

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100418\
 Data File : VU027430.D
 Acq On : 04 Oct 2018 23:53
 Operator : MD/SY
 Sample : J5165-15
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EP1-MW-E-092718

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\82U100118W.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	1.302	25	40	47	rBV	43463	113938	5.31%	0.498%
2	2.325	346	358	371	rBV	42643	68917	3.21%	0.301%
3	2.447	385	396	417	rBV2	53339	106116	4.95%	0.464%
4	4.887	1138	1155	1171	rBV	105795	241201	11.24%	1.055%
5	4.987	1172	1186	1217	rVB	143683	322377	15.03%	1.410%
6	5.312	1271	1287	1312	rBV	132157	285954	13.33%	1.251%
7	5.890	1452	1467	1493	rBV	226283	462324	21.55%	2.022%
8	7.565	1966	1988	2021	rBV	429153	796518	37.13%	3.484%
9	7.919	2086	2098	2116	rBV3	14217	26417	1.23%	0.116%
10	8.183	2166	2180	2187	rBV3	13690	29704	1.38%	0.130%
11	8.607	2299	2312	2322	rBV2	52311	98752	4.60%	0.432%
12	8.848	2380	2387	2399	rVV4	30953	65032	3.03%	0.284%
13	9.032	2434	2444	2452	rBV	24297	42705	1.99%	0.187%
14	9.090	2452	2462	2479	rVV	334934	571623	26.65%	2.500%
15	9.189	2487	2493	2501	rVB2	16274	25843	1.20%	0.113%
16	9.247	2501	2511	2525	rVB6	24410	47804	2.23%	0.209%
17	9.356	2525	2545	2550	rBV6	36320	82985	3.87%	0.363%
18	9.405	2553	2560	2567	rVV4	69958	128800	6.00%	0.563%
19	9.446	2568	2573	2600	rVB2	37263	83818	3.91%	0.367%
20	9.569	2600	2611	2623	rBV5	28648	53869	2.51%	0.236%
21	9.723	2643	2659	2671	rBV6	92496	238227	11.10%	1.042%
22	9.954	2711	2731	2741	rBV3	172590	356569	16.62%	1.560%
23	10.048	2751	2760	2768	rBV5	96508	180631	8.42%	0.790%
24	10.167	2787	2797	2806	rVB	88172	140029	6.53%	0.612%
25	10.225	2806	2815	2820	rBV6	35022	62863	2.93%	0.275%
26	10.308	2832	2841	2855	rVB2	356579	618087	28.81%	2.703%
27	10.379	2855	2863	2871	rVB3	51095	76794	3.58%	0.336%
28	10.491	2889	2898	2918	rVB5	173426	326334	15.21%	1.427%
29	10.588	2919	2928	2936	rBV	158119	251265	11.71%	1.099%
30	10.700	2955	2963	2972	rBV4	55964	92945	4.33%	0.407%
31	10.771	2972	2985	2994	rBV2	332023	611021	28.48%	2.672%
32	10.864	3009	3014	3023	rVB3	35306	52957	2.47%	0.232%
33	10.977	3040	3049	3057	rBV5	65490	137835	6.43%	0.603%
34	11.115	3071	3092	3094	rBV2	166508	311581	14.52%	1.363%

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35	11.147	3094	3102	3115	rVB	1113506	1761966	82.13%	7.707%
36	11.292	3128	3147	3151	rBV3	99062	224520	10.47%	0.982%
37	11.324	3151	3157	3171	rVV	185336	353262	16.47%	1.545%
38	11.401	3172	3181	3182	rVV2	88109	104786	4.88%	0.458%
39	11.427	3183	3189	3197	rVV	302957	462254	21.55%	2.022%
40	11.482	3197	3206	3217	rVB2	535239	843000	39.30%	3.687%
41	11.572	3224	3234	3257	rVB	1380622	2145255	100.00%	9.383%
42	11.688	3258	3270	3283	rVB	105512	210957	9.83%	0.923%
43	11.781	3285	3299	3303	rBV3	406190	736541	34.33%	3.221%
44	11.810	3304	3308	3316	rVV	464171	624253	29.10%	2.730%
45	11.868	3316	3326	3344	rVB6	478086	1165952	54.35%	5.100%
46	11.980	3351	3361	3381	rBV2	172343	320029	14.92%	1.400%
47	12.147	3403	3413	3417	rBV	297297	439022	20.46%	1.920%
48	12.176	3417	3422	3432	rVV	356809	527428	24.59%	2.307%
49	12.250	3434	3445	3453	rVV	613460	955401	44.54%	4.179%
50	12.286	3453	3456	3468	rVB3	145778	200526	9.35%	0.877%
51	12.356	3468	3478	3500	rBV6	273207	988597	46.08%	4.324%
52	12.495	3511	3521	3528	rBV5	96874	169223	7.89%	0.740%
53	12.549	3529	3538	3549	rVB3	303565	541791	25.26%	2.370%
54	12.620	3550	3560	3562	rBV4	79840	108337	5.05%	0.474%
55	12.649	3562	3569	3577	rVV	281360	417213	19.45%	1.825%
56	12.704	3577	3586	3602	rVB	483345	782719	36.49%	3.423%
57	12.787	3602	3612	3620	rBV2	108886	183880	8.57%	0.804%
58	12.880	3635	3641	3648	rVB	40847	48824	2.28%	0.214%
59	12.929	3648	3656	3661	rVV4	51520	82183	3.83%	0.359%
60	12.964	3662	3667	3674	rVB2	67334	87293	4.07%	0.382%
61	13.031	3682	3688	3695	rVB3	27087	36605	1.71%	0.160%
62	13.083	3695	3704	3707	rBV2	56502	82060	3.83%	0.359%
63	13.122	3707	3716	3735	rVB3	419172	734131	34.22%	3.211%
64	13.237	3746	3752	3759	rBV	24472	35535	1.66%	0.155%
65	13.289	3760	3768	3778	rVB2	59832	99431	4.63%	0.435%
66	13.408	3795	3805	3812	rBV6	21367	36776	1.71%	0.161%
67	13.456	3813	3820	3827	rBV2	24110	34359	1.60%	0.150%
68	13.511	3827	3837	3845	rBV4	65887	114286	5.33%	0.500%
69	13.610	3864	3868	3875	rVB	25154	28073	1.31%	0.123%
70	13.739	3898	3908	3917	rBV6	21284	39845	1.86%	0.174%
71	14.176	4028	4044	4051	rBV9	11202	25217	1.18%	0.110%

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Start Thrs: 0.2 Max Peaks: 100
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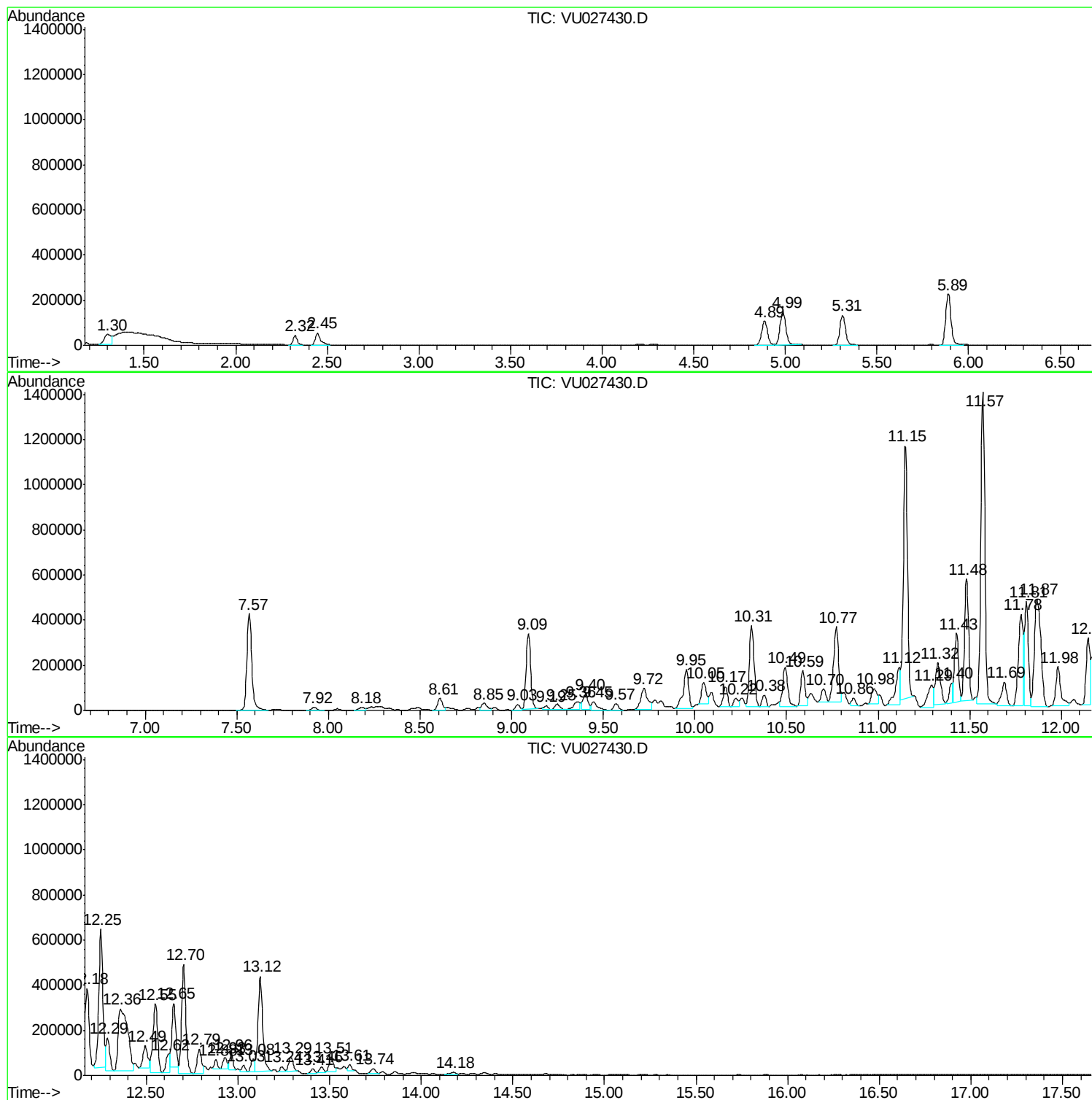
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

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



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Instrument :
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ClientSampleId :
EP1-MW-E-092718

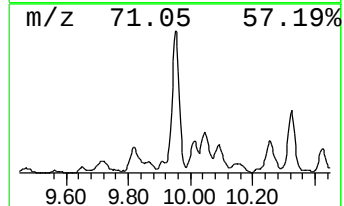
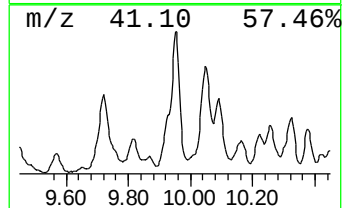
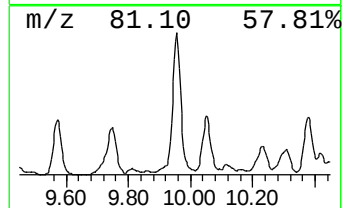
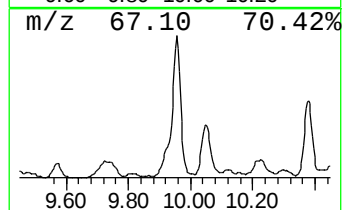
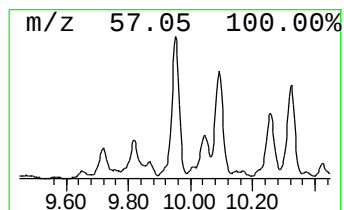
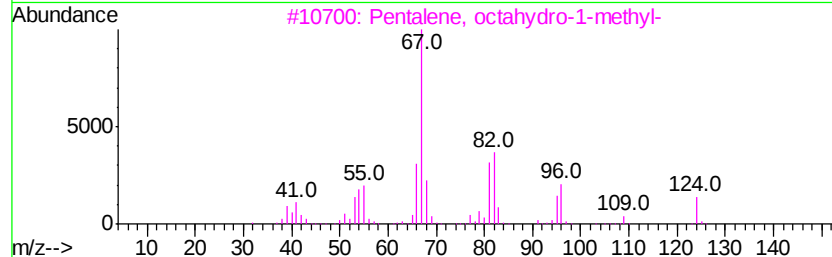
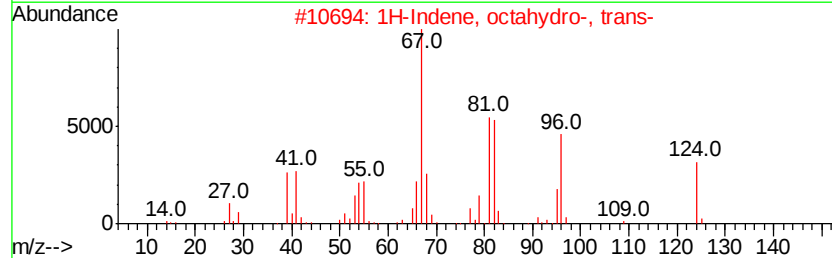
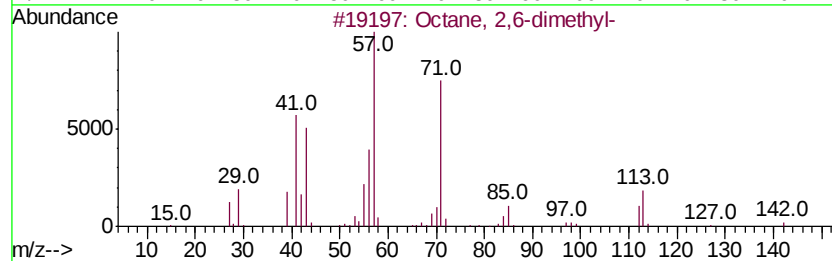
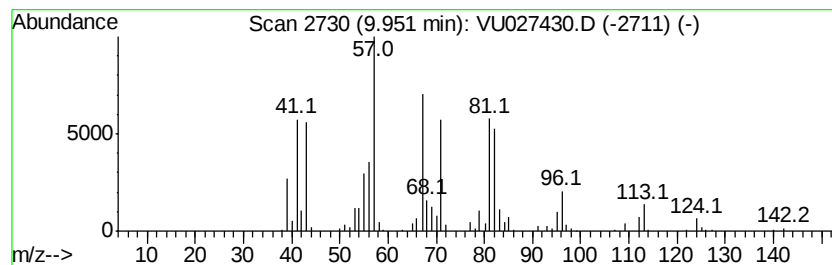
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Quant Title : SW846 8260

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

Peak Number 1 Octane, 2,6-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.95	31.19 ug/l	356569	Chlorobenzene-d5	9.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	60
2		1H-Indene, octahydro-, trans-	124	C9H16	003296-50-2	49
3		Pentalene, octahydro-1-methyl-	124	C9H16	032273-77-1	43
4		Heptanoic acid, 3-hexenyl ester,...	212	C13H24O2	061444-39-1	43
5		Bicyclo[4.1.0]heptane	96	C7H12	000286-08-8	41



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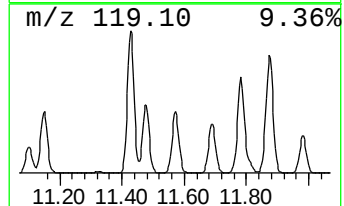
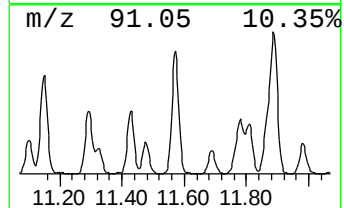
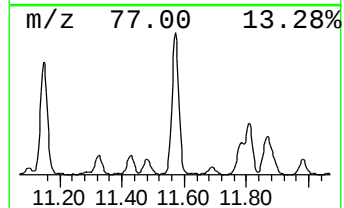
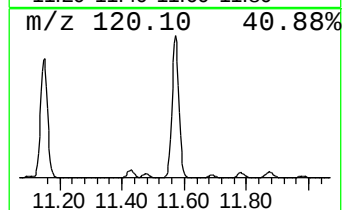
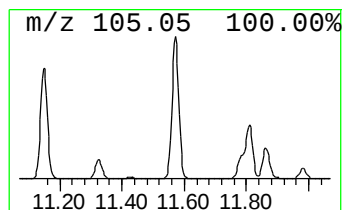
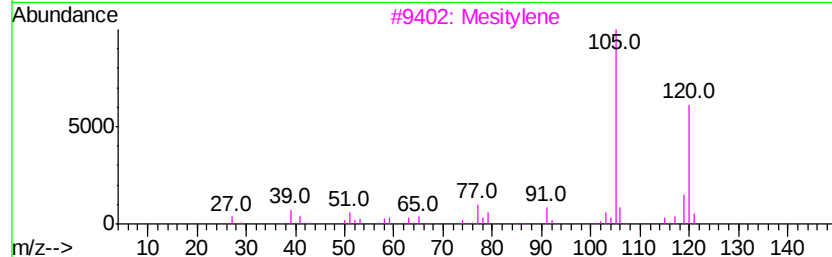
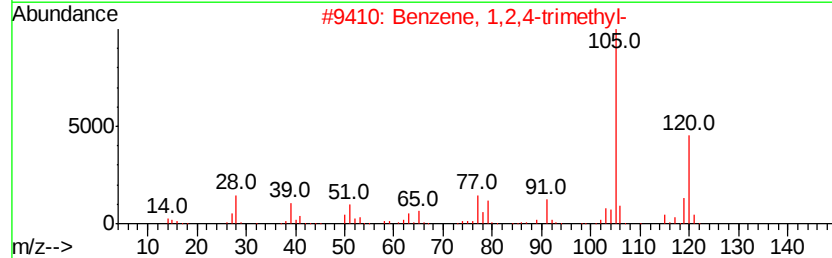
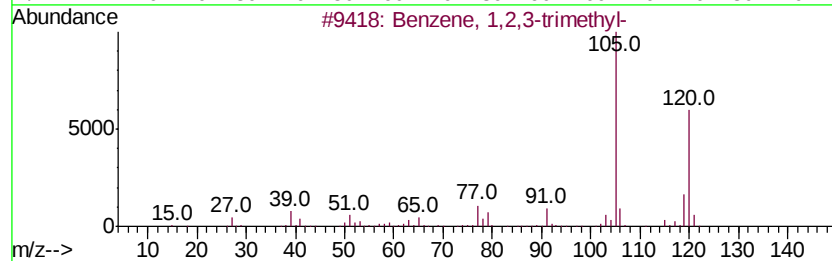
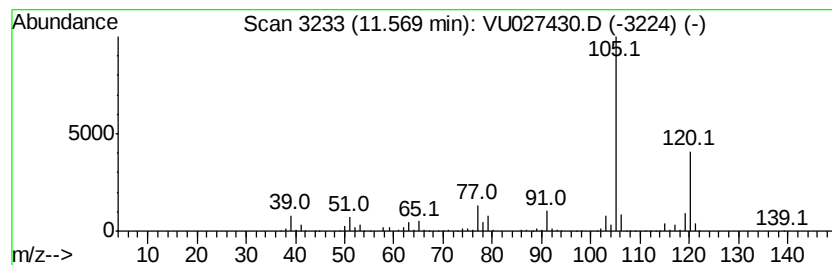
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	127.24 ug/l	2145260	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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	93
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
5		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90



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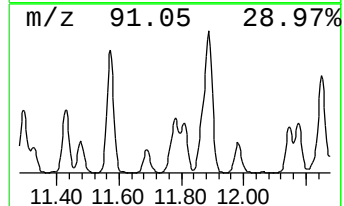
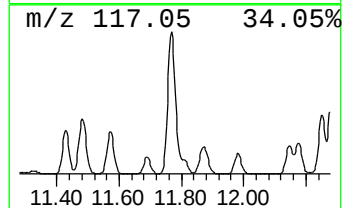
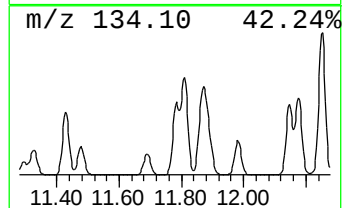
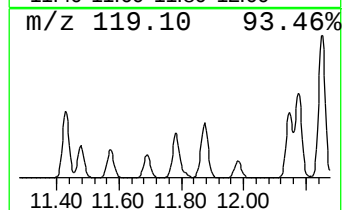
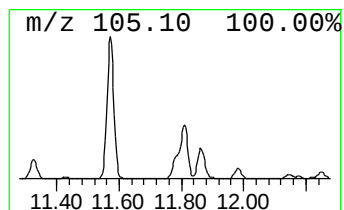
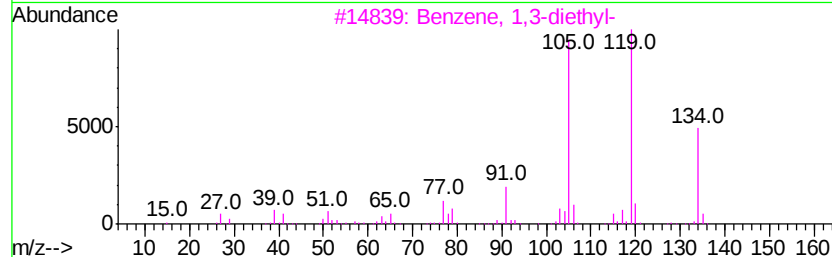
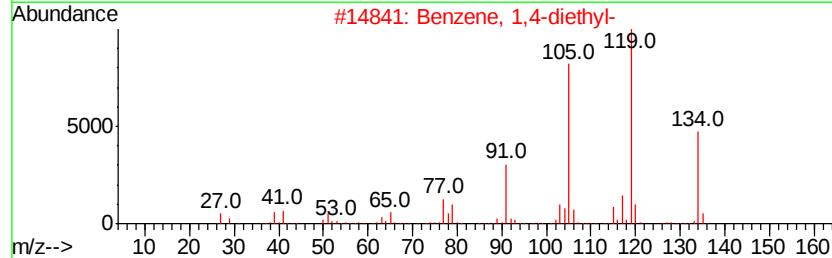
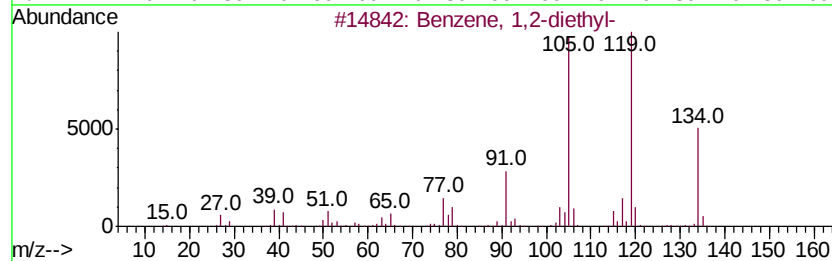
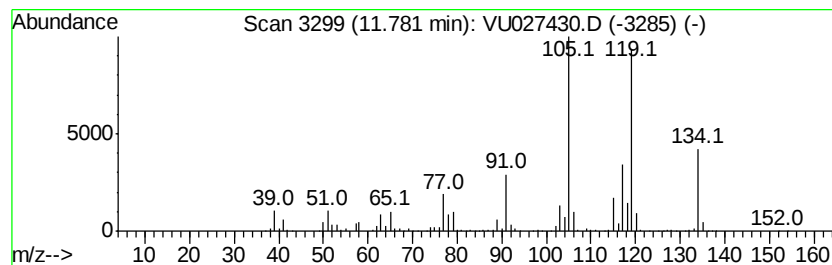
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1,2-diethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.78	43.69 ug/l	736541	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	93
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	83
5	Benzene, 1,2,3,4-tetramethyl-		134 C10H14	000488-23-3	68



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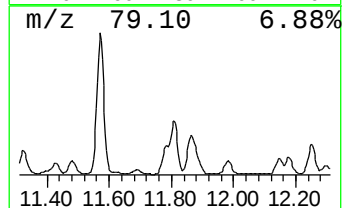
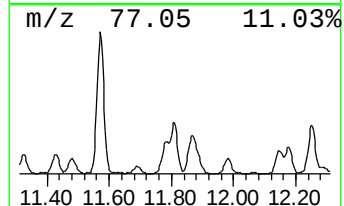
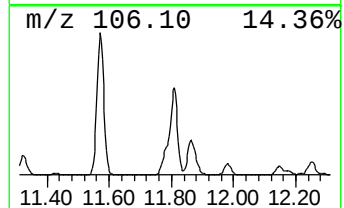
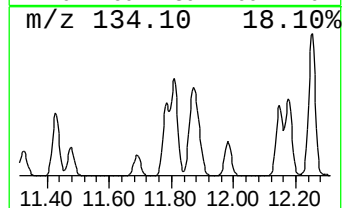
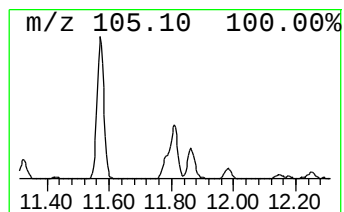
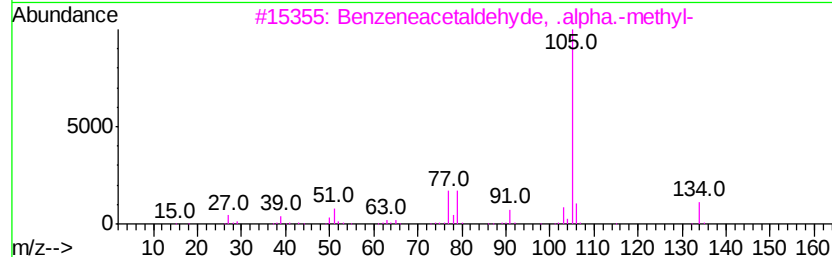
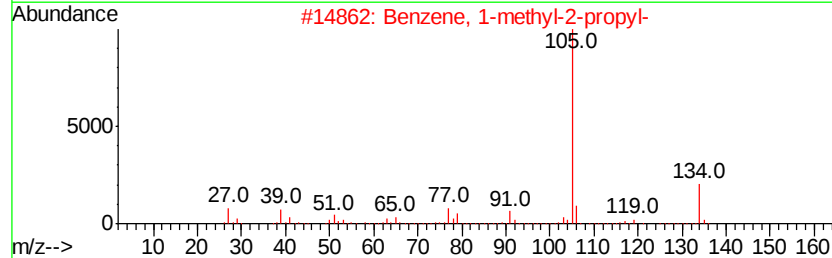
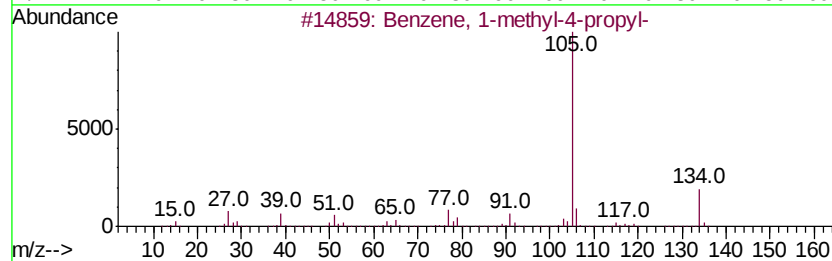
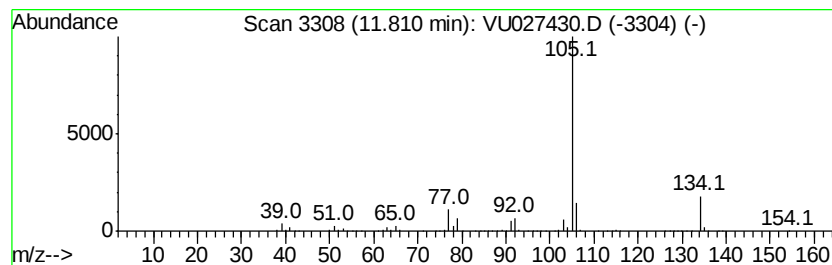
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-methyl-4-propyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.81	37.03 ug/l	624253	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	90
2		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	83
3		Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	78
4		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	72
5		Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100418\
Data File : VU027430.D
Acq On : 04 Oct 2018 23:53
Operator : MD/SY
Sample : J5165-15
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EP1-MW-E-092718

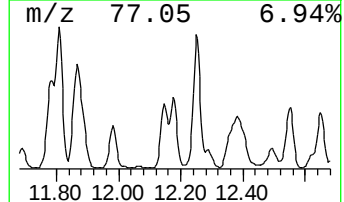
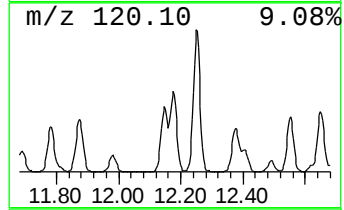
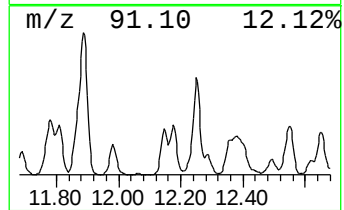
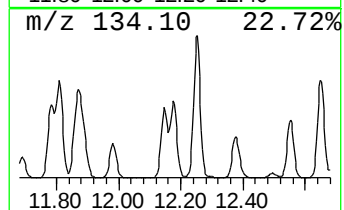
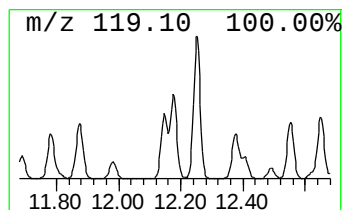
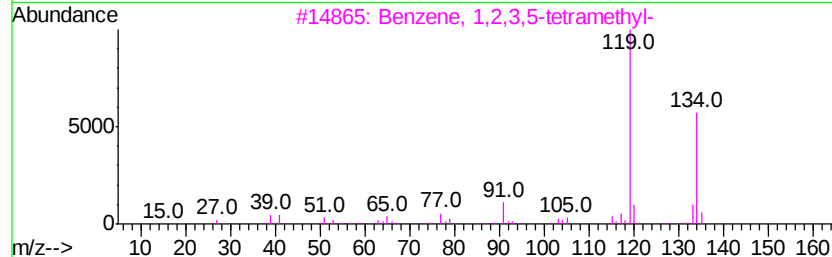
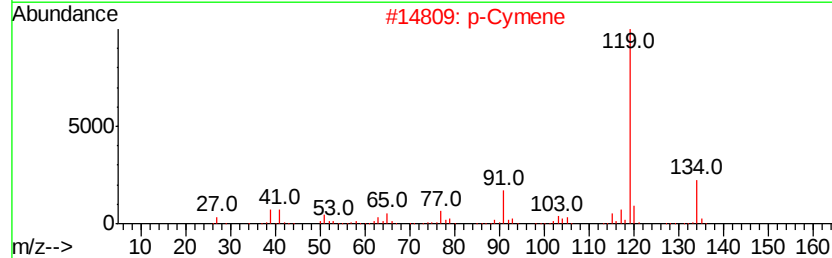
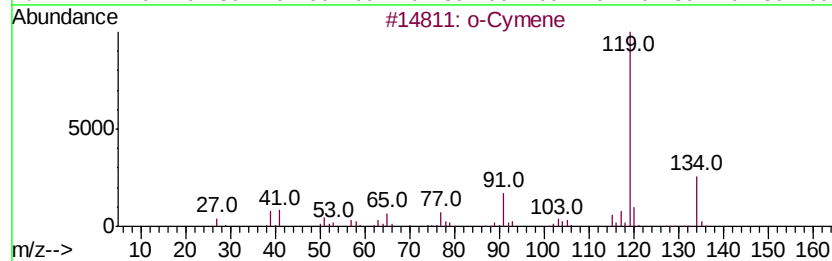
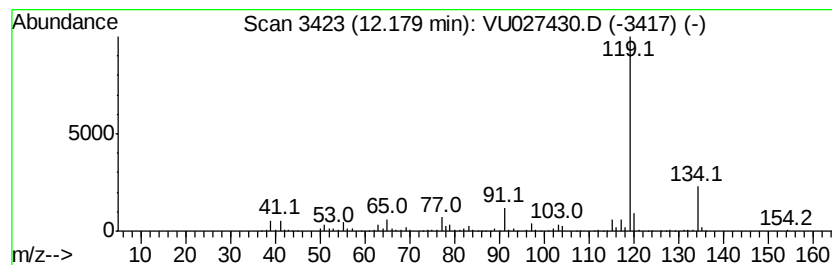
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Quant Title : SW846 8260

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

Peak Number 5 o-Cymene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.18	31.28 ug/l	527428	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, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100418\
Data File : VU027430.D
Acq On : 04 Oct 2018 23:53
Operator : MD/SY
Sample : J5165-15
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EP1-MW-E-092718

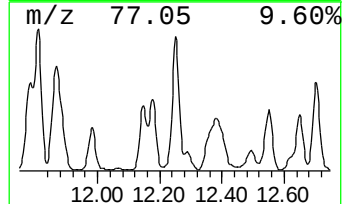
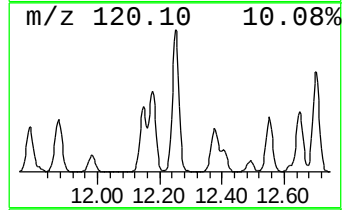
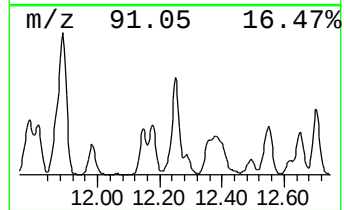
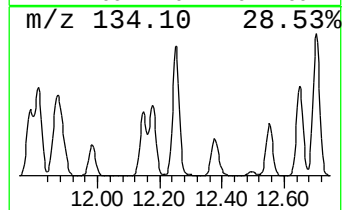
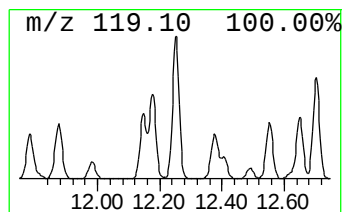
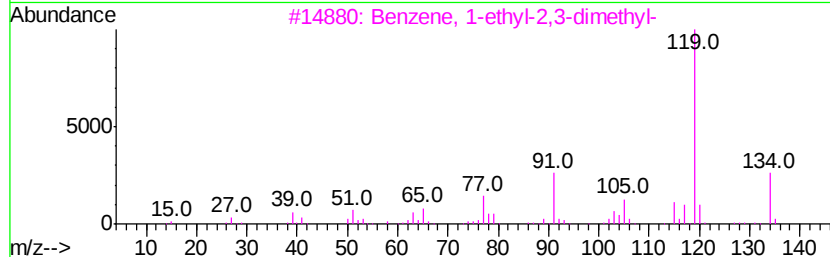
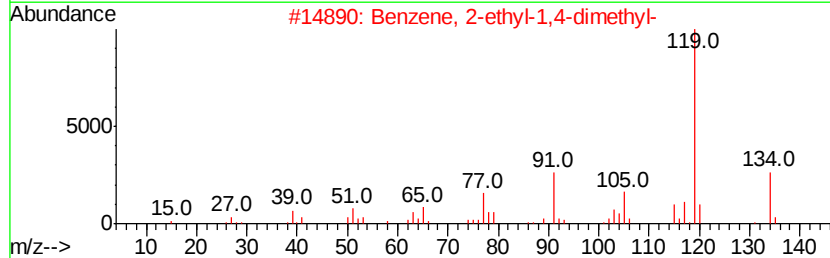
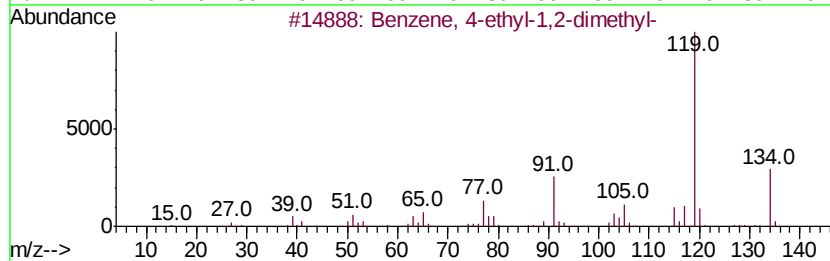
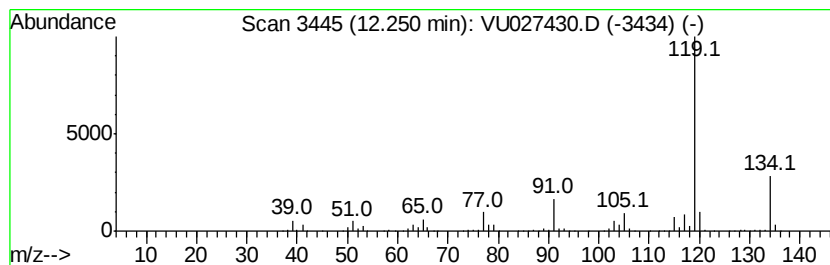
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.25	56.67 ug/l	955401	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
2		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4		o-Cymene	134	C10H14	000527-84-4	94
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100418\
Data File : VU027430.D
Acq On : 04 Oct 2018 23:53
Operator : MD/SY
Sample : J5165-15
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 41 Sample Multiplier: 1

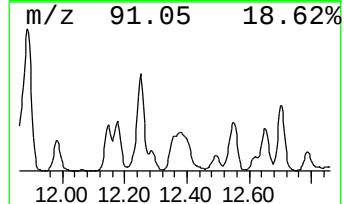
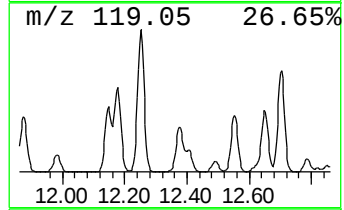
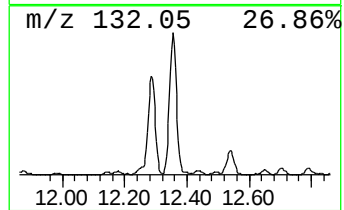
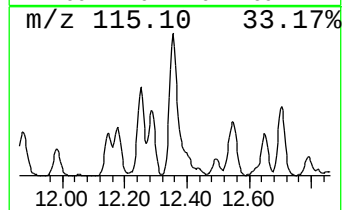
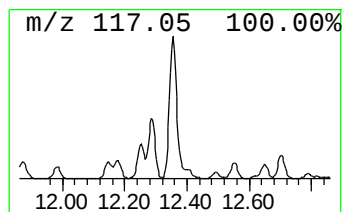
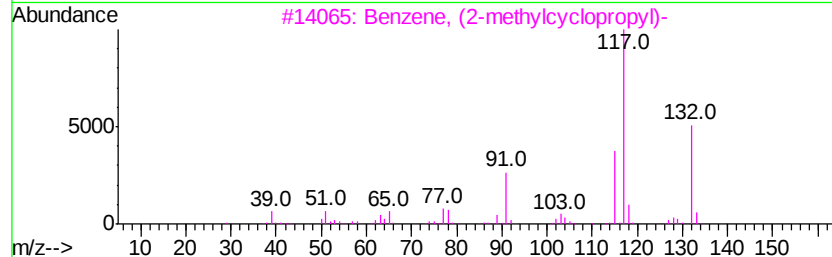
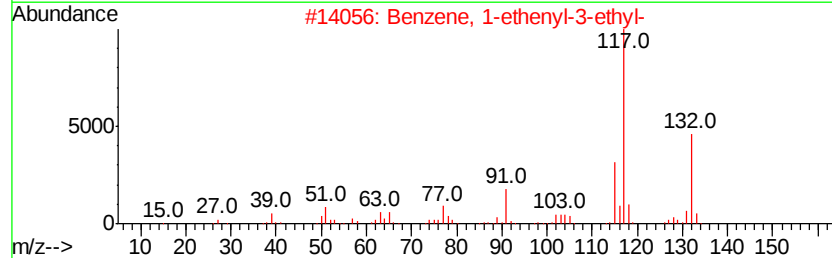
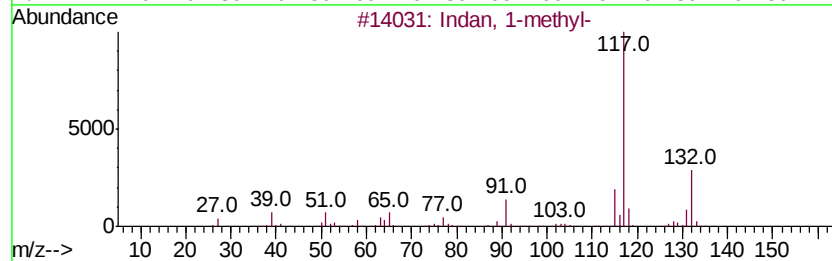
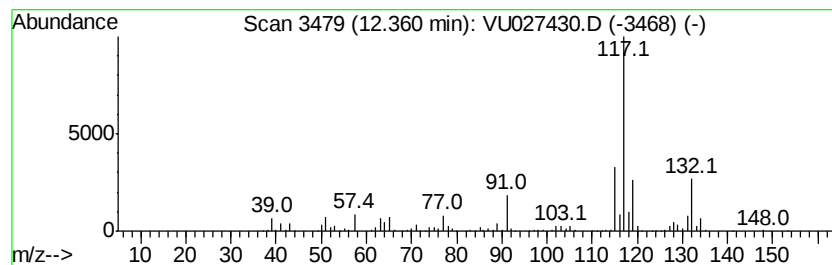
Instrument :
MSVOA_U
ClientSampleId :
EP1-MW-E-092718

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

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

Peak Number 7 Indan, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	58.64 ug/l	988597	1,4-Dichlorobenzene-d4	11.48
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indan, 1-methyl-	132 C10H12	000767-58-8	93
2	Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4	87
3	Benzene, (2-methylcyclopropyl)-	132 C10H12	1000327-39-0	87
4	Benzene, (2-methyl-1-propenyl)-	132 C10H12	000768-49-0	83
5	Benzene, 1-methyl-2-(2-propenyl)-	132 C10H12	001587-04-8	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100418\
Data File : VU027430.D
Acq On : 04 Oct 2018 23:53
Operator : MD/SY
Sample : J5165-15
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EP1-MW-E-092718

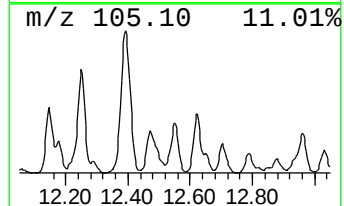
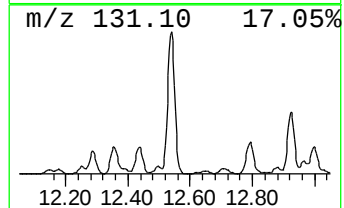
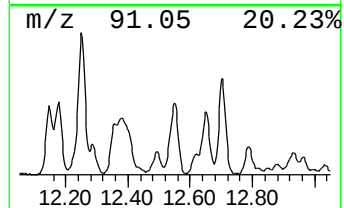
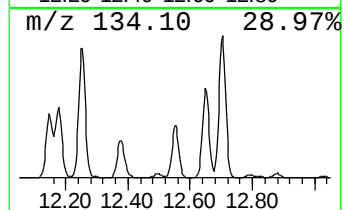
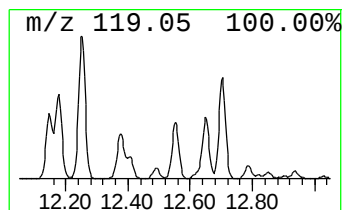
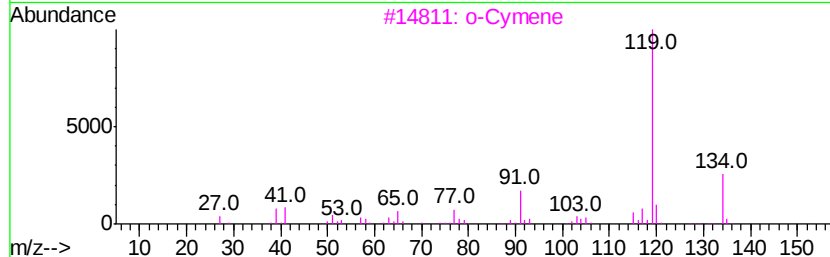
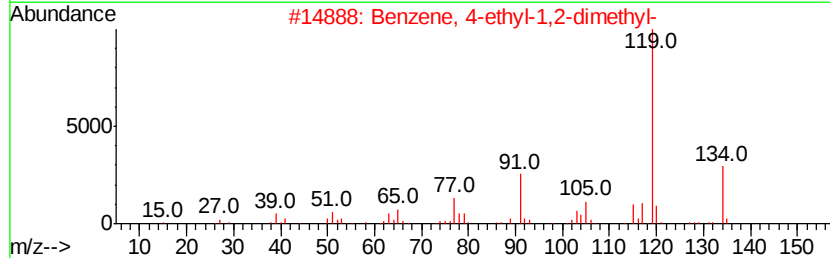
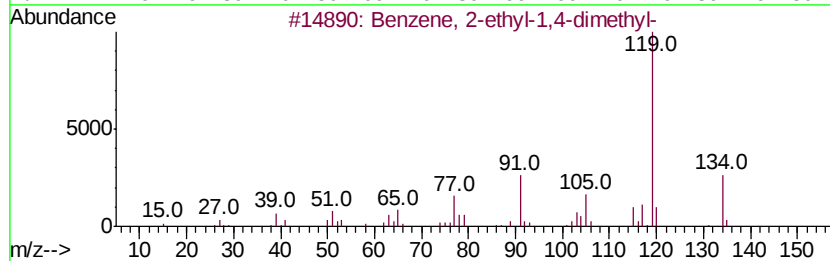
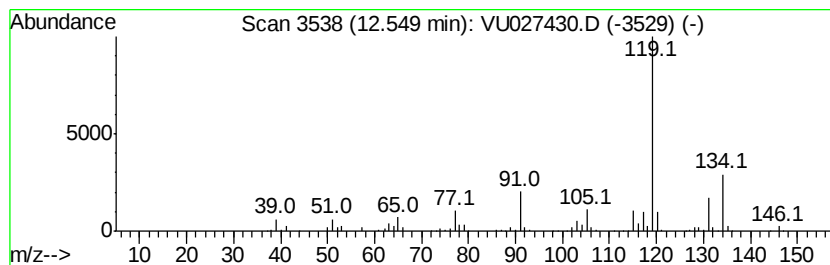
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.55	32.13 ug/l	541791	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	96
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3		o-Cymene	134	C10H14	000527-84-4	95
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5		p-Cymene	134	C10H14	000099-87-6	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100418\
Data File : VU027430.D
Acq On : 04 Oct 2018 23:53
Operator : MD/SY
Sample : J5165-15
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EP1-MW-E-092718

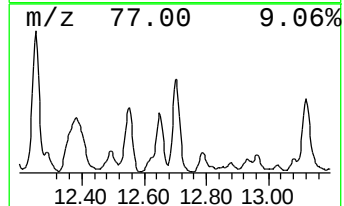
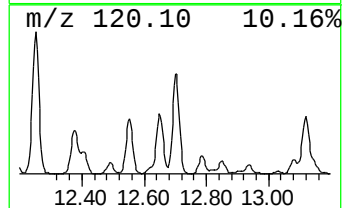
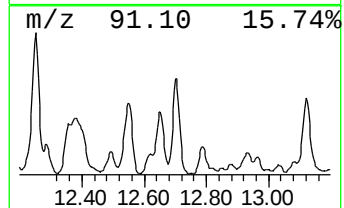
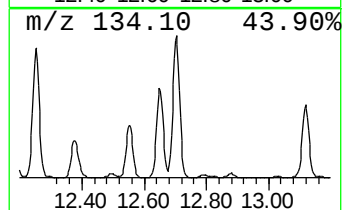
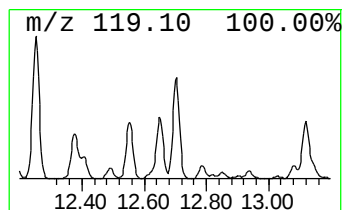
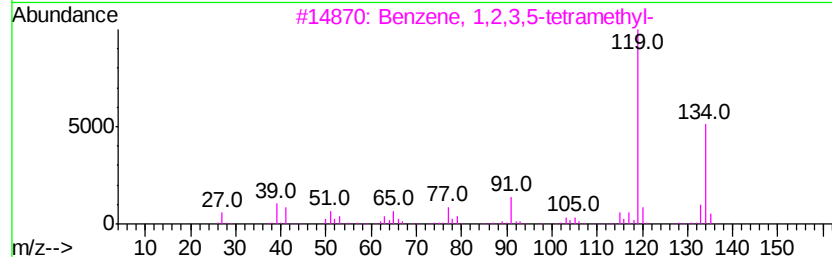
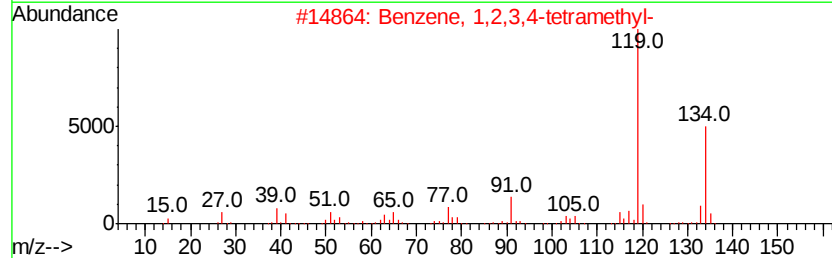
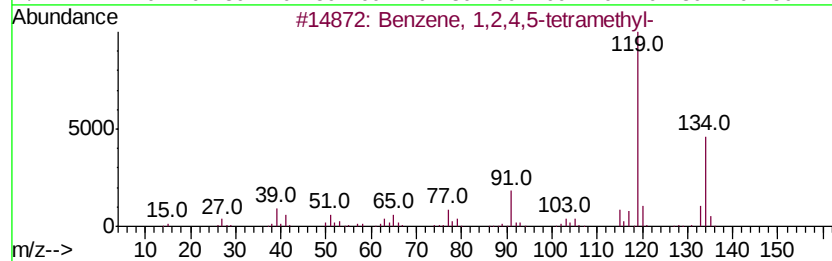
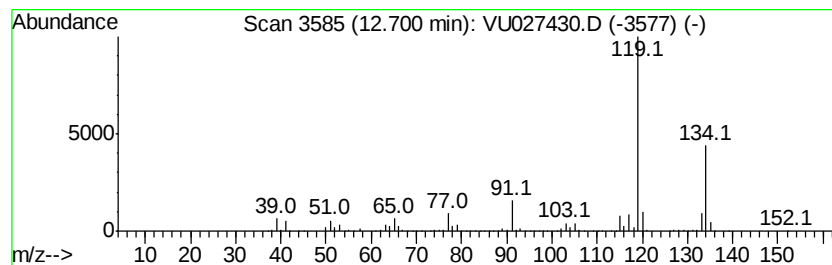
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	46.42 ug/l	782719	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	97
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
4		o-Cymene	134	C10H14	000527-84-4	94
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100418\
Data File : VU027430.D
Acq On : 04 Oct 2018 23:53
Operator : MD/SY
Sample : J5165-15
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EP1-MW-E-092718

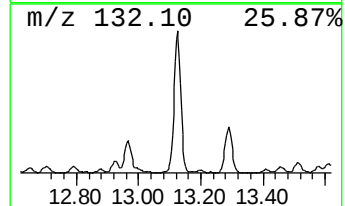
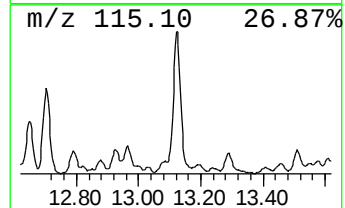
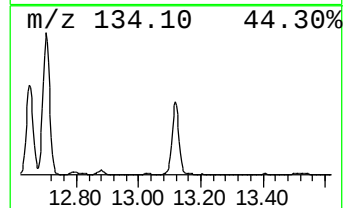
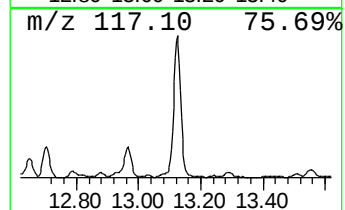
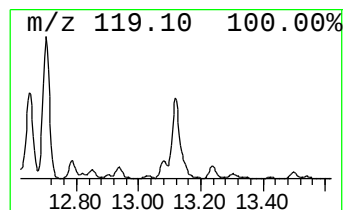
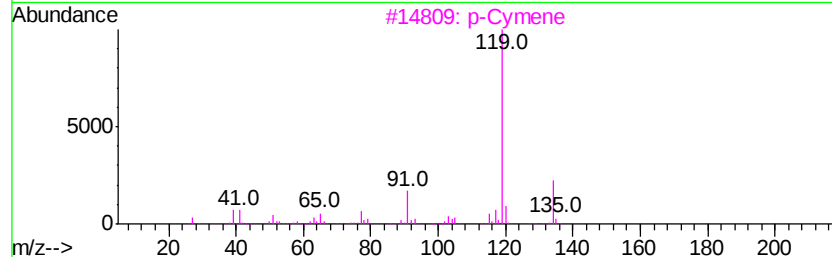
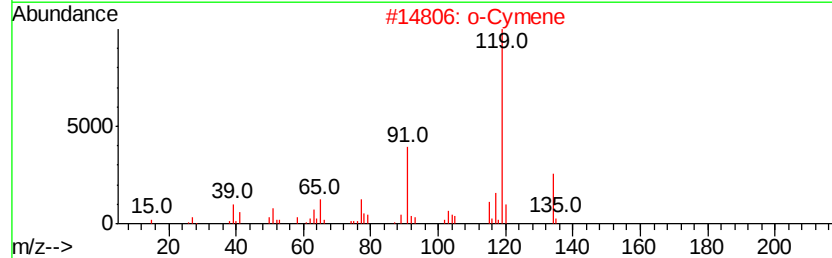
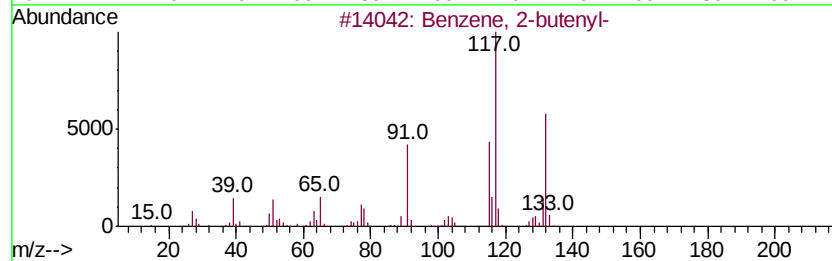
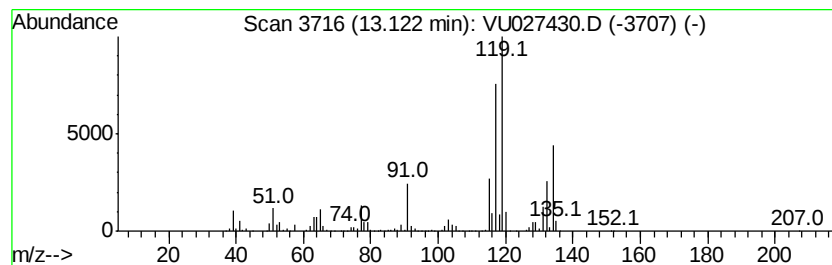
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 2-butenyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	43.54 ug/l	734131	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	64
2		o-Cymene	134	C10H14	000527-84-4	55
3		p-Cymene	134	C10H14	000099-87-6	50
4		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	50
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	50



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU100418\
 Data File : VU027430.D
 Acq On : 04 Oct 2018 23:53
 Operator : MD/SY
 Sample : J5165-15
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EP1-MW-E-092718

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.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
Octane, 2,6-dimet...	9.95	31.2	ug/l	356569	3	9.09	571623	50.0
Benzene, 1,2,3-tr...	11.57	127.2	ug/l	2145260	4	11.48	843000	50.0
Benzene, 1,2-diet...	11.78	43.7	ug/l	736541	4	11.48	843000	50.0
Benzene, 1-methyl...	11.81	37.0	ug/l	624253	4	11.48	843000	50.0
o-Cymene	12.18	31.3	ug/l	527428	4	11.48	843000	50.0
Benzene, 4-ethyl-...	12.25	56.7	ug/l	955401	4	11.48	843000	50.0
Indan, 1-methyl-	12.36	58.6	ug/l	988597	4	11.48	843000	50.0
Benzene, 2-ethyl-...	12.55	32.1	ug/l	541791	4	11.48	843000	50.0
Benzene, 1,2,4,5-...	12.70	46.4	ug/l	782719	4	11.48	843000	50.0
Benzene, 2-butenyl-	13.12	43.5	ug/l	734131	4	11.48	843000	50.0