

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU110618\
 Data File : VU027985.D
 Acq On : 06 Nov 2018 14:27
 Operator : MD/SY
 Sample : J5822-06
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MW-11

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\82U110218W.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	61	72	80	rBV	47644	64327	2.53%	0.237%
2	2.447	384	396	411	rBV	35239	64998	2.56%	0.240%
3	2.990	553	565	588	rVB3	15924	32329	1.27%	0.119%
4	4.228	929	950	989	rVV	1035707	2542580	100.00%	9.374%
5	4.887	1137	1155	1171	rBV	175421	403291	15.86%	1.487%
6	4.987	1172	1186	1208	rVB	271813	603815	23.75%	2.226%
7	5.312	1273	1287	1303	rBV	209741	446245	17.55%	1.645%
8	5.395	1304	1313	1328	rVB2	14421	31733	1.25%	0.117%
9	5.887	1451	1466	1497	rBV	451084	915597	36.01%	3.375%
10	7.057	1816	1830	1841	rBV7	12247	31992	1.26%	0.118%
11	7.565	1969	1988	2000	rBV	742582	1301795	51.20%	4.799%
12	7.639	2000	2011	2027	rVB	518144	892279	35.09%	3.290%
13	8.180	2166	2179	2188	rBV	14699	28472	1.12%	0.105%
14	9.090	2450	2462	2486	rBV	652463	1068665	42.03%	3.940%
15	9.250	2501	2512	2526	rBV	135328	217516	8.55%	0.802%
16	9.372	2538	2550	2566	rVV	859215	1422232	55.94%	5.243%
17	9.720	2645	2658	2664	rBV7	13129	27715	1.09%	0.102%
18	9.781	2665	2677	2699	rVB	558683	901617	35.46%	3.324%
19	10.044	2743	2759	2766	rBV7	19068	38241	1.50%	0.141%
20	10.308	2830	2841	2870	rVB	516026	857609	33.73%	3.162%
21	10.588	2919	2928	2941	rBV	34076	58856	2.31%	0.217%
22	10.681	2944	2957	2974	rVV	373051	843734	33.18%	3.111%
23	10.771	2974	2985	2994	rVV	520788	810437	31.87%	2.988%
24	10.813	2994	2998	3009	rVB2	76621	101041	3.97%	0.373%
25	10.970	3036	3047	3068	rVB	408072	648396	25.50%	2.390%
26	11.102	3073	3088	3093	rBV3	42566	88207	3.47%	0.325%
27	11.147	3093	3102	3115	rVB	775142	1184742	46.60%	4.368%
28	11.427	3171	3189	3197	rBV3	112683	220431	8.67%	0.813%
29	11.482	3197	3206	3218	rVV	682122	1030124	40.51%	3.798%
30	11.572	3224	3234	3255	rVB	679197	1082911	42.59%	3.992%
31	11.694	3265	3272	3281	rVB3	35482	58926	2.32%	0.217%
32	11.768	3281	3295	3302	rVV3	202044	411996	16.20%	1.519%
33	11.810	3303	3308	3316	rVV	192223	276169	10.86%	1.018%
34	11.874	3316	3328	3343	rVB2	497901	859527	33.81%	3.169%

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU110618\
 Data File : VU027985.D
 Acq On : 06 Nov 2018 14:27
 Operator : MD/SY
 Sample : J5822-06
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MW-11

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\82U110218W.M
 Title : SW846 8260

35	11.983	3353	3362	3367	rBV2	37668	59553	2.34%	0.220%
36	12.025	3368	3375	3379	rVV	83817	121141	4.76%	0.447%
37	12.054	3379	3384	3400	rVB	132616	209558	8.24%	0.773%
38	12.147	3402	3413	3417	rBV	131515	204990	8.06%	0.756%
39	12.176	3417	3422	3430	rVB	124827	172677	6.79%	0.637%
40	12.250	3436	3445	3454	rBV	424258	593516	23.34%	2.188%
41	12.356	3467	3478	3484	rBV	125359	236946	9.32%	0.874%
42	12.446	3500	3506	3513	rVB2	51566	64232	2.53%	0.237%
43	12.491	3514	3520	3529	rVB2	63139	91480	3.60%	0.337%
44	12.549	3530	3538	3546	rVB2	137566	200076	7.87%	0.738%
45	12.649	3560	3569	3578	rVV	405146	625232	24.59%	2.305%
46	12.704	3578	3586	3601	rVB	378968	612530	24.09%	2.258%
47	12.787	3603	3612	3617	rBV4	55678	87849	3.46%	0.324%
48	12.964	3659	3667	3677	rVB2	113407	159315	6.27%	0.587%
49	13.083	3696	3704	3706	rBV3	49470	61384	2.41%	0.226%
50	13.121	3707	3716	3726	rVV2	566153	931451	36.63%	3.434%
51	13.183	3726	3735	3747	rVV	469652	758171	29.82%	2.795%
52	13.289	3759	3768	3786	rVB3	136065	239256	9.41%	0.882%
53	13.385	3790	3798	3811	rVB2	85717	154336	6.07%	0.569%
54	13.456	3811	3820	3829	rBV4	125427	195190	7.68%	0.720%
55	13.511	3829	3837	3843	rBV3	57023	93369	3.67%	0.344%
56	13.549	3844	3849	3860	rVB6	28855	43911	1.73%	0.162%
57	13.678	3883	3889	3897	rBV10	15525	29132	1.15%	0.107%
58	13.742	3898	3909	3919	rBV	252170	415395	16.34%	1.531%
59	13.858	3939	3945	3950	rBV7	18732	28750	1.13%	0.106%
60	14.012	3987	3993	4006	rBV8	14813	36897	1.45%	0.136%
61	14.147	4026	4035	4041	rBV6	71627	127121	5.00%	0.469%
62	14.183	4042	4046	4061	rVB2	53270	90059	3.54%	0.332%
63	14.337	4086	4094	4105	rVB6	40884	87303	3.43%	0.322%
64	14.398	4106	4113	4123	rBV6	24155	43591	1.71%	0.161%
65	14.549	4151	4160	4174	rVB6	34528	70447	2.77%	0.260%
66	14.710	4206	4210	4217	rVB4	24135	31098	1.22%	0.115%
67	14.752	4218	4223	4239	rVB3	30138	45916	1.81%	0.169%
68	14.877	4253	4262	4276	rVB2	137216	261064	10.27%	0.962%
69	15.060	4311	4319	4336	rBV3	92649	189801	7.46%	0.700%
70	15.195	4353	4361	4370	rBV6	31409	54071	2.13%	0.199%
71	15.662	4499	4506	4515	rVB2	44921	62704	2.47%	0.231%

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU110618\
Data File : VU027985.D
Acq On : 06 Nov 2018 14:27
Operator : MD/SY
Sample : J5822-06
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-11

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

Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.M
Title : SW846 8260

72	15.977	4598	4604	4614	rVB	25361	35028	1.38%	0.129%
73	16.067	4624	4632	4638	rBV4	20797	31854	1.25%	0.117%

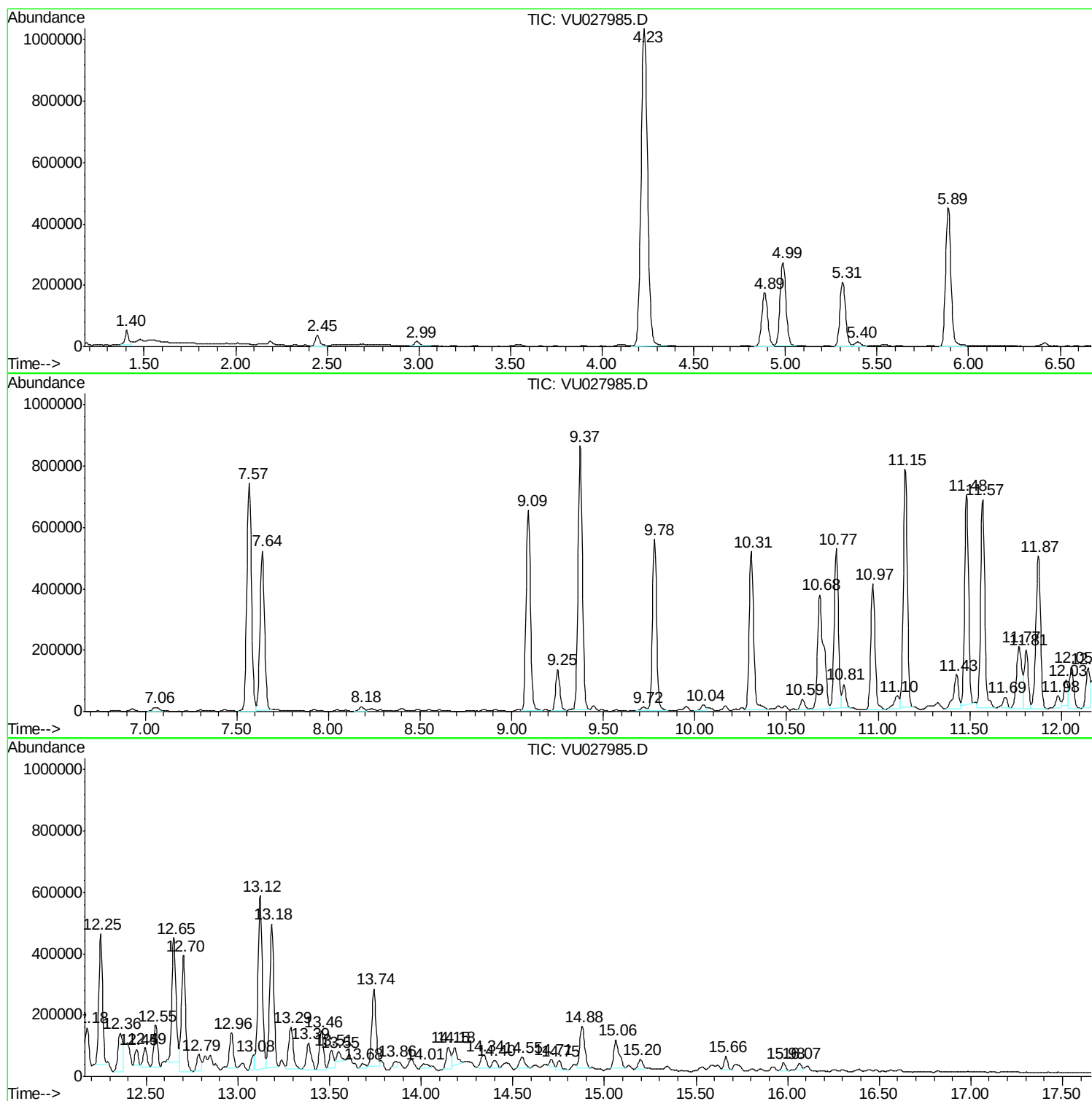
Sum of corrected areas: 27124939

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU110618\
Data File : VU027985.D
Acq On : 06 Nov 2018 14:27
Operator : MD/SY
Sample : J5822-06
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampled :
MW-11

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.M
Quant Title : SW846 8260

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU110618\
Data File : VU027985.D
Acq On : 06 Nov 2018 14:27
Operator : MD/SY
Sample : J5822-06
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-11

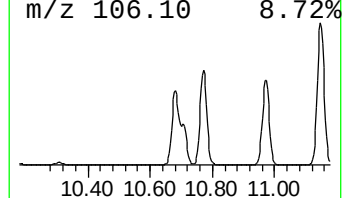
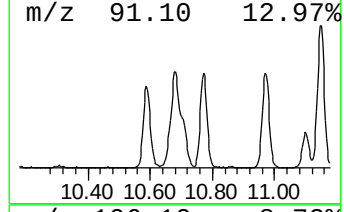
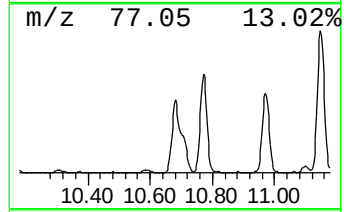
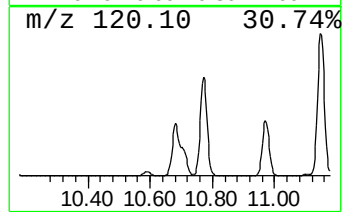
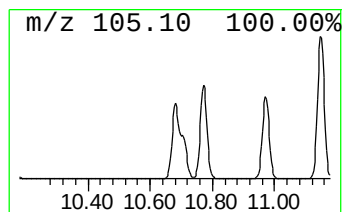
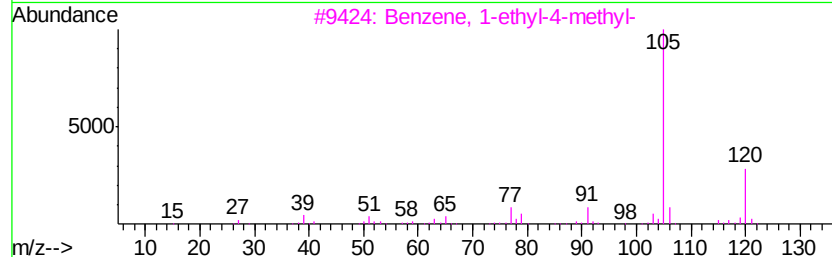
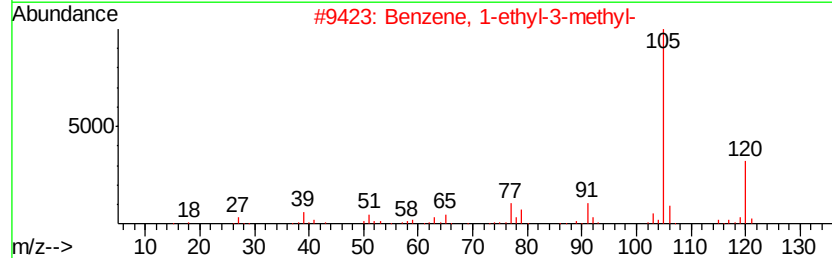
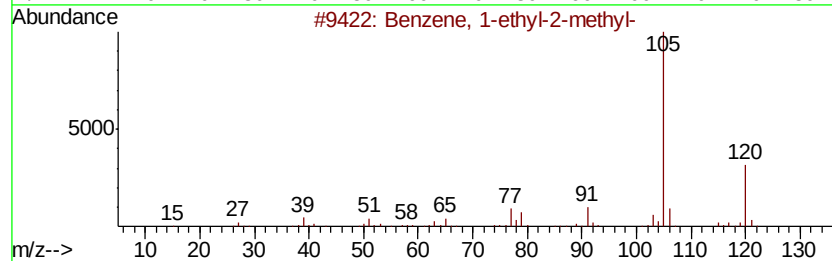
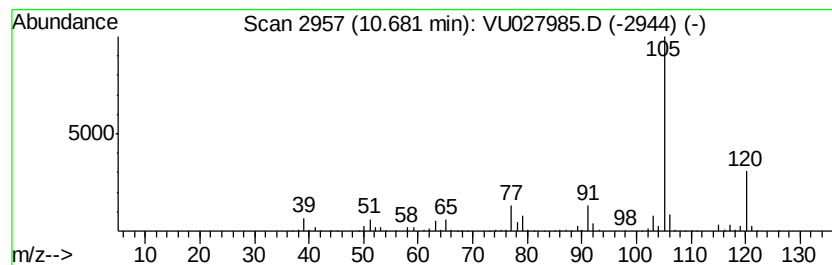
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.68	40.95 ug/l	843734	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	94
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU110618\
Data File : VU027985.D
Acq On : 06 Nov 2018 14:27
Operator : MD/SY
Sample : J5822-06
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

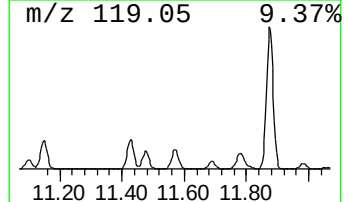
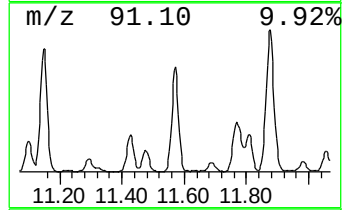
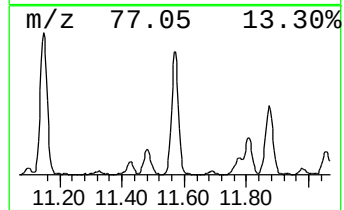
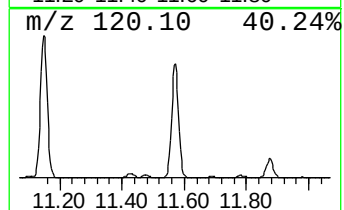
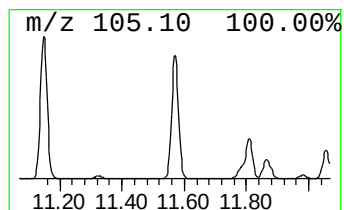
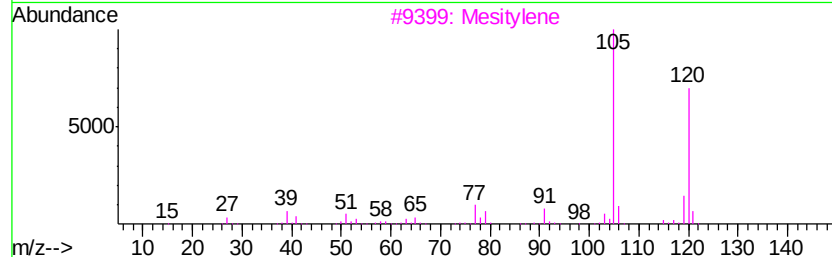
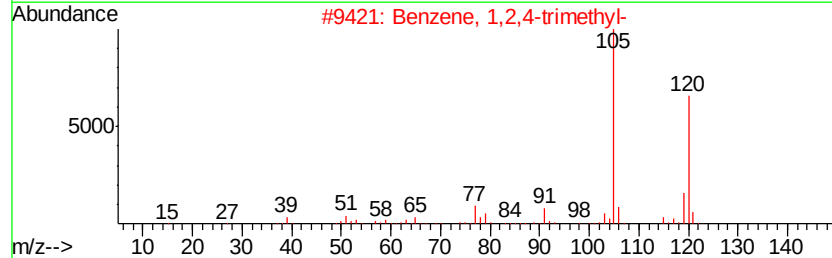
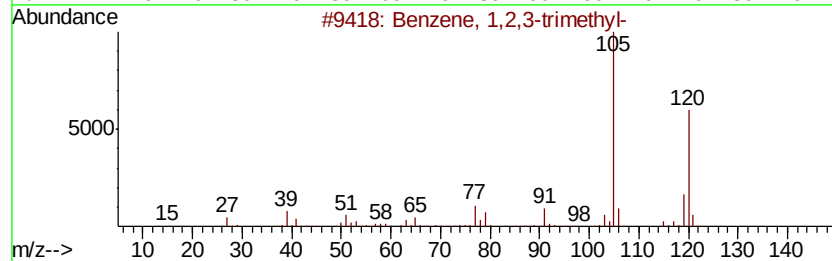
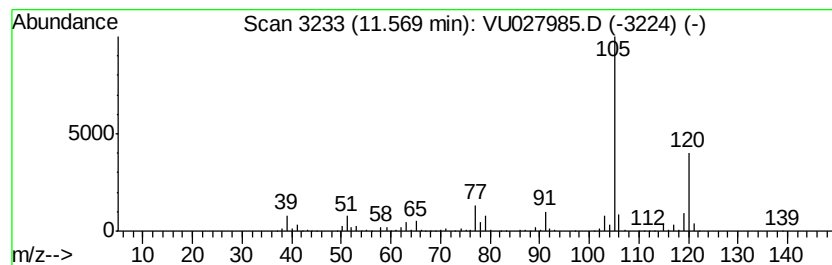
Instrument :
MSVOA_U
ClientSampled :
MW-11

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.57	52.56 ug/l	1082910	1,4-Dichlorobenzene-d4	11.48		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		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	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU110618\
Data File : VU027985.D
Acq On : 06 Nov 2018 14:27
Operator : MD/SY
Sample : J5822-06
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

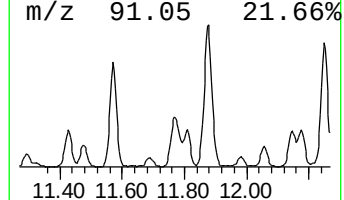
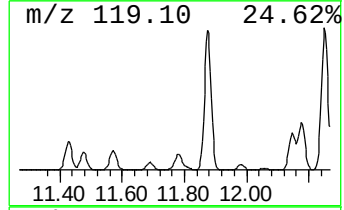
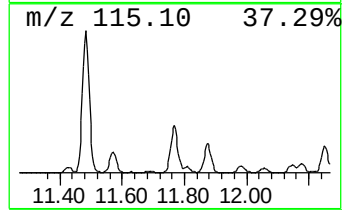
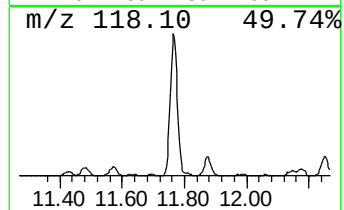
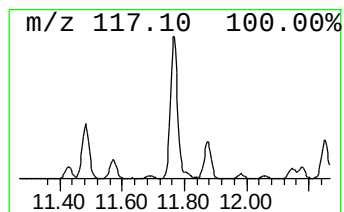
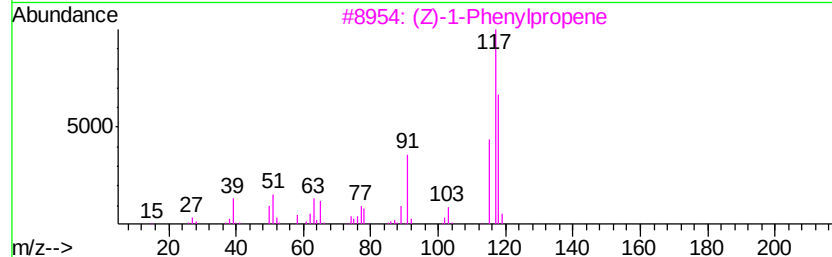
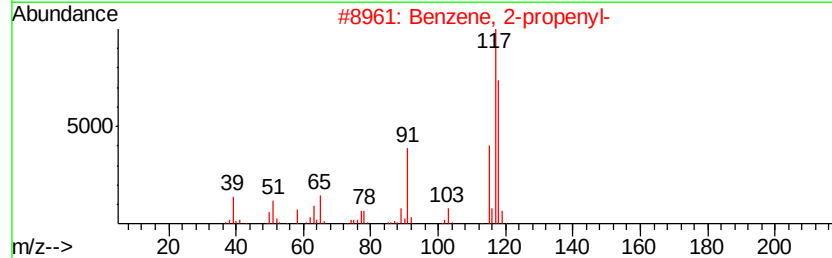
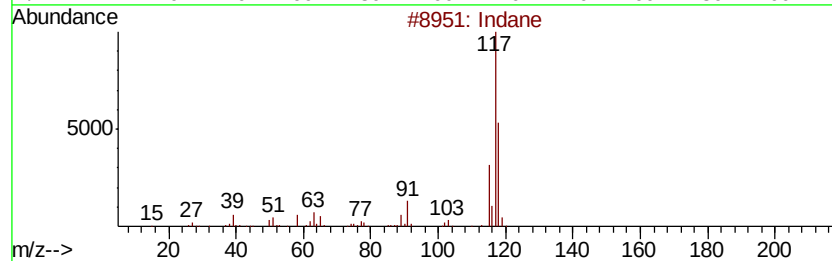
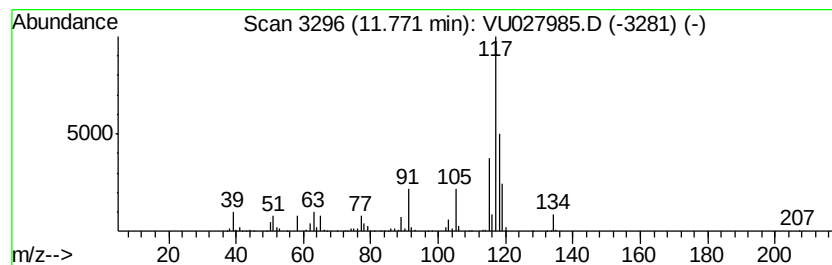
Instrument :
MSVOA_U
ClientSampleId :
MW-11

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.M
Quant Title : SW846 8260

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	20.00 ug/l	411996	1,4-Dichlorobenzene-d4	11.48
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indane		118 C9H10	000496-11-7 76
2	Benzene, 2-propenyl-		118 C9H10	000300-57-2 76
3	(Z)-1-Phenylpropene		118 C9H10	000766-90-5 76
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 58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU110618\
Data File : VU027985.D
Acq On : 06 Nov 2018 14:27
Operator : MD/SY
Sample : J5822-06
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampled :
MW-11

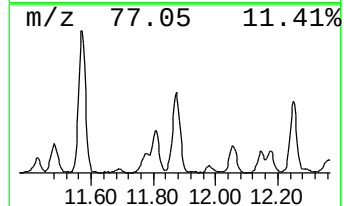
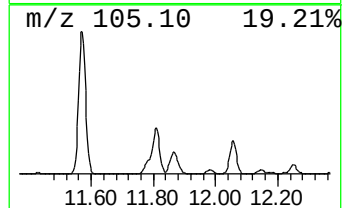
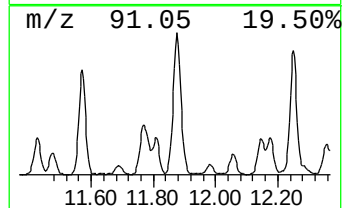
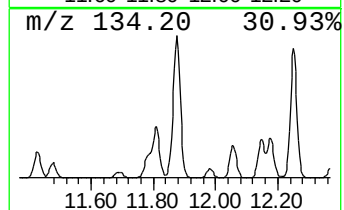
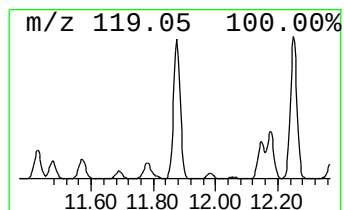
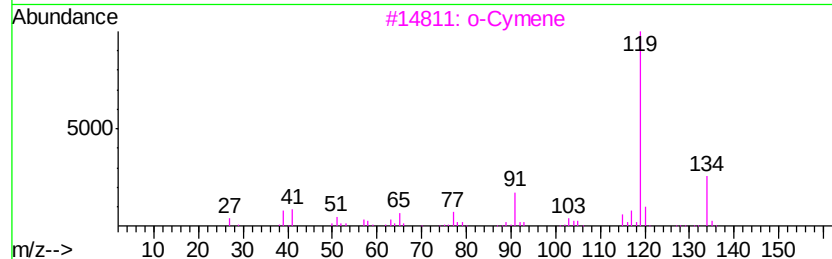
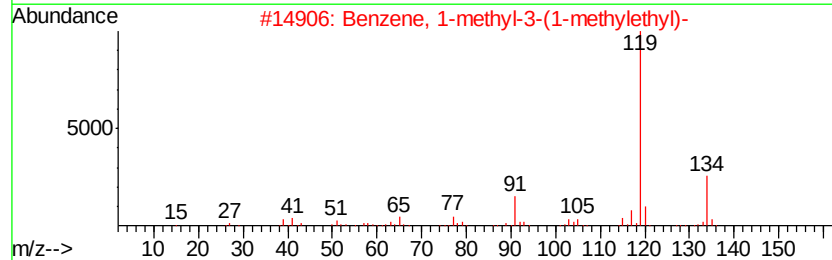
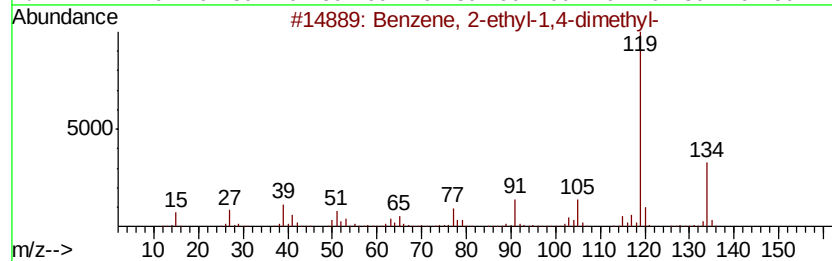
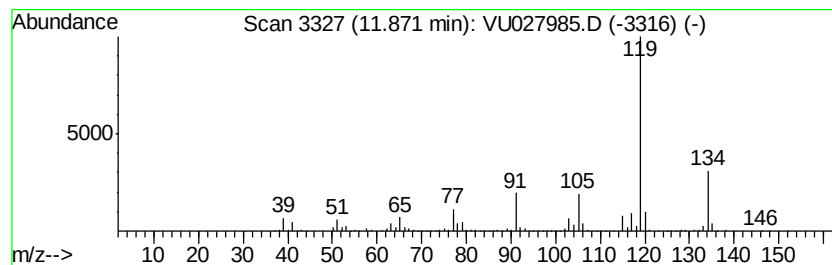
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.M
Quant Title : SW846 8260

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

Peak Number 5 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.87	41.72 ug/l	859527	1,4-Dichlorobenzene-d4	11.48

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2	Benzene, 1-methyl-3-(1-methylethyl-)	134	C10H14	000535-77-3	95
3	o-Cymene	134	C10H14	000527-84-4	95
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5	p-Cymene	134	C10H14	000099-87-6	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU110618\
Data File : VU027985.D
Acq On : 06 Nov 2018 14:27
Operator : MD/SY
Sample : J5822-06
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

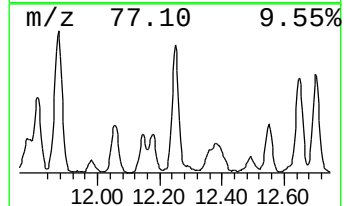
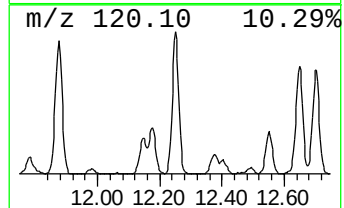
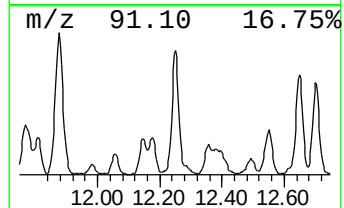
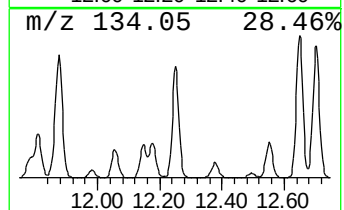
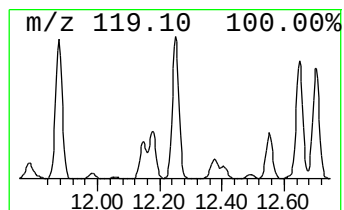
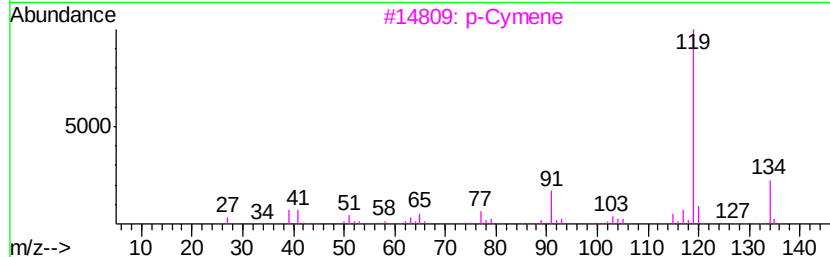
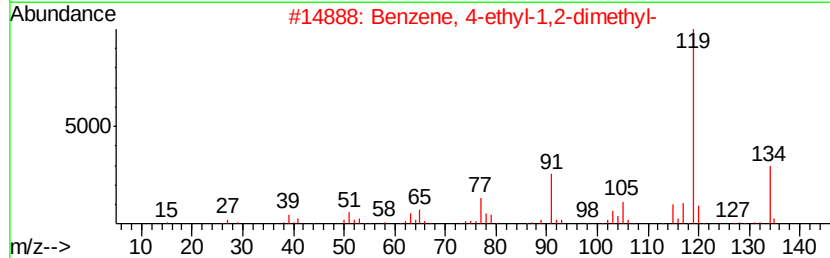
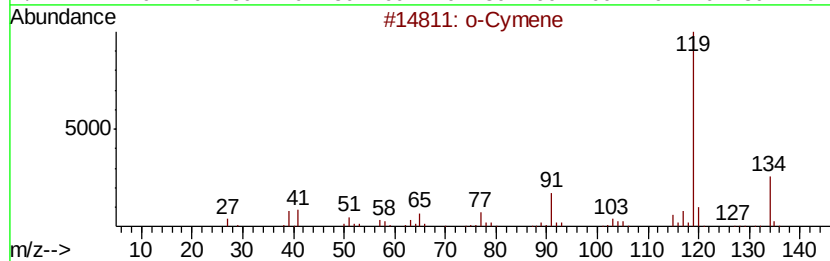
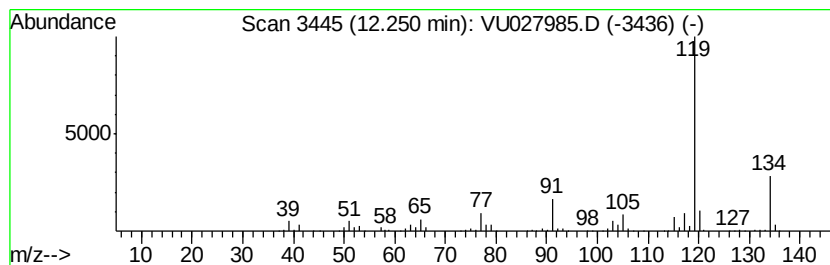
Instrument :
MSVOA_U
ClientSampleId :
MW-11

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.M
Quant Title : SW846 8260

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

Peak Number 6 o-Cymene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.25	28.81 ug/l	593516	1,4-Dichlorobenzene-d4	11.48	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	97
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3	p-Cymene	134	C10H14	000099-87-6	95
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU110618\
Data File : VU027985.D
Acq On : 06 Nov 2018 14:27
Operator : MD/SY
Sample : J5822-06
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-11

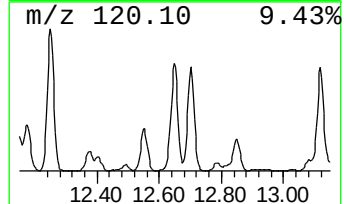
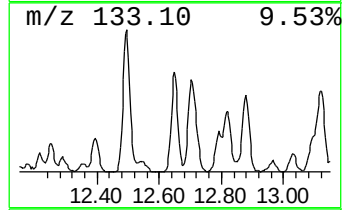
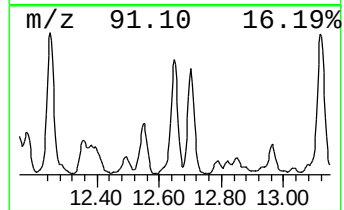
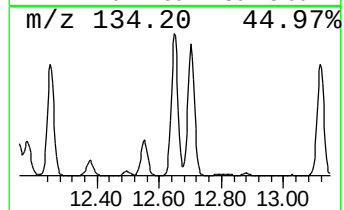
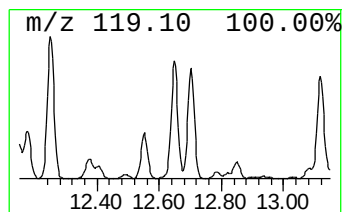
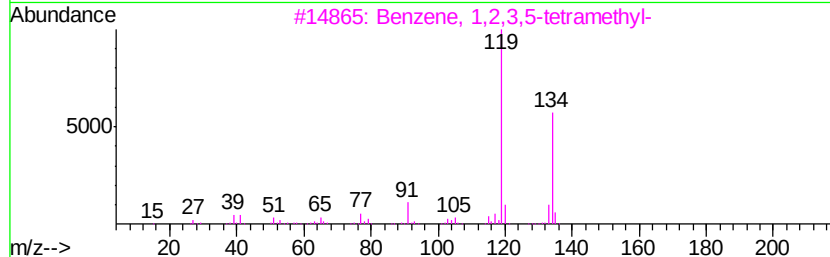
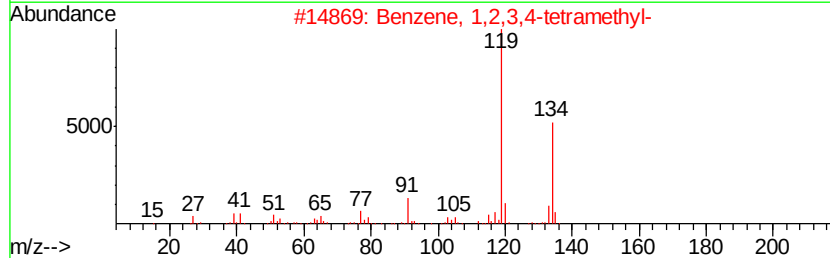
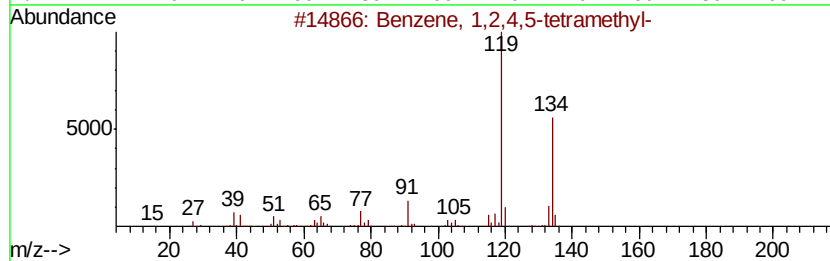
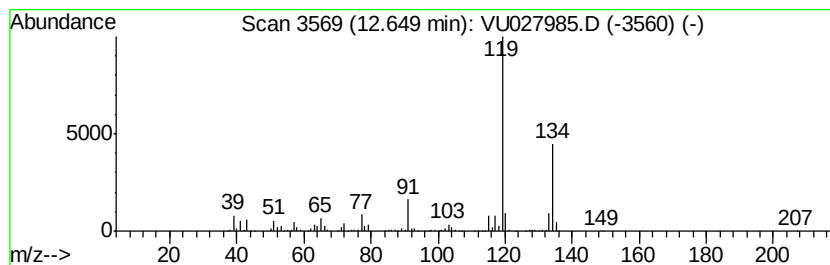
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	30.35 ug/l	625232	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, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93
5		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU110618\
Data File : VU027985.D
Acq On : 06 Nov 2018 14:27
Operator : MD/SY
Sample : J5822-06
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-11

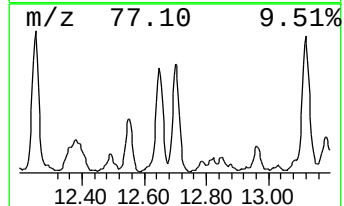
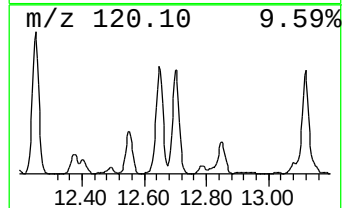
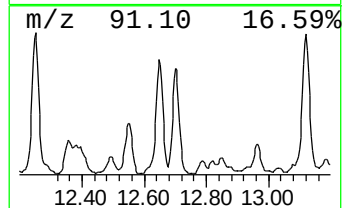
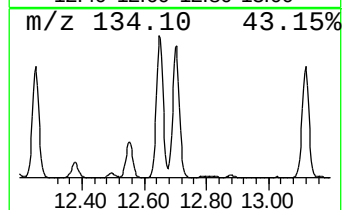
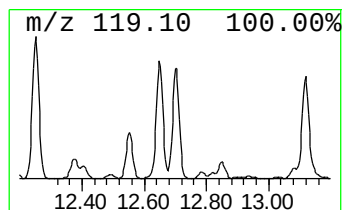
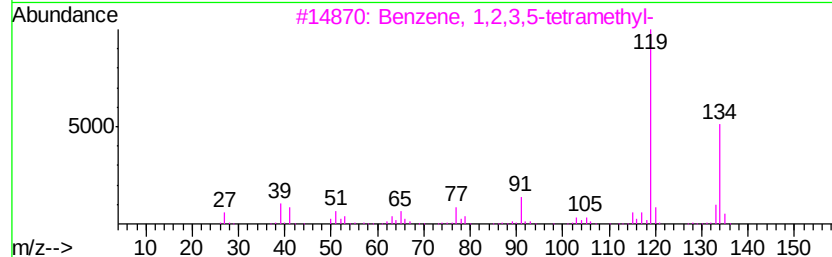
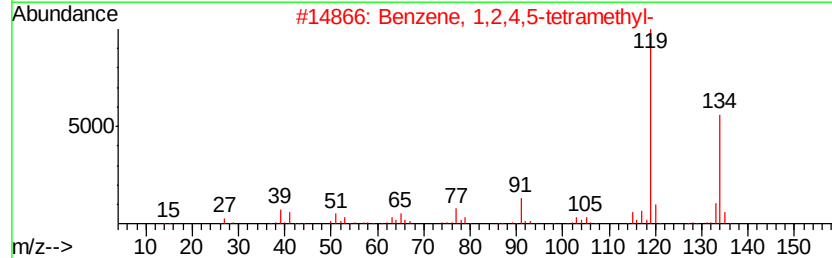
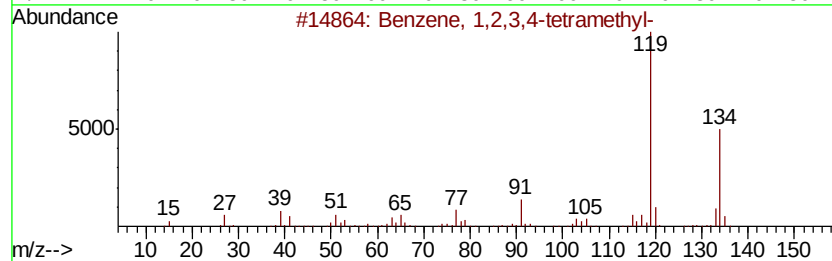
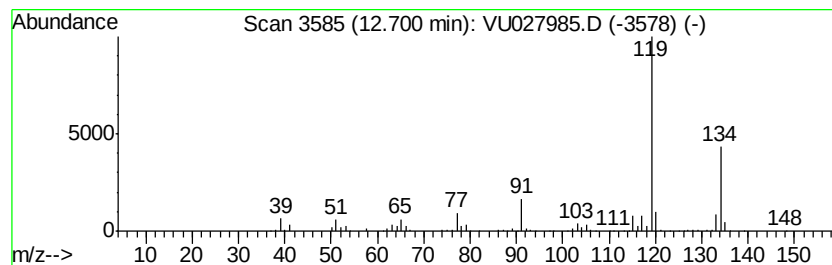
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	29.73 ug/l	612530	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU110618\
Data File : VU027985.D
Acq On : 06 Nov 2018 14:27
Operator : MD/SY
Sample : J5822-06
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-11

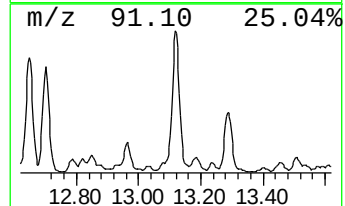
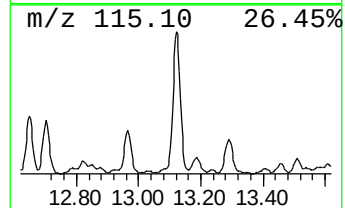
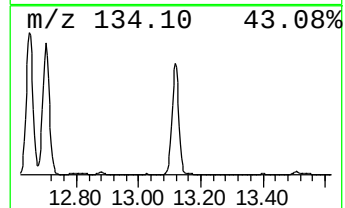
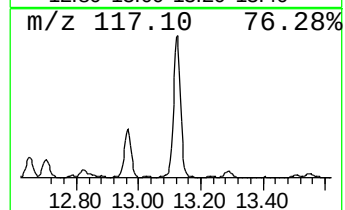
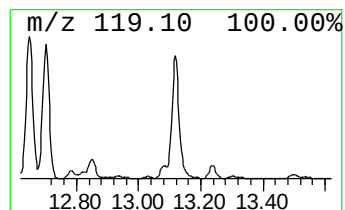
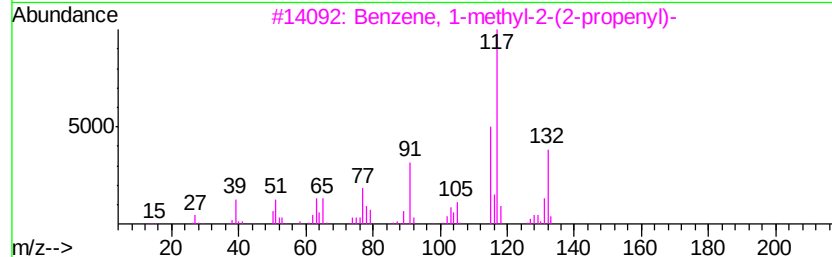
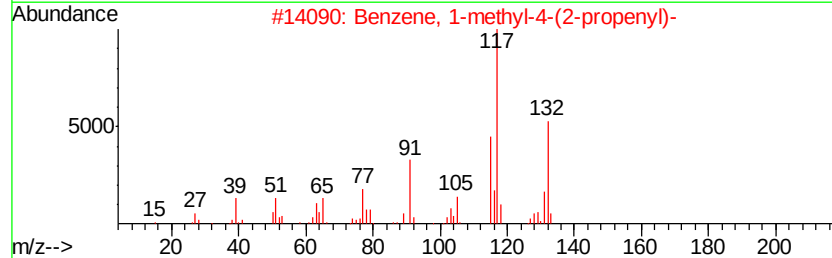
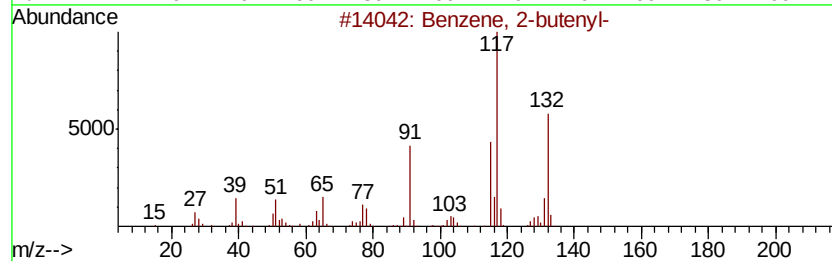
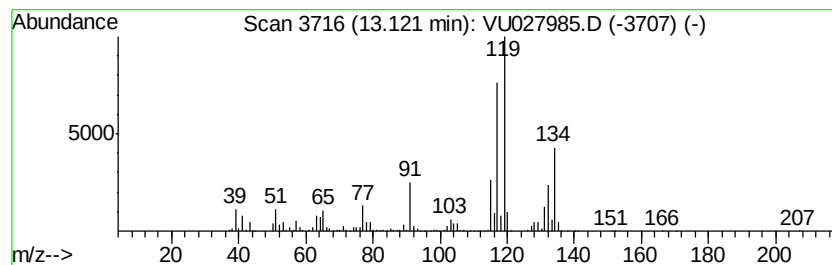
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.M
Quant Title : SW846 8260

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	45.21 ug/l	931451	1,4-Dichlorobenzene-d4	11.48

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 2-butenyl-	132	C10H12	001560-06-1	64
2		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	64
3		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	64
4		p-Cymene	134	C10H14	000099-87-6	55
5		o-Cymene	134	C10H14	000527-84-4	55



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU110618\
Data File : VU027985.D
Acq On : 06 Nov 2018 14:27
Operator : MD/SY
Sample : J5822-06
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-11

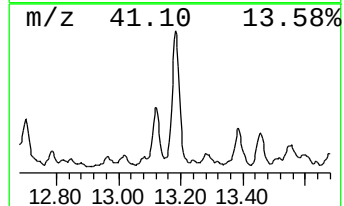
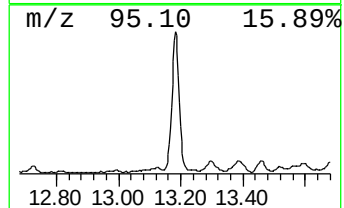
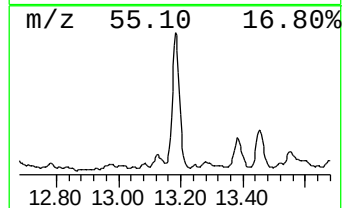
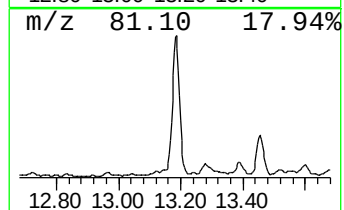
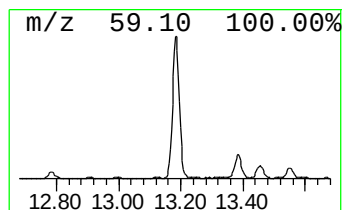
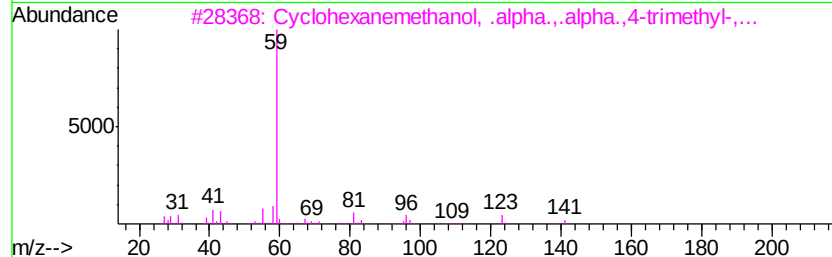
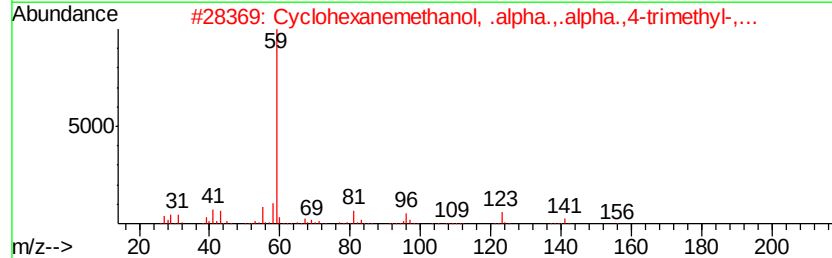
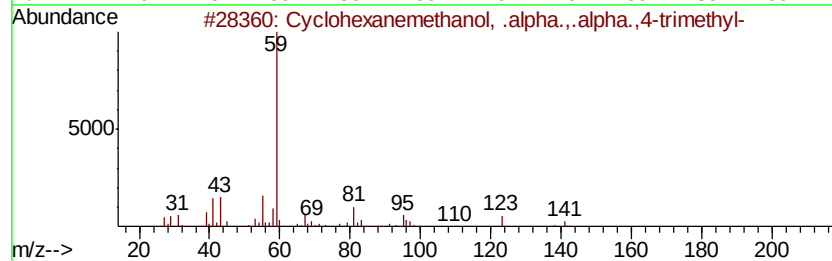
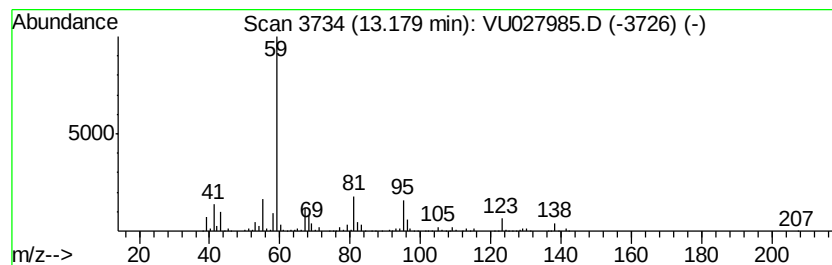
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.M
Quant Title : SW846 8260

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

Peak Number 10 Cyclohexanemethanol, .alpha... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.18	36.80 ug/l	758171	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanemethanol, .alpha...al...	156	C10H20O	000498-81-7	64
2		Cyclohexanemethanol, .alpha...al...	156	C10H20O	005114-00-1	50
3		Cyclohexanemethanol, .alpha...al...	156	C10H20O	007322-63-6	45
4		2,4-Dimethyl-4-penten-2-ol	114	C7H14O	019781-53-4	42
5		2-Cyclobutyl-2-propanol	114	C7H14O	059383-67-4	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU110618\
Data File : VU027985.D
Acq On : 06 Nov 2018 14:27
Operator : MD/SY
Sample : J5822-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-11

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.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	41.0	ug/l	843734	4	11.48	1030120	50.0
Benzene, 1,2,3-tr...	11.57	52.6	ug/l	1082910	4	11.48	1030120	50.0
Indane	11.77	20.0	ug/l	411996	4	11.48	1030120	50.0
Benzene, 2-ethyl-...	11.87	41.7	ug/l	859527	4	11.48	1030120	50.0
o-Cymene	12.25	28.8	ug/l	593516	4	11.48	1030120	50.0
Benzene, 1,2,4,5-...	12.65	30.4	ug/l	625232	4	11.48	1030120	50.0
Benzene, 1,2,3,4-...	12.70	29.7	ug/l	612530	4	11.48	1030120	50.0
Benzene, 2-butenyl-	13.12	45.2	ug/l	931451	4	11.48	1030120	50.0
Cyclohexanemethan...	13.18	36.8	ug/l	758171	4	11.48	1030120	50.0