

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\U112918\
 Data File : VU028405.D
 Acq On : 29 Nov 2018 19:39
 Operator : MD/SY
 Sample : J6134-07
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 DUP-1

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\82U112818W.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	1.405	62	72	78	rBV	69915	93090	2.73%	0.260%
2	2.444	385	395	417	rBV	38239	74472	2.18%	0.208%
3	2.987	552	564	581	rBV3	23642	46719	1.37%	0.130%
4	4.096	891	909	924	rBV6	12141	37181	1.09%	0.104%
5	4.225	928	949	988	rVV	1419304	3411723	100.00%	9.525%
6	4.884	1136	1154	1170	rBV2	235640	530686	15.55%	1.482%
7	4.984	1170	1185	1203	rVV	353326	783563	22.97%	2.188%
8	5.308	1271	1286	1303	rBV	298299	623809	18.28%	1.742%
9	5.392	1303	1312	1325	rVB3	18106	40982	1.20%	0.114%
10	5.887	1451	1466	1493	rBV	617704	1242584	36.42%	3.469%
11	6.418	1613	1631	1644	rBV7	13995	37844	1.11%	0.106%
12	7.054	1815	1829	1842	rBV9	18164	46529	1.36%	0.130%
13	7.566	1968	1988	2000	rBV	1012766	1766939	51.79%	4.933%
14	7.636	2000	2010	2042	rVB	774238	1369127	40.13%	3.822%
15	9.090	2449	2462	2487	rBV	903098	1511875	44.31%	4.221%
16	9.250	2500	2512	2528	rBV	174965	288724	8.46%	0.806%
17	9.373	2537	2550	2566	rBV	1118157	1842083	53.99%	5.143%
18	9.778	2664	2676	2703	rVB	809232	1299978	38.10%	3.629%
19	10.045	2750	2759	2776	rBV10	17426	42844	1.26%	0.120%
20	10.305	2829	2840	2858	rBV	748215	1202340	35.24%	3.357%
21	10.585	2918	2927	2937	rBV	41763	62932	1.84%	0.176%
22	10.681	2942	2957	2974	rBV	442420	1087736	31.88%	3.037%
23	10.771	2974	2985	2994	rBV	674671	1013876	29.72%	2.831%
24	10.971	3034	3047	3068	rBV	526530	850015	24.91%	2.373%
25	11.099	3072	3087	3092	rBV2	48331	92050	2.70%	0.257%
26	11.147	3092	3102	3115	rVB	885229	1349873	39.57%	3.769%
27	11.321	3152	3156	3171	rVB2	26138	41794	1.23%	0.117%
28	11.424	3171	3188	3196	rBV2	146283	261735	7.67%	0.731%
29	11.482	3196	3206	3221	rVV	1029342	1648947	48.33%	4.604%
30	11.569	3222	3233	3253	rVB	884449	1370032	40.16%	3.825%
31	11.688	3264	3270	3281	rVB3	38294	62334	1.83%	0.174%
32	11.768	3281	3295	3302	rVV3	263379	551034	16.15%	1.538%
33	11.807	3302	3307	3316	rVV	245141	365156	10.70%	1.019%
34	11.874	3316	3328	3342	rVB2	639959	1088219	31.90%	3.038%

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35	11.980	3352	3361	3368	rBV2	53718	85774	2.51%	0.239%
36	12.028	3369	3376	3377	rVV	45045	53058	1.56%	0.148%
37	12.054	3377	3384	3398	rVB	175953	280125	8.21%	0.782%
38	12.144	3398	3412	3416	rBV	154419	237199	6.95%	0.662%
39	12.173	3417	3421	3429	rVB	140618	184925	5.42%	0.516%
40	12.250	3436	3445	3454	rBV	543397	770962	22.60%	2.152%
41	12.353	3467	3477	3483	rBV	151134	276244	8.10%	0.771%
42	12.446	3499	3506	3513	rBV3	71822	96491	2.83%	0.269%
43	12.491	3514	3520	3528	rVB2	80960	112244	3.29%	0.313%
44	12.549	3529	3538	3546	rBV	173191	255675	7.49%	0.714%
45	12.649	3560	3569	3577	rBV	480592	724975	21.25%	2.024%
46	12.700	3578	3585	3600	rVB	473178	742246	21.76%	2.072%
47	12.784	3602	3611	3617	rBV5	68103	112061	3.28%	0.313%
48	12.819	3617	3622	3627	rVV4	34293	42704	1.25%	0.119%
49	12.848	3628	3631	3637	rVB	39600	36461	1.07%	0.102%
50	12.961	3658	3666	3676	rVB2	139405	199091	5.84%	0.556%
51	13.080	3695	3703	3706	rBV3	59335	81589	2.39%	0.228%
52	13.118	3706	3715	3725	rVV2	693765	1149853	33.70%	3.210%
53	13.183	3725	3735	3746	rVV	669596	1089922	31.95%	3.043%
54	13.241	3747	3753	3759	rVV3	47159	77040	2.26%	0.215%
55	13.286	3759	3767	3784	rVB2	164144	278937	8.18%	0.779%
56	13.385	3789	3798	3810	rVB	104964	186622	5.47%	0.521%
57	13.453	3810	3819	3828	rBV4	141803	216763	6.35%	0.605%
58	13.507	3828	3836	3842	rBV4	74660	121453	3.56%	0.339%
59	13.546	3843	3848	3861	rVB3	57226	87681	2.57%	0.245%
60	13.739	3898	3908	3931	rVB	322457	605617	17.75%	1.691%
61	13.855	3936	3944	3952	rBV8	22882	46535	1.36%	0.130%
62	13.945	3965	3972	3984	rVB10	35736	71381	2.09%	0.199%
63	14.147	4026	4035	4040	rBV4	97925	160768	4.71%	0.449%
64	14.179	4041	4045	4060	rVB	93376	149666	4.39%	0.418%
65	14.337	4087	4094	4105	rVB7	42404	90564	2.65%	0.253%
66	14.395	4106	4112	4122	rVB7	26788	49084	1.44%	0.137%
67	14.453	4125	4130	4146	rVB7	24950	65773	1.93%	0.184%
68	14.549	4150	4160	4172	rBV2	74115	132566	3.89%	0.370%
69	14.707	4204	4209	4218	rVB6	40517	59931	1.76%	0.167%
70	14.874	4251	4261	4274	rVB	177880	304356	8.92%	0.850%
71	15.060	4310	4319	4334	rBV	121300	198128	5.81%	0.553%

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72	15.195	4353	4361	4372	rBV4	41053	67890	1.99%	0.190%
73	15.340	4401	4406	4422	rVB5	30881	48426	1.42%	0.135%
74	15.588	4475	4483	4490	rBV4	23080	40929	1.20%	0.114%
75	15.726	4518	4526	4537	rVB9	24654	46218	1.35%	0.129%
76	16.060	4623	4630	4637	rBV4	26689	40958	1.20%	0.114%

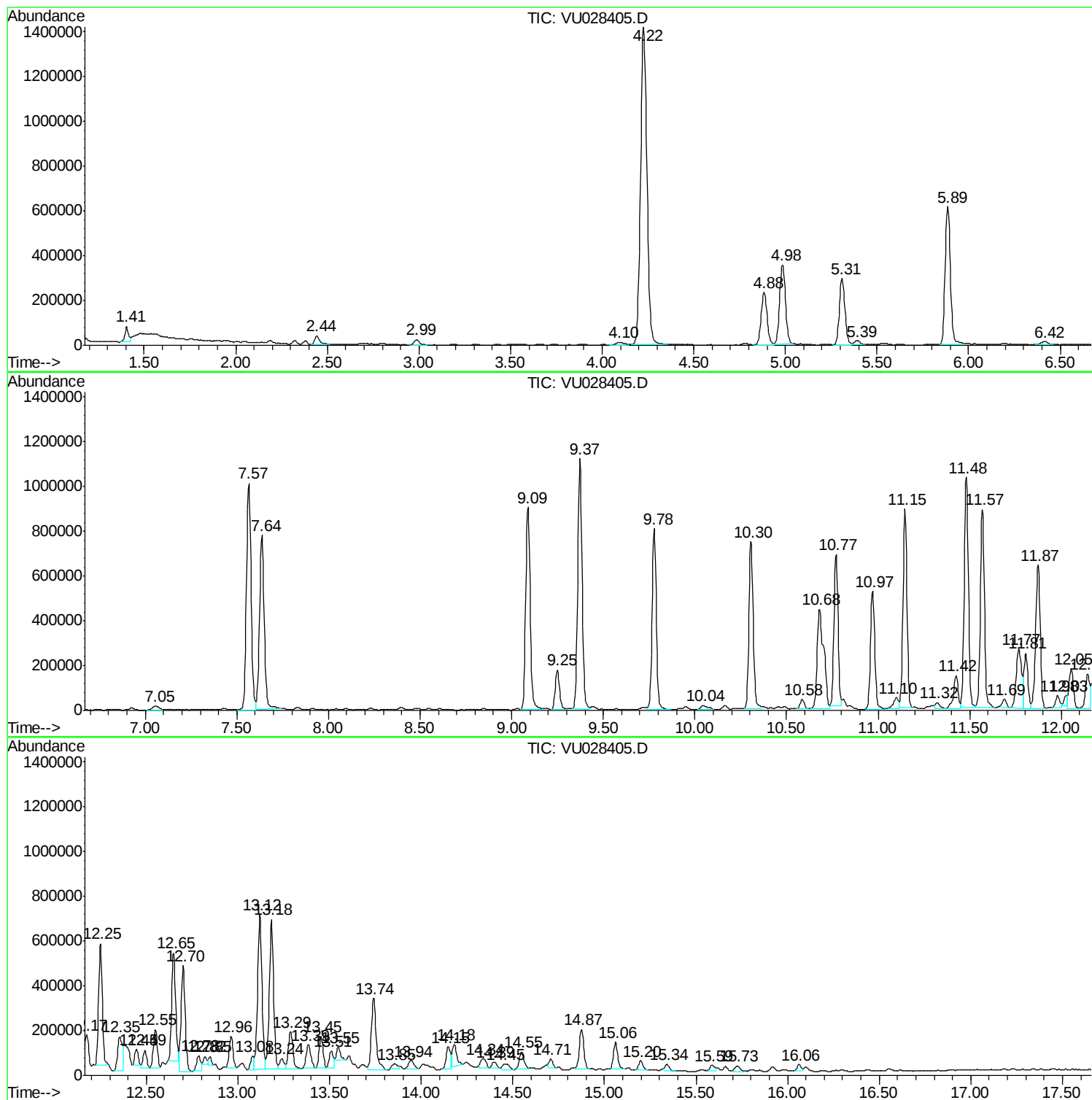
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Instrument :
MSVOA_U
ClientSampled :
DUP-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U112818W.M
Quant Title : SW846 8260

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



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Instrument :
MSVOA_U
ClientSampleId :
DUP-1

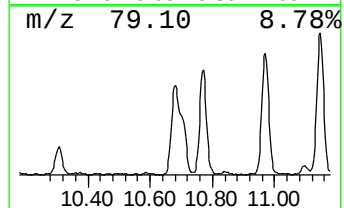
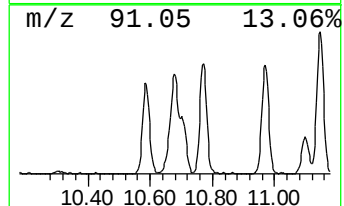
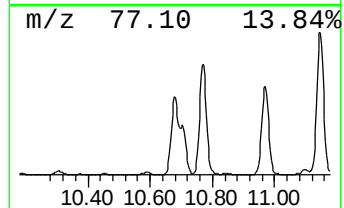
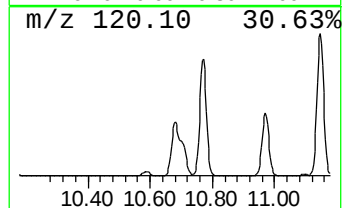
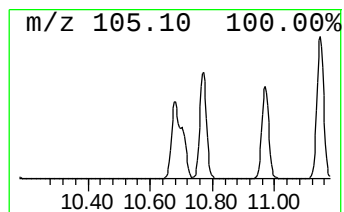
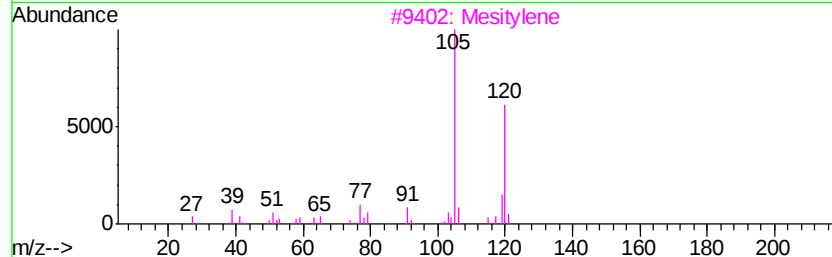
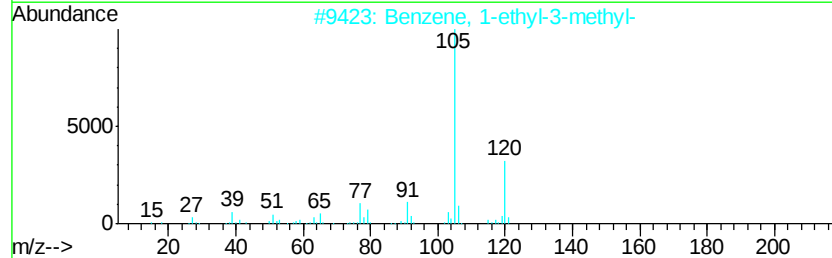
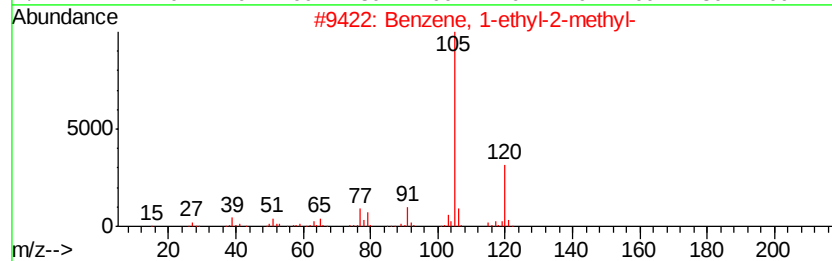
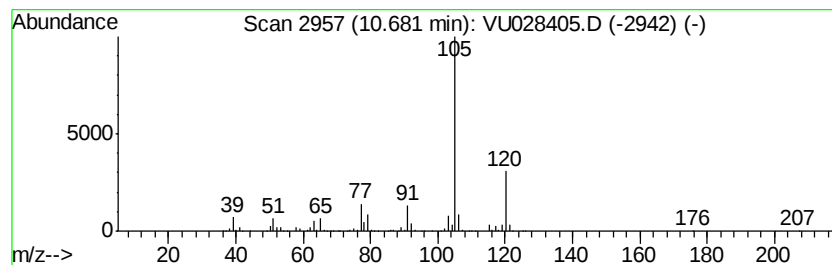
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U112818W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.68	32.98 ug/l	1087740	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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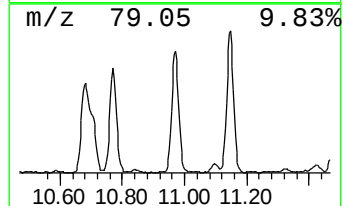
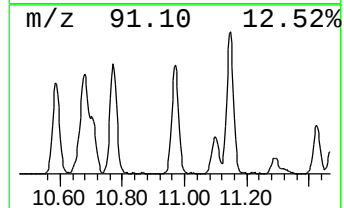
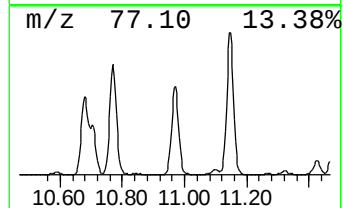
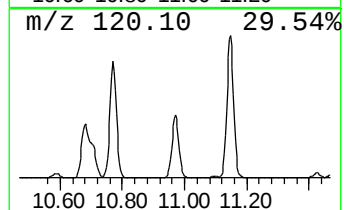
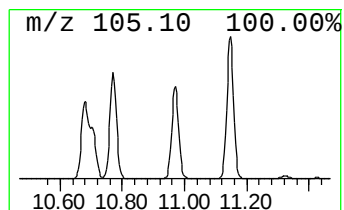
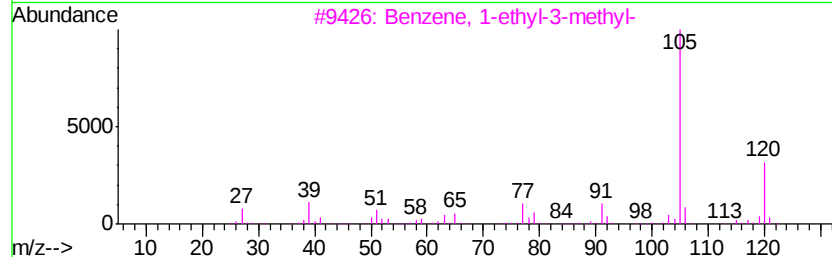
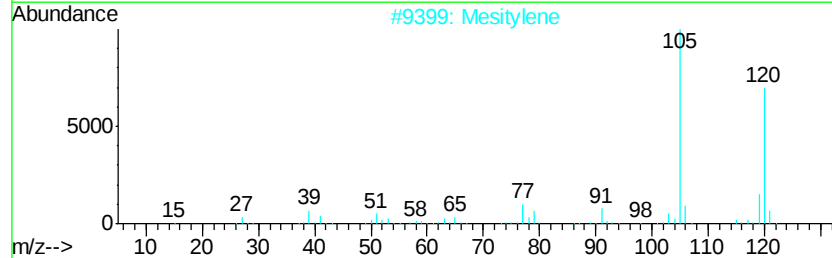
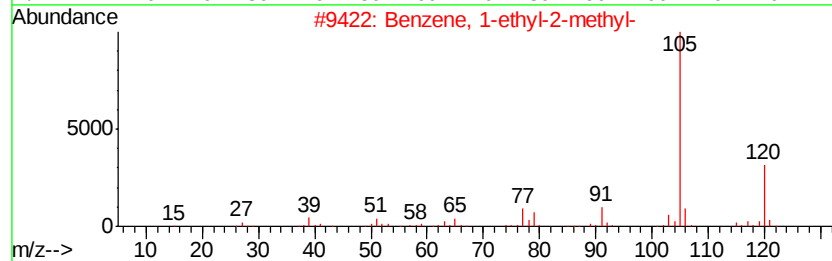
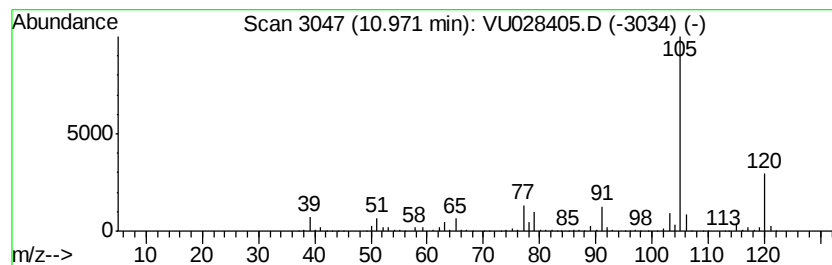
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1-ethyl-3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.97	25.77 ug/l	850015	1,4-Dichlorobenzene-d4	11.48

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



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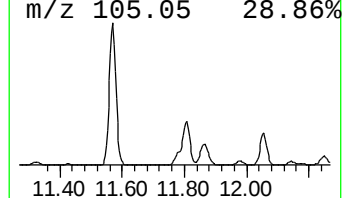
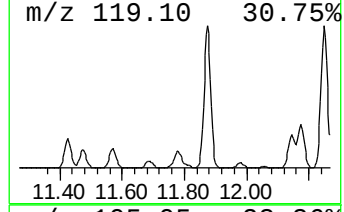
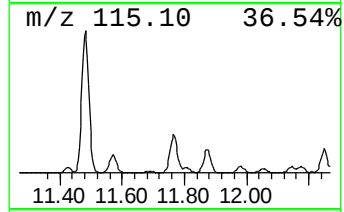
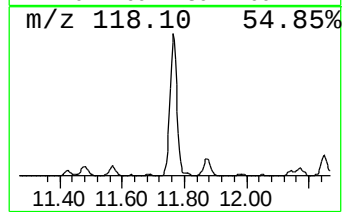
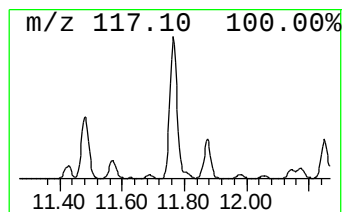
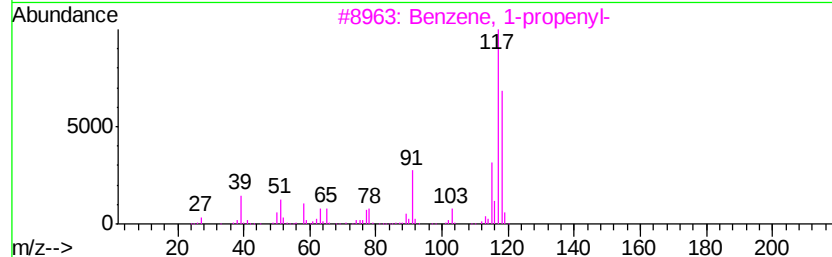
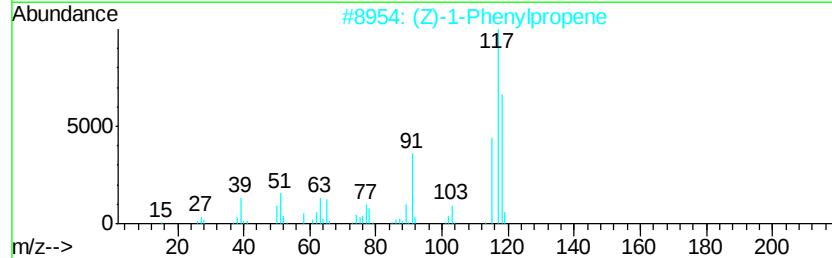
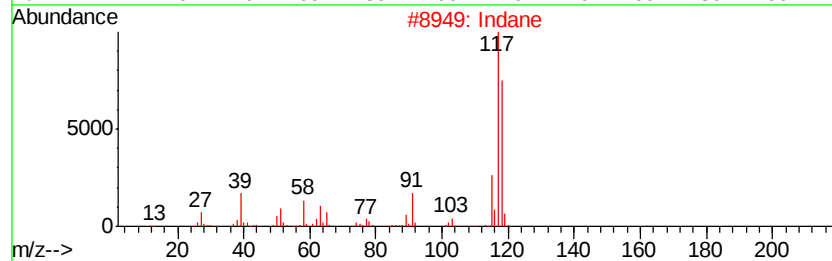
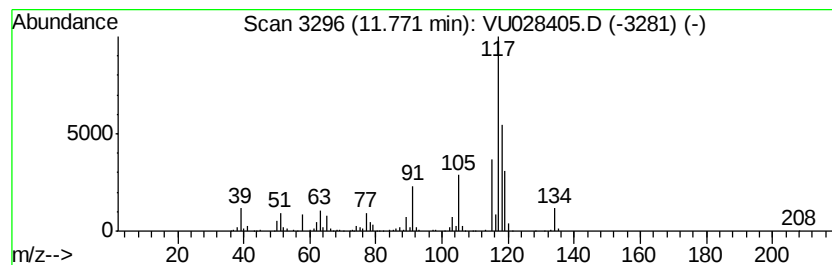
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Indane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	16.71 ug/l	551034	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	76
2		(Z)-1-Phenylpropene	118	C9H10	000766-90-5	70
3		Benzene, 1-propenyl-	118	C9H10	000637-50-3	70
4		Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
5		Benzene, 2-propenyl-	118	C9H10	000300-57-2	68



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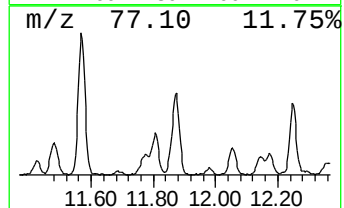
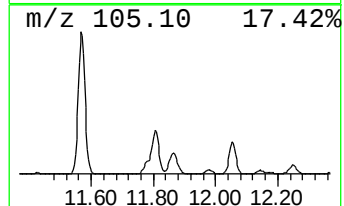
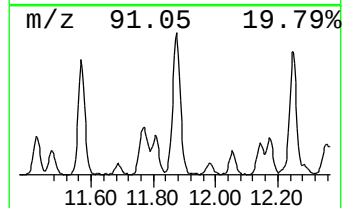
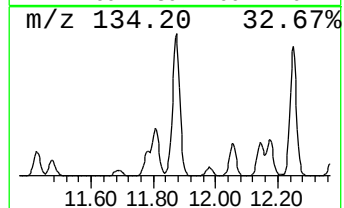
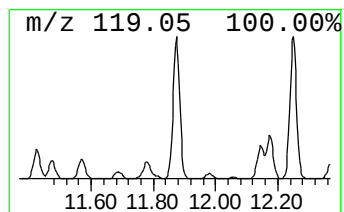
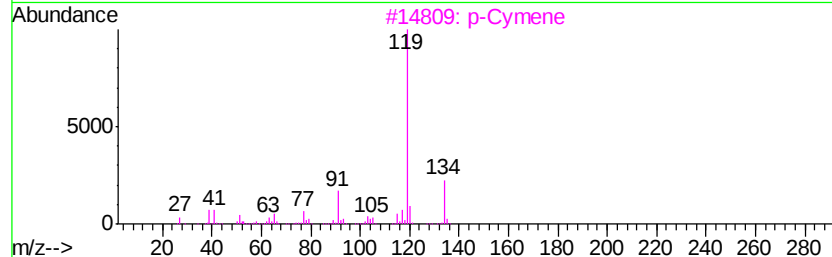
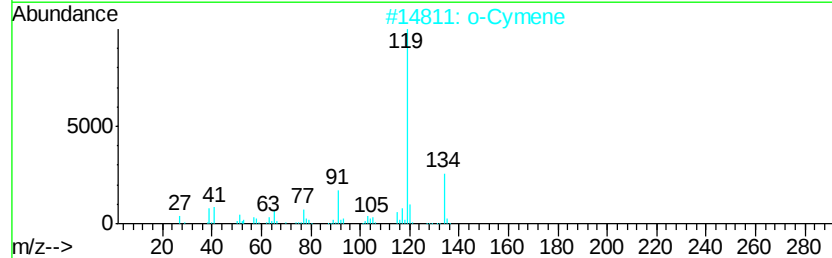
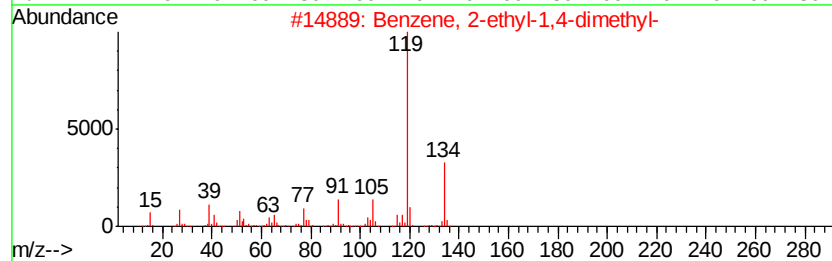
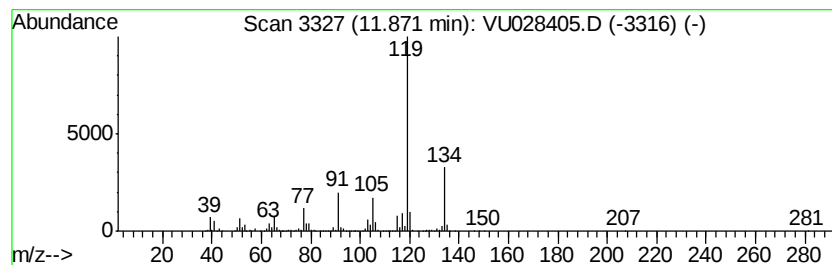
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.87	33.00 ug/l	1088220	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2		o-Cymene	134	C10H14	000527-84-4	95
3		p-Cymene	134	C10H14	000099-87-6	95
4		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU112918\
Data File : VU028405.D
Acq On : 29 Nov 2018 19:39
Operator : MD/SY
Sample : J6134-07
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DUP-1

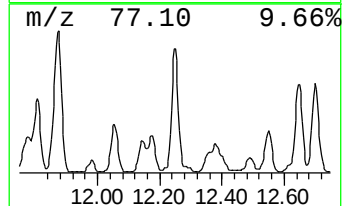
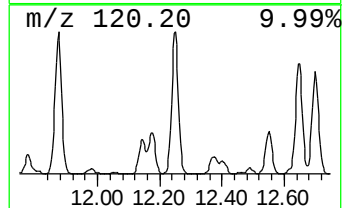
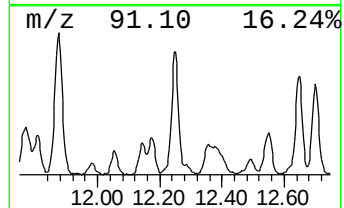
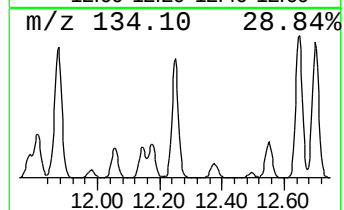
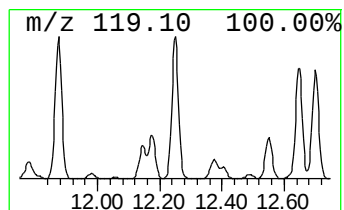
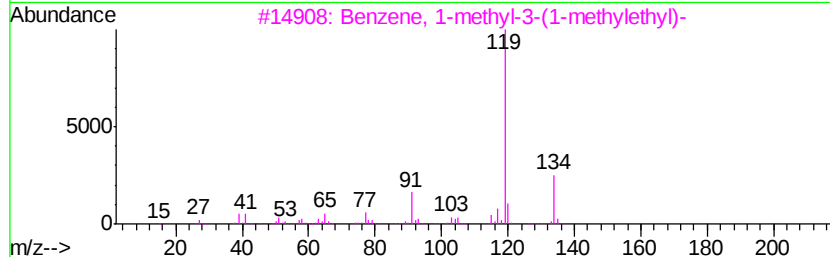
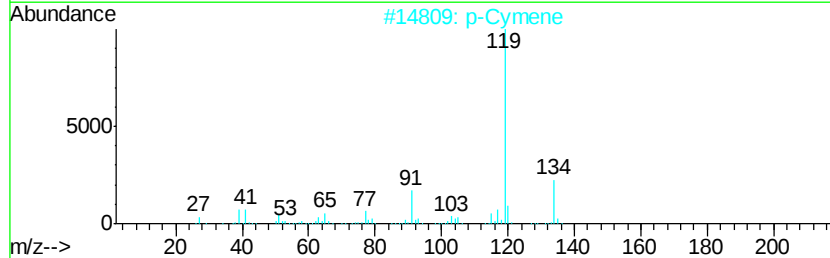
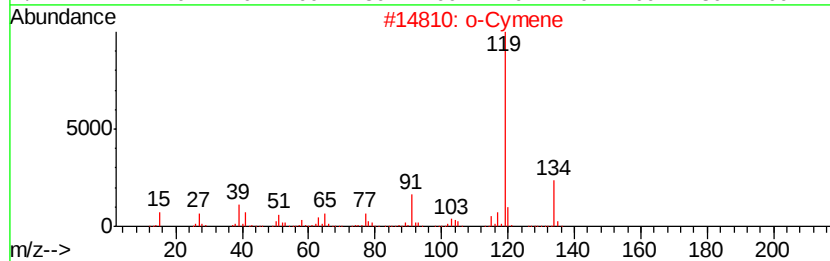
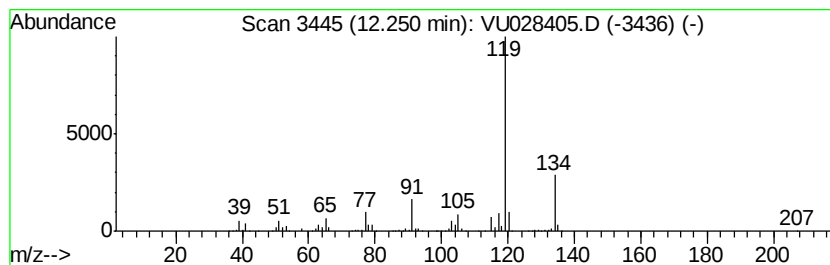
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Quant Title : SW846 8260

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

Peak Number 6 o-Cymene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.25	23.38 ug/l	770962	1,4-Dichlorobenzene-d4	11.48

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		o-Cymene	134	C10H14	000527-84-4	95
2		p-Cymene	134	C10H14	000099-87-6	95
3		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93
5		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU112918\
Data File : VU028405.D
Acq On : 29 Nov 2018 19:39
Operator : MD/SY
Sample : J6134-07
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DUP-1

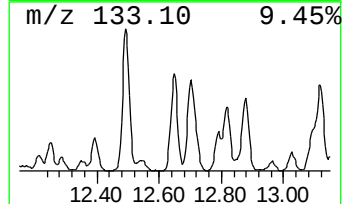
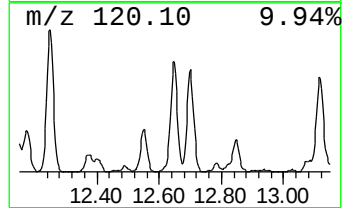
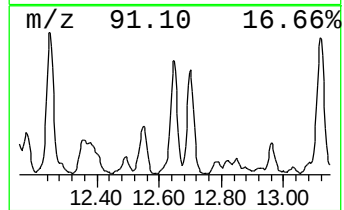
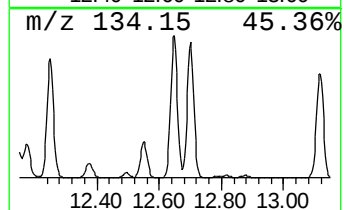
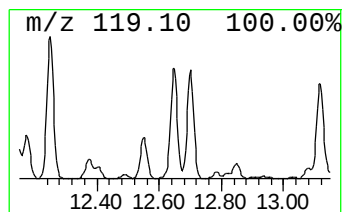
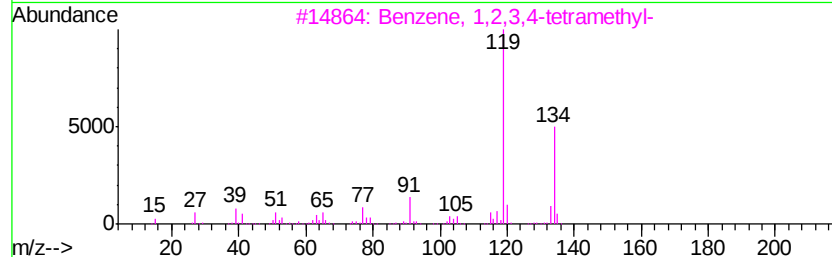
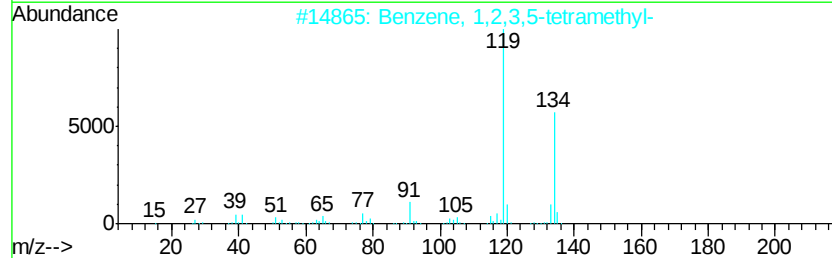
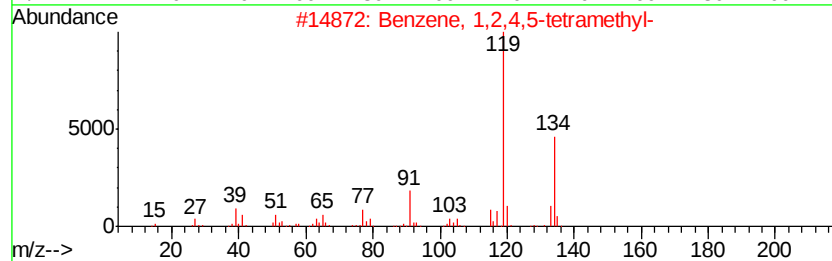
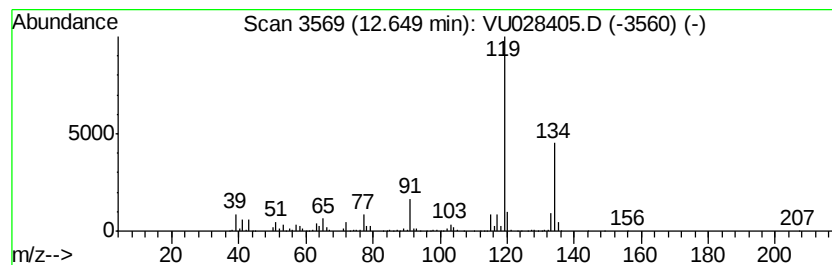
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	21.98 ug/l	724975	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	95
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
4		o-Cymene	134	C10H14	000527-84-4	93
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU112918\
Data File : VU028405.D
Acq On : 29 Nov 2018 19:39
Operator : MD/SY
Sample : J6134-07
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DUP-1

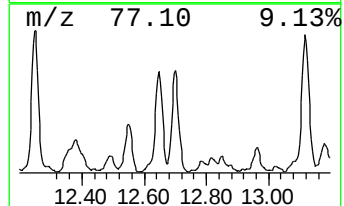
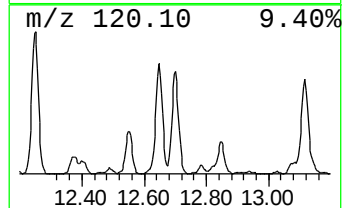
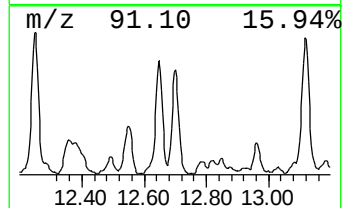
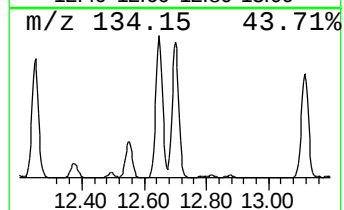
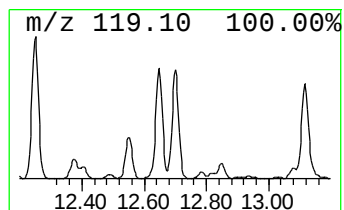
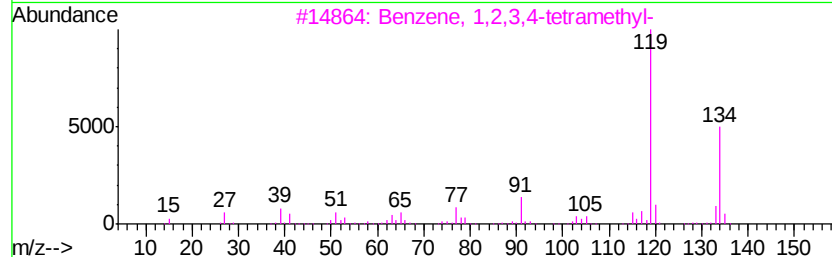
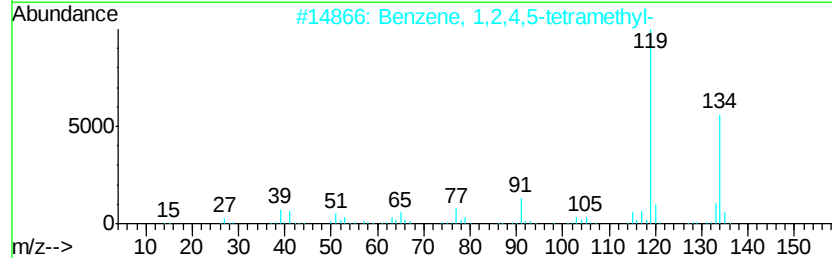
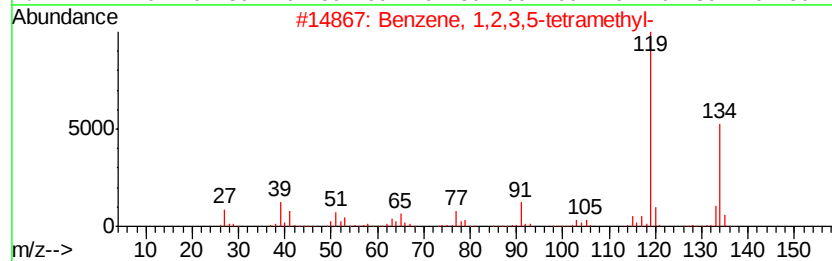
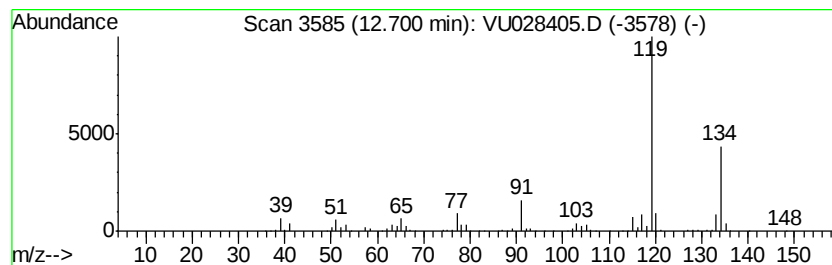
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	22.51 ug/l	742246	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
5		o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU112918\
Data File : VU028405.D
Acq On : 29 Nov 2018 19:39
Operator : MD/SY
Sample : J6134-07
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DUP-1

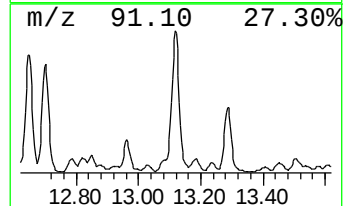
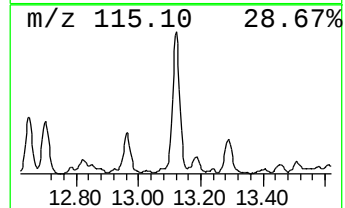
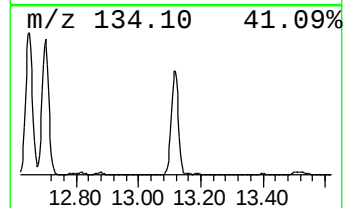
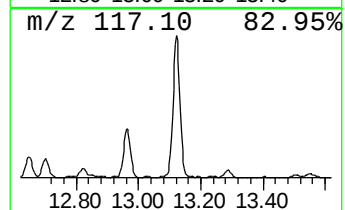
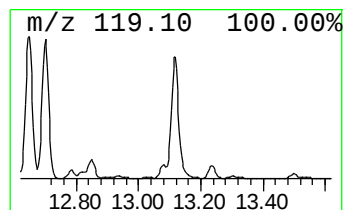
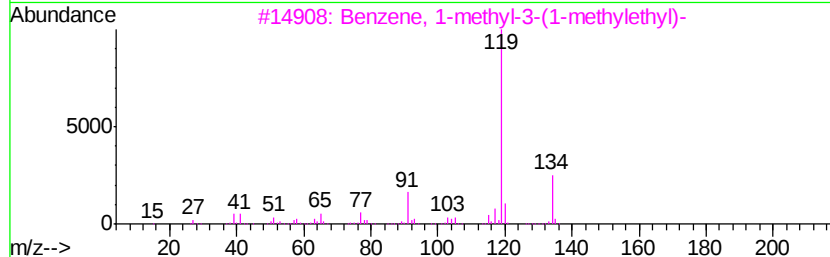
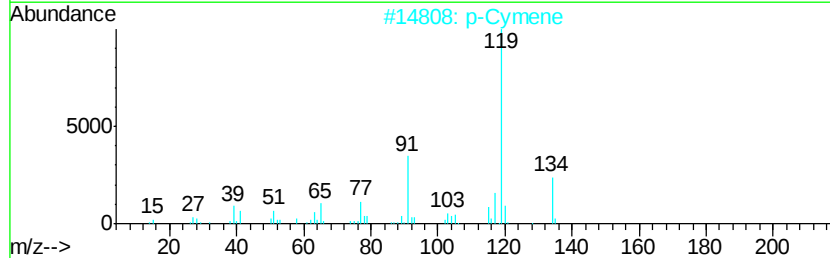
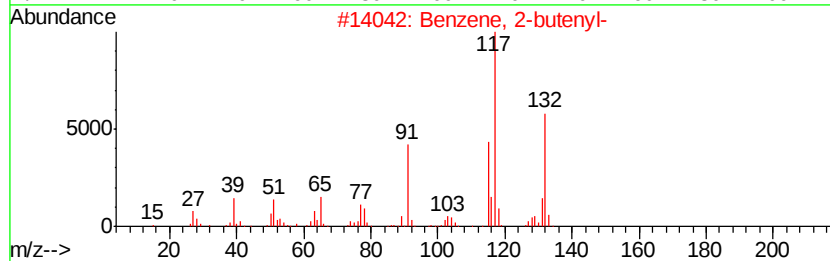
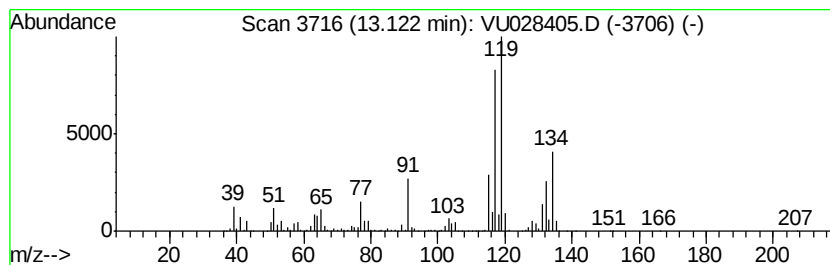
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 2-butenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	34.87 ug/l	1149850	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	83
2		p-Cymene	134	C10H14	000099-87-6	60
3		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	60
4		o-Cymene	134	C10H14	000527-84-4	60
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU112918\
Data File : VU028405.D
Acq On : 29 Nov 2018 19:39
Operator : MD/SY
Sample : J6134-07
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DUP-1

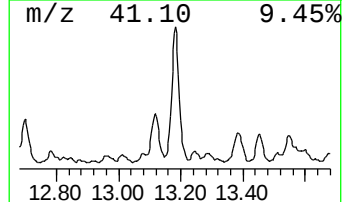
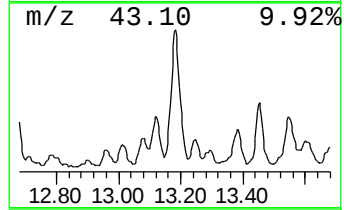
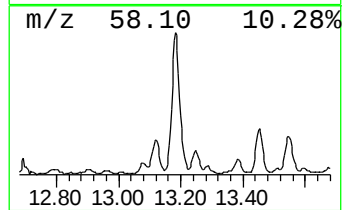
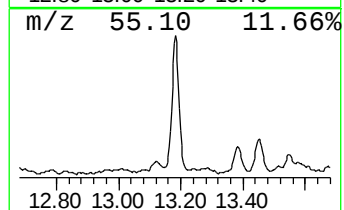
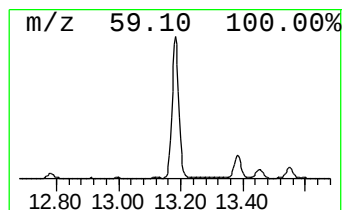
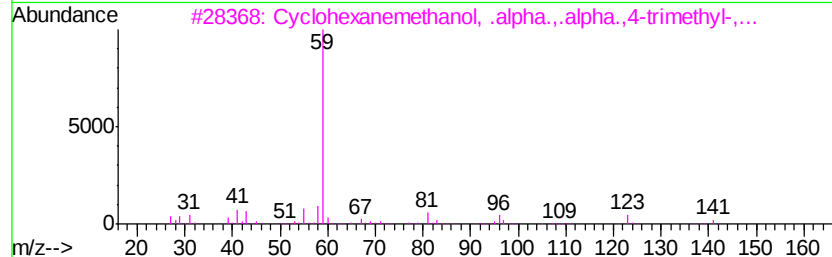
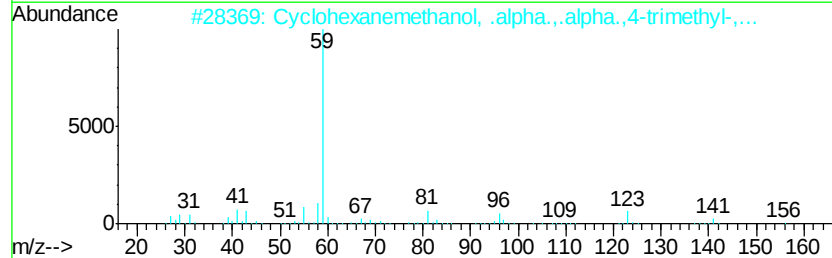
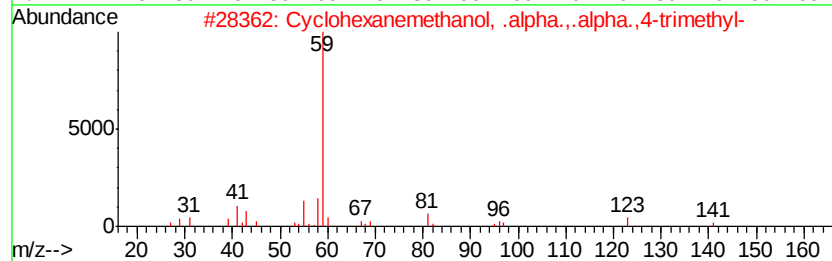
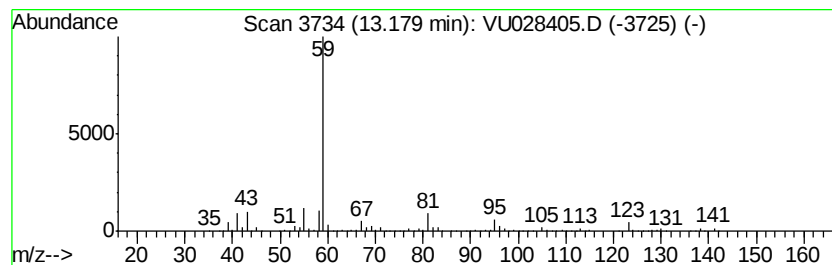
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U112818W.M
Quant Title : SW846 8260

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

Peak Number 10 Cyclohexanemethanol, .alpha... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.18	33.05 ug/l	1089920	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	78
2		Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	005114-00-1	72
3		Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	007322-63-6	72
4		1-Cyclohexylethanol, methyl ether	142	C9H18O	1000365-12-8	72
5		2,4-Dimethyl-4-penten-2-ol	114	C7H14O	019781-53-4	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU112918\
Data File : VU028405.D
Acq On : 29 Nov 2018 19:39
Operator : MD/SY
Sample : J6134-07
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DUP-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U112818W.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
Benzene, 1-ethyl-...	10.68	33.0	ug/l	1087740	4	11.48	1648950	50.0
Benzene, 1-ethyl-...	10.97	25.8	ug/l	850015	4	11.48	1648950	50.0
Indane	11.77	16.7	ug/l	551034	4	11.48	1648950	50.0
Benzene, 2-ethyl-...	11.87	33.0	ug/l	1088220	4	11.48	1648950	50.0
o-Cymene	12.25	23.4	ug/l	770962	4	11.48	1648950	50.0
Benzene, 1,2,4,5-...	12.65	22.0	ug/l	724975	4	11.48	1648950	50.0
Benzene, 1,2,3,5-...	12.70	22.5	ug/l	742246	4	11.48	1648950	50.0
Benzene, 2-butenyl-	13.12	34.9	ug/l	1149850	4	11.48	1648950	50.0
Cyclohexanemethan...	13.18	33.0	ug/l	1089920	4	11.48	1648950	50.0