

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU121118\
 Data File : VU028592.D
 Acq On : 11 Dec 2018 08:29
 Operator : MD/SY
 Sample : J6305-08
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 P001-SW005-01

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_U\METHOD\82U120618W.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	1.276	27	32	43	rVB3	7804	9393	1.63%	0.162%
2	1.386	57	66	73	rBV2	5591	6752	1.17%	0.116%
3	1.472	77	93	99	rBV2	20410	36983	6.43%	0.637%
4	1.514	99	106	113	rBV	61642	61659	10.72%	1.062%
5	1.556	114	119	123	rBV3	7492	9062	1.58%	0.156%
6	2.244	324	333	344	rVB	8123	13475	2.34%	0.232%
7	2.321	345	357	368	rVV	307682	482289	83.82%	8.309%
8	2.373	369	373	384	rVV	14978	22914	3.98%	0.395%
9	2.450	387	397	416	rVV	50860	101671	17.67%	1.752%
10	2.553	418	429	446	rVV	51118	96042	16.69%	1.655%
11	2.836	505	517	536	rVB2	26256	61003	10.60%	1.051%
12	2.942	542	550	564	rBV2	3405	6532	1.14%	0.113%
13	4.199	923	941	951	rBV2	28308	76197	13.24%	1.313%
14	4.263	952	961	976	rVB	38228	86561	15.04%	1.491%
15	4.575	1042	1058	1066	rBV3	8190	20607	3.58%	0.355%
16	4.640	1067	1078	1106	rVB3	40407	98974	17.20%	1.705%
17	4.881	1134	1153	1170	rBV	77859	175348	30.48%	3.021%
18	4.980	1170	1184	1207	rVB	119682	257946	44.83%	4.444%
19	5.305	1271	1285	1299	rBV	96281	199575	34.69%	3.438%
20	5.389	1299	1311	1329	rVB	46358	99171	17.24%	1.708%
21	5.787	1417	1435	1451	rBV3	10014	22267	3.87%	0.384%
22	5.884	1451	1465	1492	rVV	202179	393627	68.41%	6.781%
23	6.421	1615	1632	1645	rBV2	19811	40466	7.03%	0.697%
24	6.614	1685	1692	1713	rVB2	11560	24108	4.19%	0.415%
25	7.459	1942	1955	1969	rBV	36786	64790	11.26%	1.116%
26	7.562	1969	1987	2001	rVV	320893	575363	100.00%	9.912%
27	7.636	2001	2010	2023	rVV	66955	123498	21.46%	2.128%
28	7.704	2023	2031	2044	rVB6	4379	9669	1.68%	0.167%
29	7.910	2084	2095	2113	rBV2	13625	26187	4.55%	0.451%
30	8.231	2184	2195	2203	rVV4	5628	10128	1.76%	0.174%
31	8.286	2203	2212	2226	rVV2	12323	24006	4.17%	0.414%
32	8.369	2226	2238	2253	rVV5	5942	12874	2.24%	0.222%
33	8.459	2256	2266	2289	rVV9	8065	26724	4.64%	0.460%
34	8.575	2289	2302	2333	rVB	99192	178336	31.00%	3.072%

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Integration Parameters: RTEINT.P

Integrator: RTE
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 Sampling : 1
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 Stop Thrs : 0

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35	8.791	2359	2369	2385	rVB2	11290	19806	3.44%	0.341%
36	9.086	2450	2461	2482	rBV	280209	466047	81.00%	8.029%
37	9.250	2501	2512	2523	rBV	30732	50345	8.75%	0.867%
38	9.324	2523	2535	2543	rVV	53038	88406	15.37%	1.523%
39	9.372	2543	2550	2565	rVB	31640	52272	9.09%	0.901%
40	9.495	2578	2588	2598	rVB2	12431	21124	3.67%	0.364%
41	9.604	2610	2622	2632	rBV	15190	24436	4.25%	0.421%
42	9.790	2666	2680	2703	rBV2	115139	199044	34.59%	3.429%
43	9.897	2703	2713	2717	rBV5	7690	12362	2.15%	0.213%
44	10.080	2759	2770	2781	rVB5	10808	19495	3.39%	0.336%
45	10.163	2784	2796	2808	rVB7	6085	12192	2.12%	0.210%
46	10.234	2808	2818	2825	rBV6	3818	6564	1.14%	0.113%
47	10.305	2828	2840	2857	rBV	221009	346277	60.18%	5.966%
48	10.581	2916	2926	2940	rVB8	4567	10297	1.79%	0.177%
49	10.736	2968	2974	2981	rVB5	9699	13143	2.28%	0.226%
50	10.996	3035	3055	3073	rBV7	13293	29715	5.16%	0.512%
51	11.147	3093	3102	3113	rVB2	14931	23764	4.13%	0.409%
52	11.398	3174	3180	3190	rBV5	4116	6099	1.06%	0.105%
53	11.482	3195	3206	3223	rBV	269771	418192	72.68%	7.205%
54	11.597	3223	3242	3265	rVB	46563	89426	15.54%	1.541%
55	11.742	3277	3287	3303	rBV	40224	71358	12.40%	1.229%
56	11.823	3303	3312	3322	rVB8	2912	6821	1.19%	0.118%
57	11.983	3350	3362	3375	rBV4	27873	50096	8.71%	0.863%
58	12.475	3501	3515	3530	rBV	48354	89811	15.61%	1.547%
59	12.671	3567	3576	3580	rBV7	4478	7405	1.29%	0.128%
60	13.089	3695	3706	3713	rBV5	11127	19595	3.41%	0.338%
61	13.186	3726	3736	3748	rBV	32176	48878	8.50%	0.842%
62	13.324	3773	3779	3792	rVV8	5500	11104	1.93%	0.191%
63	13.395	3794	3801	3808	rVB8	3706	6068	1.05%	0.105%
64	13.453	3811	3819	3829	rBV3	6304	9751	1.69%	0.168%
65	13.527	3831	3842	3849	rBV8	3280	6275	1.09%	0.108%
66	13.742	3895	3909	3923	rBV	46408	83523	14.52%	1.439%
67	14.388	4105	4110	4123	rVB8	4297	6723	1.17%	0.116%
68	14.880	4255	4263	4272	rVB3	4625	7457	1.30%	0.128%
69	15.064	4311	4320	4328	rBV2	4451	6457	1.12%	0.111%
70	15.610	4477	4490	4499	rBV7	5307	10054	1.75%	0.173%

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ClientSampleId :
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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

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

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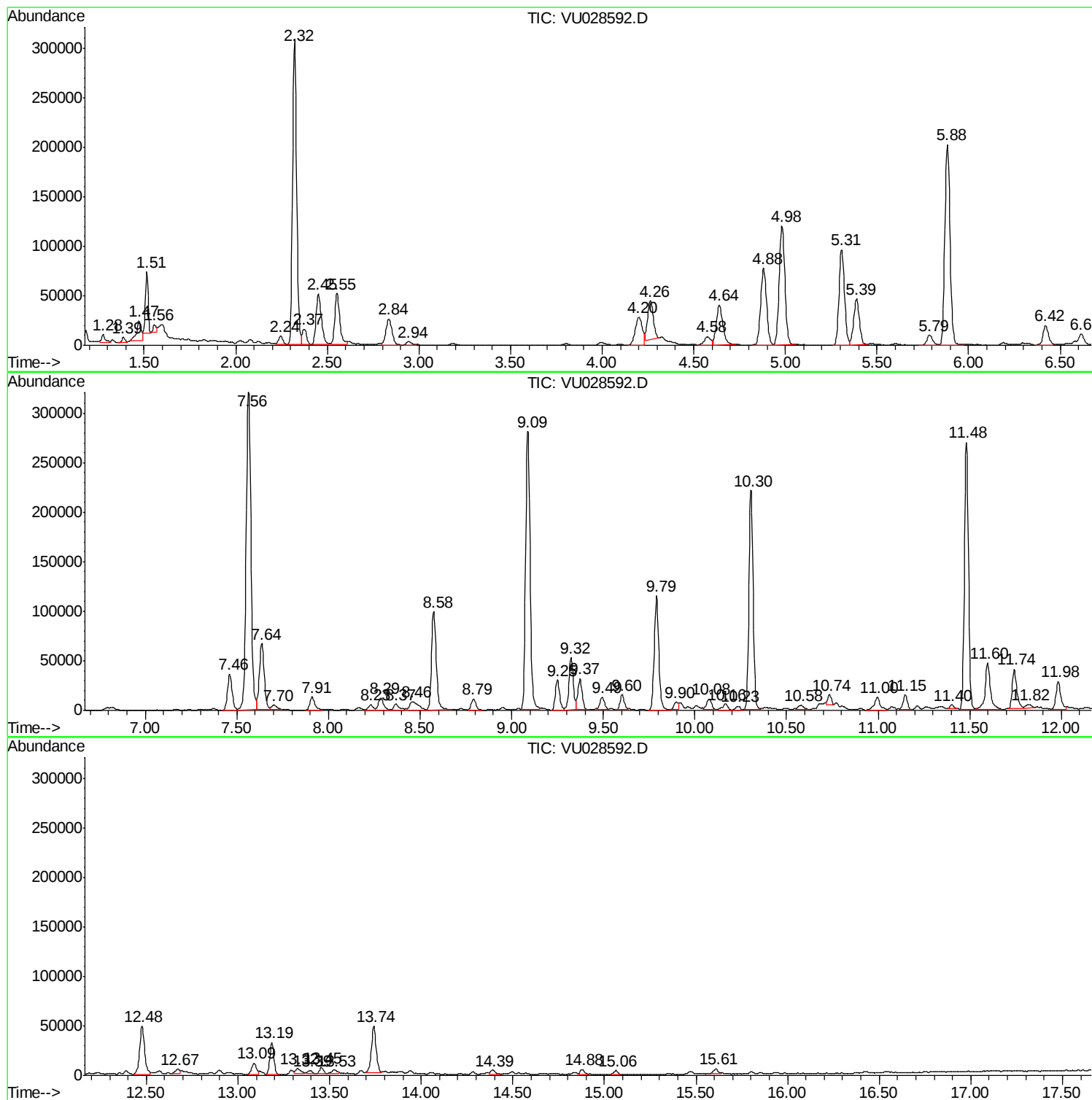
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U120618W.M
Quant Title : SW846 8260

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



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Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 49 Sample Multiplier: 1

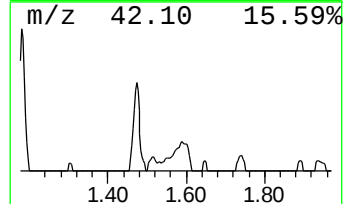
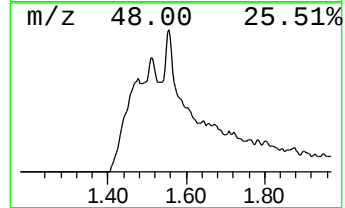
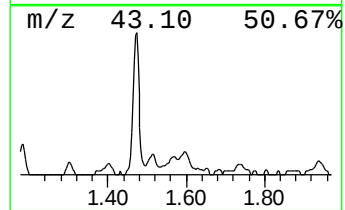
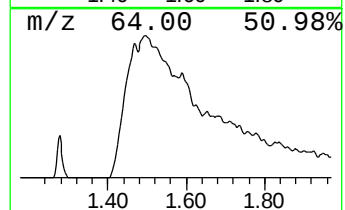
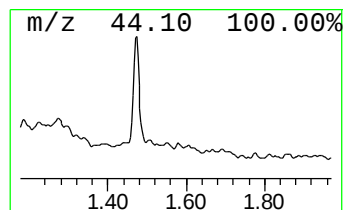
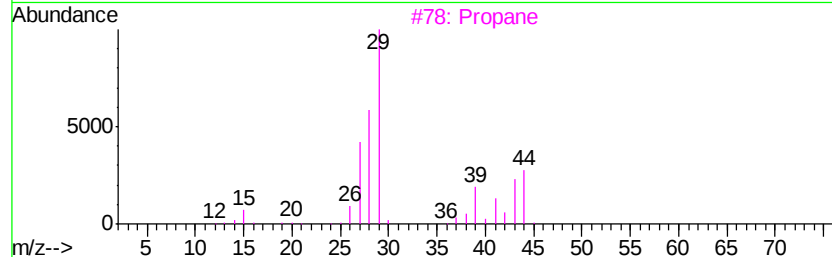
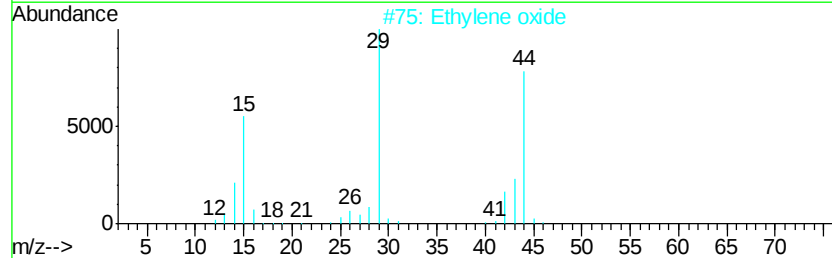
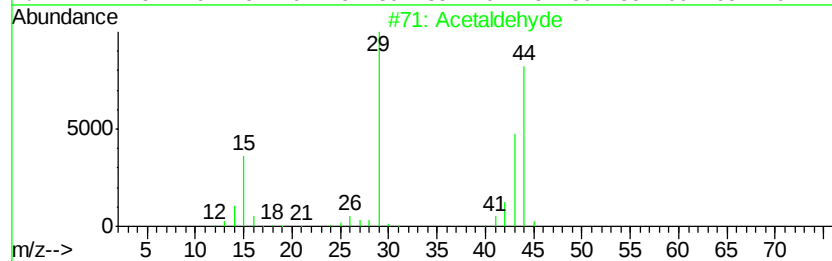
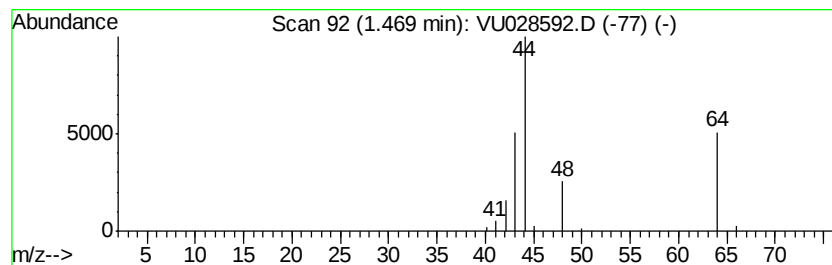
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Quant Title : SW846 8260

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

Peak Number 1 unknown1.47 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
1.47	7.17 ug/l	36983	Pentafluorobenzene	4.98		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Acetaldehyde	44	C2H4O	000075-07-0	12
2		Ethylene oxide	44	C2H4O	000075-21-8	5
3		Propane	44	C3H8	000074-98-6	4
4		1,2-Propanediamine	74	C3H10N2	000078-90-0	4
5		Silane, methyl-	46	CH6Si	000992-94-9	4



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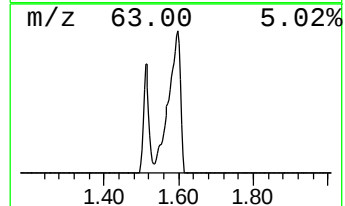
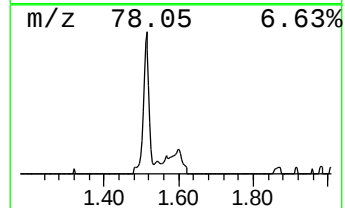
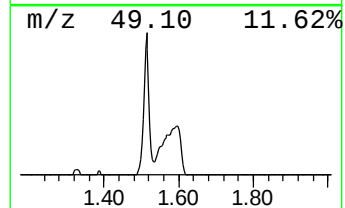
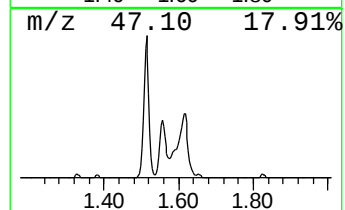
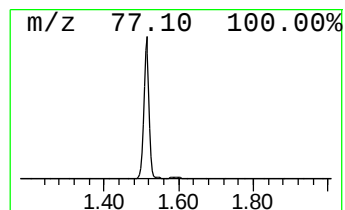
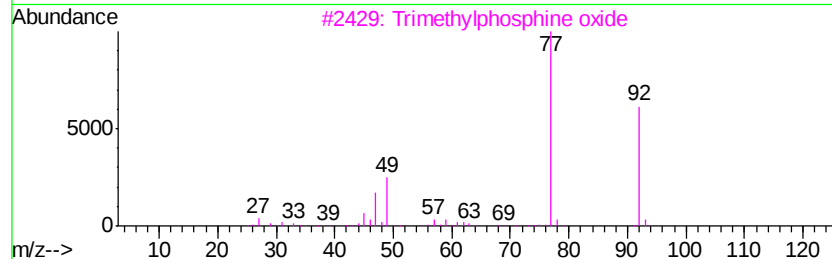
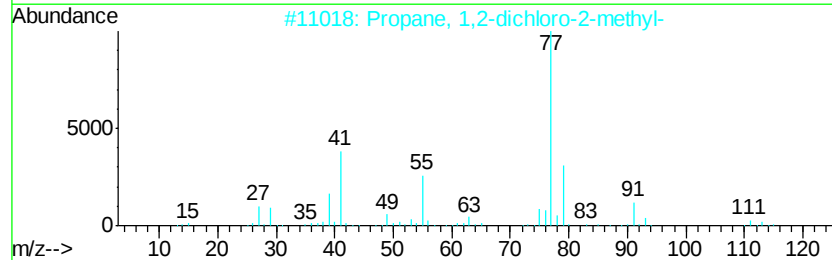
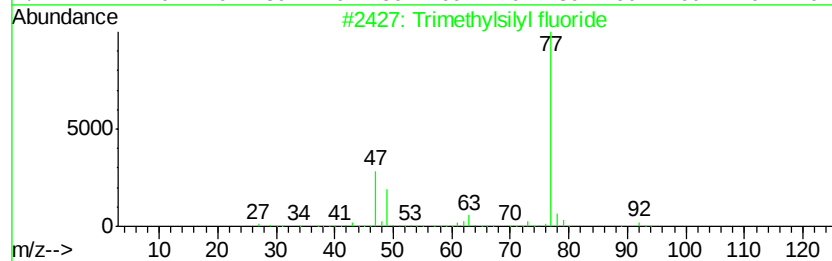
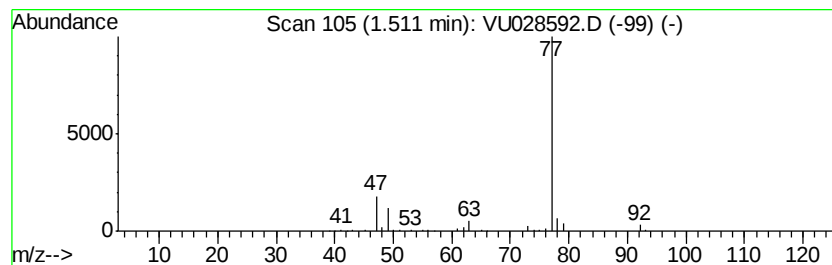
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U120618W.M
Quant Title : SW846 8260

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

Peak Number 2 Trimethylsilyl fluoride Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.51	11.95 ug/l	61659	Pentafluorobenzene	4.98

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trimethylsilyl fluoride	92	C3H9FSi	000420-56-4	91
2			Propane, 1,2-dichloro-2-methyl-	126	C4H8Cl2	000594-37-6	39
3			Trimethylphosphine oxide	92	C3H9OP	000676-96-0	38
4			Propane, 2-chloro-2-nitro-	123	C3H6ClNO2	000594-71-8	4
5			2-Propanone, 1,1,3-trichloro-	160	C3H3Cl3O	000921-03-9	2



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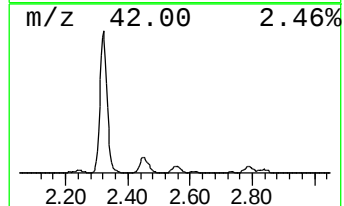
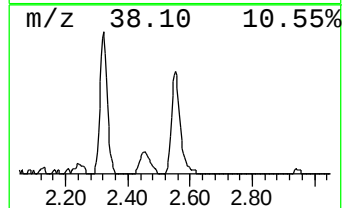
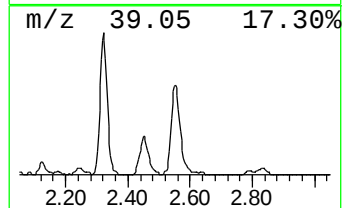
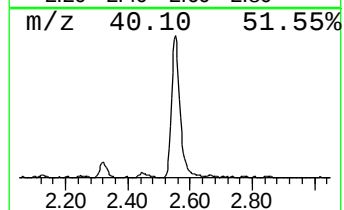
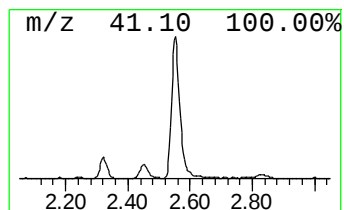
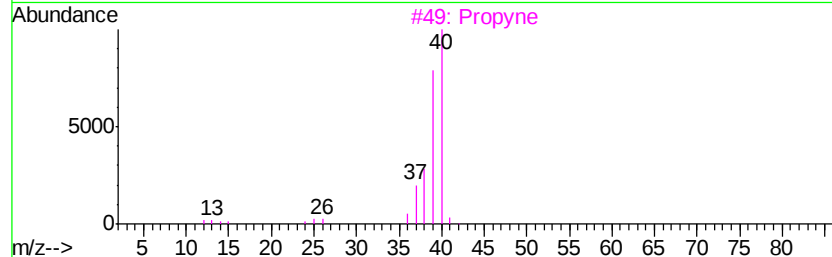
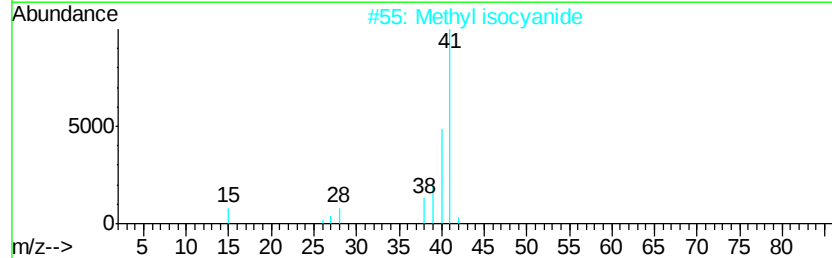
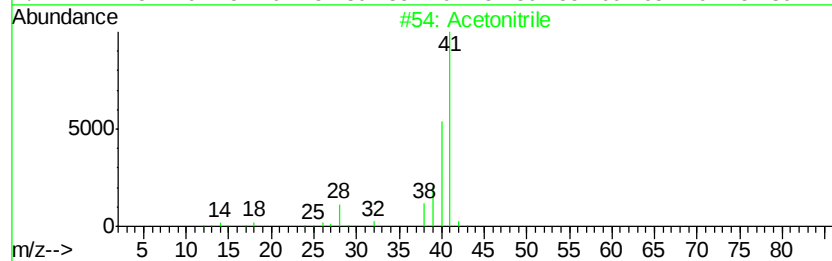
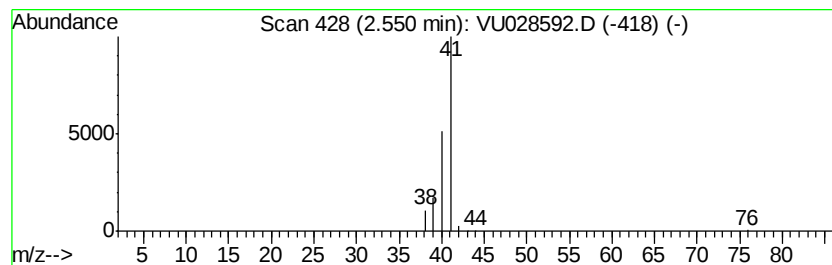
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Acetonitrile Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.55	18.62 ug/l	96042	Pentafluorobenzene	4.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Acetonitrile	41	C2H3N	000075-05-8	56
2		Methyl isocyanide	41	C2H3N	000593-75-9	56
3		Propyne	40	C3H4	000074-99-7	4
4		2,5-Furandione, dihydro-3-methyl...	112	C5H4O3	002170-03-8	2
5		2-Azirane carbonitrile, 1-formyl	96	C4H4N2O	1000196-87-5	2



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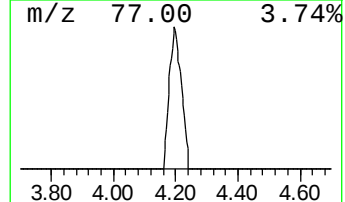
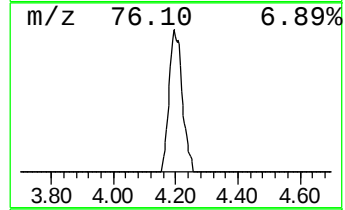
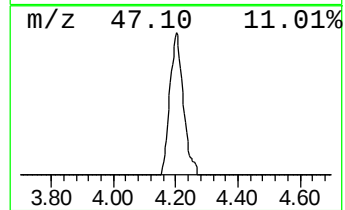
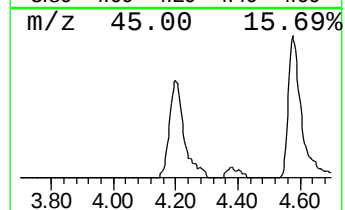
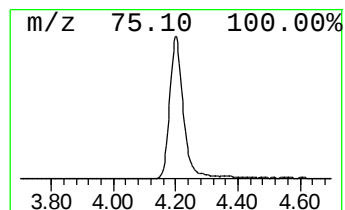
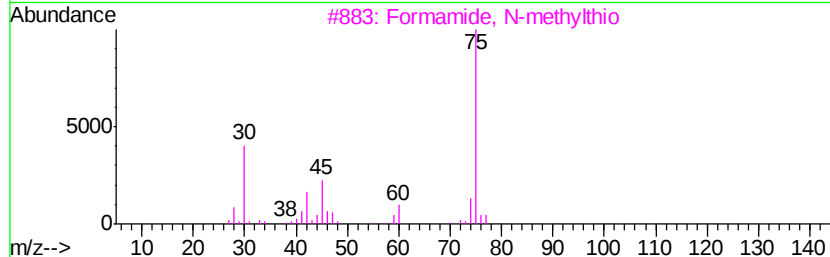
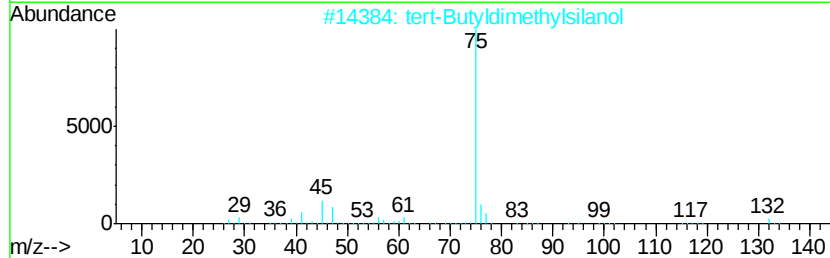
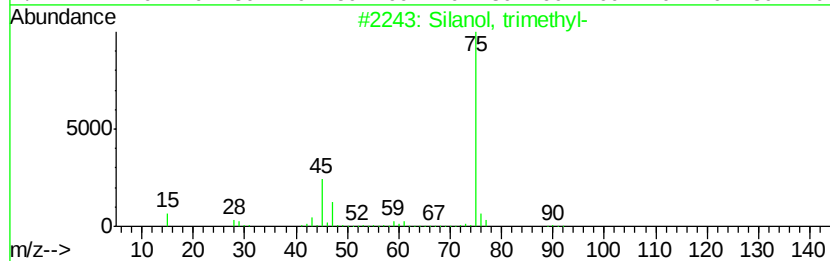
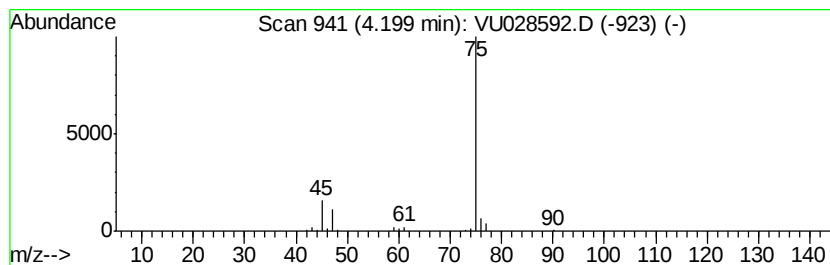
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U120618W.M
Quant Title : SW846 8260

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

Peak Number 4 Silanol, trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.20	14.77 ug/l	76197	Pentafluorobenzene	4.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Silanol, trimethyl-	90	C3H10OSi	001066-40-6	86
2		tert-Butyldimethylsilanol	132	C6H16OSi	018173-64-3	83
3		Formamide, N-methylthio	75	C2H5NS	018952-41-5	72
4		Ethanol, 2-(trimethylsilyl)-	118	C5H14OSi	002916-68-9	43
5		2-Propanol, 1,1-dimethoxy-	120	C5H12O3	042919-42-6	39



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU121118\
Data File : VU028592.D
Acq On : 11 Dec 2018 08:29
Operator : MD/SY
Sample : J6305-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
P001-SW005-01

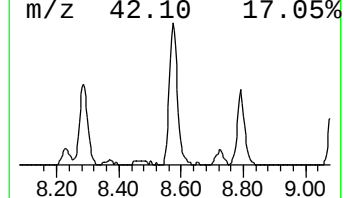
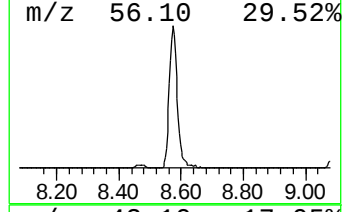
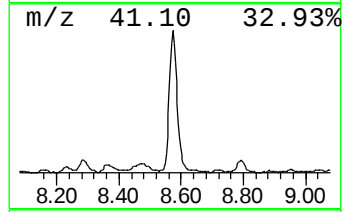
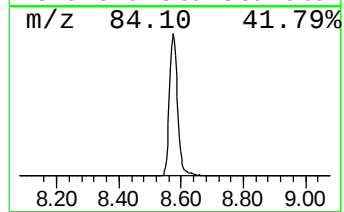
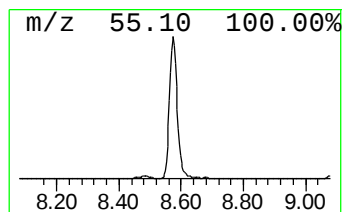
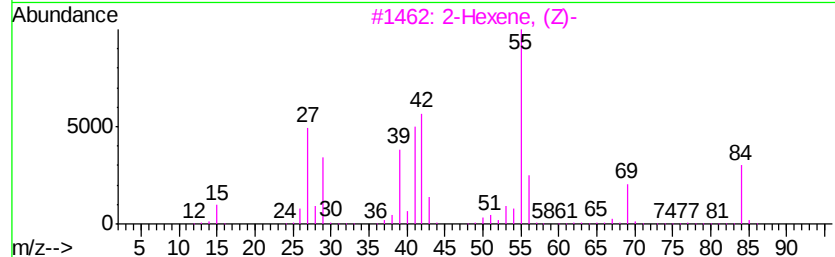
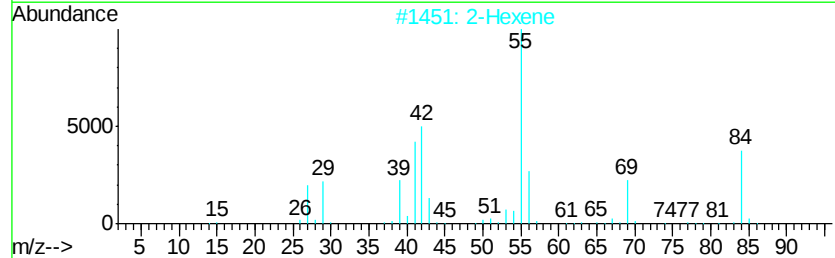
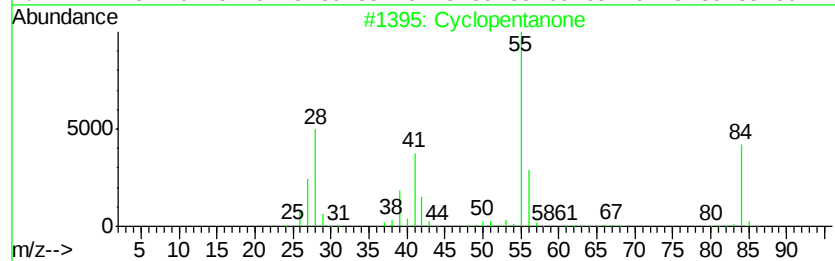
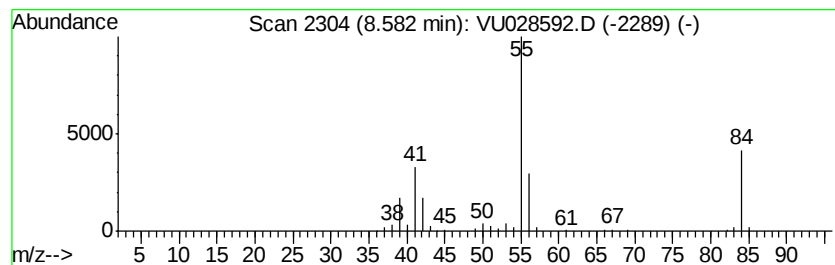
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Quant Title : SW846 8260

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

Peak Number 5 Cyclopentanone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.58	19.13 ug/l	178336	Chlorobenzene-d5	9.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentanone	84	C5H8O	000120-92-3	91
2		2-Hexene	84	C6H12	000592-43-8	59
3		2-Hexene, (Z)-	84	C6H12	007688-21-3	59
4		1H-Tetrazole, 1-methyl-	84	C2H4N4	016681-77-9	59
5		2-Hexene, (E)-	84	C6H12	004050-45-7	45



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU121118\
Data File : VU028592.D
Acq On : 11 Dec 2018 08:29
Operator : MD/SY
Sample : J6305-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
P001-SW005-01

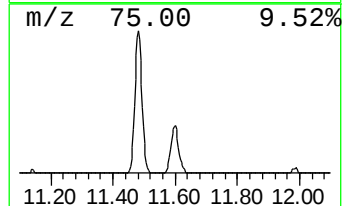
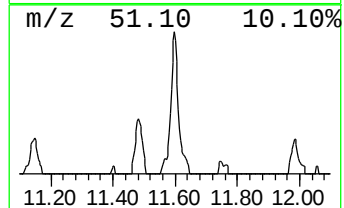
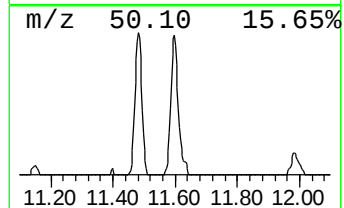
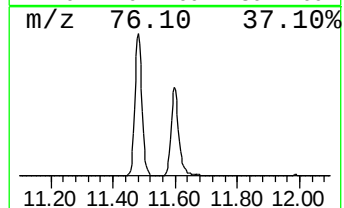
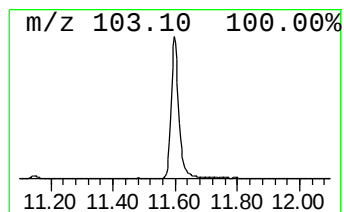
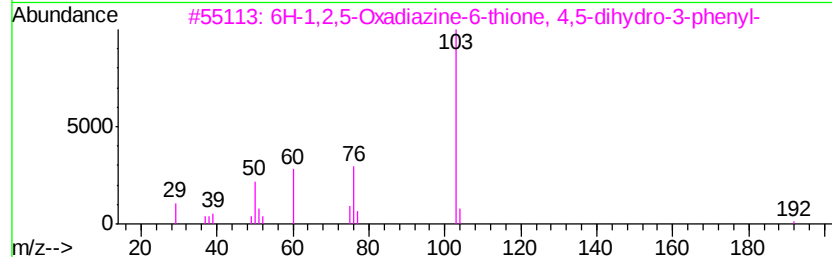
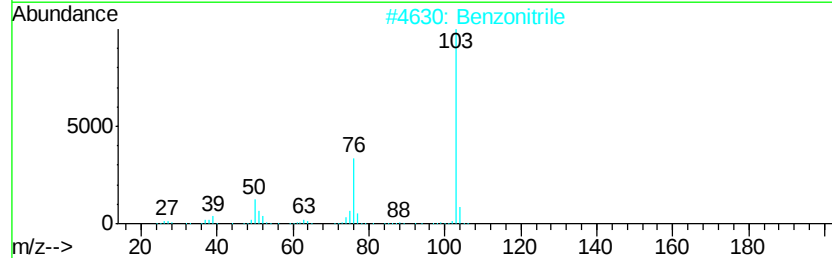
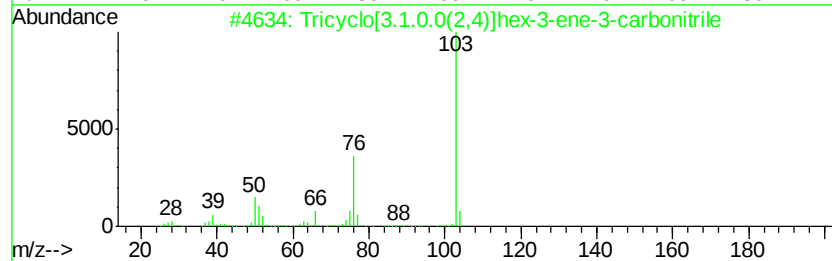
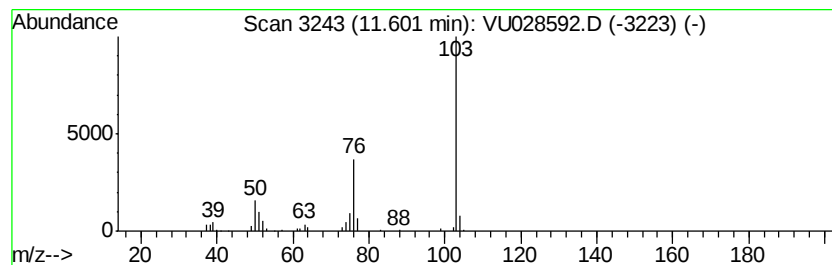
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Quant Title : SW846 8260

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

Peak Number 6 Tricyclo[3.1.0.0(2,4)]hex-3... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.60	10.69 ug/l	89426	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricyclo[3.1.0.0(2,4)]hex-3-ene-...	103	C7H5N	103495-51-8	97
2		Benzonitrile	103	C7H5N	000100-47-0	95
3		6H-1,2,5-Oxadiazine-6-thione, 4,...	192	C9H8N2OS	033406-07-4	83
4		Bicyclo[2.2.2]octa-2,5-diene-2-c...	131	C9H9N	039863-23-5	72
5		Ethanethioamide, N,N-dimethyl-	103	C4H9NS	000631-67-4	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU121118\
Data File : VU028592.D
Acq On : 11 Dec 2018 08:29
Operator : MD/SY
Sample : J6305-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
P001-SW005-01

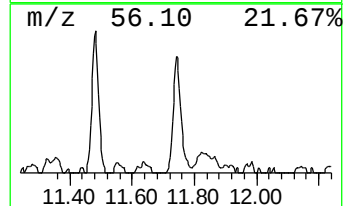
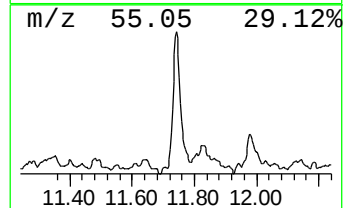
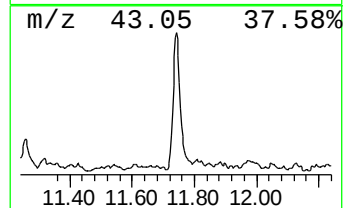
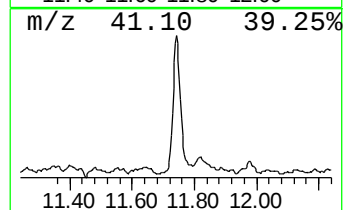
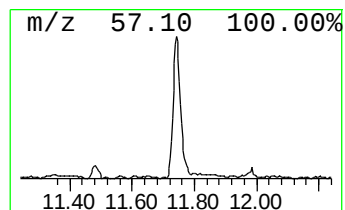
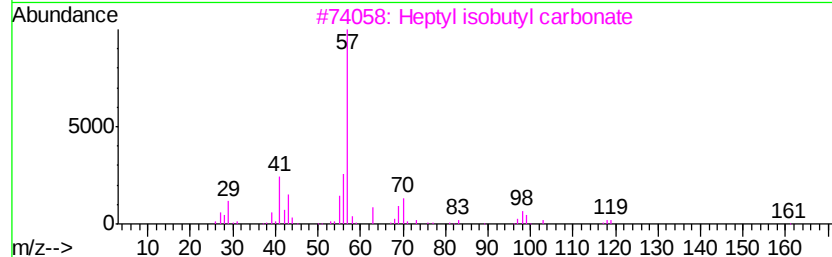
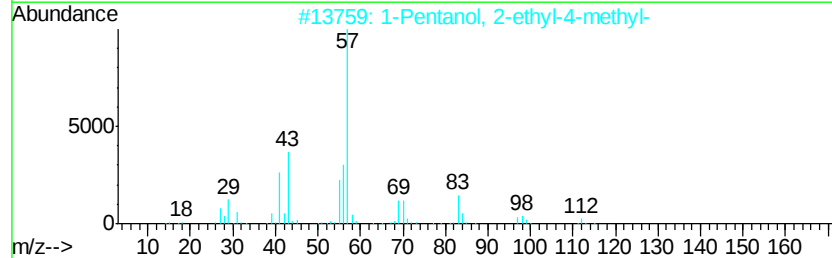
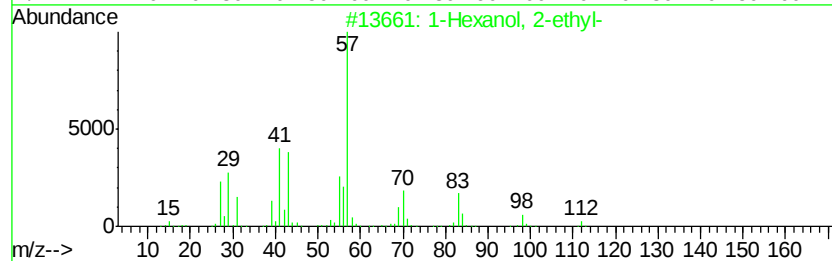
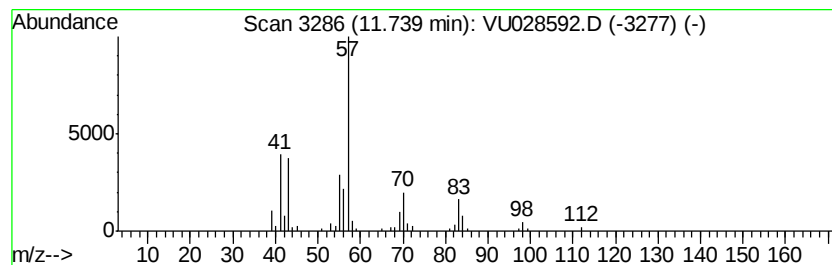
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.74	8.53 ug/l	71358	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2		1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	72
3		Heptyl isobutyl carbonate	216	C12H24O3	959068-08-7	53
4		2-Ethyl-1-hexanol, methyl ether	144	C9H20O	1000365-10-2	53
5		2-Propyl-1-pentanol	130	C8H18O	058175-57-8	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU121118\
Data File : VU028592.D
Acq On : 11 Dec 2018 08:29
Operator : MD/SY
Sample : J6305-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
P001-SW005-01

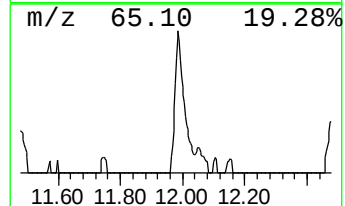
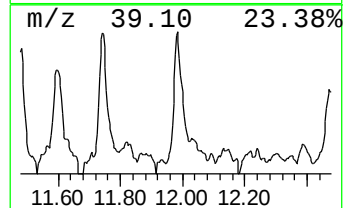
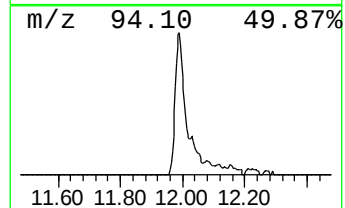
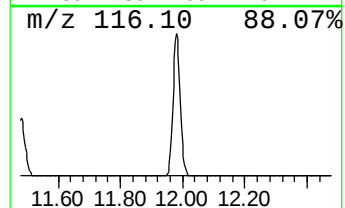
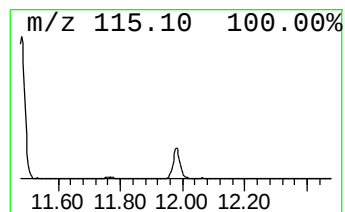
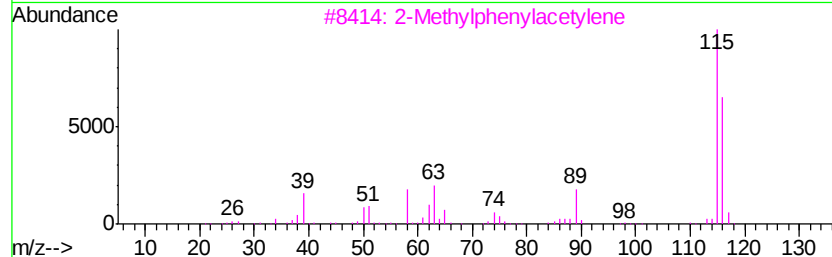
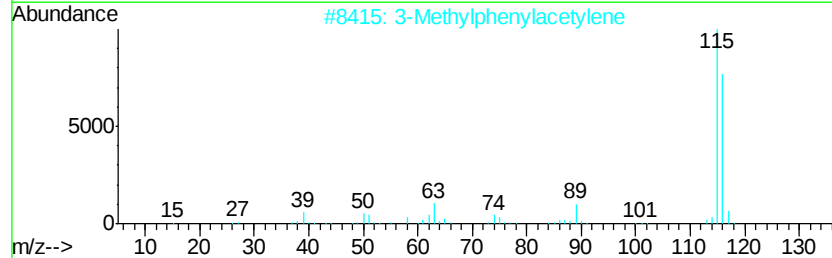
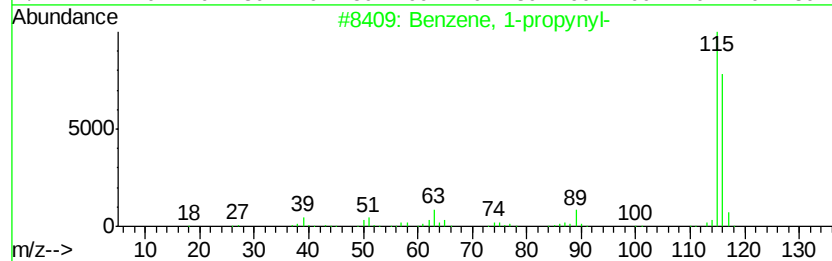
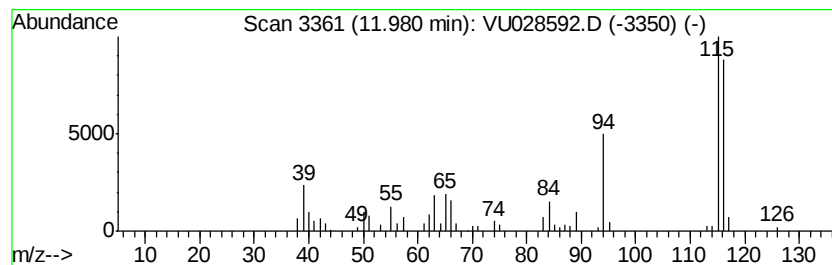
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-propynyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.98	5.99 ug/l	50096	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-propynyl-	116	C9H8	000673-32-5	60
2		3-Methylphenylacetylene	116	C9H8	000766-82-5	60
3		2-Methylphenylacetylene	116	C9H8	000766-47-2	55
4		Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	53
5		Indene	116	C9H8	000095-13-6	41



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU121118\
Data File : VU028592.D
Acq On : 11 Dec 2018 08:29
Operator : MD/SY
Sample : J6305-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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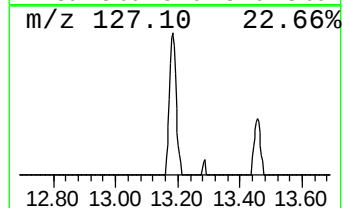
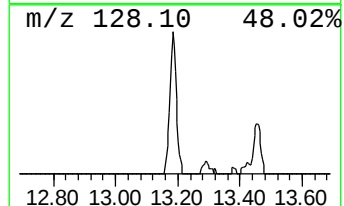
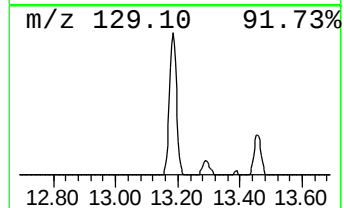
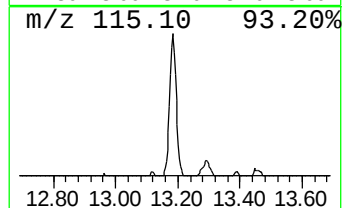
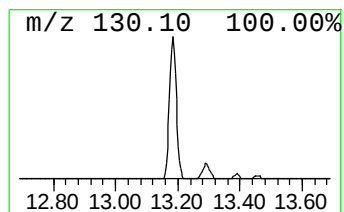
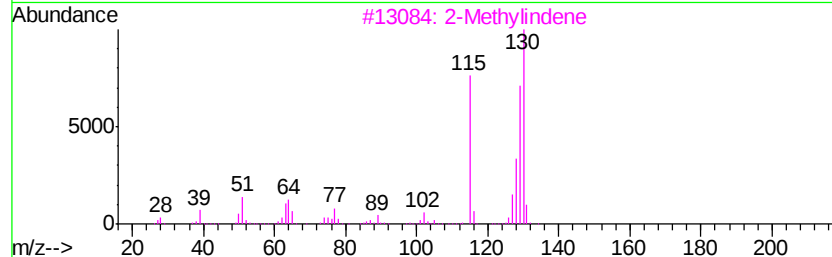
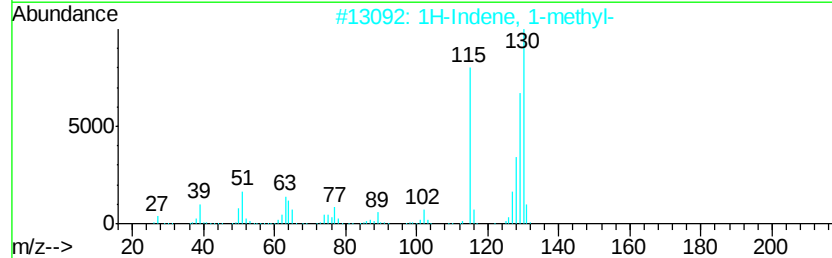
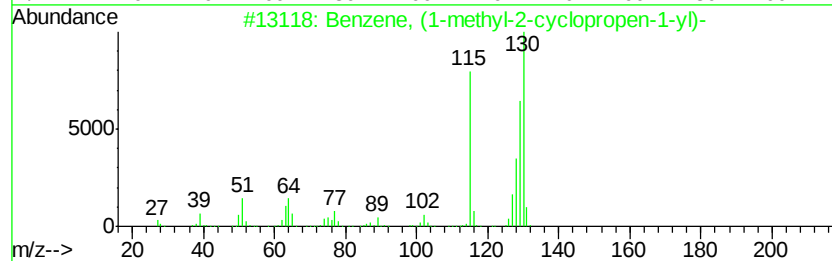
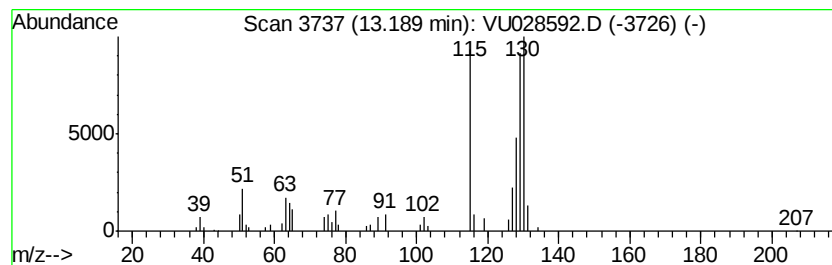
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, (1-methyl-2-cyclop... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.19	5.84 ug/l	48878	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	96
2		1H-Indene, 1-methyl-	130	C10H10	000767-59-9	96
3		2-Methylindene	130	C10H10	002177-47-1	95
4		Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	95
5		Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU121118\
Data File : VU028592.D
Acq On : 11 Dec 2018 08:29
Operator : MD/SY
Sample : J6305-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
P001-SW005-01

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U120618W.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
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Trimethylsilyl fl...	1.51	11.9	ug/l	61659	1	4.98	257946	50.0
Acetonitrile	2.55	18.6	ug/l	96042	1	4.98	257946	50.0
Silanol, trimethyl-	4.20	14.8	ug/l	76197	1	4.98	257946	50.0
Cyclopentanone	8.58	19.1	ug/l	178336	3	9.09	466047	50.0
Tricyclo[3.1.0.0(...	11.60	10.7	ug/l	89426	4	11.48	418192	50.0
1-Hexanol, 2-ethyl-	11.74	8.5	ug/l	71358	4	11.48	418192	50.0
Benzene, 1-propynyl-	11.98	6.0	ug/l	50096	4	11.48	418192	50.0
Benzene, (1-methy...	13.19	5.8	ug/l	48878	4	11.48	418192	50.0