

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX122718\
 Data File : VX006845.D
 Acq On : 27 Dec 2018 16:29
 Operator : JC/MD
 Sample : J6571-02
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 10827

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % 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:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M

Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.489	51	55	64	rBV	43563	57238	1.33%	0.102%
2	1.647	67	81	85	rVB6	52555	188842	4.39%	0.336%
3	2.446	206	212	223	rBV	436920	721372	16.77%	1.284%
4	2.861	275	280	287	rVV	23916	44385	1.03%	0.079%
5	5.500	700	713	726	rBV	351138	1000021	23.25%	1.780%
6	5.665	728	740	758	rVB	635306	1795353	41.74%	3.196%
7	6.067	794	806	818	rBV	358943	946093	22.00%	1.684%
8	6.458	862	870	880	rBV2	26597	72126	1.68%	0.128%
9	6.860	925	936	953	rBV	938509	2211231	51.41%	3.936%
10	7.549	1044	1049	1054	rBV	29530	55238	1.28%	0.098%
11	7.732	1074	1079	1083	rVV	25090	44239	1.03%	0.079%
12	7.878	1096	1103	1124	rBV	2488691	4301066	100.00%	7.656%
13	8.616	1214	1224	1234	rBV	420264	774409	18.01%	1.379%
14	8.713	1234	1240	1246	rVV	1941732	3048403	70.88%	5.426%
15	8.786	1246	1252	1263	rVB	1301229	2026857	47.12%	3.608%
16	9.018	1286	1290	1298	rBV	47493	78888	1.83%	0.140%
17	9.103	1298	1304	1310	rVB	26671	43539	1.01%	0.078%
18	9.311	1330	1338	1344	rBV	184071	269514	6.27%	0.480%
19	9.372	1344	1348	1350	rVV2	36307	51309	1.19%	0.091%
20	9.408	1350	1354	1357	rVV	109835	154609	3.59%	0.275%
21	9.439	1357	1359	1365	rVB	34540	45457	1.06%	0.081%
22	9.542	1371	1376	1381	rBV	1911678	2439663	56.72%	4.343%
23	9.579	1381	1382	1388	rVV3	96052	125115	2.91%	0.223%
24	9.640	1388	1392	1401	rVB	149403	209392	4.87%	0.373%
25	9.969	1440	1446	1456	rBV	1065169	1338968	31.13%	2.384%
26	10.109	1464	1469	1476	rBV	1970269	2642457	61.44%	4.704%
27	10.250	1485	1492	1498	rBV	395726	516104	12.00%	0.919%
28	10.359	1503	1510	1516	rVV	1505454	2126047	49.43%	3.785%
29	10.414	1516	1519	1528	rVB2	40174	57577	1.34%	0.102%
30	10.695	1560	1565	1571	rBV	1199587	1579316	36.72%	2.811%
31	10.829	1580	1587	1592	rBV4	25680	43897	1.02%	0.078%
32	10.914	1596	1601	1605	rBV2	43194	59387	1.38%	0.106%
33	11.018	1612	1618	1622	rBV	92751	118538	2.76%	0.211%
34	11.134	1632	1637	1648	rVB	1692611	2227559	51.79%	3.965%

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Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
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 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

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 Title : SW846 8260

35	11.359	1667	1674	1681	rBV	207625	261506	6.08%	0.466%
36	11.432	1681	1686	1693	rBV	789775	1354783	31.50%	2.412%
37	11.505	1693	1698	1701	rBV	316560	388719	9.04%	0.692%
38	11.664	1720	1724	1731	rVB	550963	712914	16.58%	1.269%
39	11.804	1743	1747	1757	rVB	1590969	1905217	44.30%	3.391%
40	11.944	1763	1770	1774	rBV	130506	188505	4.38%	0.336%
41	11.999	1774	1779	1780	rVV4	39266	70060	1.63%	0.125%
42	12.024	1780	1783	1786	rVB	114591	130113	3.03%	0.232%
43	12.079	1786	1792	1797	rBV	2215905	2731721	63.51%	4.863%
44	12.139	1797	1802	1809	rVB	750873	979611	22.78%	1.744%
45	12.274	1819	1824	1826	rBV	1421979	1896227	44.09%	3.375%
46	12.298	1826	1828	1836	rVV2	1086288	1609931	37.43%	2.866%
47	12.371	1836	1840	1845	rVB3	389054	595575	13.85%	1.060%
48	12.475	1851	1857	1860	rVB2	144635	221042	5.14%	0.393%
49	12.517	1860	1864	1870	rVB	202010	252927	5.88%	0.450%
50	12.585	1870	1875	1877	rBV	254453	323560	7.52%	0.576%
51	12.609	1877	1879	1882	rVV2	239110	259359	6.03%	0.462%
52	12.664	1885	1888	1892	rVV	310247	398888	9.27%	0.710%
53	12.700	1892	1894	1898	rVB	134136	133655	3.11%	0.238%
54	12.755	1898	1903	1911	rBV2	381409	651007	15.14%	1.159%
55	12.902	1924	1927	1931	rVB	166688	217044	5.05%	0.386%
56	12.975	1931	1939	1942	rBV	187654	246856	5.74%	0.439%
57	13.017	1942	1946	1953	rVB	330685	421859	9.81%	0.751%
58	13.097	1953	1959	1962	rBV5	65036	133616	3.11%	0.238%
59	13.194	1972	1975	1976	rBV	114256	116615	2.71%	0.208%
60	13.225	1976	1980	1984	rVB	460829	563575	13.10%	1.003%
61	13.267	1984	1987	1990	rVB2	52111	58307	1.36%	0.104%
62	13.316	1990	1995	1996	rBV3	100321	139718	3.25%	0.249%
63	13.347	1996	2000	2005	rVB2	999746	1305405	30.35%	2.324%
64	13.481	2017	2022	2026	rBV	674392	867193	20.16%	1.544%
65	13.560	2031	2035	2038	rBV2	54475	62783	1.46%	0.112%
66	13.603	2038	2042	2044	rBV	92109	116138	2.70%	0.207%
67	13.645	2044	2049	2053	rBV4	188354	274529	6.38%	0.489%
68	13.718	2058	2061	2068	rVV2	126445	190025	4.42%	0.338%
69	13.804	2070	2075	2077	rVV	1125502	1370432	31.86%	2.440%
70	13.828	2077	2079	2088	rVV2	953872	1404843	32.66%	2.501%
71	13.907	2088	2092	2097	rVV	230914	336498	7.82%	0.599%

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Integration Parameters: RTEINT.P

Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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72	13.981	2098	2104	2109	rVV2	292549	449479	10.45%	0.800%
73	14.023	2109	2111	2116	rVV	54992	68792	1.60%	0.122%
74	14.133	2125	2129	2132	rBV	95732	112930	2.63%	0.201%
75	14.285	2147	2154	2161	rVB2	245821	421155	9.79%	0.750%
76	14.432	2174	2178	2182	rVB2	69472	87951	2.04%	0.157%
77	14.535	2191	2195	2205	rVB2	142469	200882	4.67%	0.358%
78	14.694	2214	2221	2225	rBV	256434	380029	8.84%	0.676%
79	14.840	2241	2245	2249	rVB	118061	150267	3.49%	0.267%

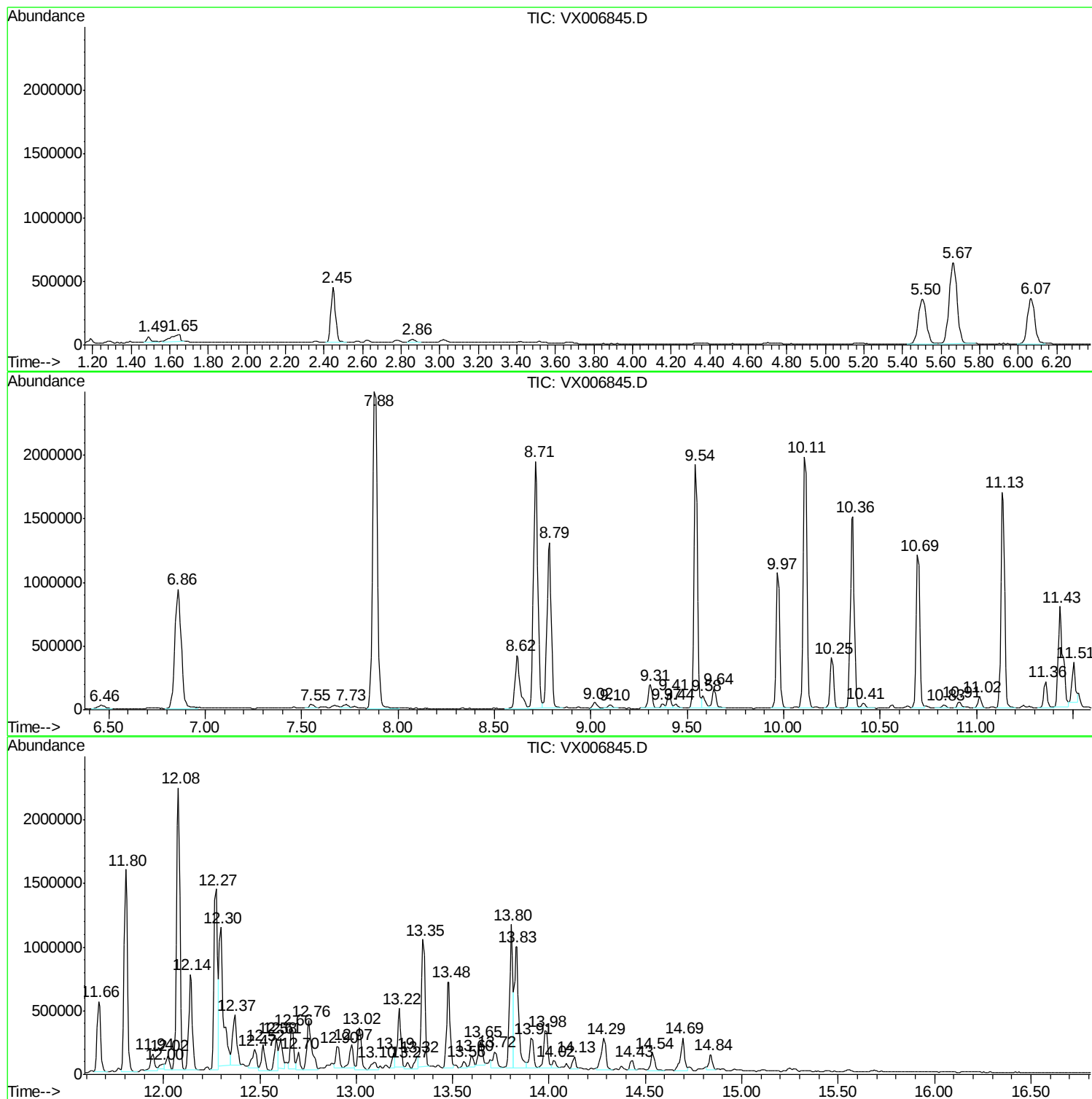
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Instrument :
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ClientSampled :
10827

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260

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



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Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
10827

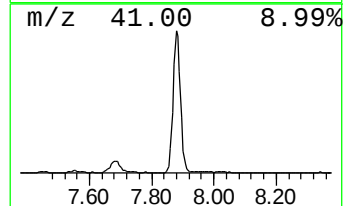
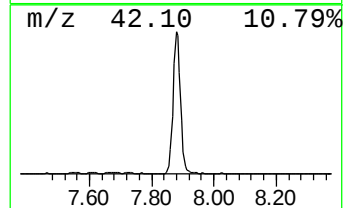
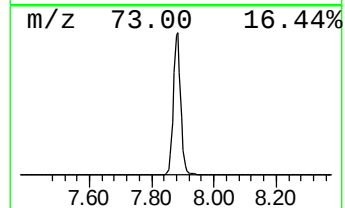
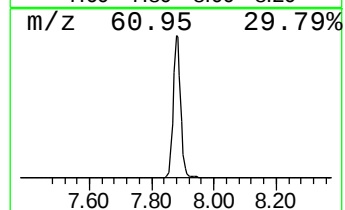
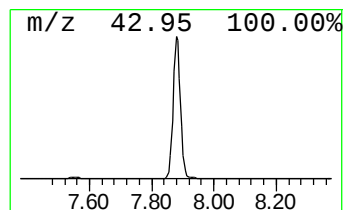
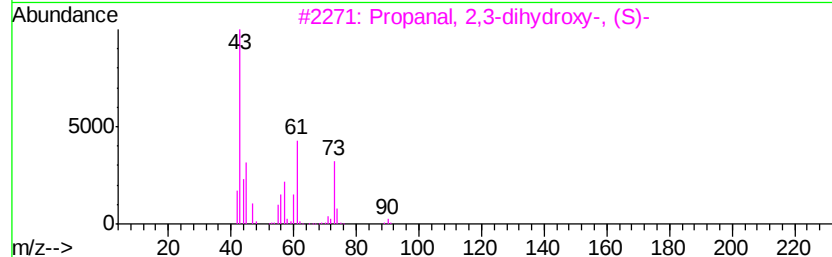
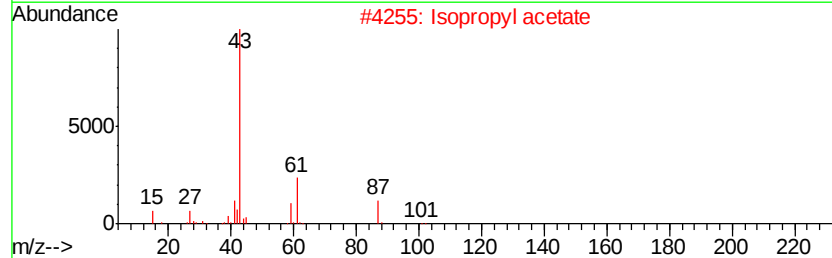
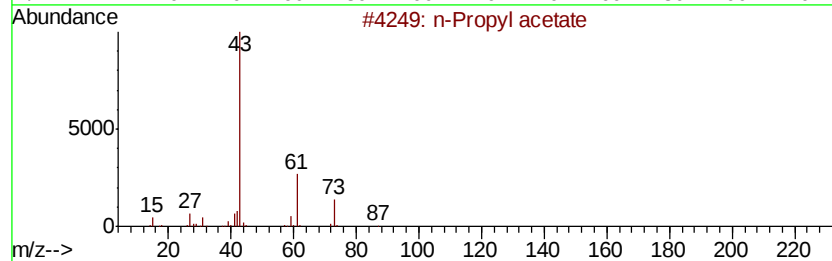
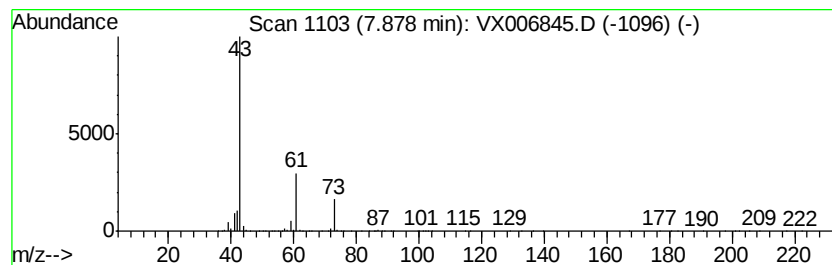
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260

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

Peak Number 1 n-Propyl acetate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.88	97.26 ug/l	4301070	1,4-Difluorobenzene	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Propyl acetate	102	C5H10O2	000109-60-4	83
2		Isopropyl acetate	102	C5H10O2	000108-21-4	9
3		Propanal, 2,3-dihydroxy-, (S)-	90	C3H6O3	000497-09-6	9
4		1-Butanamine	73	C4H11N	000109-73-9	7
5		Isobutylamine	73	C4H11N	000078-81-9	4



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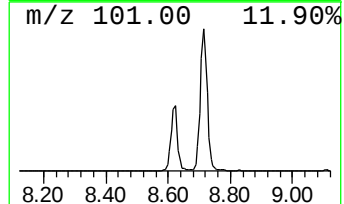
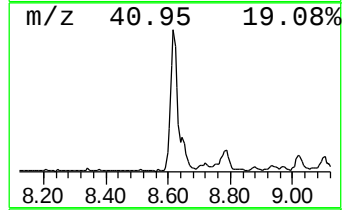
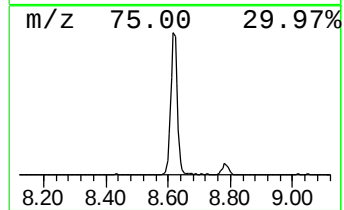
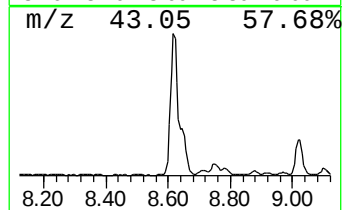
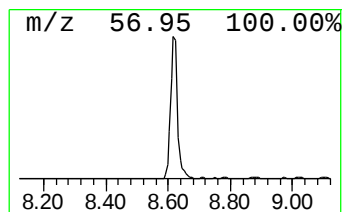
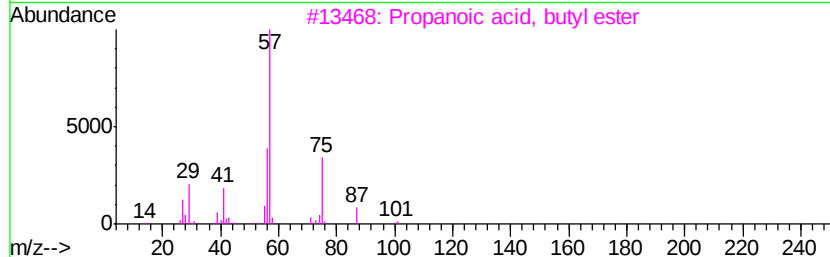
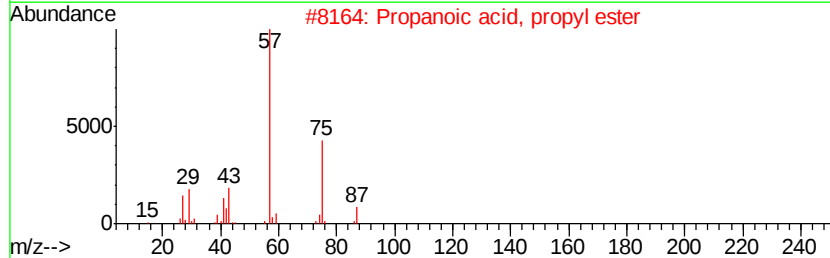
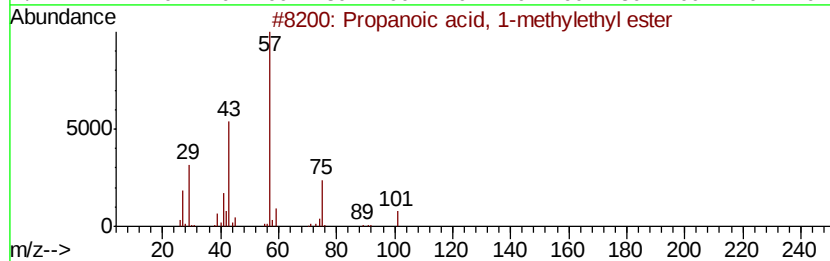
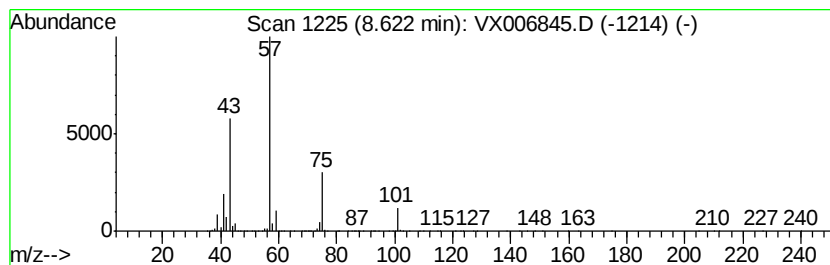
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Propanoic acid, 1-methyleth... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.62	14.65 ug/l	774409	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	90
2		Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	50
3		Propanoic acid, butyl ester	130	C7H14O2	000590-01-2	25
4		Propane, 1-(1-methylethoxy)-	102	C6H14O	000627-08-7	23
5		Butanoic acid, methyl ester	102	C5H10O2	000623-42-7	9



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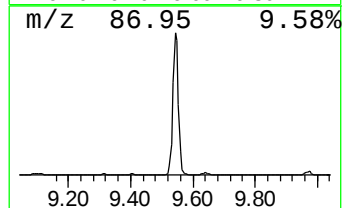
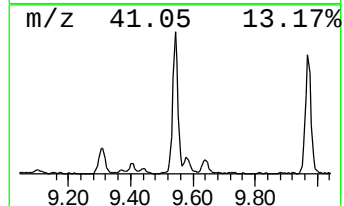
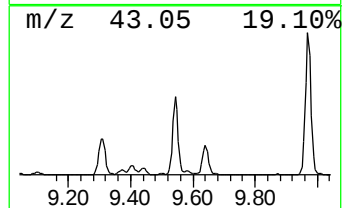
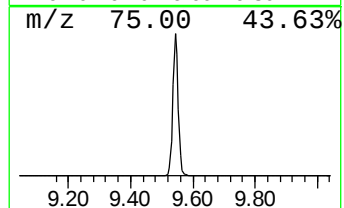
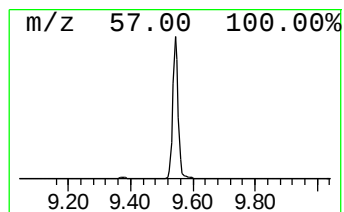
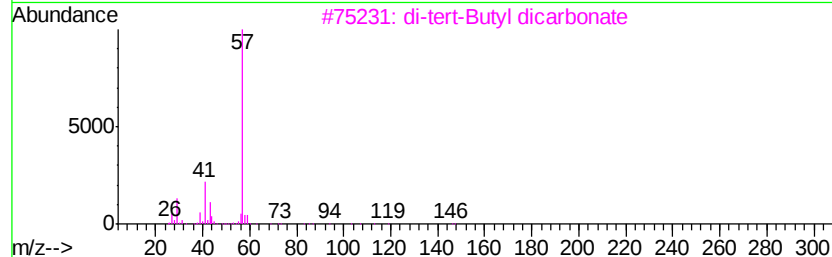
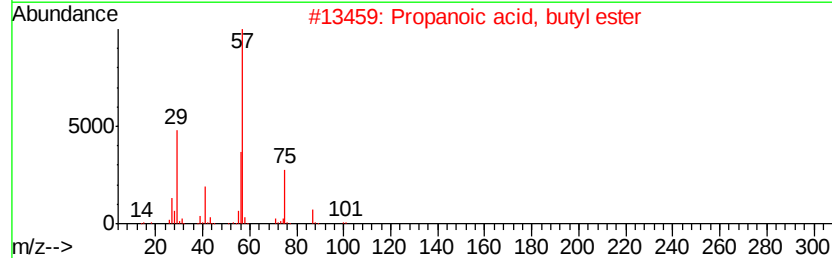
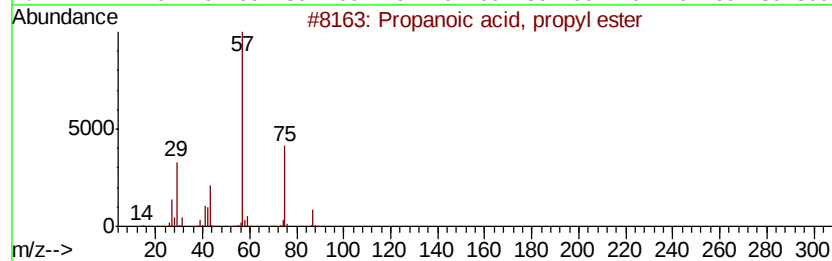
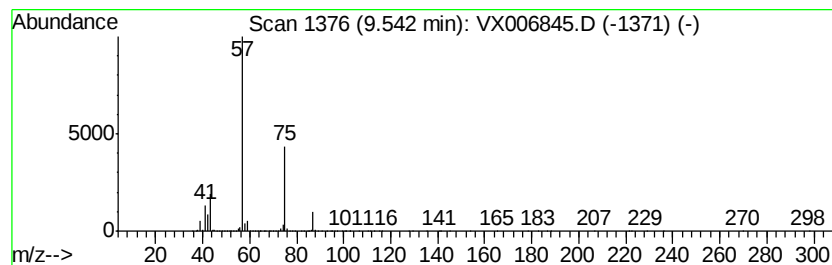
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.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.
9.54	46.16 ug/l	2439660	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
2		Propanoic acid, butyl ester	130	C7H14O2	000590-01-2	28
3		di-tert-Butyl dicarbonate	218	C10H18O5	024424-99-5	9
4		Acetaldehyde, O-ethyloxime-	87	C4H9NO	1000288-34-6	7
5		1-Butanamine, 2-methyl-	87	C5H13N	000096-15-1	7



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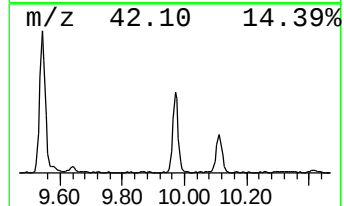
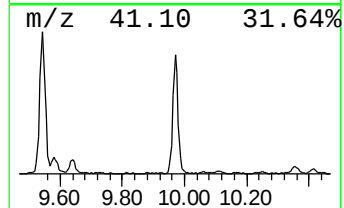
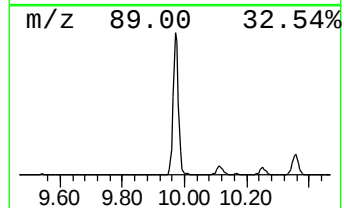
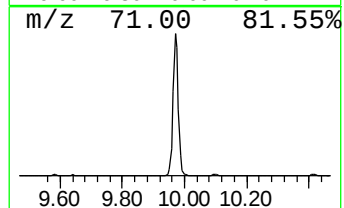
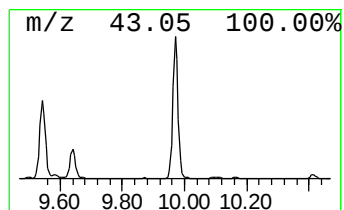
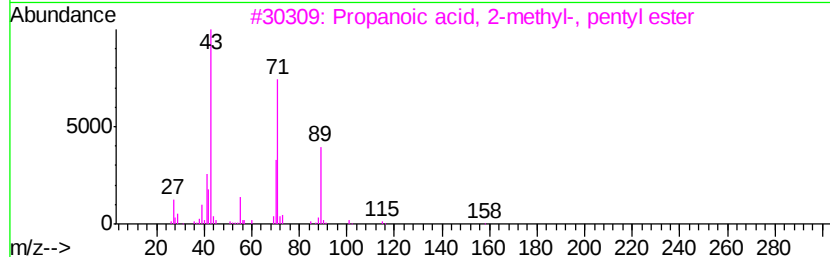
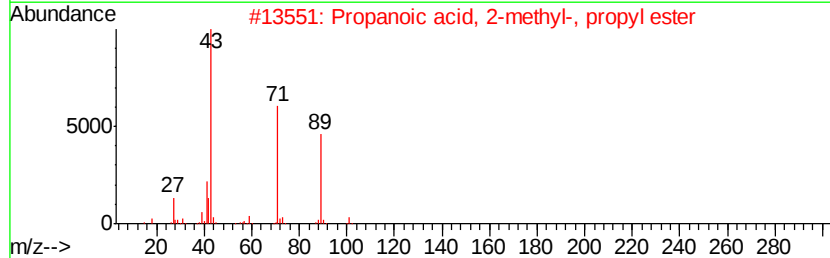
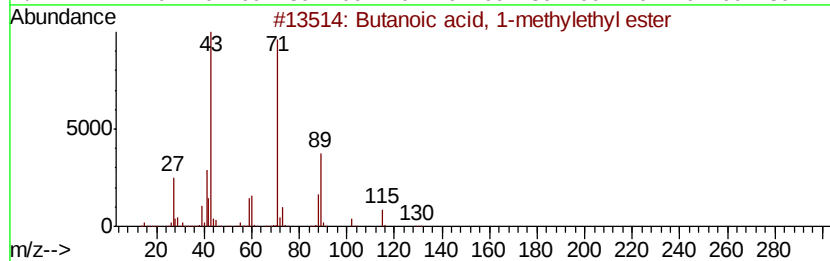
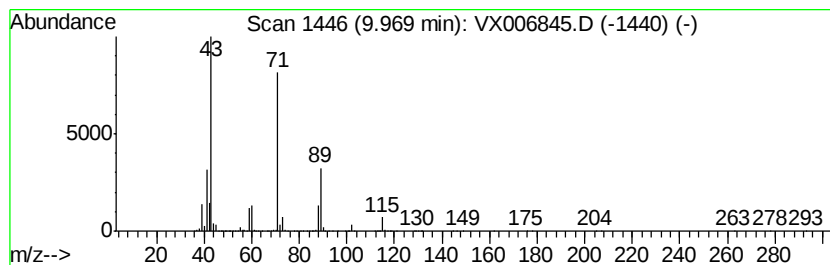
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Quant Title : SW846 8260

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

Peak Number 4 Butanoic acid, 1-methylethy... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.97	25.34 ug/l	1338970	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	91
2		Propanoic acid, 2-methyl-, propyl...	130	C7H14O2	000644-49-5	72
3		Propanoic acid, 2-methyl-, pentyl...	158	C9H18O2	002445-72-9	64
4		Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	56
5		Ethane-1,1-diol dibutanoate	202	C10H18O4	025572-25-2	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122718\
Data File : VX006845.D
Acq On : 27 Dec 2018 16:29
Operator : JC/MD
Sample : J6571-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
10827

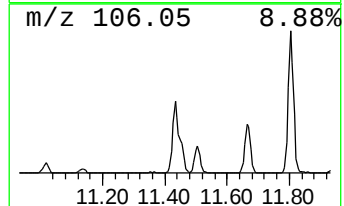
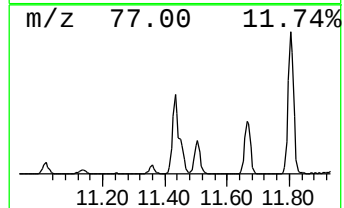
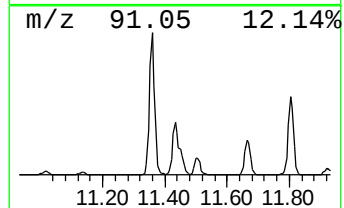
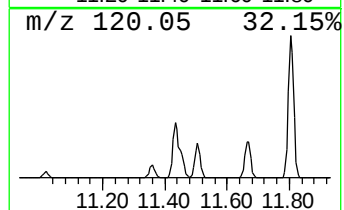
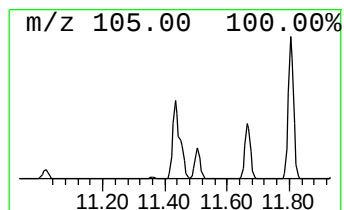
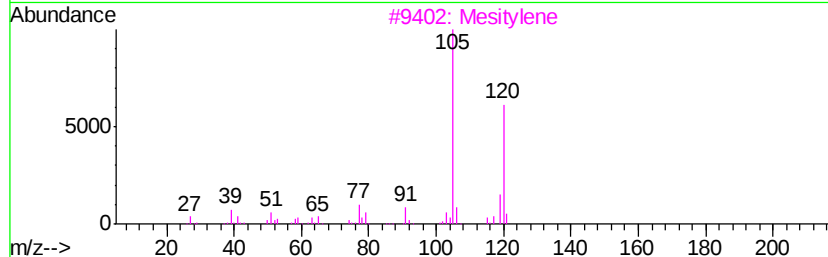
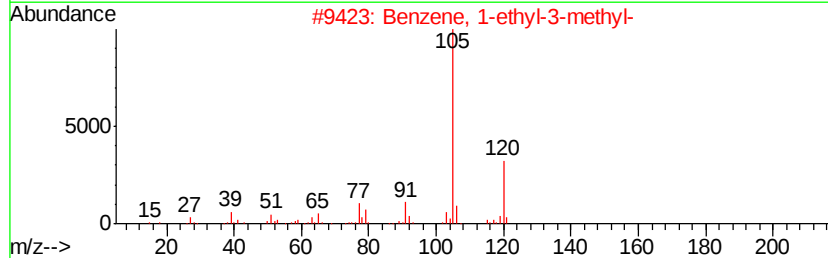
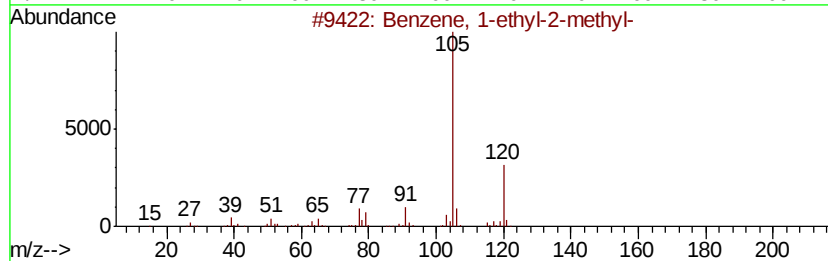
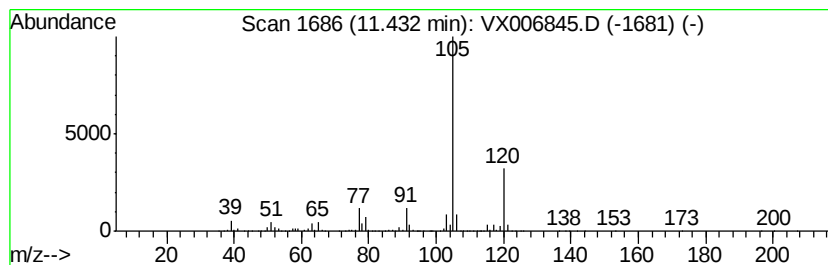
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.43	24.80 ug/l	1354780	1,4-Dichlorobenzene-d4	12.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3		Mesitylene	120	C9H12	000108-67-8	91
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122718\
Data File : VX006845.D
Acq On : 27 Dec 2018 16:29
Operator : JC/MD
Sample : J6571-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
10827

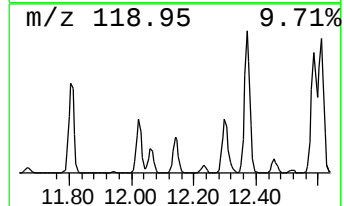
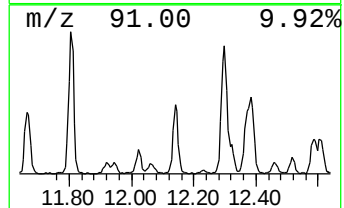
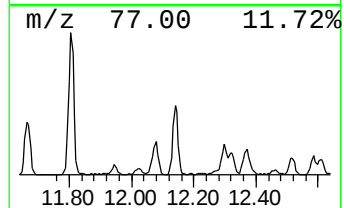
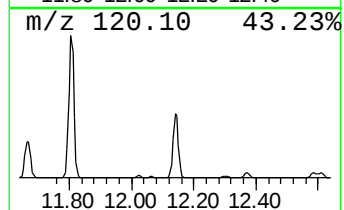
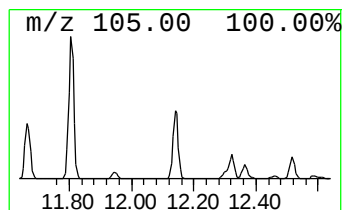
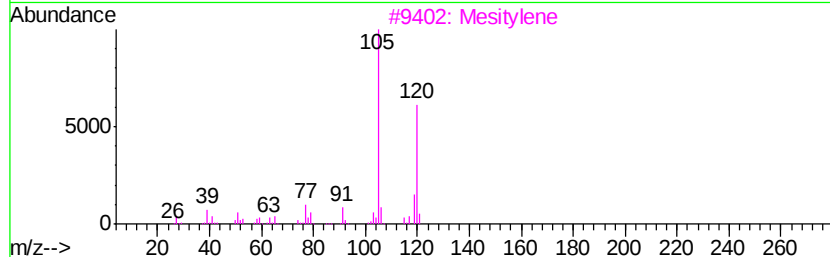
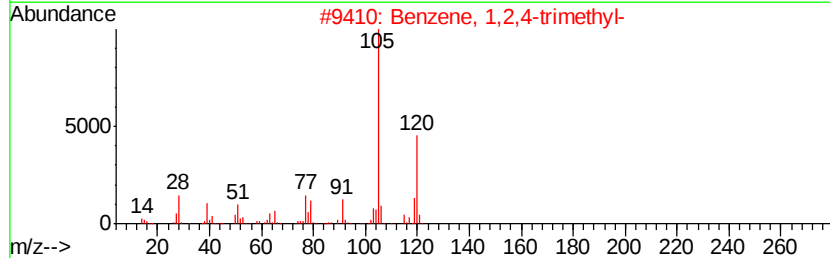
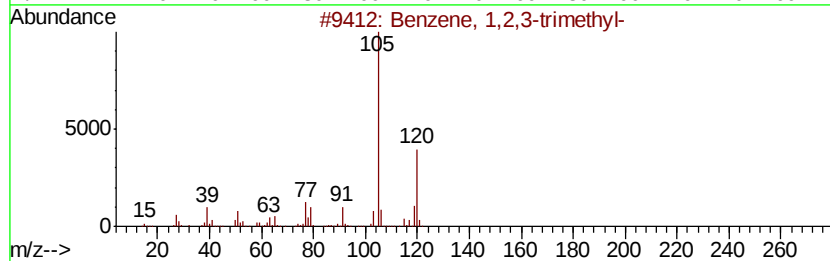
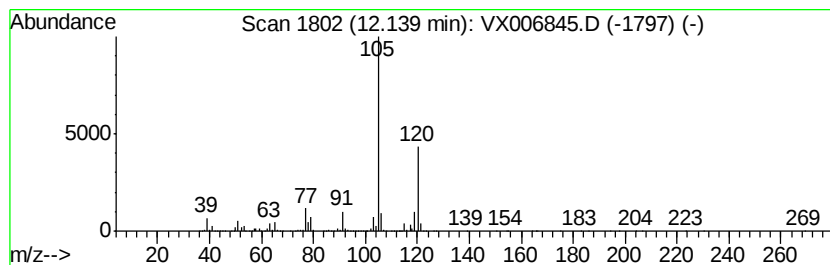
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	17.93 ug/l	979611	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3		Mesitylene	120	C9H12	000108-67-8	93
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX122718\
Data File : VX006845.D
Acq On : 27 Dec 2018 16:29
Operator : JC/MD
Sample : J6571-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
10827

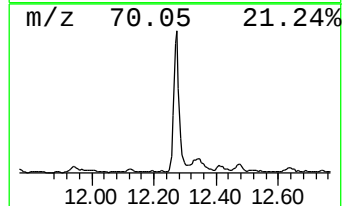
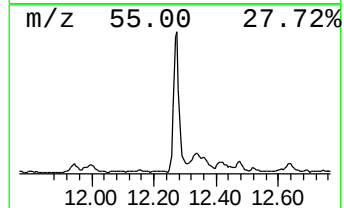
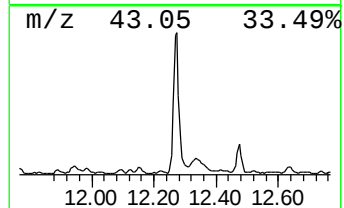
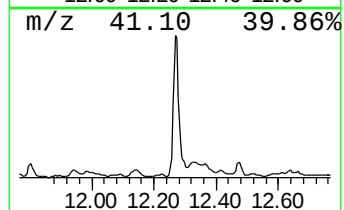
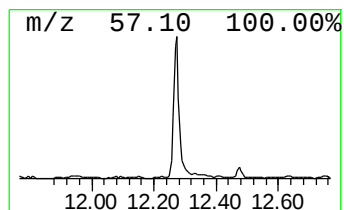
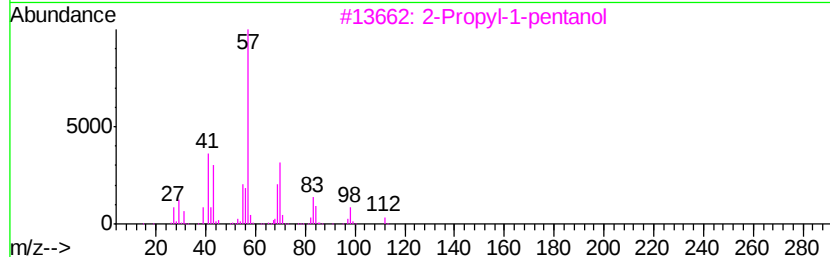
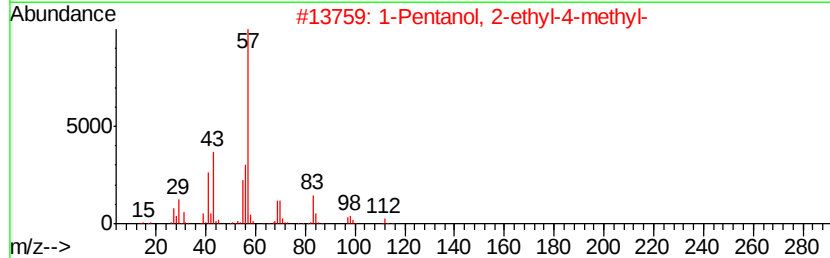
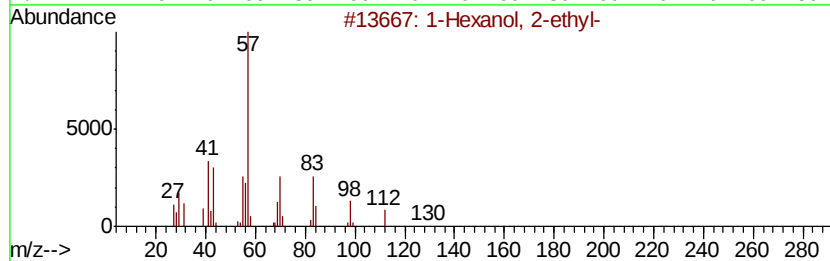
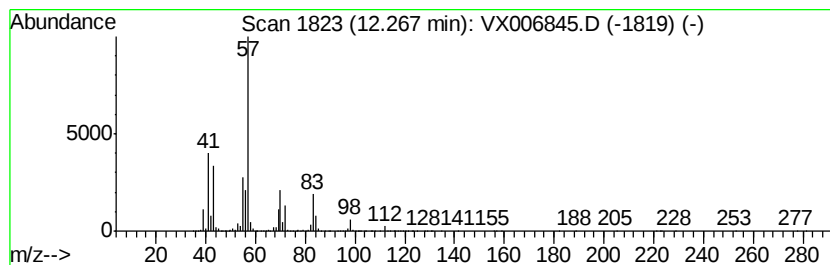
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Quant Title : SW846 8260

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

Peak Number 7 1-Hexanol, 2-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.27	34.71 ug/l	1896230	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
2		1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	53
3		2-Propyl-1-pentanol	130	C8H18O	058175-57-8	47
4		Heptyl isobutyl carbonate	216	C12H24O3	959068-08-7	43
5		Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122718\
Data File : VX006845.D
Acq On : 27 Dec 2018 16:29
Operator : JC/MD
Sample : J6571-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

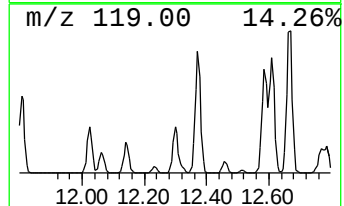
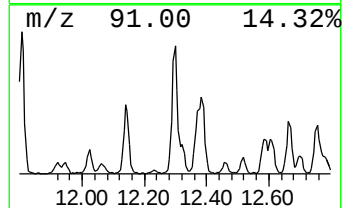
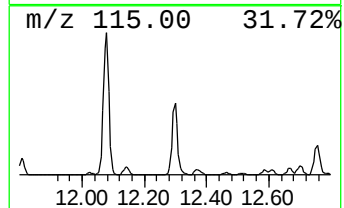
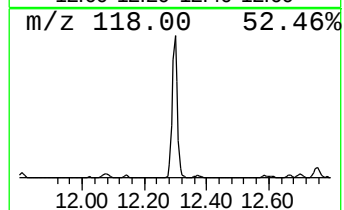
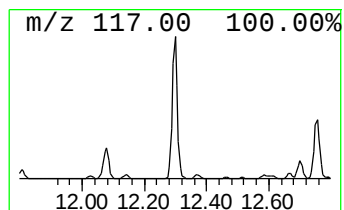
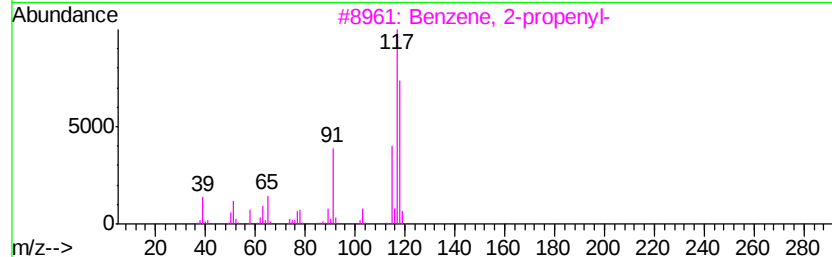
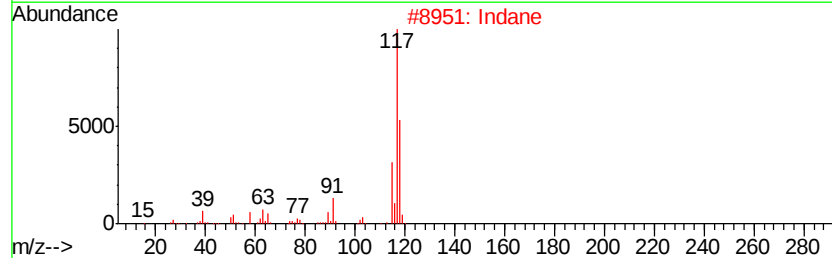
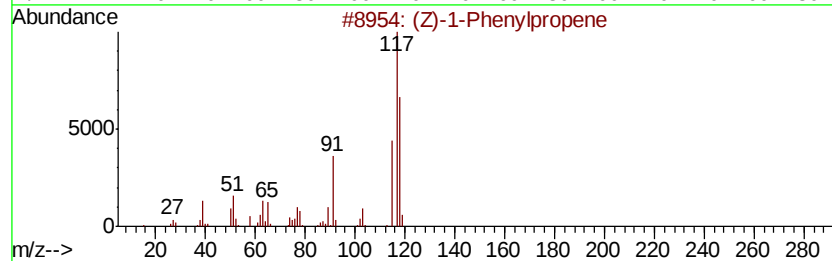
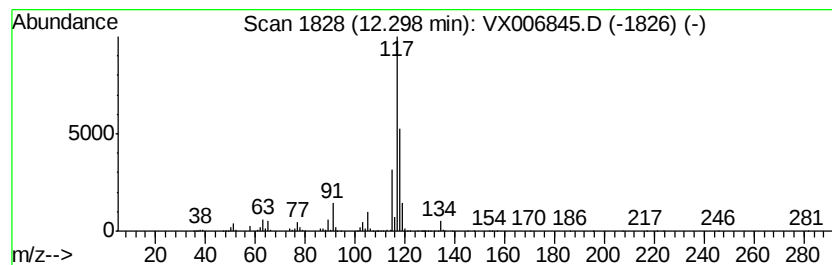
Instrument :
MSVOA_X
ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260

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

Peak Number 8 (Z)-1-Phenylpropene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	29.47 ug/l	1609930	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		(Z)-1-Phenylpropene	118 C9H10	000766-90-5 76
2		Indane	118 C9H10	000496-11-7 68
3		Benzene, 2-propenyl-	118 C9H10	000300-57-2 64
4		Benzene, cyclopropyl-	118 C9H10	000873-49-4 64
5		Benzeneethanol, .beta.-ethenyl-	148 C10H12O	006052-63-7 50



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX122718\
Data File : VX006845.D
Acq On : 27 Dec 2018 16:29
Operator : JC/MD
Sample : J6571-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
10827

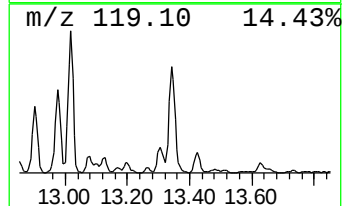
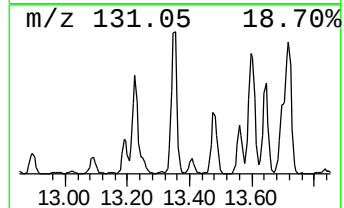
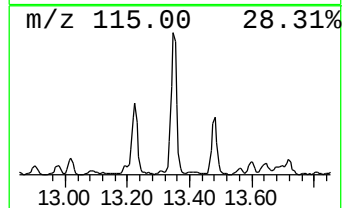
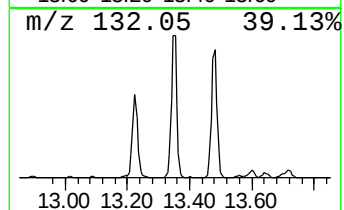
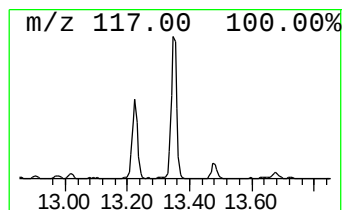
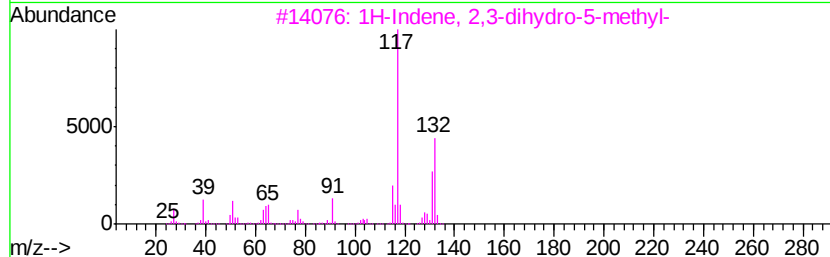
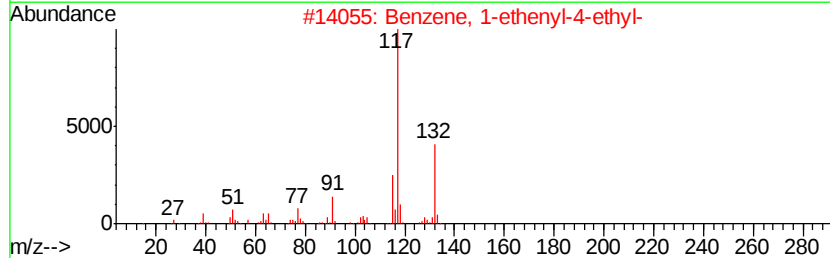
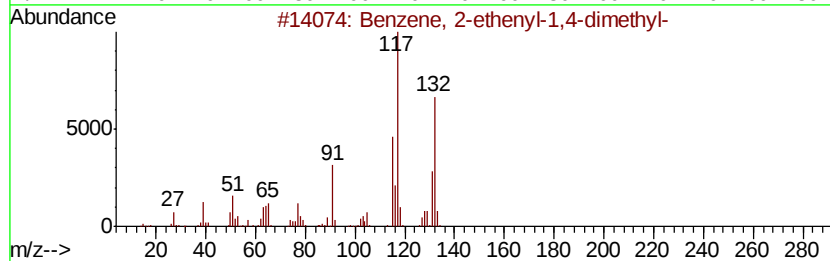
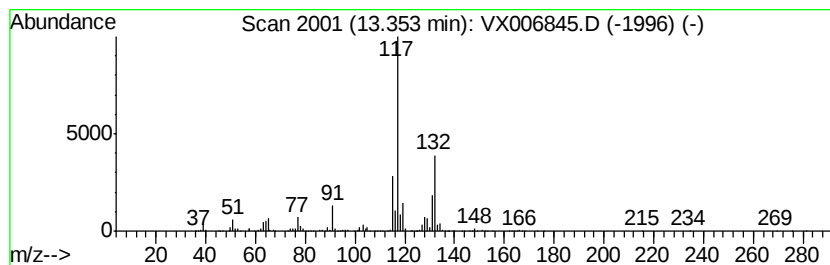
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	23.89 ug/l	1305410	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	90
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	89
3		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	89
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122718\
Data File : VX006845.D
Acq On : 27 Dec 2018 16:29
Operator : JC/MD
Sample : J6571-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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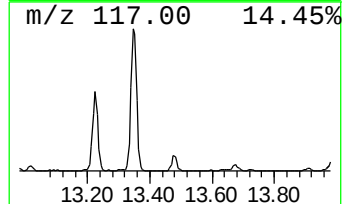
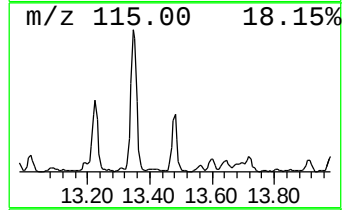
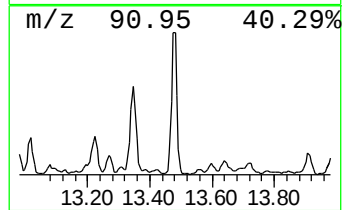
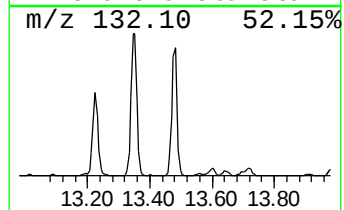
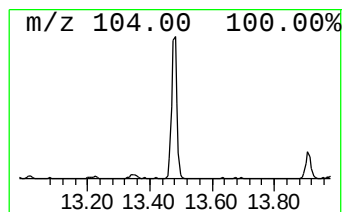
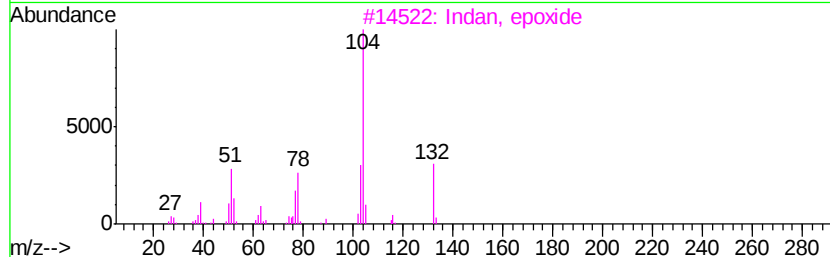
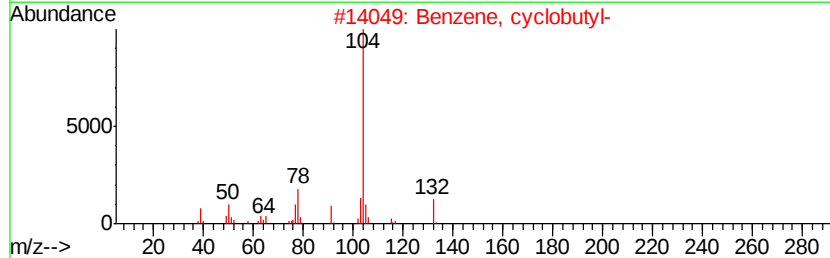
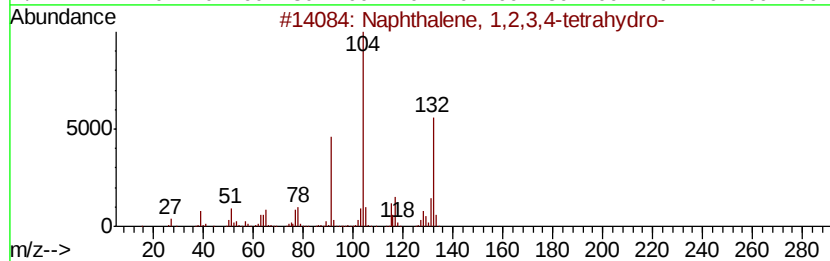
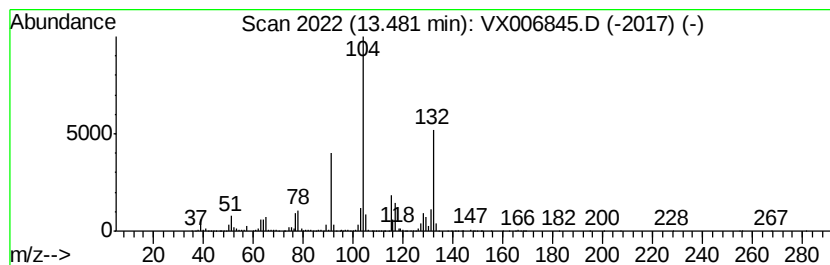
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetrahydro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.48	15.87 ug/l	867193	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	93
2		Benzene, cyclobutyl-	132	C10H12	004392-30-7	53
3		Indan, epoxide	132	C9H8O	000768-22-9	53
4		Benzene, 3-cyclohexen-1-yl-	158	C12H14	004994-16-5	47
5		2-Propenoic acid, 2-phenylethyl ...	176	C11H12O2	003530-36-7	38



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX122718\
Data File : VX006845.D
Acq On : 27 Dec 2018 16:29
Operator : JC/MD
Sample : J6571-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
10827

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
n-Propyl acetate	7.88	97.3	ug/l	4301070	2	6.86	2211230	50.0
Propanoic acid, 1...	8.62	14.7	ug/l	774409	3	10.11	2642460	50.0
Propanoic acid, p...	9.54	46.2	ug/l	2439660	3	10.11	2642460	50.0
Butanoic acid, 1-...	9.97	25.3	ug/l	1338970	3	10.11	2642460	50.0
Benzene, 1-ethyl-...	11.43	24.8	ug/l	1354780	4	12.08	2731720	50.0
Benzene, 1,2,3-tr...	12.14	17.9	ug/l	979611	4	12.08	2731720	50.0
1-Hexanol, 2-ethyl-	12.27	34.7	ug/l	1896230	4	12.08	2731720	50.0
(Z)-1-Phenylpropene	12.30	29.5	ug/l	1609930	4	12.08	2731720	50.0
Benzene, 2-etheny...	13.35	23.9	ug/l	1305410	4	12.08	2731720	50.0
Naphthalene, 1,2,...	13.48	15.9	ug/l	867193	4	12.08	2731720	50.0