

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU080618\
 Data File : VU025923.D
 Acq On : 07 Aug 2018 06:35
 Operator : MD/SY
 Sample : J4344-12
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 FL-0531

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\82U072718W.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	1.088	137	155	173	rBV	250523	295414	2.12%	0.313%
2	2.747	644	671	684	rBV	181444	467677	3.36%	0.495%
3	3.002	733	750	773	rBV	186518	395349	2.84%	0.418%
4	3.310	830	846	863	rBV	95613	203107	1.46%	0.215%
5	4.101	1073	1092	1115	rVB	323467	864151	6.20%	0.915%
6	4.889	1320	1337	1353	rBV	204562	462784	3.32%	0.490%
7	4.995	1353	1370	1387	rVV3	669283	1687705	12.12%	1.786%
8	5.085	1387	1398	1422	rVB2	76861	192124	1.38%	0.203%
9	5.223	1423	1441	1459	rBV4	115189	380930	2.74%	0.403%
10	5.313	1460	1469	1498	rVB	197215	430686	3.09%	0.456%
11	5.487	1507	1523	1533	rBV2	115697	260769	1.87%	0.276%
12	5.558	1534	1545	1556	rVV3	112054	267658	1.92%	0.283%
13	5.628	1556	1567	1588	rVB2	165473	361645	2.60%	0.383%
14	5.892	1631	1649	1680	rBV	402103	800656	5.75%	0.847%
15	6.419	1783	1813	1829	rBV	1107773	2389059	17.15%	2.528%
16	6.644	1872	1883	1898	rVB	101402	194323	1.40%	0.206%
17	6.911	1952	1966	1989	rVB6	59579	166431	1.20%	0.176%
18	7.535	2145	2160	2161	rBV	166500	218845	1.57%	0.232%
19	7.570	2162	2171	2201	rVB	802253	1476394	10.60%	1.562%
20	7.712	2202	2215	2229	rBV3	83697	204903	1.47%	0.217%
21	7.921	2269	2280	2296	rVB2	181909	332485	2.39%	0.352%
22	8.050	2301	2320	2331	rBV	84024	155846	1.12%	0.165%
23	8.548	2466	2475	2485	rVV	151754	241360	1.73%	0.255%
24	8.609	2485	2494	2504	rVV	105112	171635	1.23%	0.182%
25	9.095	2633	2645	2662	rBV	589797	985993	7.08%	1.043%
26	9.252	2684	2694	2711	rVB	451084	725475	5.21%	0.768%
27	9.381	2718	2734	2750	rBV	1045418	1821892	13.08%	1.928%
28	9.956	2889	2913	2924	rBV4	64113	141261	1.01%	0.149%
29	10.168	2968	2979	2991	rBV	693570	1070601	7.69%	1.133%
30	10.310	3011	3023	3036	rBV	635118	1000656	7.19%	1.059%
31	10.590	3093	3110	3127	rBV	1113267	1710989	12.29%	1.811%
32	10.686	3127	3140	3145	rBV	3741937	6456556	46.36%	6.833%
33	10.776	3158	3168	3189	rVB	2995589	4577504	32.87%	4.844%
34	10.975	3216	3230	3251	rBV	3271125	5041646	36.20%	5.336%

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Integration Parameters: RTEINT.P

Integrator: RTE
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 Sampling : 1
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 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\82U072718W.M
 Title : SW846 8260

35	11.152	3273	3285	3312	rVB	9296339	13926902	100.00%	14.739%
36	11.294	3312	3329	3333	rBV	105989	184749	1.33%	0.196%
37	11.326	3333	3339	3354	rVB	280878	439036	3.15%	0.465%
38	11.429	3355	3371	3379	rBV	608988	945598	6.79%	1.001%
39	11.483	3379	3388	3403	rVB3	1120732	1826945	13.12%	1.933%
40	11.573	3403	3416	3440	rVB	3871601	5907625	42.42%	6.252%
41	11.692	3441	3453	3464	rVB	149592	235830	1.69%	0.250%
42	11.779	3465	3480	3485	rBV4	690784	1385773	9.95%	1.467%
43	11.811	3485	3490	3499	rVV	884584	1253504	9.00%	1.327%
44	11.876	3499	3510	3531	rVB3	1206191	2357554	16.93%	2.495%
45	11.985	3531	3544	3555	rBV	356202	556993	4.00%	0.589%
46	12.059	3555	3567	3580	rBV	951862	1457041	10.46%	1.542%
47	12.149	3583	3595	3599	rBV	943633	1392231	10.00%	1.473%
48	12.178	3599	3604	3614	rVB	1166923	1682673	12.08%	1.781%
49	12.255	3617	3628	3635	rBV	1134716	1697695	12.19%	1.797%
50	12.291	3635	3639	3650	rVB2	310708	420366	3.02%	0.445%
51	12.358	3650	3660	3691	rBV4	658039	2016822	14.48%	2.134%
52	12.496	3691	3703	3710	rBV3	178317	307796	2.21%	0.326%
53	12.551	3710	3720	3733	rVB2	839140	1467914	10.54%	1.553%
54	12.651	3733	3751	3759	rBV	980120	1633965	11.73%	1.729%
55	12.705	3759	3768	3783	rVB	1614849	2486263	17.85%	2.631%
56	12.792	3784	3795	3801	rBV3	215303	349465	2.51%	0.370%
57	12.930	3830	3838	3843	rBV4	109282	179728	1.29%	0.190%
58	12.966	3843	3849	3857	rVB2	300596	421542	3.03%	0.446%
59	13.085	3877	3886	3887	rBV2	151239	169062	1.21%	0.179%
60	13.123	3888	3898	3913	rVB2	2284129	3721845	26.72%	3.939%
61	13.290	3940	3950	3974	rVB	763295	1302734	9.35%	1.379%
62	13.409	3976	3987	3994	rBV3	126387	221035	1.59%	0.234%
63	13.458	3995	4002	4009	rVB	113603	159902	1.15%	0.169%
64	13.512	4009	4019	4026	rBV3	384161	657302	4.72%	0.696%
65	13.612	4045	4050	4056	rVB	132763	153369	1.10%	0.162%
66	13.744	4072	4091	4100	rBV	739553	1247550	8.96%	1.320%
67	13.789	4100	4105	4117	rVB	142159	213200	1.53%	0.226%
68	13.856	4117	4126	4141	rBV	106800	182439	1.31%	0.193%
69	13.959	4141	4158	4169	rVV4	139352	346991	2.49%	0.367%
70	14.155	4208	4219	4224	rBV	82902	141382	1.02%	0.150%
71	14.352	4265	4280	4300	rVB5	230576	558278	4.01%	0.591%

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Integration Parameters: RTEINT.P

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

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72	14.544	4331	4340	4352	rVB6	78033	144225	1.04%	0.153%
73	14.679	4371	4382	4401	rBV2	208894	384324	2.76%	0.407%
74	14.879	4432	4444	4455	rBV	1344430	2134272	15.32%	2.259%
75	15.062	4490	4501	4516	rBV	877508	1448405	10.40%	1.533%
76	15.917	4754	4767	4783	rBV	110460	212796	1.53%	0.225%
77	16.062	4801	4812	4819	rBV	158024	257194	1.85%	0.272%
78	16.101	4819	4824	4840	rVB2	94492	146503	1.05%	0.155%

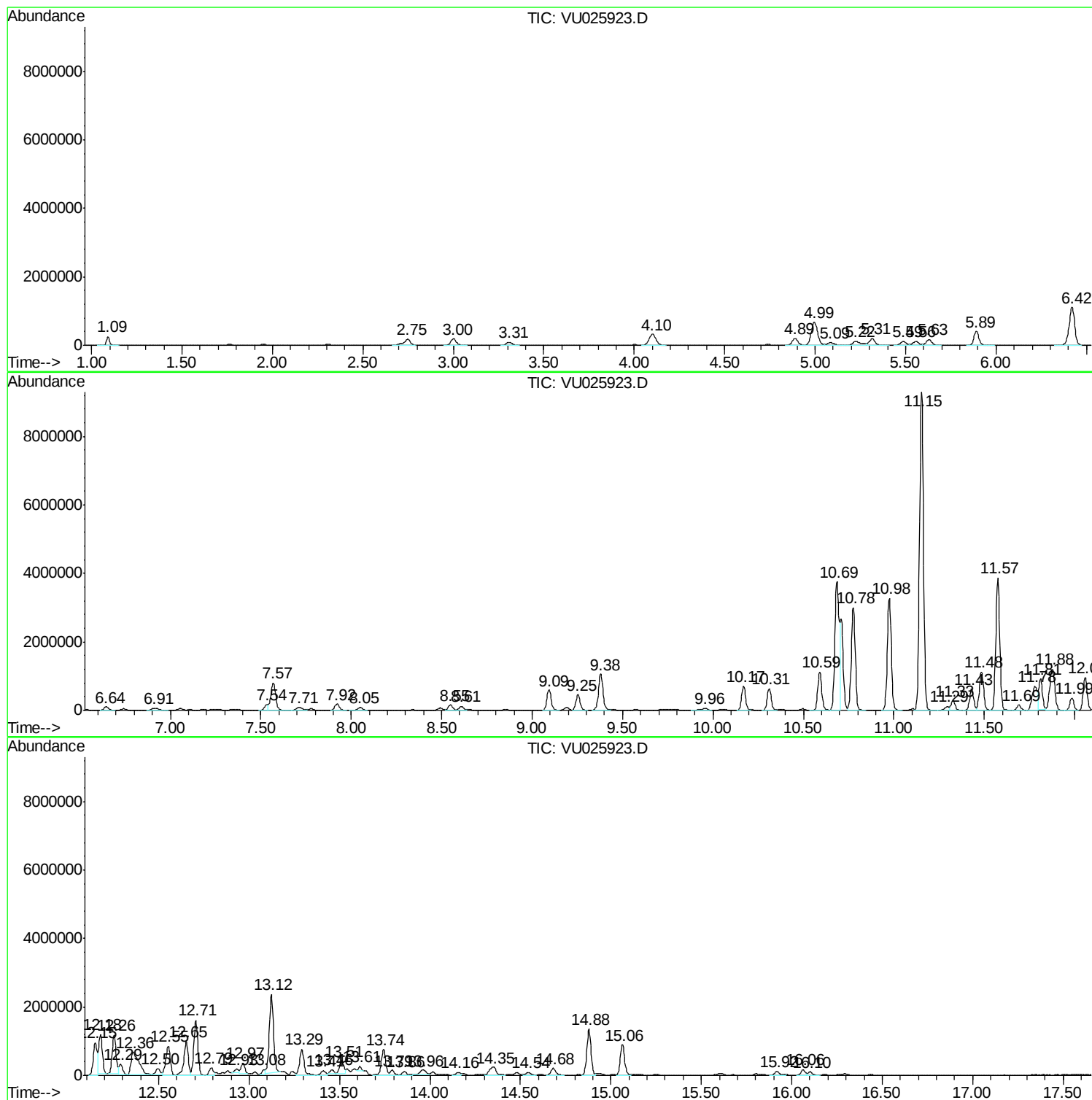
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.M
Quant Title : SW846 8260

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



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Instrument :
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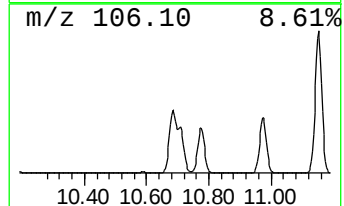
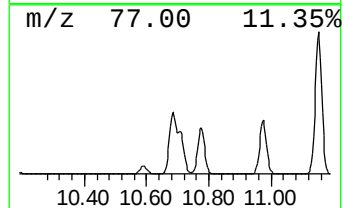
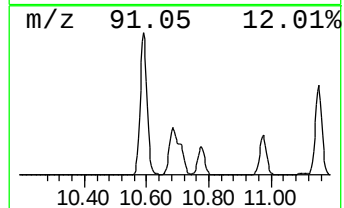
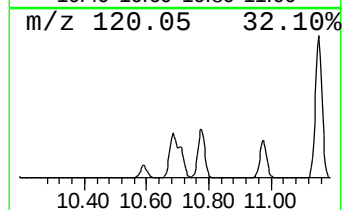
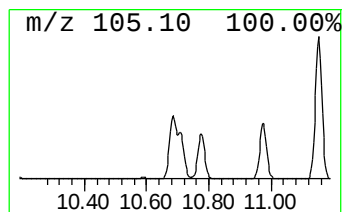
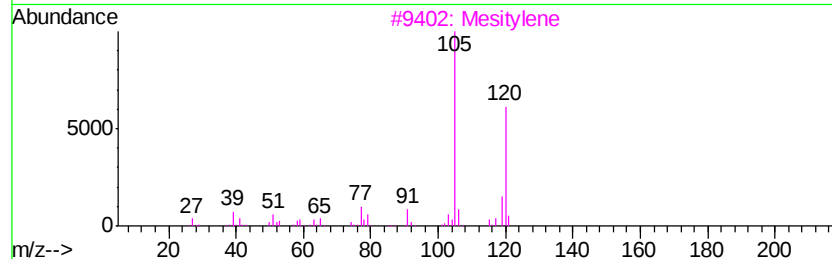
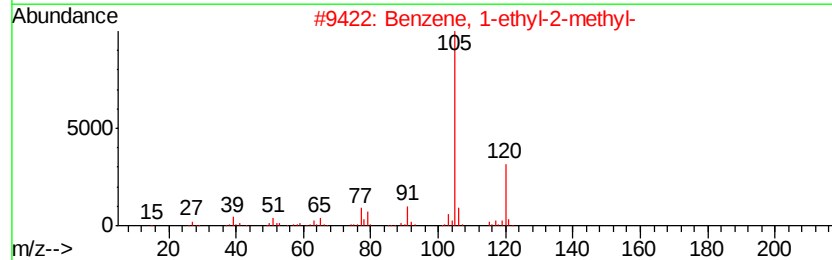
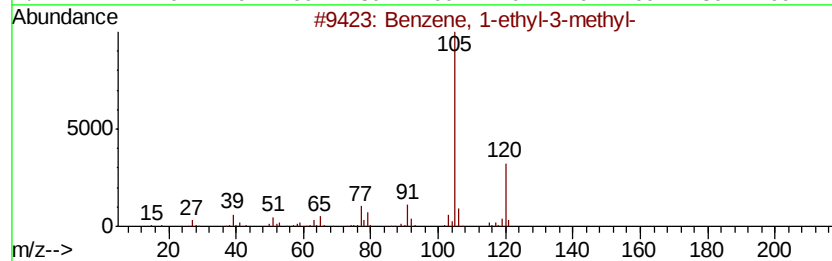
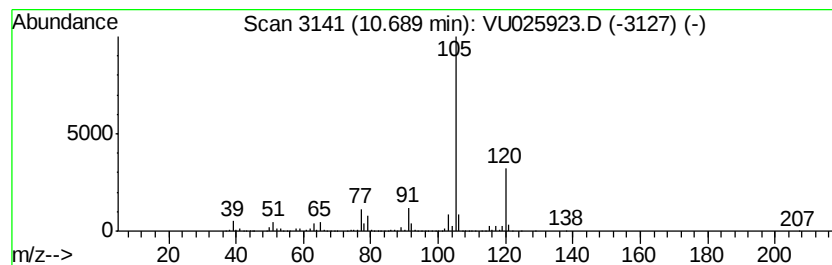
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.69	176.70 ug/l	6456560	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3		Mesitylene	120	C9H12	000108-67-8	91
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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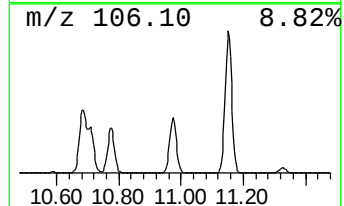
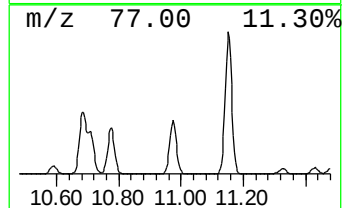
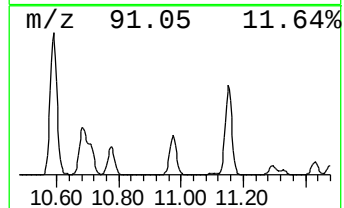
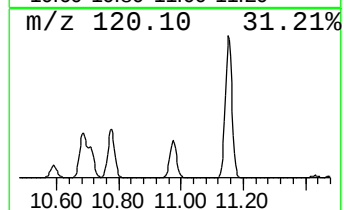
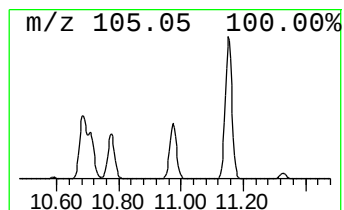
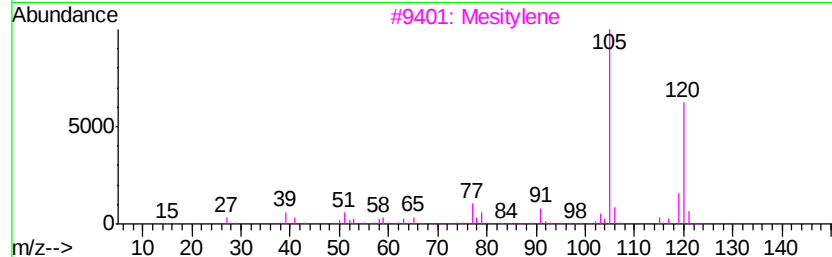
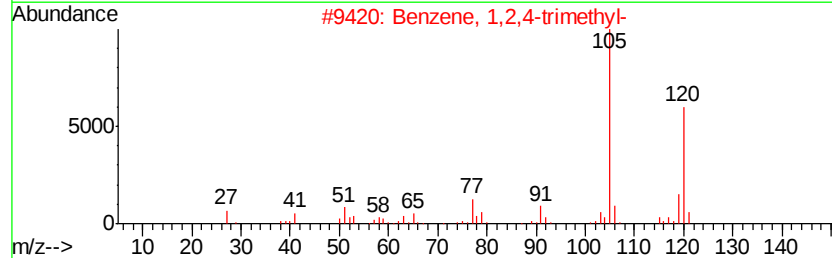
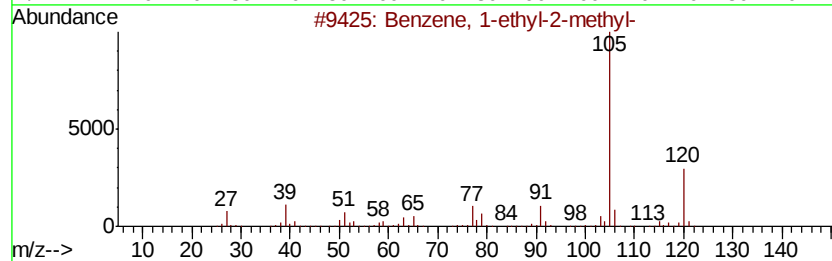
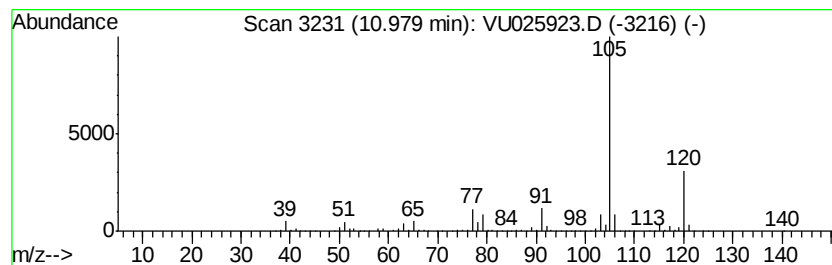
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.98	137.98 ug/l	5041650	1,4-Dichlorobenzene-d4	11.49

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



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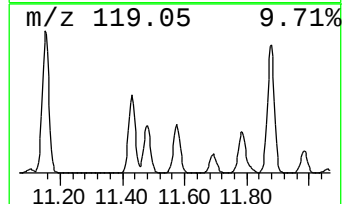
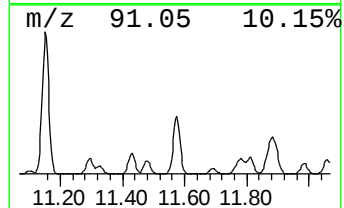
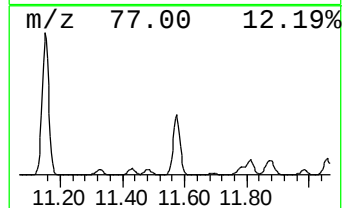
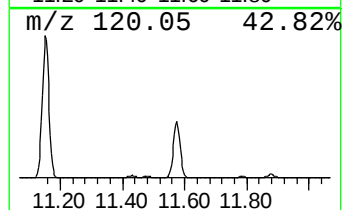
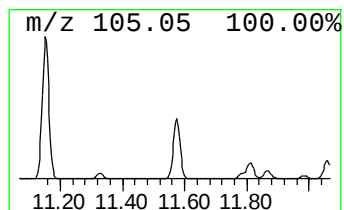
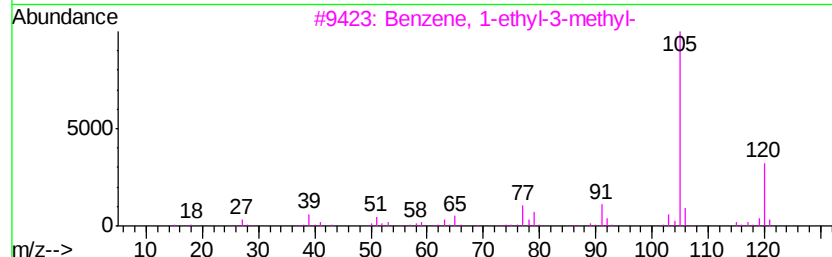
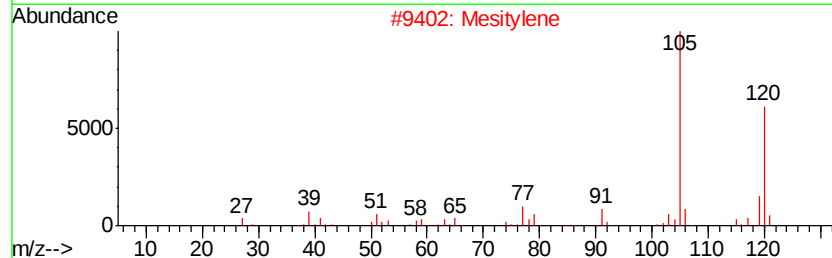
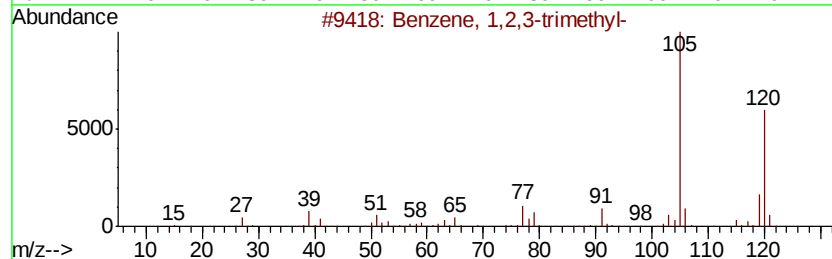
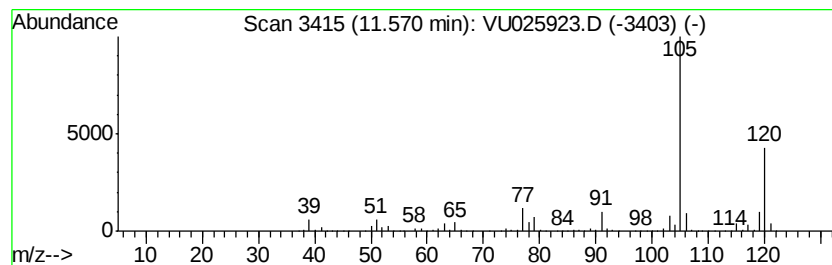
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	161.68 ug/l	5907630	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Mesitylene	120	C9H12	000108-67-8	93
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-ethyl-4-methyl-	120	C9H12	000622-96-8	90



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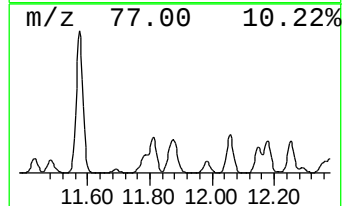
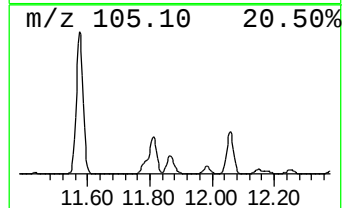
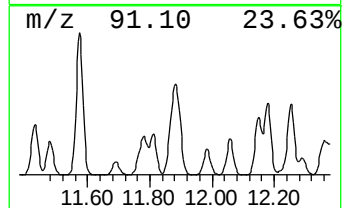
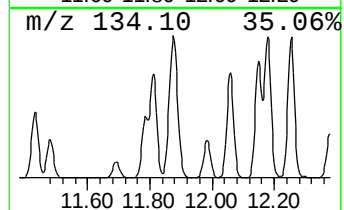
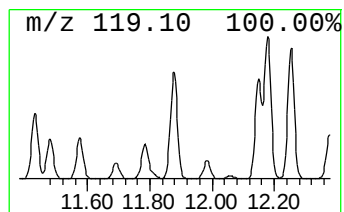
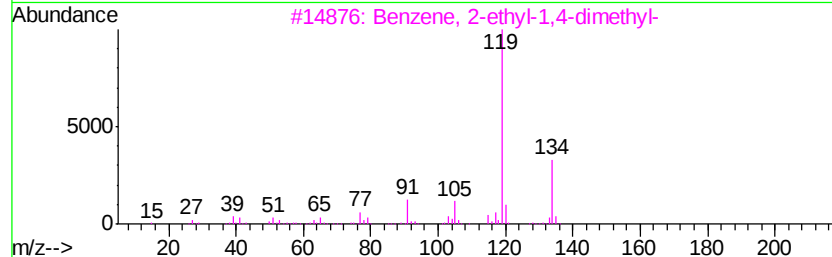
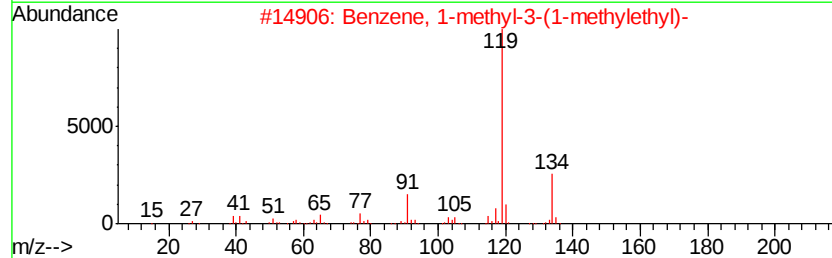
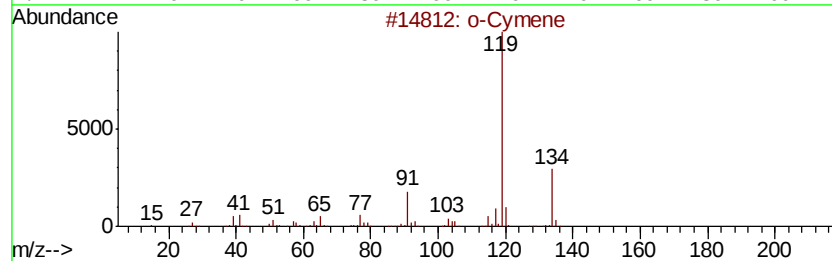
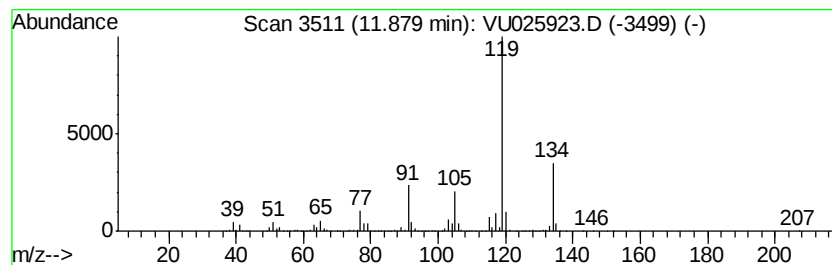
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Quant Title : SW846 8260

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

Peak Number 4 o-Cymene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.88	64.52 ug/l	2357550	1,4-Dichlorobenzene-d4	11.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		o-Cymene	134	C10H14	000527-84-4	95
2		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	95
3		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
4		p-Cymene	134	C10H14	000099-87-6	93
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU080618\
Data File : VU025923.D
Acq On : 07 Aug 2018 06:35
Operator : MD/SY
Sample : J4344-12
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
FL-0531

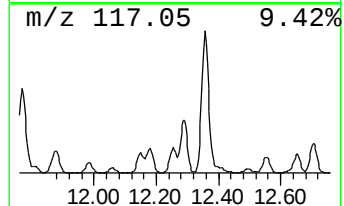
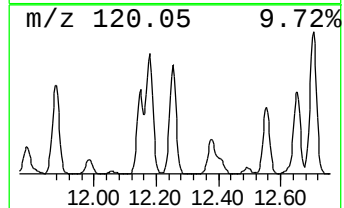
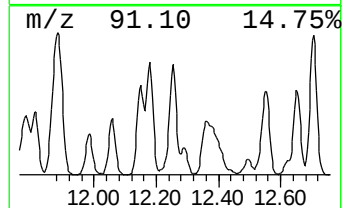
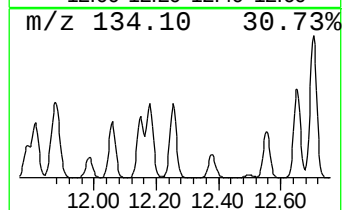
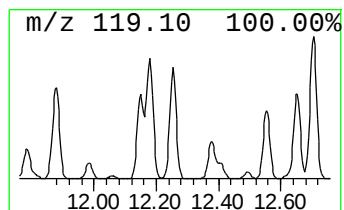
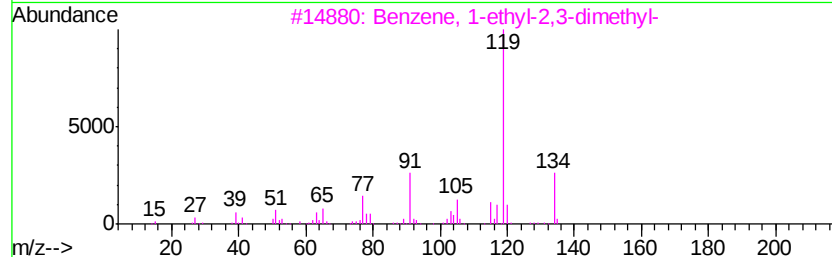
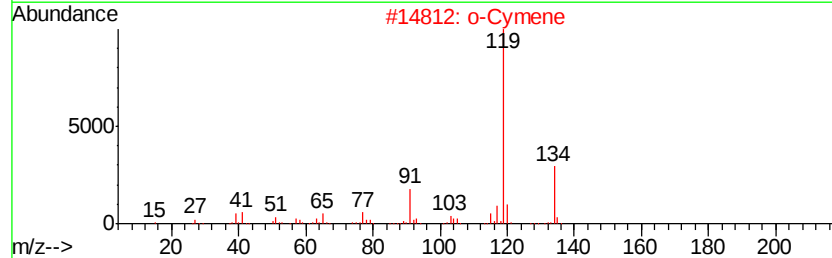
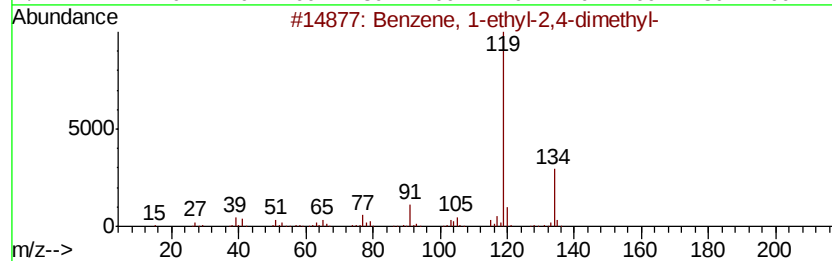
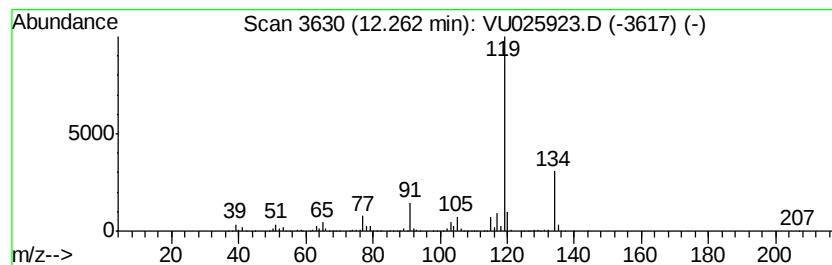
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.26	46.46 ug/l	1697700	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
2		o-Cymene	134	C10H14	000527-84-4	95
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
5		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU080618\
Data File : VU025923.D
Acq On : 07 Aug 2018 06:35
Operator : MD/SY
Sample : J4344-12
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
FL-0531

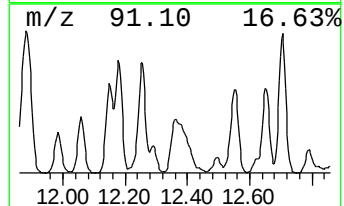
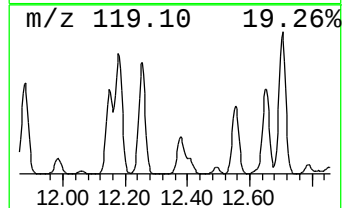
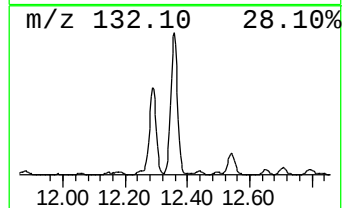
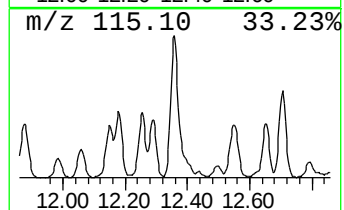
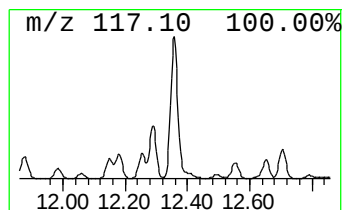
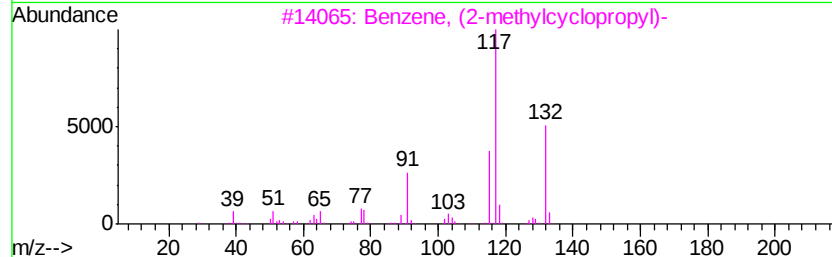
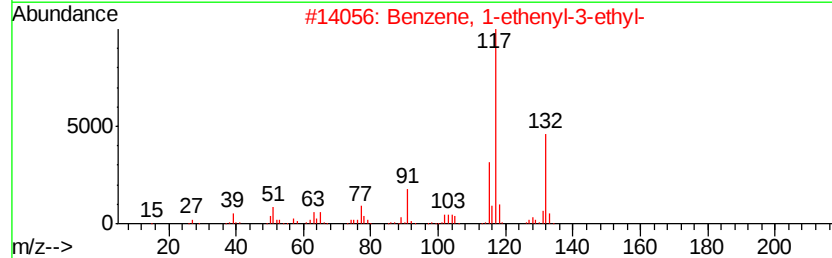
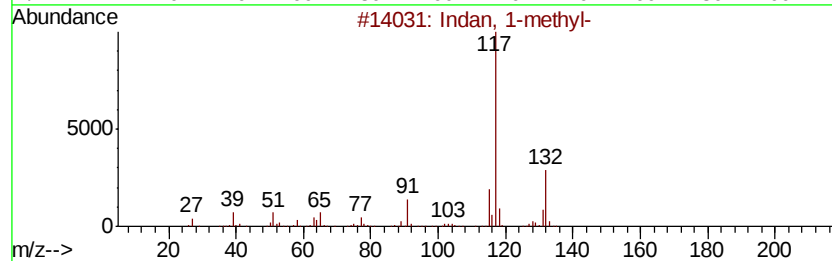
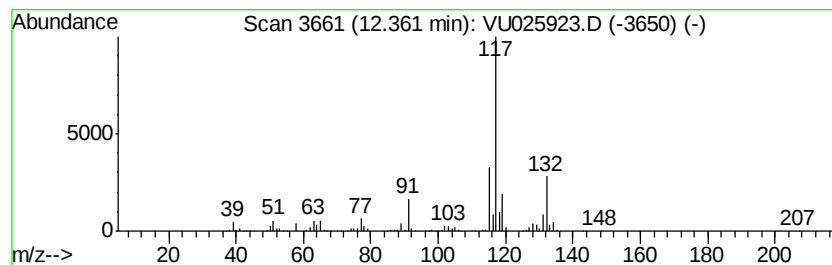
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Quant Title : SW846 8260

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

Peak Number 7 Indan, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	55.20 ug/l	2016820	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indan, 1-methyl-	132	C10H12	000767-58-8	94
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
3		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	87
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	86
5		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU080618\
Data File : VU025923.D
Acq On : 07 Aug 2018 06:35
Operator : MD/SY
Sample : J4344-12
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
FL-0531

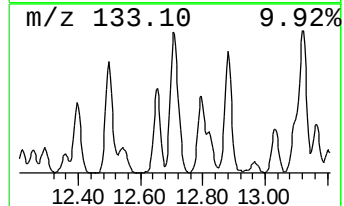
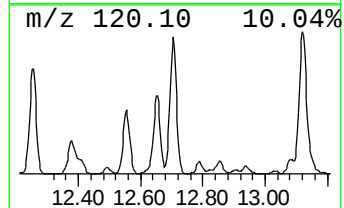
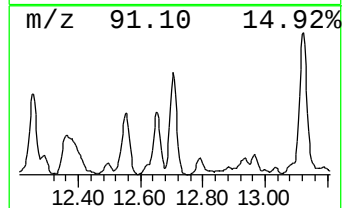
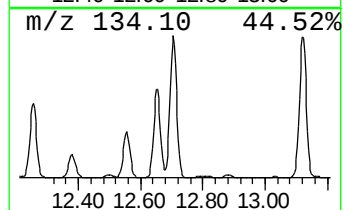
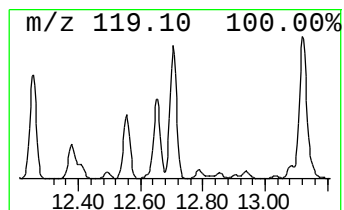
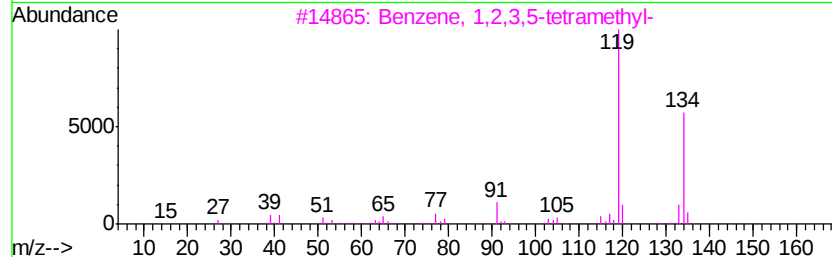
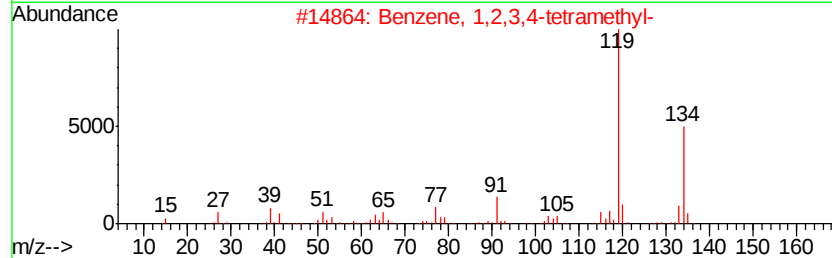
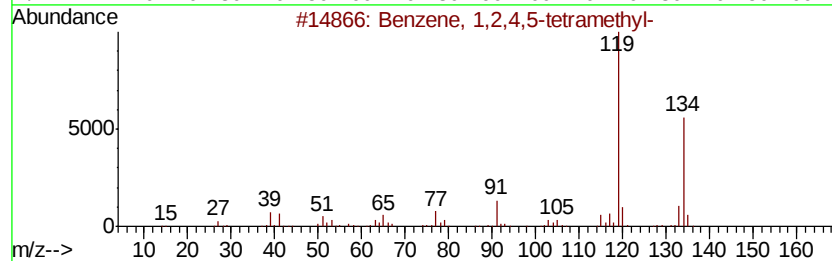
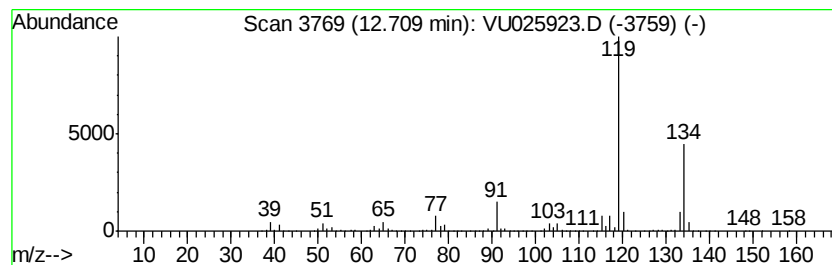
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.71	68.04 ug/l	2486260	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
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	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU080618\
Data File : VU025923.D
Acq On : 07 Aug 2018 06:35
Operator : MD/SY
Sample : J4344-12
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
FL-0531

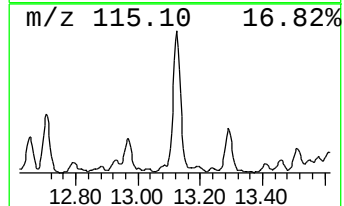
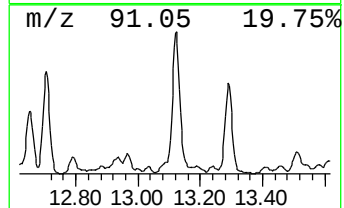
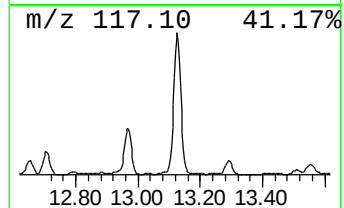
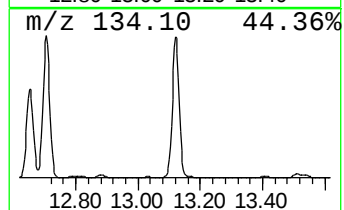
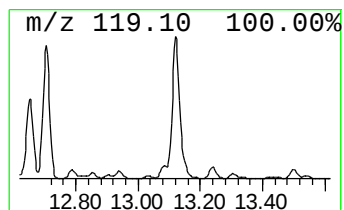
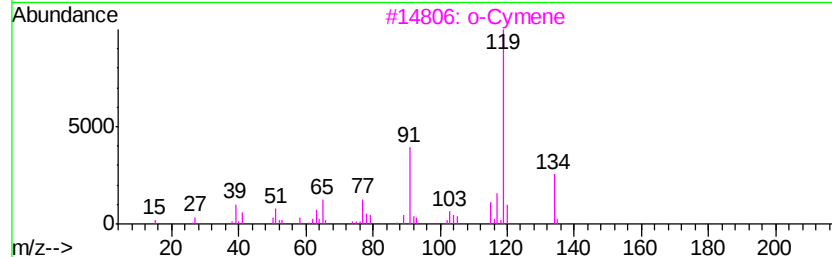
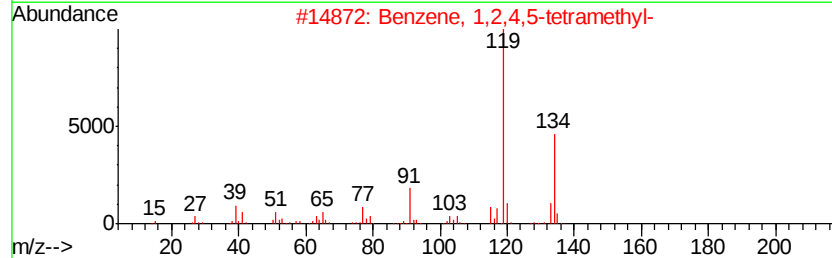
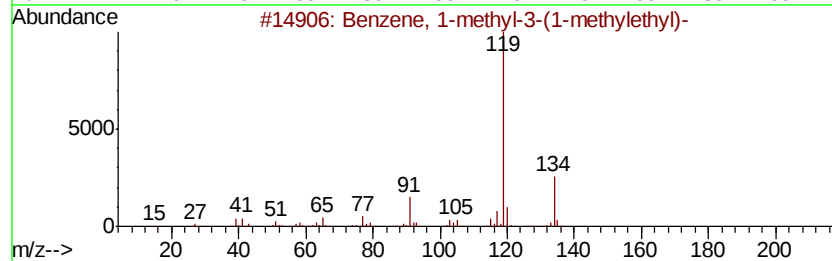
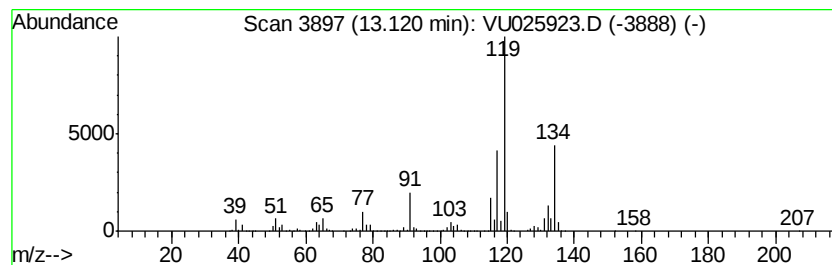
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-methyl-3-(1-meth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	101.86 ug/l	3721850	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	70
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	70
3		o-Cymene	134	C10H14	000527-84-4	70
4		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	62
5		p-Cymene	134	C10H14	000099-87-6	62



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU080618\
Data File : VU025923.D
Acq On : 07 Aug 2018 06:35
Operator : MD/SY
Sample : J4344-12
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
FL-0531

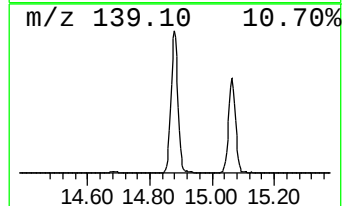
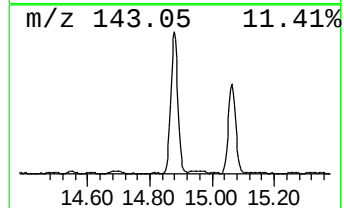
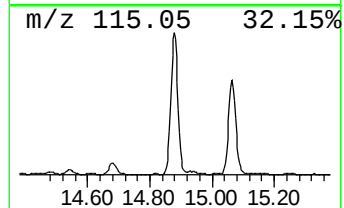
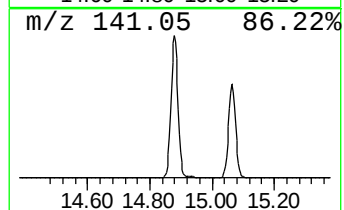
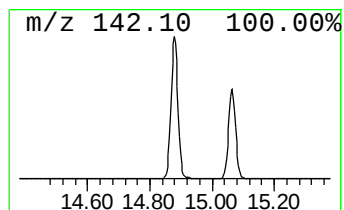
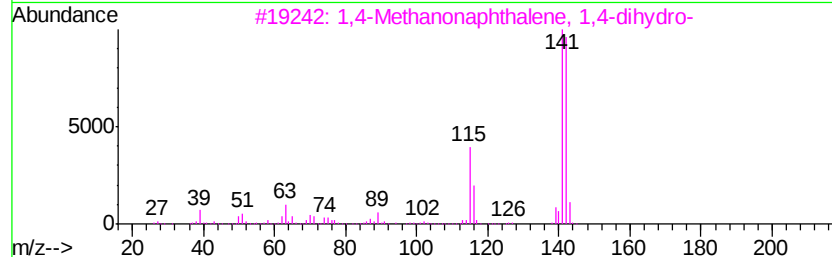
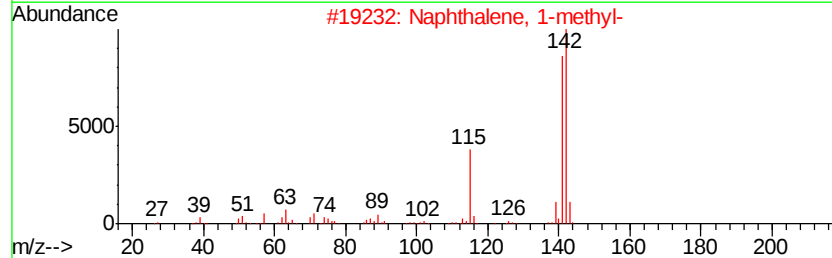
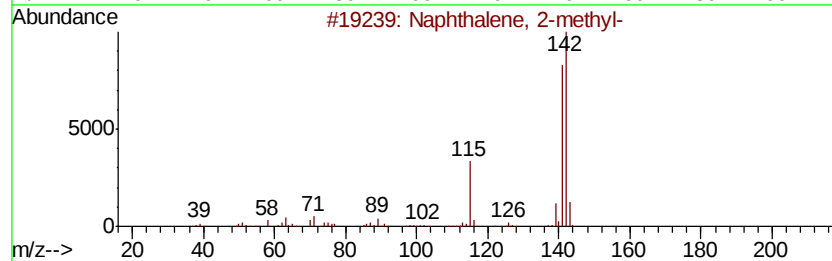
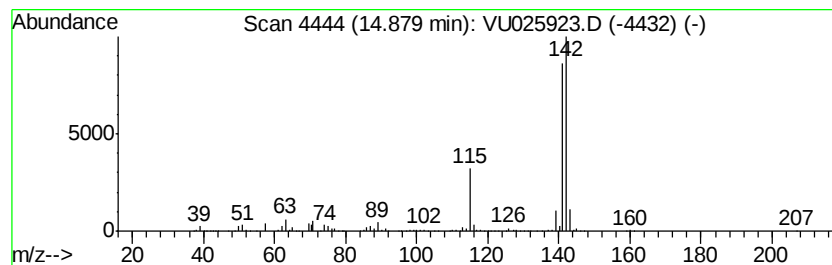
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.88	58.41 ug/l	2134270	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3		1,4-Methanonaphthalene, 1,4-dihv...	142	C11H10	004453-90-1	91
4		Benzocycloheptatriene	142	C11H10	000264-09-5	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU080618\
Data File : VU025923.D
Acq On : 07 Aug 2018 06:35
Operator : MD/SY
Sample : J4344-12
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
FL-0531

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.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-...	10.69	176.7	ug/l	6456560	4	11.49	1826950	50.0
Benzene, 1-ethyl-...	10.98	138.0	ug/l	5041650	4	11.49	1826950	50.0
Benzene, 1,2,3-tr...	11.57	161.7	ug/l	5907630	4	11.49	1826950	50.0
o-Cymene	11.88	64.5	ug/l	2357550	4	11.49	1826950	50.0
Benzene, 1-ethyl-...	12.26	46.5	ug/l	1697700	4	11.49	1826950	50.0
Indan, 1-methyl-	12.36	55.2	ug/l	2016820	4	11.49	1826950	50.0
Benzene, 1,2,4,5-...	12.71	68.0	ug/l	2486260	4	11.49	1826950	50.0
Benzene, 1-methyl...	13.12	101.9	ug/l	3721850	4	11.49	1826950	50.0
Naphthalene, 1-me...	14.88	58.4	ug/l	2134270	4	11.49	1826950	50.0