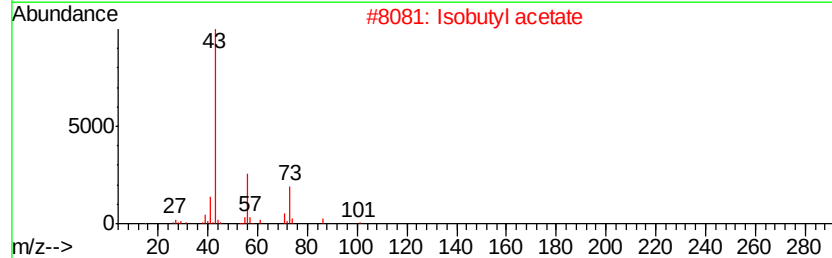
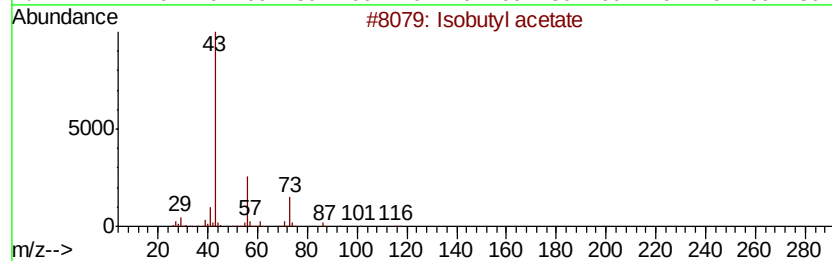
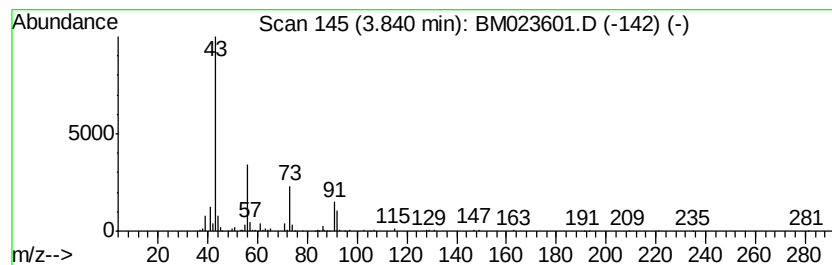
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Isobutyl acetate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.84	2.48 ng/ul	272919	1,4-Dichlorobenzene-d4	7.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Isobutyl acetate	116	C6H12O2	000110-19-0	64
2		Isobutyl acetate	116	C6H12O2	000110-19-0	59
3		Isobutyl acetate	116	C6H12O2	000110-19-0	53
4		Isobutyl acetate	116	C6H12O2	000110-19-0	50
5		Isobutyl acetate	116	C6H12O2	000110-19-0	42



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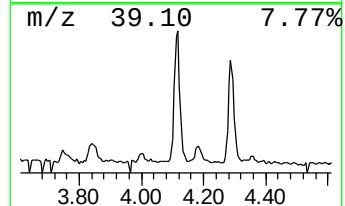
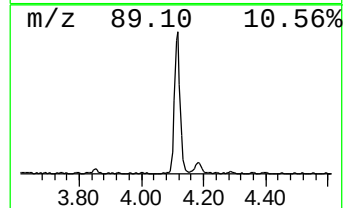
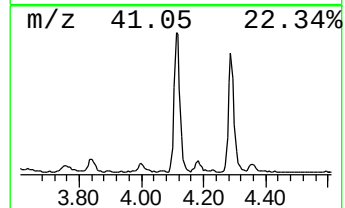
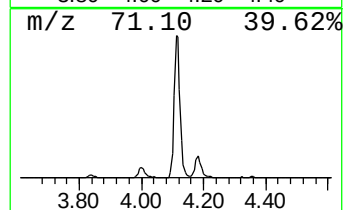
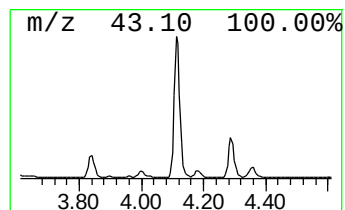
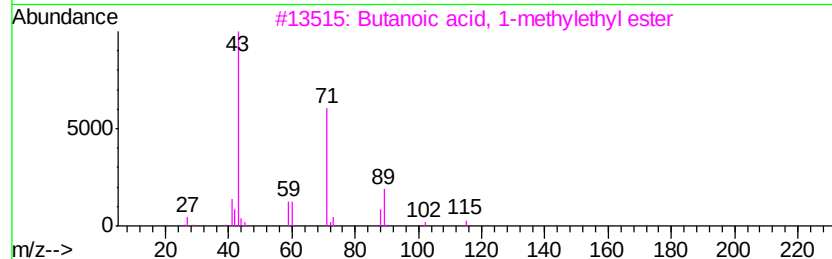
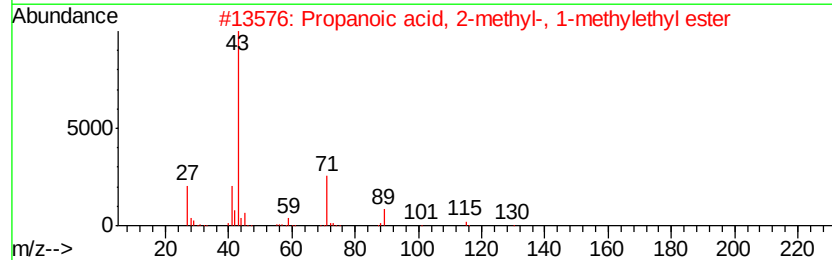
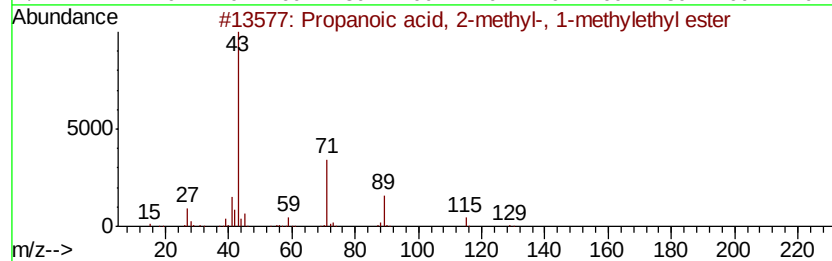
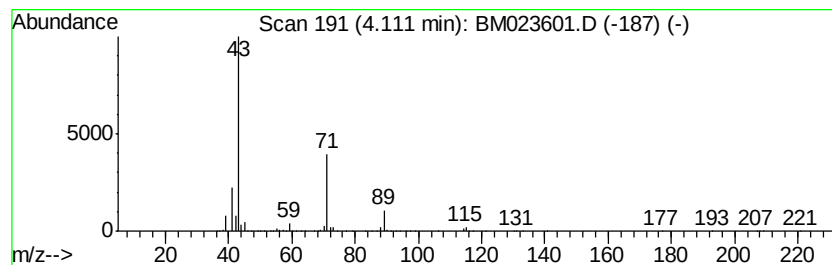
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Peak Number 3 Propanoic acid, 2-methyl-, ... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.11	11.75 ng/ul	1294240	1,4-Dichlorobenzene-d4	7.60

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	83
3		Butanoic acid, 1-methyl ethyl ester	130	C7H14O2	000638-11-9	38
4		Propanoic acid, 2-methyl-, 3-met...	158	C9H18O2	002050-01-3	38
5		Pentane, 3-ethyl-	100	C7H16	000617-78-7	36



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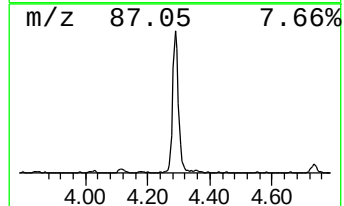
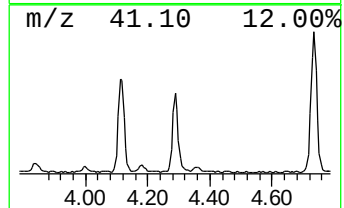
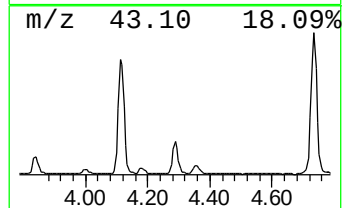
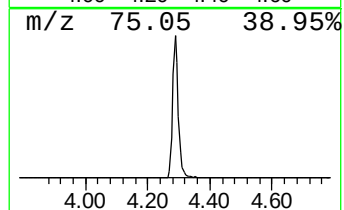
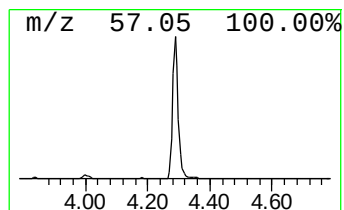
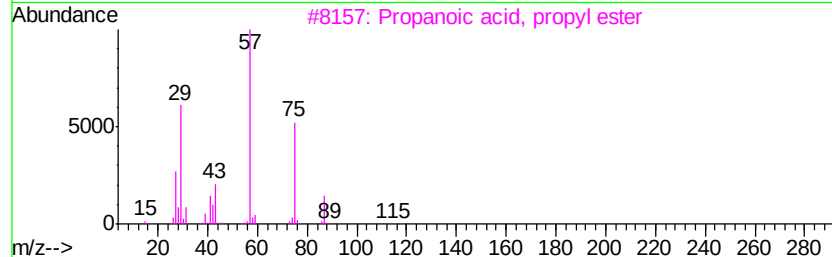
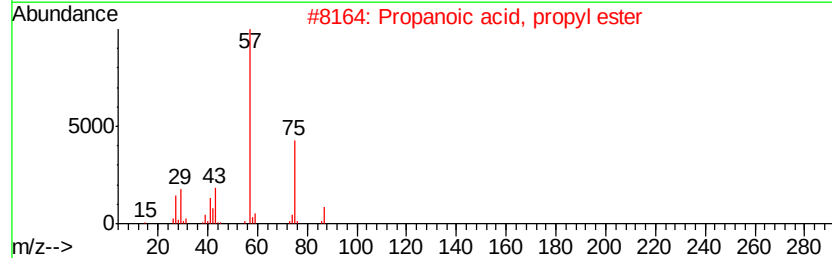
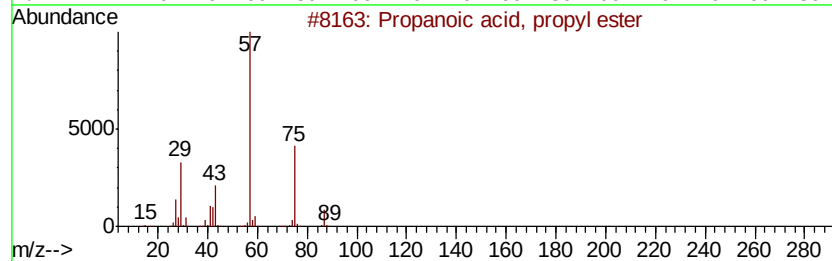
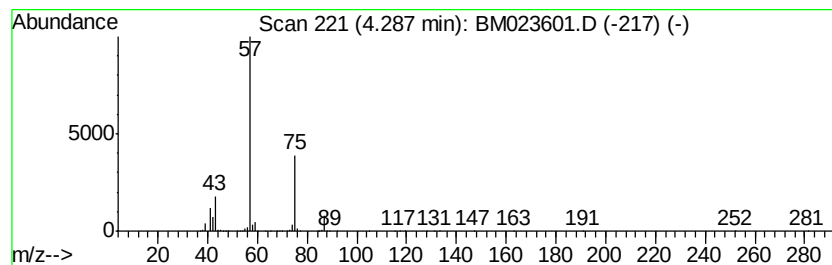
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Propanoic acid, propyl ester Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.29	17.15 ng/ul	1887880	1,4-Dichlorobenzene-d4	7.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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	83
5		Propanoic acid, butyl ester	130	C7H14O2	000590-01-2	28



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM110519\
Data File : BM023601.D
Acq On : 05 Nov 2019 11:13
Operator : JU
Sample : K5536-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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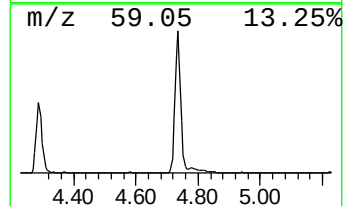
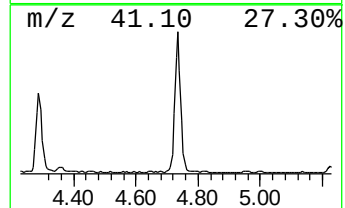
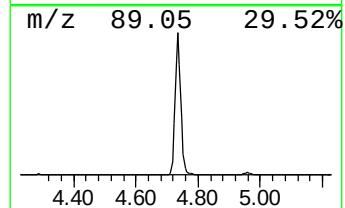
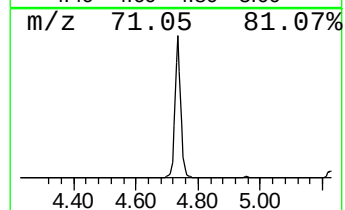
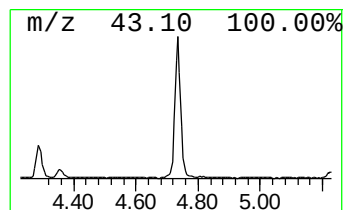
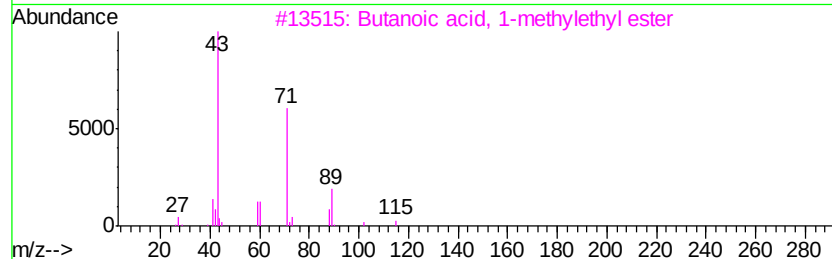
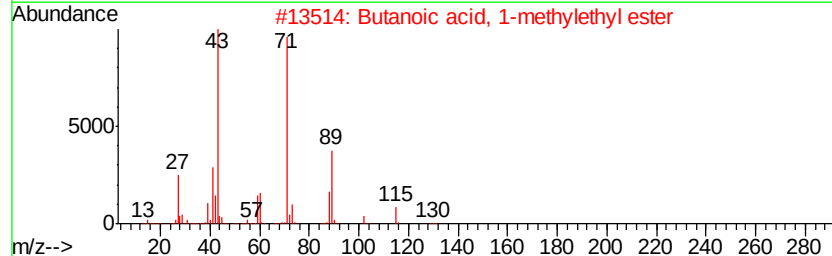
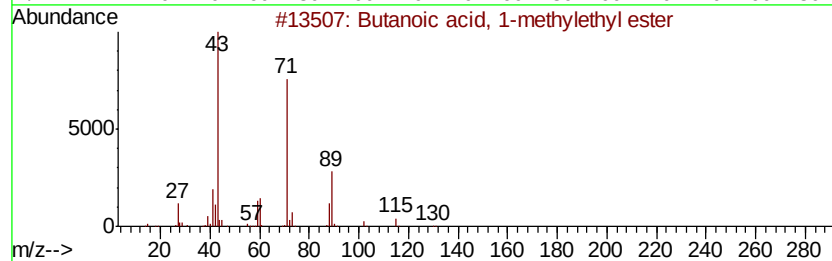
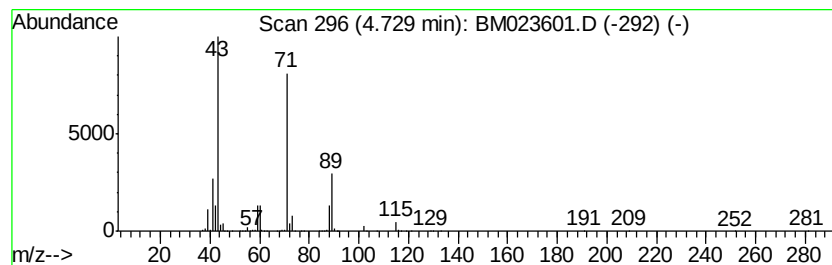
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Butanoic acid, 1-methylethy... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.73	21.85 ng/ul	2405660	1,4-Dichlorobenzene-d4	7.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	95
2		Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	90
3		Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	83
4		Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	64
5		Butanoic acid, pentyl ester	158	C9H18O2	000540-18-1	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM110519\
Data File : BM023601.D
Acq On : 05 Nov 2019 11:13
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Sample : K5536-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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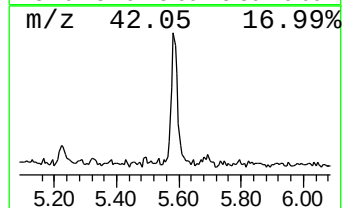
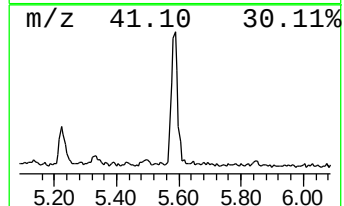
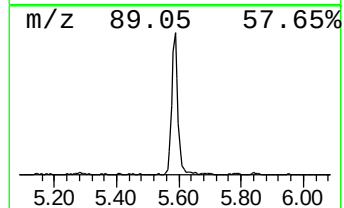
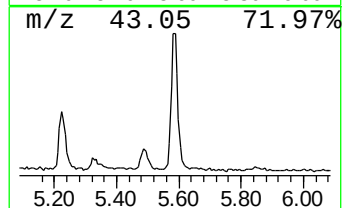
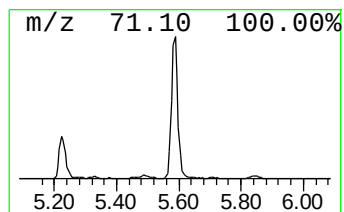
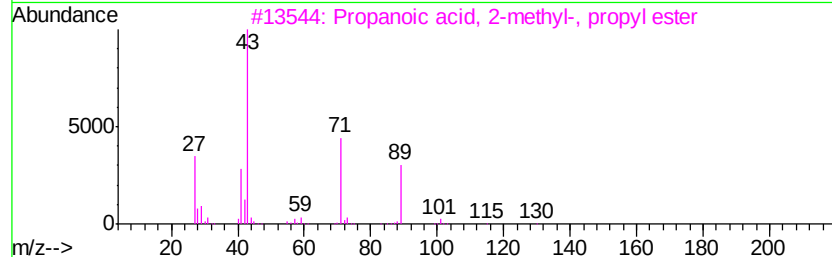
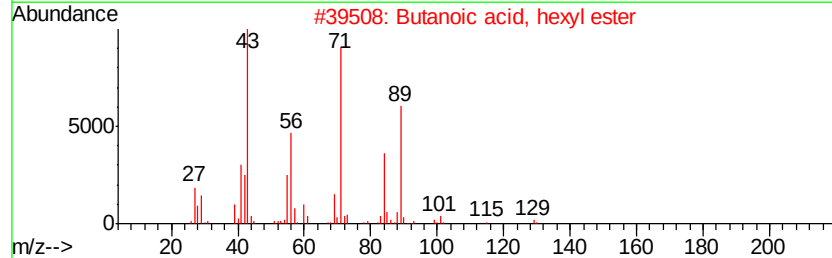
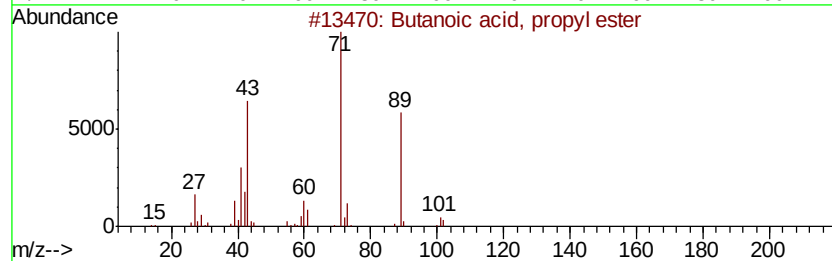
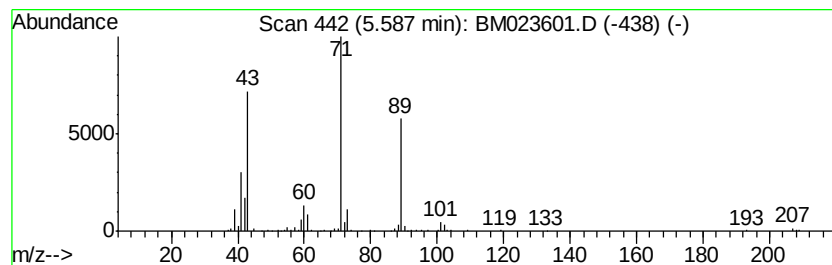
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Butanoic acid, propyl ester Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.59	3.36 ng/ul	369874	1,4-Dichlorobenzene-d4	7.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	90
2		Butanoic acid, hexyl ester	172	C10H20O2	002639-63-6	78
3		Propanoic acid, 2-methyl-, propyl...	130	C7H14O2	000644-49-5	78
4		Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	74
5		Propanoic acid, 2-methyl-, propyl...	130	C7H14O2	000644-49-5	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM110519\
Data File : BM023601.D
Acq On : 05 Nov 2019 11:13
Operator : JU
Sample : K5536-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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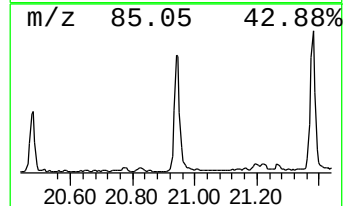
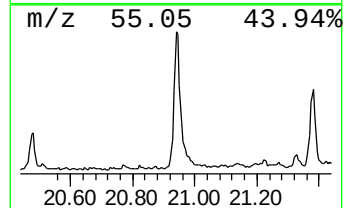
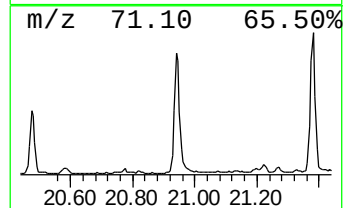
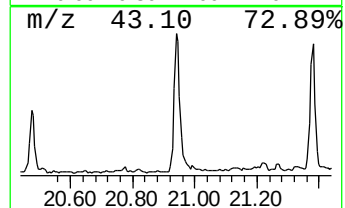
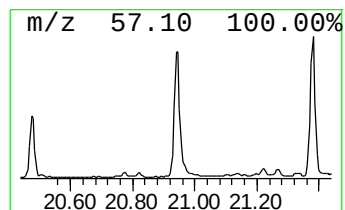
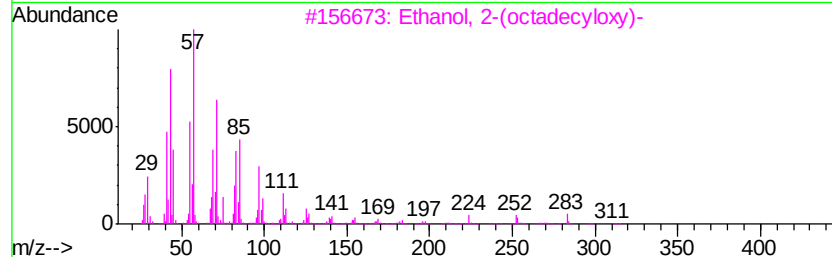
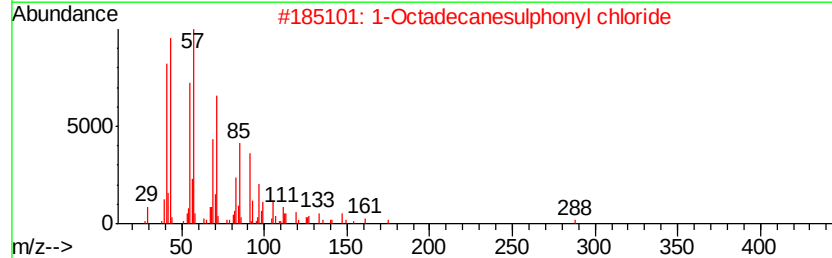
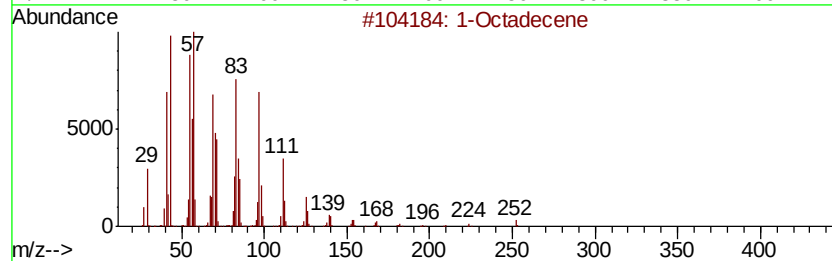
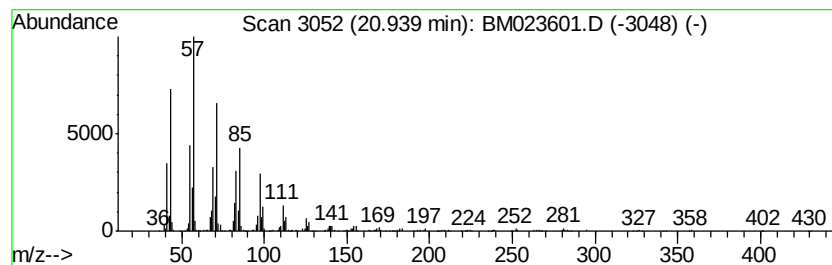
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Cyclic20.94 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.94	3.70 ng/ul	1089590	Chrysene-d12	21.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Octadecene	252	C18H36	000112-88-9	95
2		1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	91
3		Ethanol, 2-(octadecyloxy)-	314	C20H42O2	002136-72-3	90
4		Octatriacontyl pentafluoropropio...	697	C41H77F5O2	1000351-89-1	83
5		Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM110519\
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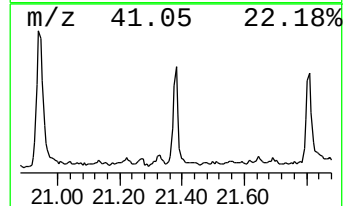
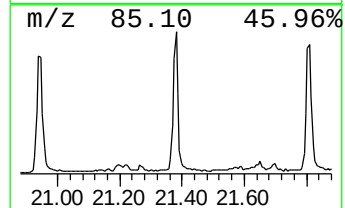
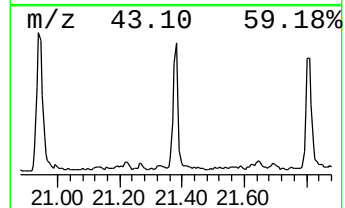
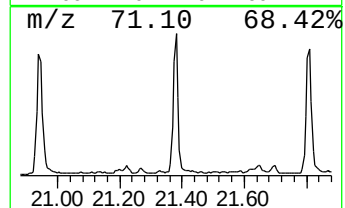
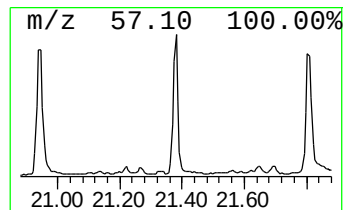
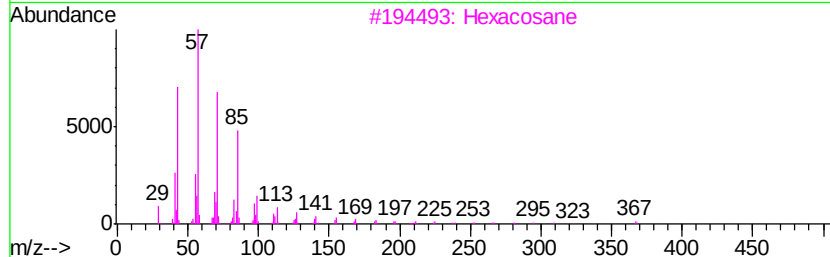
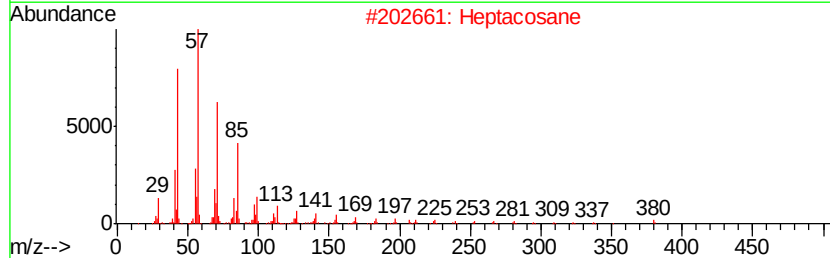
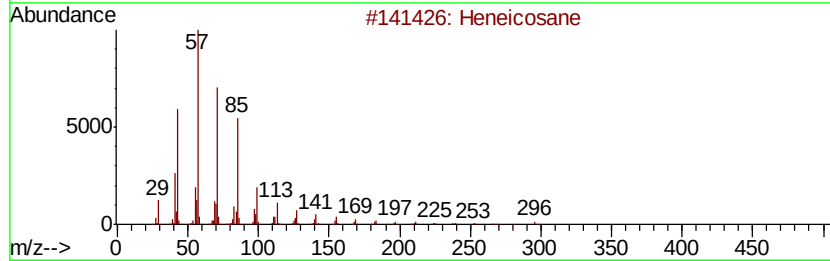
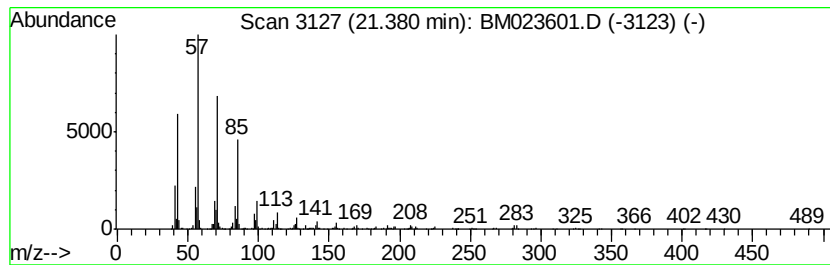
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
21.38	2.31 ng/ul	677998	Chrysene-d12		21.20
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	94
2	Heptacosane	380	C27H56	000593-49-7	90
3	Hexacosane	366	C26H54	000630-01-3	90
4	Octacosane	394	C28H58	000630-02-4	87
5	Heptadecane	240	C17H36	000629-78-7	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM110519\
Data File : BM023601.D
Acq On : 05 Nov 2019 11:13
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Sample : K5536-02
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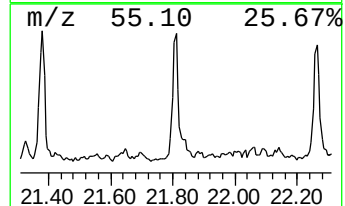
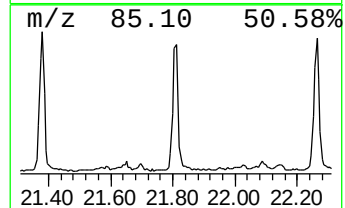
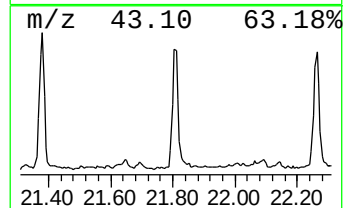
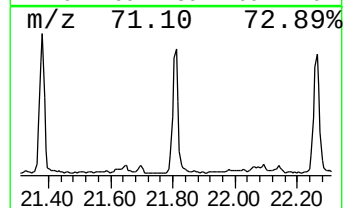
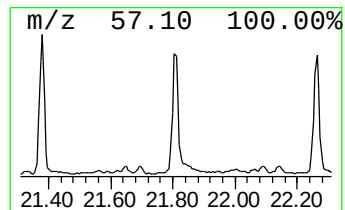
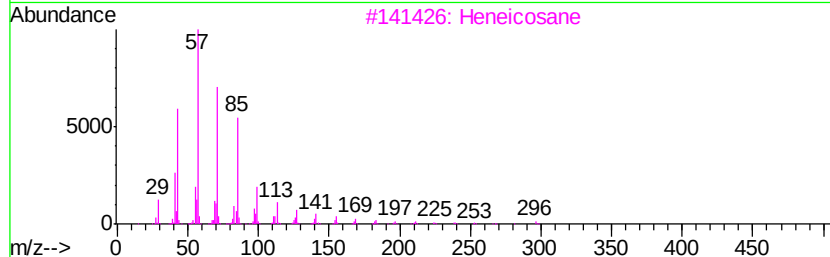
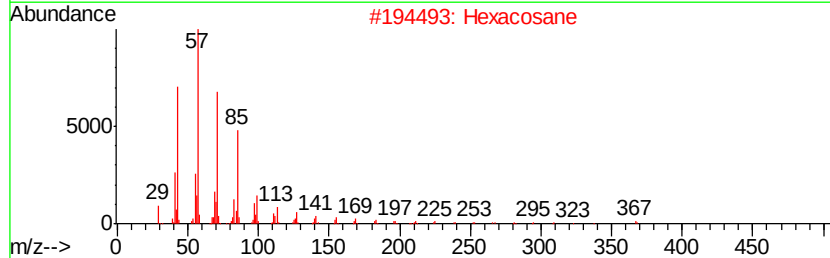
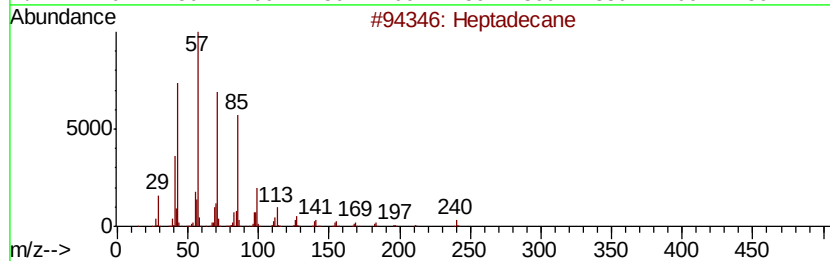
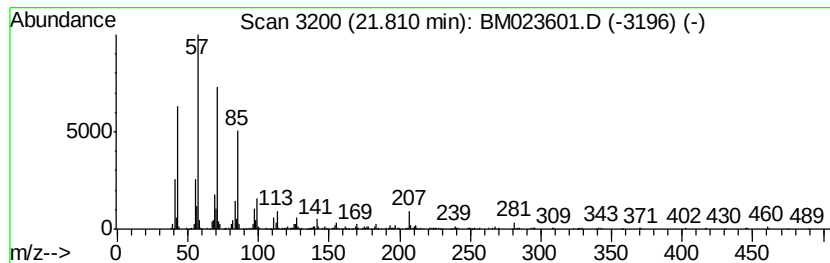
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.81	3.19 ng/ul	938130	Chrysene-d12	21.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	93
2		Hexacosane	366	C26H54	000630-01-3	87
3		Heneicosane	296	C21H44	000629-94-7	87
4		Octacosane	394	C28H58	000630-02-4	87
5		Tetratetracontane	619	C44H90	007098-22-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM110519\
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Acq On : 05 Nov 2019 11:13
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ALS Vial : 4 Sample Multiplier: 1

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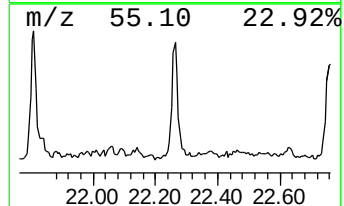
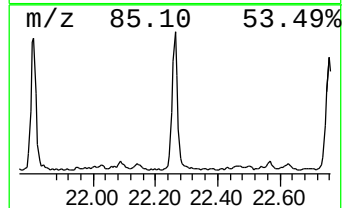
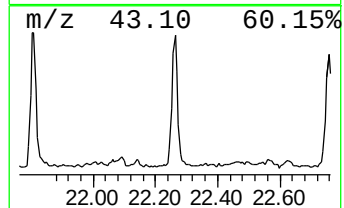
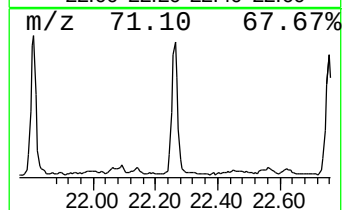
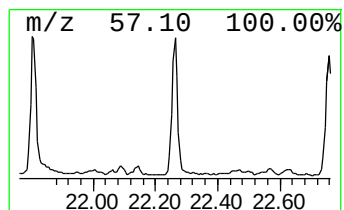
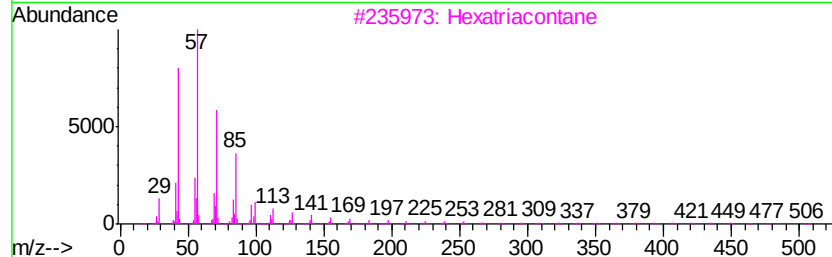
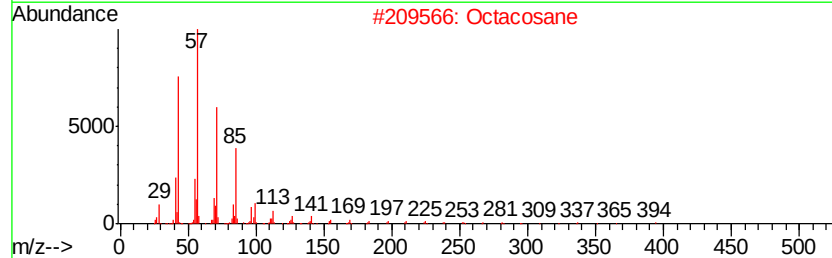
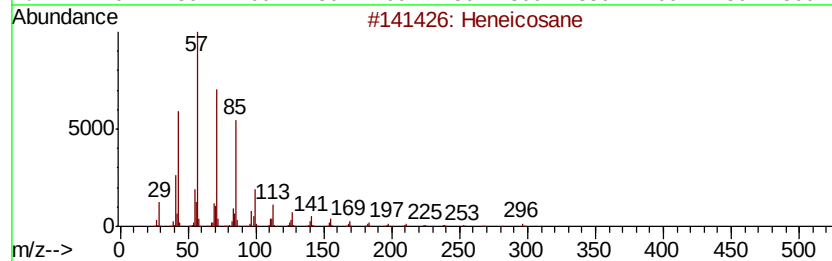
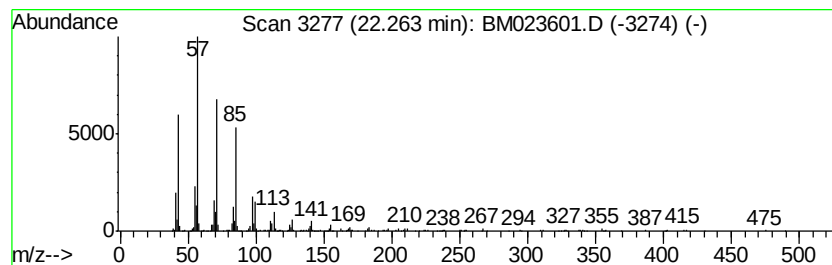
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.26	2.22 ng/ul	652756	Chrysene-d12	21.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	90
2		Octacosane	394	C28H58	000630-02-4	87
3		Hexatriacontane	507	C36H74	000630-06-8	87
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
5		Octadecane	254	C18H38	000593-45-3	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM110519\
Data File : BM023601.D
Acq On : 05 Nov 2019 11:13
Operator : JU
Sample : K5536-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propanoic acid, 1...	3.59	5.1	ng/ul	559579	1	7.60	2202160	20.0
Isobutyl acetate	3.84	2.5	ng/ul	272919	1	7.60	2202160	20.0
Propanoic acid, 2...	4.11	11.8	ng/ul	1294240	1	7.60	2202160	20.0
Propanoic acid, p...	4.29	17.1	ng/ul	1887880	1	7.60	2202160	20.0
Butanoic acid, 1-...	4.73	21.9	ng/ul	2405660	1	7.60	2202160	20.0
Butanoic acid, pr...	5.59	3.4	ng/ul	369874	1	7.60	2202160	20.0
(DEL) Alkane: Cyc...	20.94	3.7	ng/ul	1089590	5	21.20	5881840	20.0
(DEL) Alkane: Str...	21.38	2.3	ng/ul	677998	5	21.20	5881840	20.0
(DEL) Alkane: Str...	21.81	3.2	ng/ul	938130	5	21.20	5881840	20.0
(DEL) Alkane: Str...	22.26	2.2	ng/ul	652756	5	21.20	5881840	20.0