

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU121518\
 Data File : VU028701.D
 Acq On : 14 Dec 2018 22:29
 Operator : MD/SY
 Sample : J6395-03
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MW-2A_R1

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	1.562	113	121	128	rVB6	3922	5674	1.18%	0.151%
2	4.884	1139	1154	1170	rBV	68525	150541	31.32%	4.011%
3	4.983	1171	1185	1206	rVB	101420	225343	46.88%	6.005%
4	5.308	1273	1286	1305	rBV	81277	169241	35.21%	4.510%
5	5.887	1447	1466	1485	rBV	168295	329241	68.50%	8.773%
6	6.414	1612	1630	1641	rVB5	7433	17258	3.59%	0.460%
7	7.565	1964	1988	2018	rBV	270361	480635	100.00%	12.807%
8	7.707	2023	2032	2049	rVB3	6037	11572	2.41%	0.308%
9	7.919	2087	2098	2107	rBV3	11183	19336	4.02%	0.515%
10	8.044	2129	2137	2147	rBV3	5711	9469	1.97%	0.252%
11	8.488	2265	2275	2282	rVB3	5188	8379	1.74%	0.223%
12	8.607	2301	2312	2321	rBV3	14748	24871	5.17%	0.663%
13	8.845	2378	2386	2397	rVB2	7072	11394	2.37%	0.304%
14	9.089	2449	2462	2480	rBV	233199	385579	80.22%	10.274%
15	9.186	2484	2492	2501	rVV4	10924	19024	3.96%	0.507%
16	9.247	2501	2511	2522	rVB5	7141	13096	2.72%	0.349%
17	9.363	2535	2547	2549	rBV5	12398	17540	3.65%	0.467%
18	9.443	2567	2572	2582	rVB3	5833	8449	1.76%	0.225%
19	9.569	2601	2611	2622	rVB	16141	26312	5.47%	0.701%
20	9.723	2641	2659	2663	rBV6	22750	51091	10.63%	1.361%
21	9.803	2677	2684	2698	rVB2	7739	12890	2.68%	0.343%
22	9.919	2705	2720	2723	rBV7	11153	20148	4.19%	0.537%
23	9.951	2724	2730	2742	rVB2	28510	45215	9.41%	1.205%
24	10.051	2748	2761	2777	rBV10	20692	49755	10.35%	1.326%
25	10.118	2777	2782	2788	rVB7	4392	5475	1.14%	0.146%
26	10.167	2788	2797	2804	rBV	15034	21790	4.53%	0.581%
27	10.218	2804	2813	2830	rVB10	7929	17283	3.60%	0.461%
28	10.308	2830	2841	2853	rBV	171579	271934	56.58%	7.246%
29	10.376	2853	2862	2870	rVB4	26058	41142	8.56%	1.096%
30	10.449	2879	2885	2888	rBV4	4361	5502	1.14%	0.147%
31	10.491	2889	2898	2910	rVB4	38686	64893	13.50%	1.729%
32	10.588	2919	2928	2935	rBV	22459	32340	6.73%	0.862%
33	10.697	2953	2962	2973	rVV7	12487	25136	5.23%	0.670%
34	10.755	2974	2980	2990	rVV8	5028	10029	2.09%	0.267%

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35	10.810	2991	2997	3004	rVV5	4768	6545	1.36%	0.174%
36	10.864	3004	3014	3025	rVB5	9762	17568	3.66%	0.468%
37	10.980	3036	3050	3065	rBV4	17386	37282	7.76%	0.993%
38	11.099	3078	3087	3095	rBV2	15438	22487	4.68%	0.599%
39	11.147	3095	3102	3114	rVB9	4097	7488	1.56%	0.200%
40	11.289	3131	3146	3153	rBV	60370	108089	22.49%	2.880%
41	11.398	3172	3180	3189	rVB6	3184	4943	1.03%	0.132%
42	11.482	3195	3206	3221	rBV	206826	327580	68.16%	8.729%
43	11.687	3258	3270	3284	rVB	31125	49538	10.31%	1.320%
44	11.768	3284	3295	3308	rBV2	32485	61890	12.88%	1.649%
45	11.861	3313	3324	3347	rVB7	24962	57807	12.03%	1.540%
46	11.980	3351	3361	3372	rBV2	33429	51963	10.81%	1.385%
47	12.244	3431	3443	3448	rBV5	6185	12872	2.68%	0.343%
48	12.285	3448	3456	3467	rVB2	40095	67008	13.94%	1.786%
49	12.353	3468	3477	3485	rBV	55691	89997	18.72%	2.398%
50	12.433	3496	3502	3509	rVB2	8697	11064	2.30%	0.295%
51	12.536	3526	3534	3550	rVB	29305	45792	9.53%	1.220%
52	12.620	3550	3560	3562	rBV3	15370	18810	3.91%	0.501%
53	12.649	3562	3569	3581	rVV	46372	73468	15.29%	1.958%
54	12.720	3584	3591	3604	rVV6	3641	5363	1.12%	0.143%
55	12.787	3604	3612	3622	rVB4	5525	7425	1.54%	0.198%
56	12.925	3644	3655	3669	rBV4	11352	21819	4.54%	0.581%
57	12.996	3670	3677	3687	rVB2	3870	5631	1.17%	0.150%
58	13.125	3709	3717	3725	rBV3	7149	10804	2.25%	0.288%
59	13.507	3827	3836	3843	rBV5	6595	9712	2.02%	0.259%
60	13.578	3843	3858	3865	rBV6	5572	12273	2.55%	0.327%

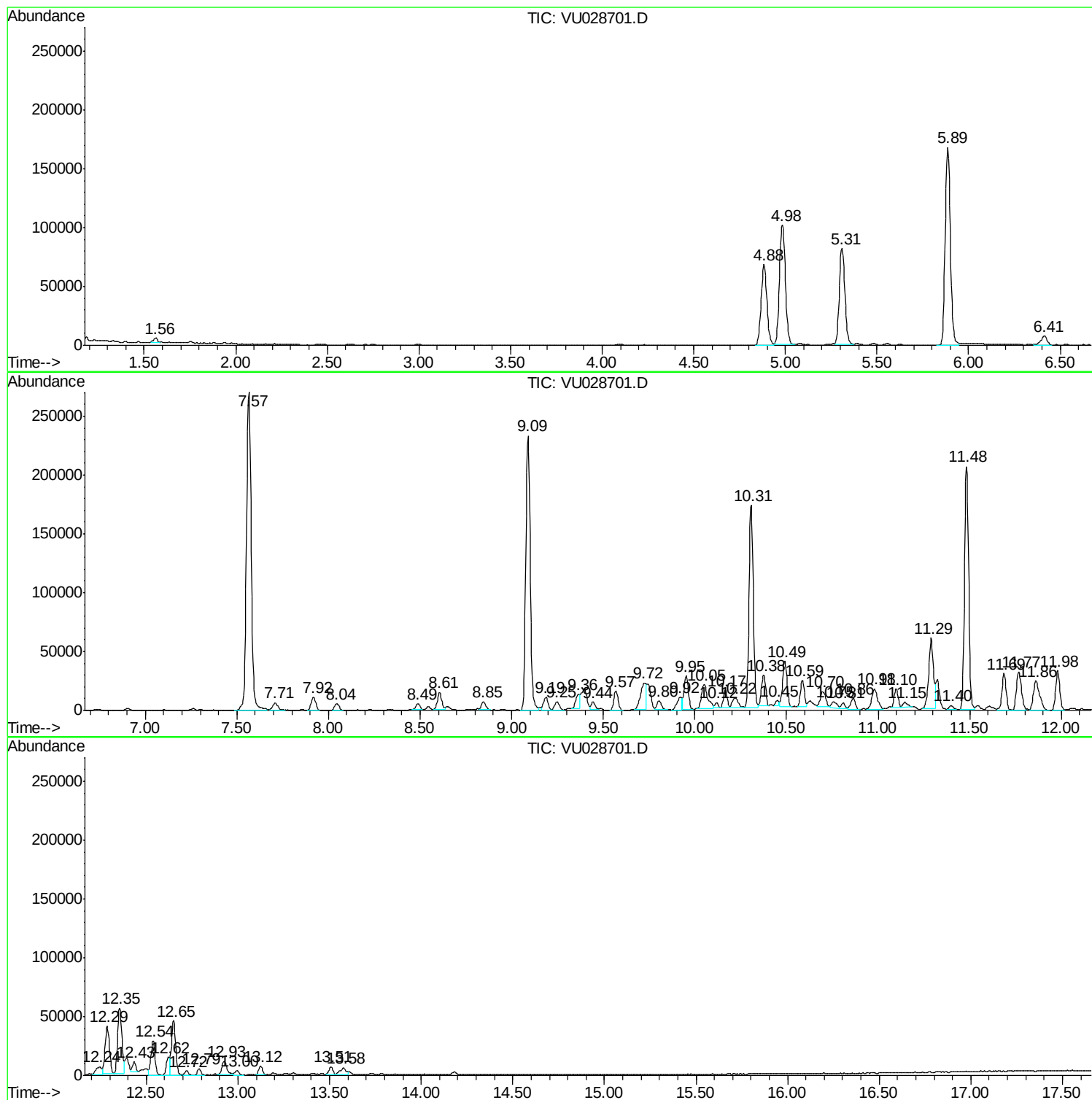
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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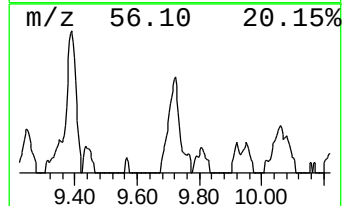
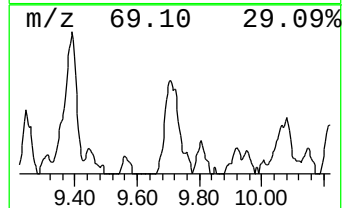
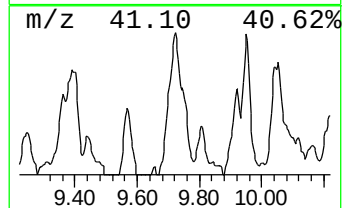
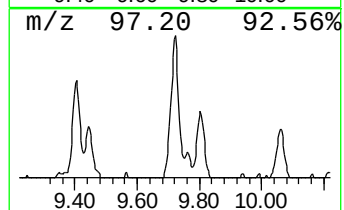
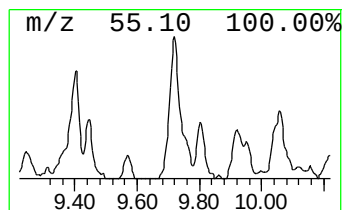
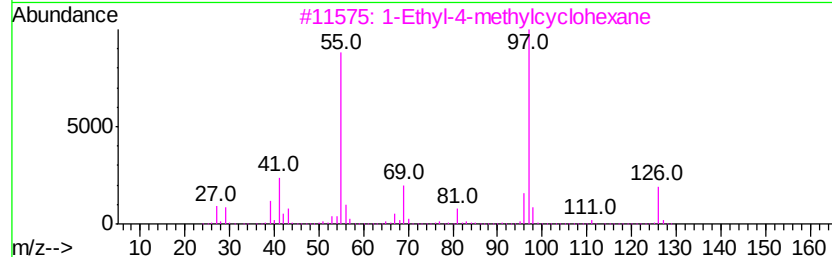
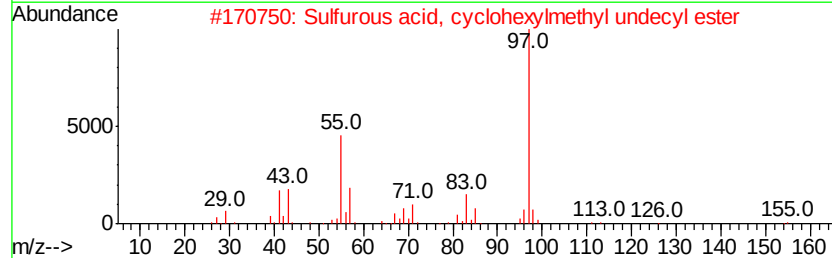
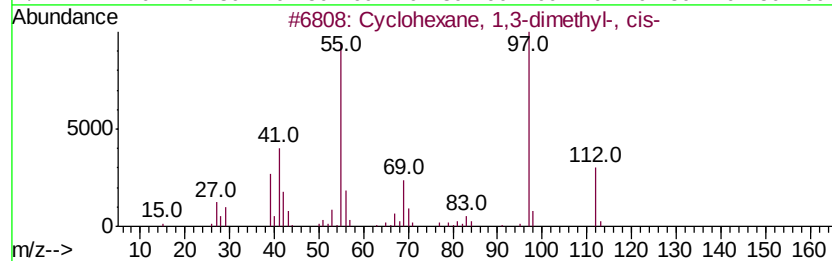
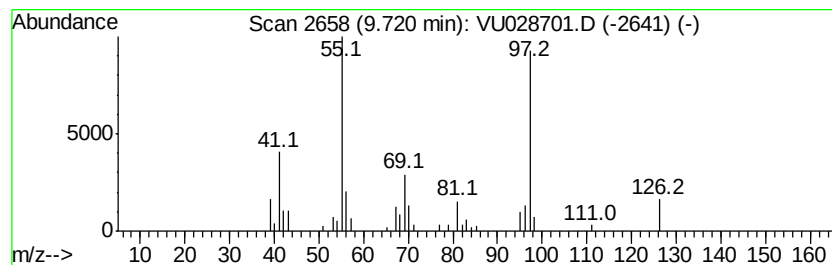
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U120618W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclohexane, 1,3-dimethyl-,... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.72	6.63 ug/l	51091	Chlorobenzene-d5	9.09

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	72
2			Sulfurous acid, cyclohexylmethyl...	332	C18H36O3S	1000309-21-9	64
3			1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	64
4			Oxalic acid, cyclohexylmethyl is...	270	C15H26O4	1000309-68-3	64
5			3,5-Dimethyl-3-heptene	126	C9H18	059643-68-4	59



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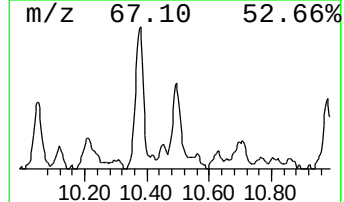
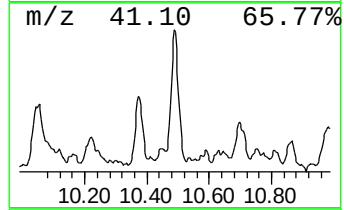
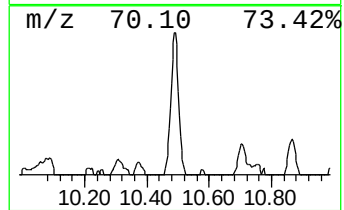
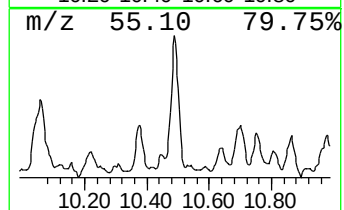
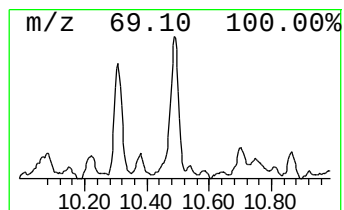
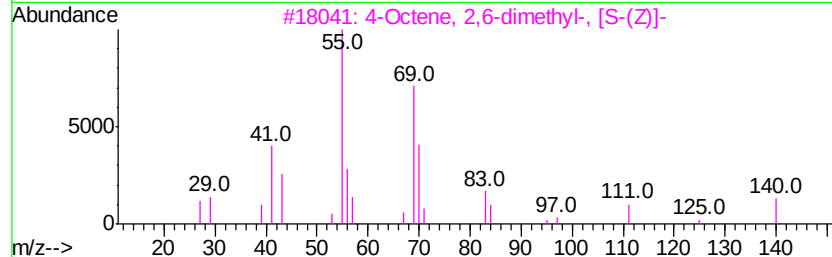
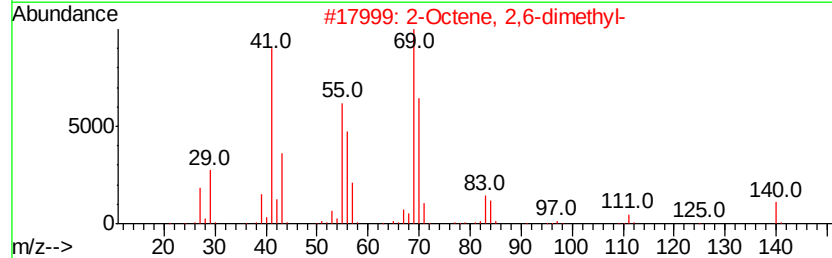
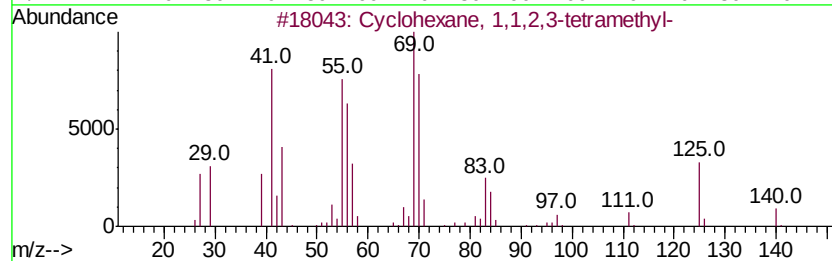
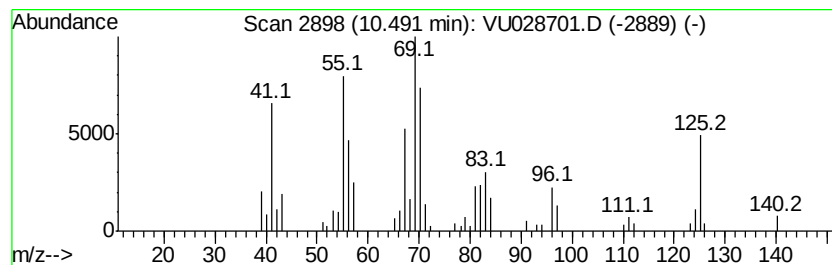
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclohexane, 1,1,2,3-tetram... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.49	9.90 ug/l	64893	1,4-Dichlorobenzene-d4	11.48	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,1,2,3-tetramethyl-	140	C10H20	006783-92-2	70
2	2-Octene, 2,6-dimethyl-	140	C10H20	004057-42-5	53
3	4-Octene, 2,6-dimethyl-, [S-(Z)]-	140	C10H20	062960-77-4	47
4	Cyclohexane, 1,2,3-trimethyl-, (...)	126	C9H18	001678-81-5	45
5	1-Hexene, 3,3,5-trimethyl-	126	C9H18	013427-43-5	38



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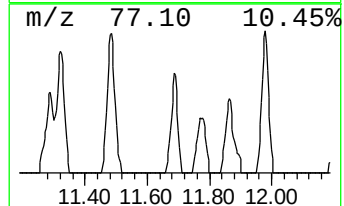
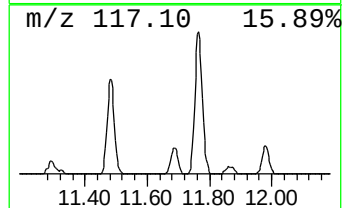
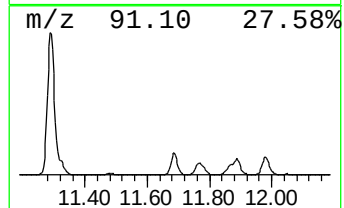
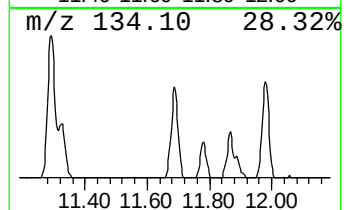
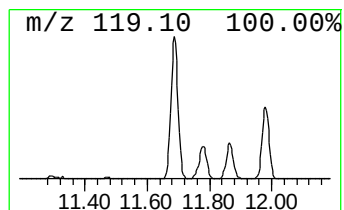
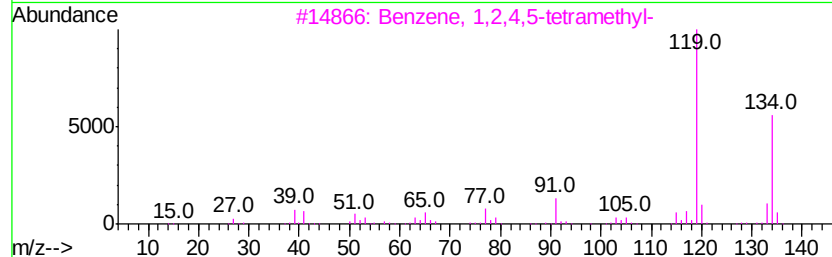
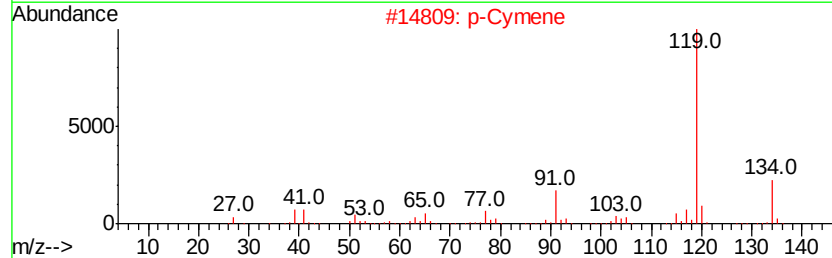
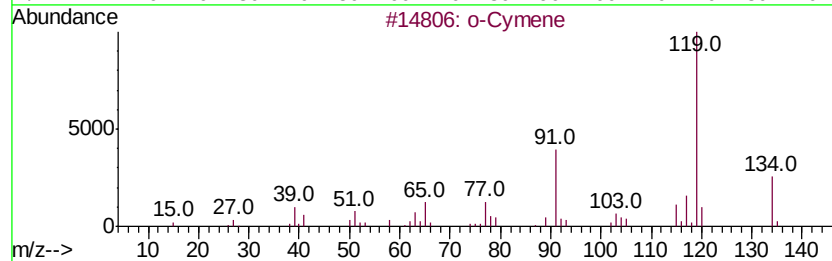
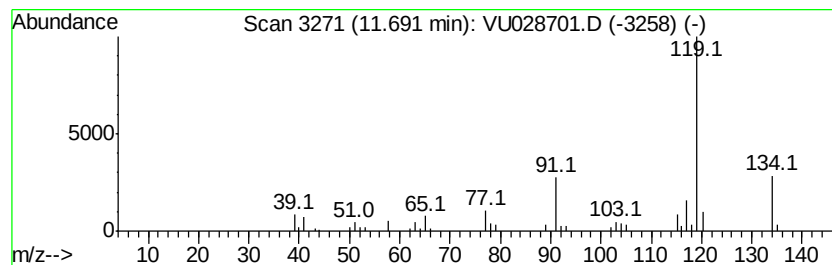
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 o-Cymene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.69	7.56 ug/l	49538	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	96
2		p-Cymene	134	C10H14	000099-87-6	94
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91
4		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91
5		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	91



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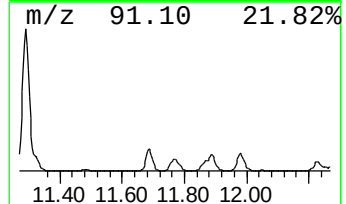
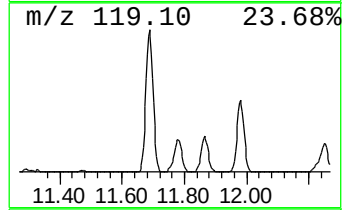
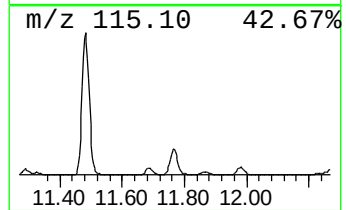
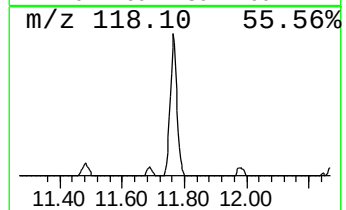
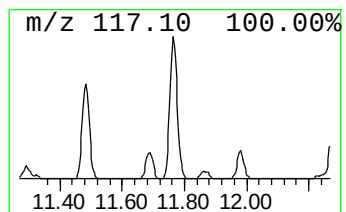
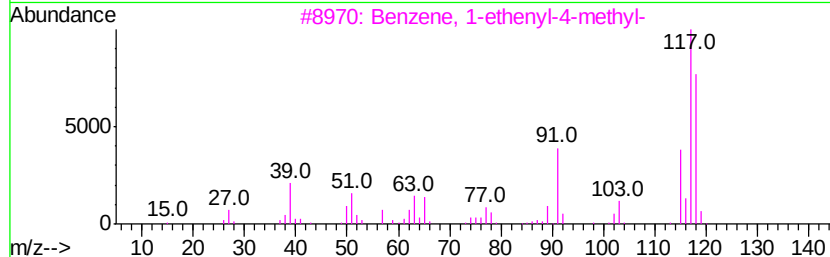
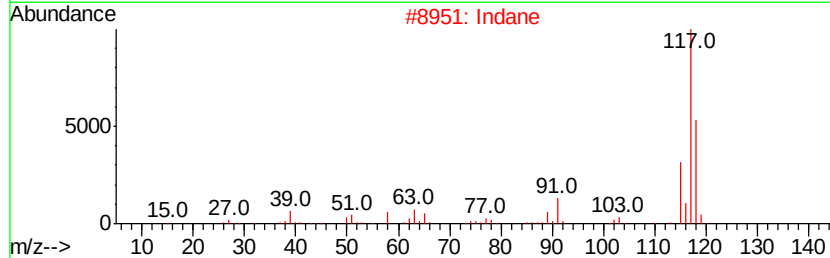
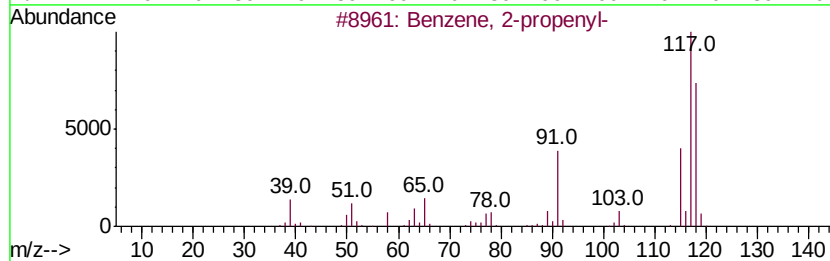
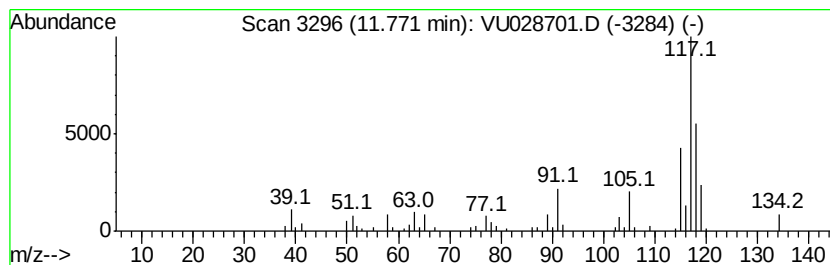
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 2-propenyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	9.45 ug/l	61890	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
2		Indane	118	C9H10	000496-11-7	70
3		Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	70
4		Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
5		Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	62



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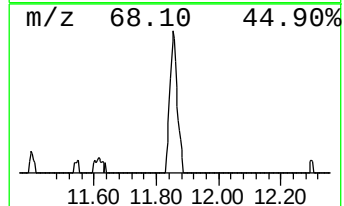
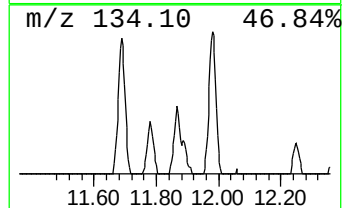
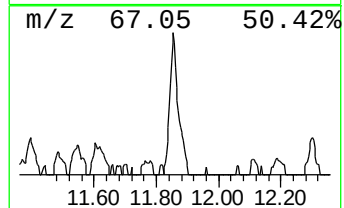
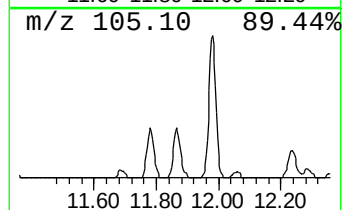
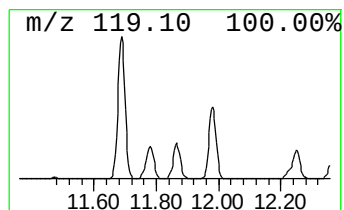
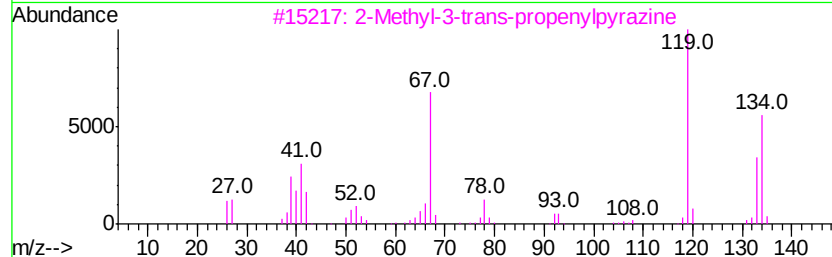
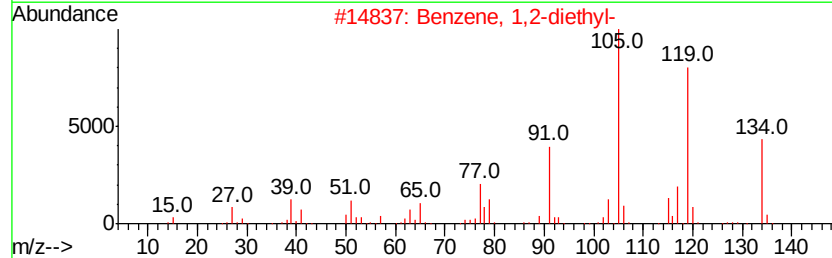
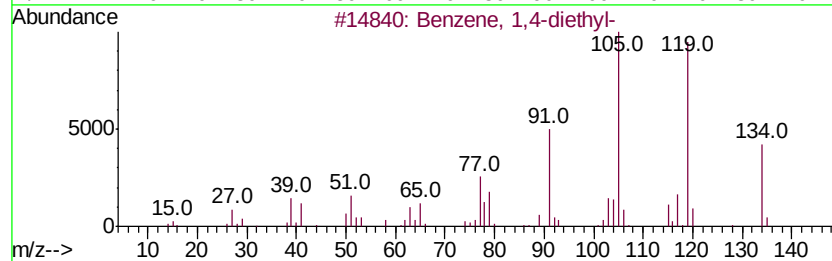
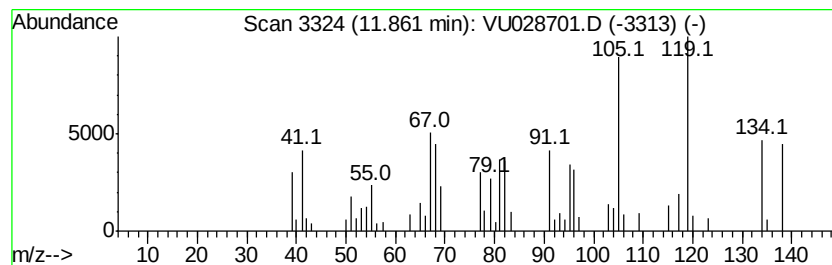
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MSVOA_U
ClientSampleId :
MW-2A_R1

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 5 Benzene, 1,4-diethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.86	8.82 ug/l	57807	1,4-Dichlorobenzene-d4	11.48
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,4-diethyl-	134 C10H14	000105-05-5 94
2		Benzene, 1,2-diethyl-	134 C10H14	000135-01-3 83
3		2-Methyl-3-trans-propenylpyrazine	134 C8H10N2	232255-41-3 46
4		Pyrazine, 2-methyl-3-(1-propenyl-...	134 C8H10N2	055138-65-3 46
5		Benzene, 1,2,4,5-tetramethyl-	134 C10H14	000095-93-2 30



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU121518\
Data File : VU028701.D
Acq On : 14 Dec 2018 22:29
Operator : MD/SY
Sample : J6395-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 23 Sample Multiplier: 1

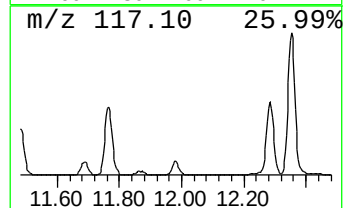
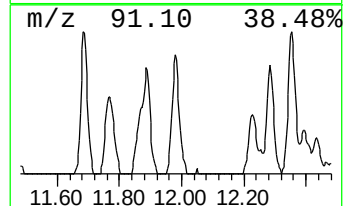
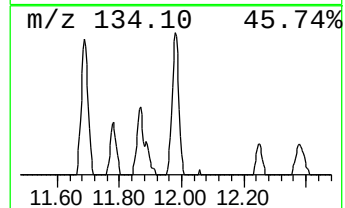
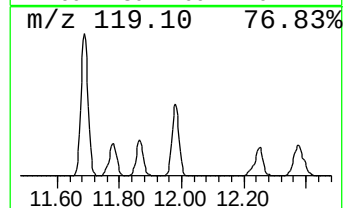
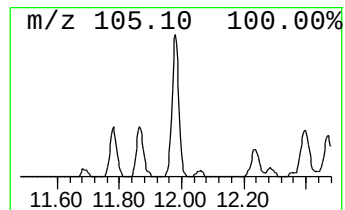
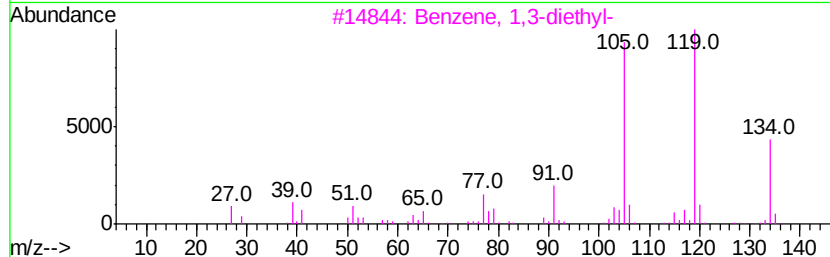
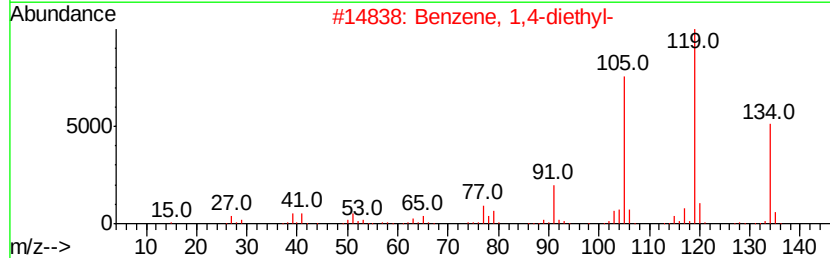
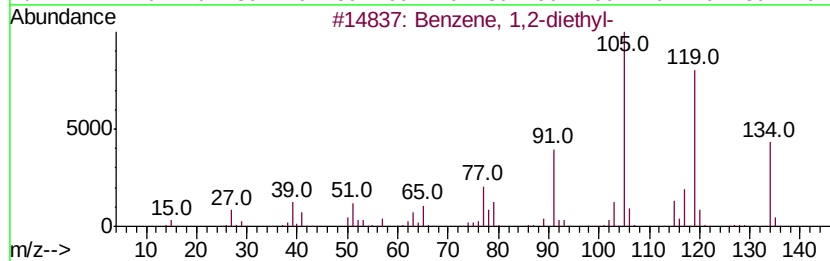
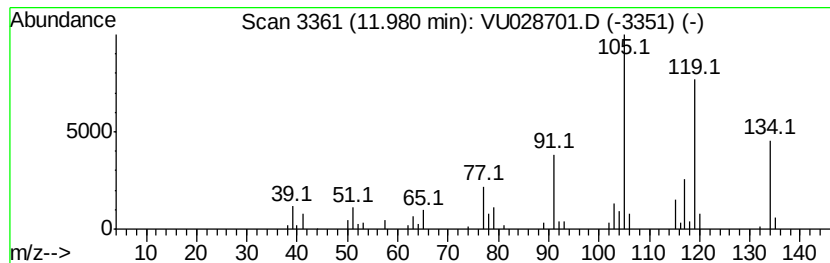
Instrument :
MSVOA_U
ClientSampled :
MW-2A_R1

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 6 Benzene, 1,2-diethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.98	7.93 ug/l	51963	1,4-Dichlorobenzene-d4	11.48
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2-diethyl-	134 C10H14	000135-01-3 96
2		Benzene, 1,4-diethyl-	134 C10H14	000105-05-5 90
3		Benzene, 1,3-diethyl-	134 C10H14	000141-93-5 87
4		Benzene, 1-ethyl-3,5-dimethyl-	134 C10H14	000934-74-7 68
5		Benzene, 4-ethyl-1,2-dimethyl-	134 C10H14	000934-80-5 64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU121518\
Data File : VU028701.D
Acq On : 14 Dec 2018 22:29
Operator : MD/SY
Sample : J6395-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-2A_R1

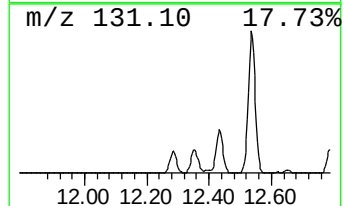
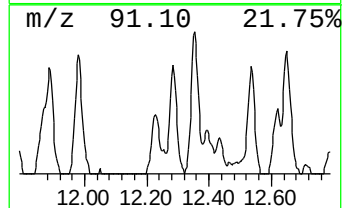
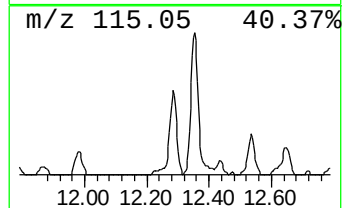
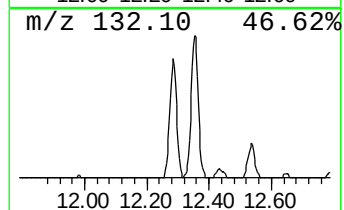
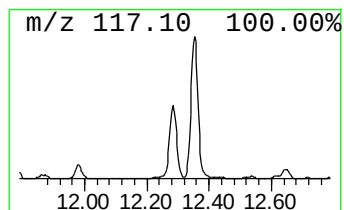
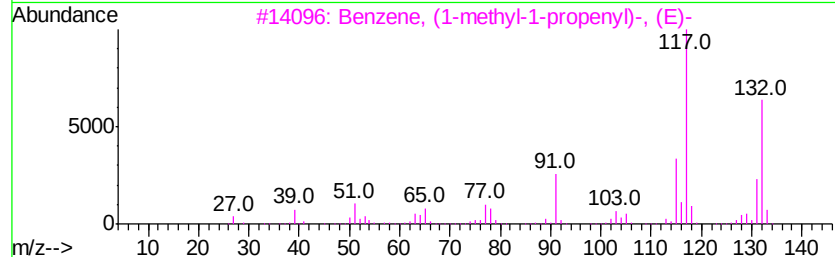
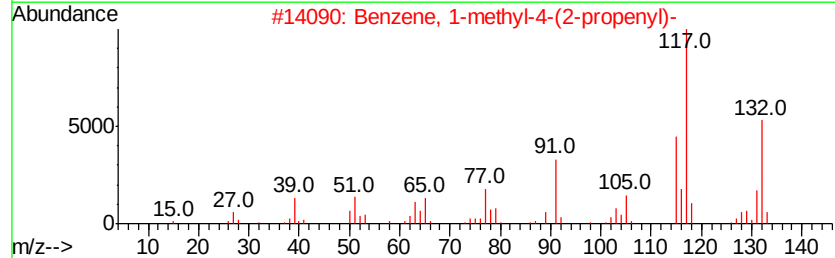
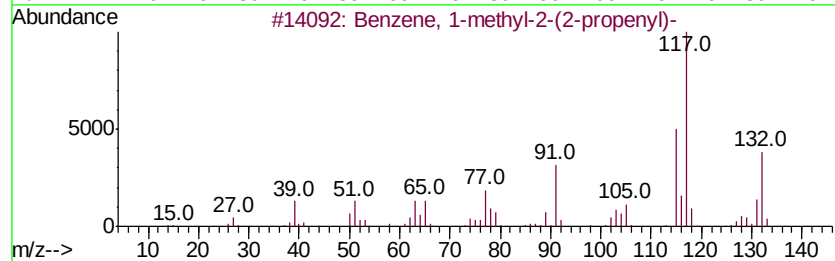
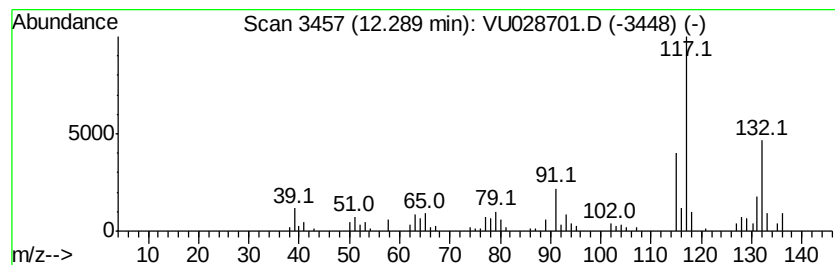
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-methyl-2-(2-prop... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	10.23 ug/l	67008	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	94
2		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	93
3		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	91
4		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	91
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU121518\
Data File : VU028701.D
Acq On : 14 Dec 2018 22:29
Operator : MD/SY
Sample : J6395-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 23 Sample Multiplier: 1

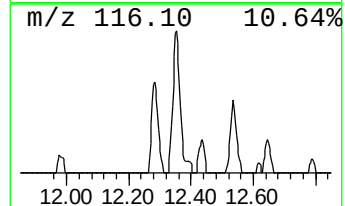
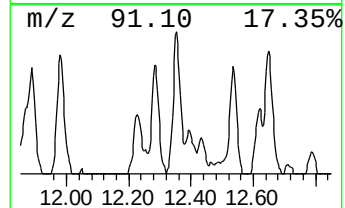
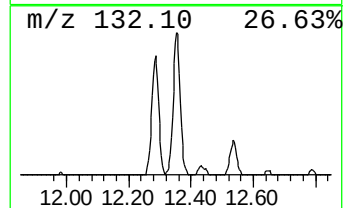
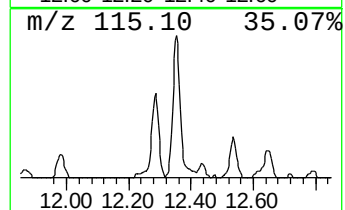
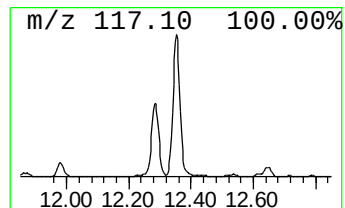
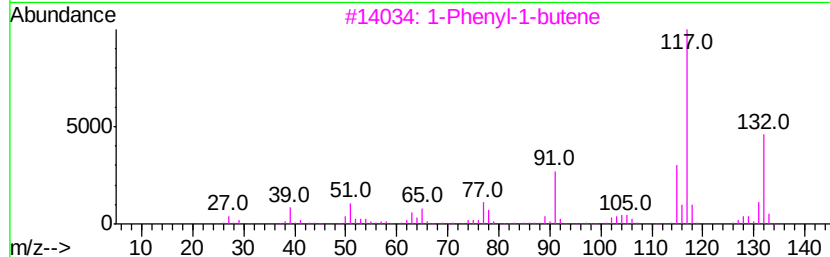
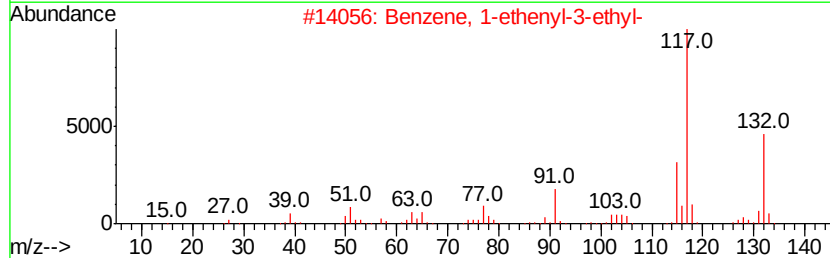
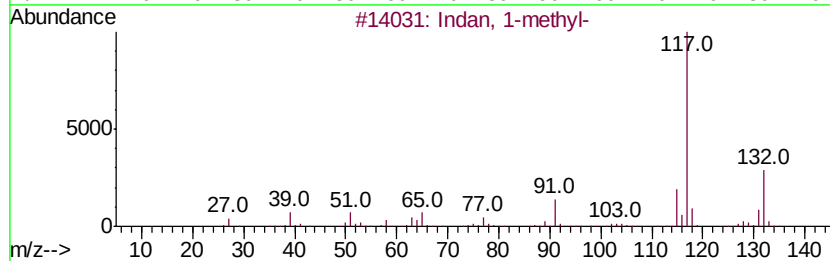
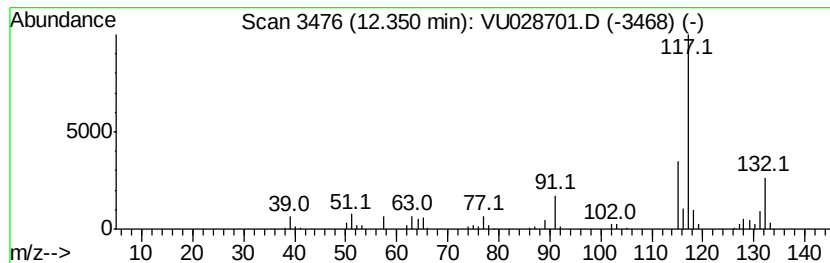
Instrument :
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ClientSampleId :
MW-2A_R1

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 8 Indan, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.35	13.74 ug/l	89997	1,4-Dichlorobenzene-d4	11.48
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indan, 1-methyl-	132 C10H12	000767-58-8	91
2	Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4	90
3	1-Phenyl-1-butene	132 C10H12	000824-90-8	87
4	Benzene, (2-methylcyclopropyl)-	132 C10H12	1000327-39-0	87
5	1H-Indene, 2,3-dihydro-2-methyl-	132 C10H12	000824-63-5	86



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU121518\
Data File : VU028701.D
Acq On : 14 Dec 2018 22:29
Operator : MD/SY
Sample : J6395-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-2A_R1

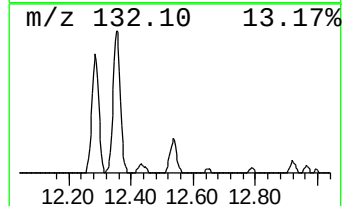
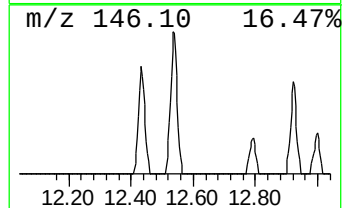
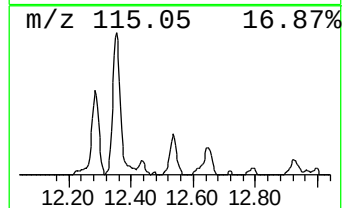
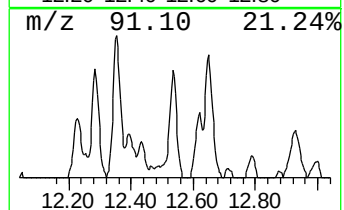
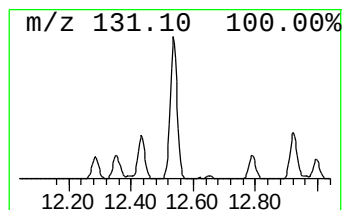
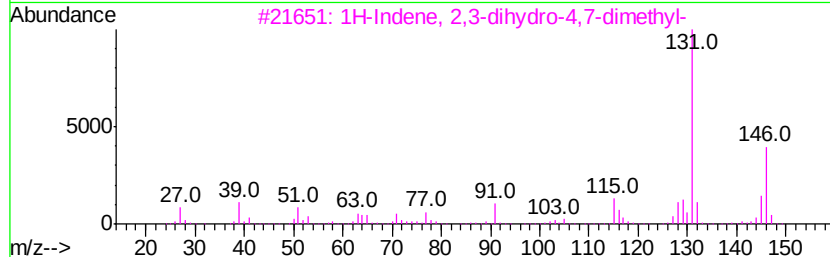
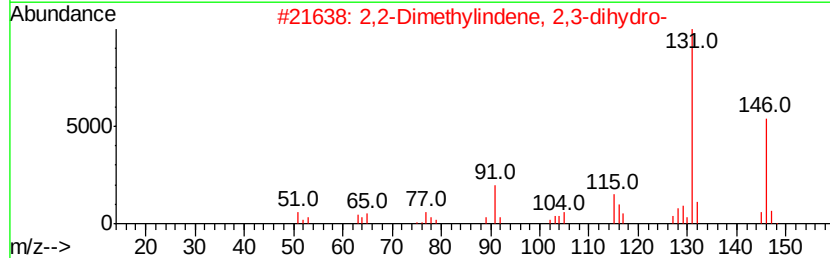
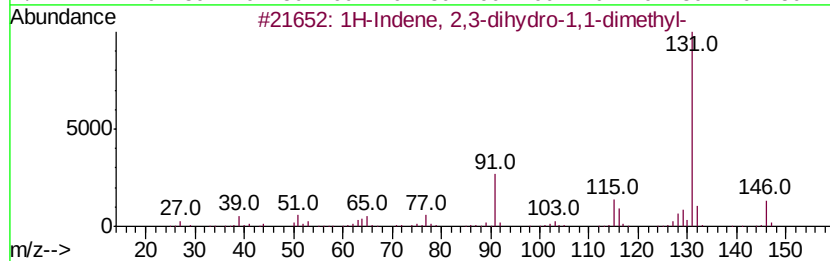
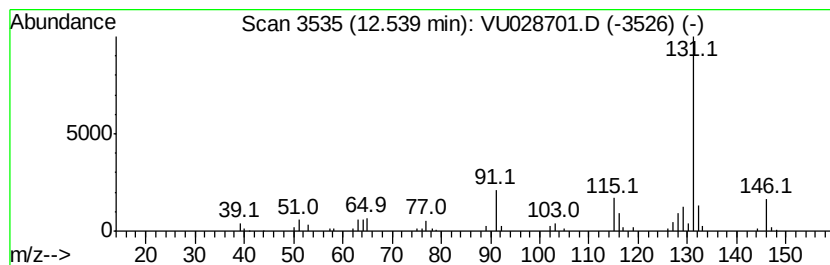
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Quant Title : SW846 8260

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

Peak Number 9 1H-Indene, 2,3-dihydro-1,1-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	6.99 ug/l	45792	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	95
2		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91
3		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	90
4		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90
5		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU121518\
Data File : VU028701.D
Acq On : 14 Dec 2018 22:29
Operator : MD/SY
Sample : J6395-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-2A_R1

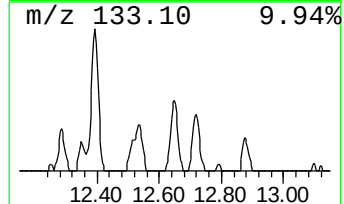
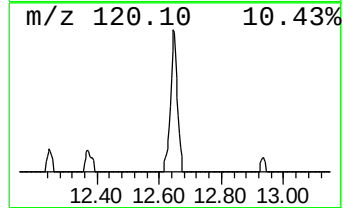
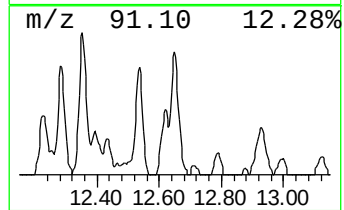
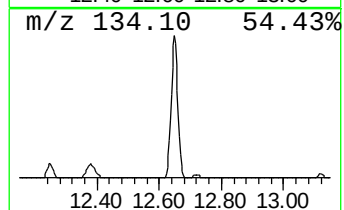
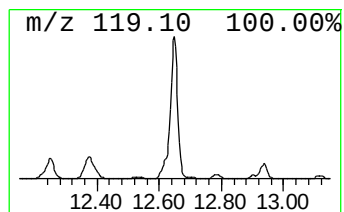
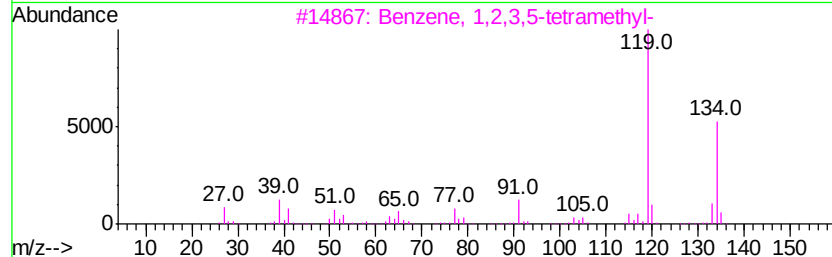
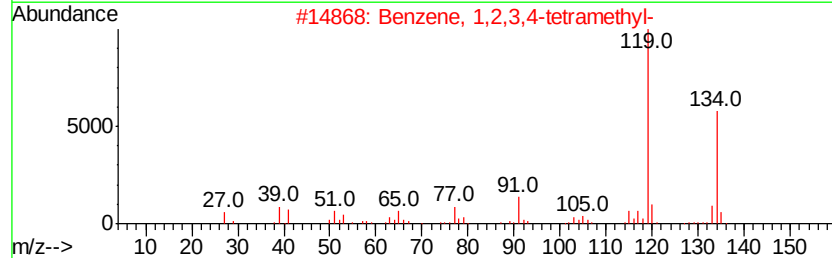
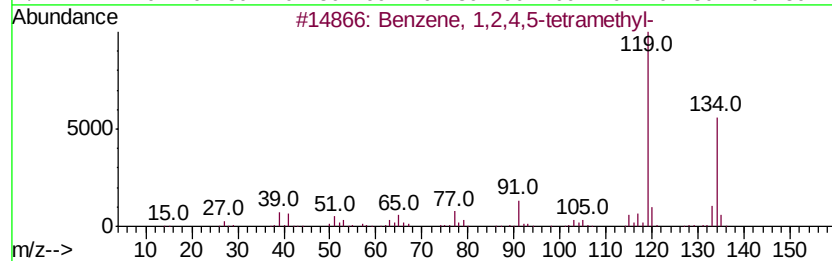
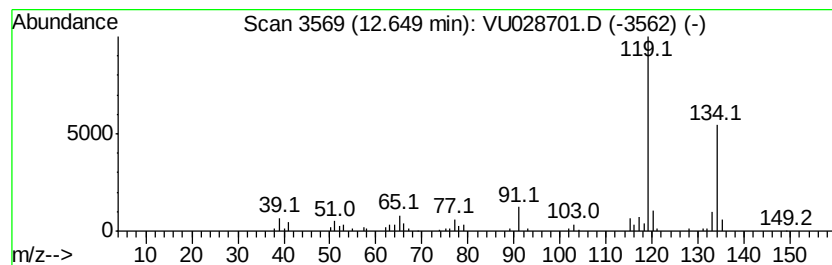
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U120618W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	11.21 ug/l	73468	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
5		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU121518\
Data File : VU028701.D
Acq On : 14 Dec 2018 22:29
Operator : MD/SY
Sample : J6395-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-2A_R1

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
Cyclohexane, 1,3-...	9.72	6.6	ug/l	51091	3	9.09	385579	50.0
Cyclohexane, 1,1,...	10.49	9.9	ug/l	64893	4	11.48	327580	50.0
o-Cymene	11.69	7.6	ug/l	49538	4	11.48	327580	50.0
Benzene, 2-propenyl-	11.77	9.4	ug/l	61890	4	11.48	327580	50.0
Benzene, 1,4-diet...	11.86	8.8	ug/l	57807	4	11.48	327580	50.0
Benzene, 1,2-diet...	11.98	7.9	ug/l	51963	4	11.48	327580	50.0
Benzene, 1-methyl...	12.29	10.2	ug/l	67008	4	11.48	327580	50.0
Indan, 1-methyl-	12.35	13.7	ug/l	89997	4	11.48	327580	50.0
1H-Indene, 2,3-di...	12.54	7.0	ug/l	45792	4	11.48	327580	50.0
Benzene, 1,2,4,5-...	12.65	11.2	ug/l	73468	4	11.48	327580	50.0