

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU072718\
 Data File : VU025707.D
 Acq On : 28 Jul 2018 06:55
 Operator : MD/SY
 Sample : J4208-04
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 FL-0514

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	1.089	136	155	175	rBV	389244	435047	10.21%	1.321%
2	1.761	356	364	377	rBV	60400	78686	1.85%	0.239%
3	1.950	415	423	434	rVB	31884	43119	1.01%	0.131%
4	2.748	665	671	681	rVB3	27664	48463	1.14%	0.147%
5	3.002	736	750	769	rBV2	34629	74394	1.75%	0.226%
6	4.101	1075	1092	1111	rVB	68382	180081	4.23%	0.547%
7	4.889	1320	1337	1353	rBV	168752	387068	9.09%	1.175%
8	4.992	1353	1369	1386	rVV2	320303	766366	17.99%	2.327%
9	5.085	1386	1398	1418	rVB3	41954	102854	2.41%	0.312%
10	5.223	1427	1441	1456	rBV2	38784	103547	2.43%	0.314%
11	5.313	1457	1469	1487	rVB	171909	359586	8.44%	1.092%
12	5.487	1508	1523	1529	rBV4	26943	55817	1.31%	0.169%
13	5.542	1530	1540	1556	rVV	89942	231415	5.43%	0.703%
14	5.629	1556	1567	1583	rVB3	30121	67543	1.59%	0.205%
15	5.889	1633	1648	1680	rBV	334264	662522	15.56%	2.011%
16	6.420	1792	1813	1825	rBV2	172304	394593	9.26%	1.198%
17	6.741	1899	1913	1927	rVB4	21925	47950	1.13%	0.146%
18	6.928	1951	1971	1989	rBV4	93514	222607	5.23%	0.676%
19	7.053	1994	2010	2021	rBV2	120724	246080	5.78%	0.747%
20	7.532	2145	2159	2161	rBV3	116946	162545	3.82%	0.494%
21	7.567	2162	2170	2188	rVB	668377	1182385	27.76%	3.590%
22	7.712	2202	2215	2229	rBV4	40504	104647	2.46%	0.318%
23	7.780	2230	2236	2248	rVB3	32017	52815	1.24%	0.160%
24	7.921	2268	2280	2297	rVB	96090	174096	4.09%	0.529%
25	8.050	2303	2320	2331	rBV2	47766	89750	2.11%	0.272%
26	8.493	2438	2458	2466	rBV5	41736	89019	2.09%	0.270%
27	8.548	2466	2475	2484	rVB	95602	151311	3.55%	0.459%
28	8.609	2484	2494	2504	rBV	112141	185134	4.35%	0.562%
29	9.091	2633	2644	2659	rVV	500808	817809	19.20%	2.483%
30	9.191	2666	2675	2683	rVB	33343	53554	1.26%	0.163%
31	9.252	2683	2694	2710	rBV	1044025	1633194	38.35%	4.959%
32	9.378	2717	2733	2750	rBV	2578063	4259156	100.00%	12.931%
33	9.448	2751	2755	2780	rVB7	20212	49247	1.16%	0.150%
34	9.725	2826	2841	2855	rBV9	34485	103248	2.42%	0.313%

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Integrator: RTE
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35	9.956	2892	2913	2927	rBV5	51465	111763	2.62%	0.339%
36	10.050	2932	2942	2968	rVB7	35310	80346	1.89%	0.244%
37	10.169	2968	2979	2990	rBV	215089	328843	7.72%	0.998%
38	10.310	3012	3023	3037	rBV	534157	844900	19.84%	2.565%
39	10.497	3073	3081	3098	rVB4	34841	63568	1.49%	0.193%
40	10.590	3099	3110	3121	rBV	250328	386696	9.08%	1.174%
41	10.683	3127	3139	3144	rBV	1312700	2148483	50.44%	6.523%
42	10.773	3157	3167	3188	rVB	1006921	1551723	36.43%	4.711%
43	10.972	3210	3229	3250	rBV	937958	1489392	34.97%	4.522%
44	11.152	3273	3285	3309	rVB	2140185	3247418	76.25%	9.859%
45	11.326	3332	3339	3354	rVB	41111	66783	1.57%	0.203%
46	11.429	3355	3371	3378	rBV	120649	191291	4.49%	0.581%
47	11.487	3378	3389	3403	rVV2	741521	1197571	28.12%	3.636%
48	11.574	3405	3416	3429	rVB	393344	593814	13.94%	1.803%
49	11.693	3441	3453	3463	rVB	31773	52465	1.23%	0.159%
50	11.770	3463	3477	3485	rBV3	305372	573647	13.47%	1.742%
51	11.812	3486	3490	3498	rVV	134867	181957	4.27%	0.552%
52	11.873	3498	3509	3527	rVB2	207486	426604	10.02%	1.295%
53	11.982	3531	3543	3555	rBV	70263	116637	2.74%	0.354%
54	12.059	3555	3567	3580	rVB	172135	268406	6.30%	0.815%
55	12.149	3580	3595	3599	rBV	166379	256106	6.01%	0.778%
56	12.178	3599	3604	3614	rVB	199297	285570	6.70%	0.867%
57	12.255	3617	3628	3634	rVV	178804	269472	6.33%	0.818%
58	12.287	3634	3638	3649	rVB	74390	103015	2.42%	0.313%
59	12.358	3650	3660	3683	rVV3	157249	395141	9.28%	1.200%
60	12.496	3692	3703	3710	rBV4	33337	57405	1.35%	0.174%
61	12.551	3711	3720	3732	rVB2	89052	162135	3.81%	0.492%
62	12.651	3732	3751	3759	rBV	136246	240374	5.64%	0.730%
63	12.705	3759	3768	3783	rVB	242253	373360	8.77%	1.134%
64	12.792	3783	3795	3802	rBV2	37396	64631	1.52%	0.196%
65	12.966	3843	3849	3864	rVB2	72304	107071	2.51%	0.325%
66	13.123	3888	3898	3913	rVB3	460696	728968	17.12%	2.213%
67	13.291	3940	3950	3960	rBV	228428	348891	8.19%	1.059%
68	13.512	4009	4019	4026	rBV3	54534	92650	2.18%	0.281%
69	13.744	4074	4091	4102	rBV	611931	976675	22.93%	2.965%
70	13.860	4117	4127	4143	rVB3	26382	49026	1.15%	0.149%
71	13.956	4147	4157	4168	rBV6	28258	60523	1.42%	0.184%

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Integration Parameters: RTEINT.P

Integrator: RTE
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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	14.352	4266	4280	4292	rVB6	44931	105279	2.47%	0.320%
73	14.680	4372	4382	4400	rBV2	43492	85517	2.01%	0.260%
74	14.879	4432	4444	4456	rBV	225657	374362	8.79%	1.137%
75	15.065	4491	4502	4516	rBV	156103	262906	6.17%	0.798%

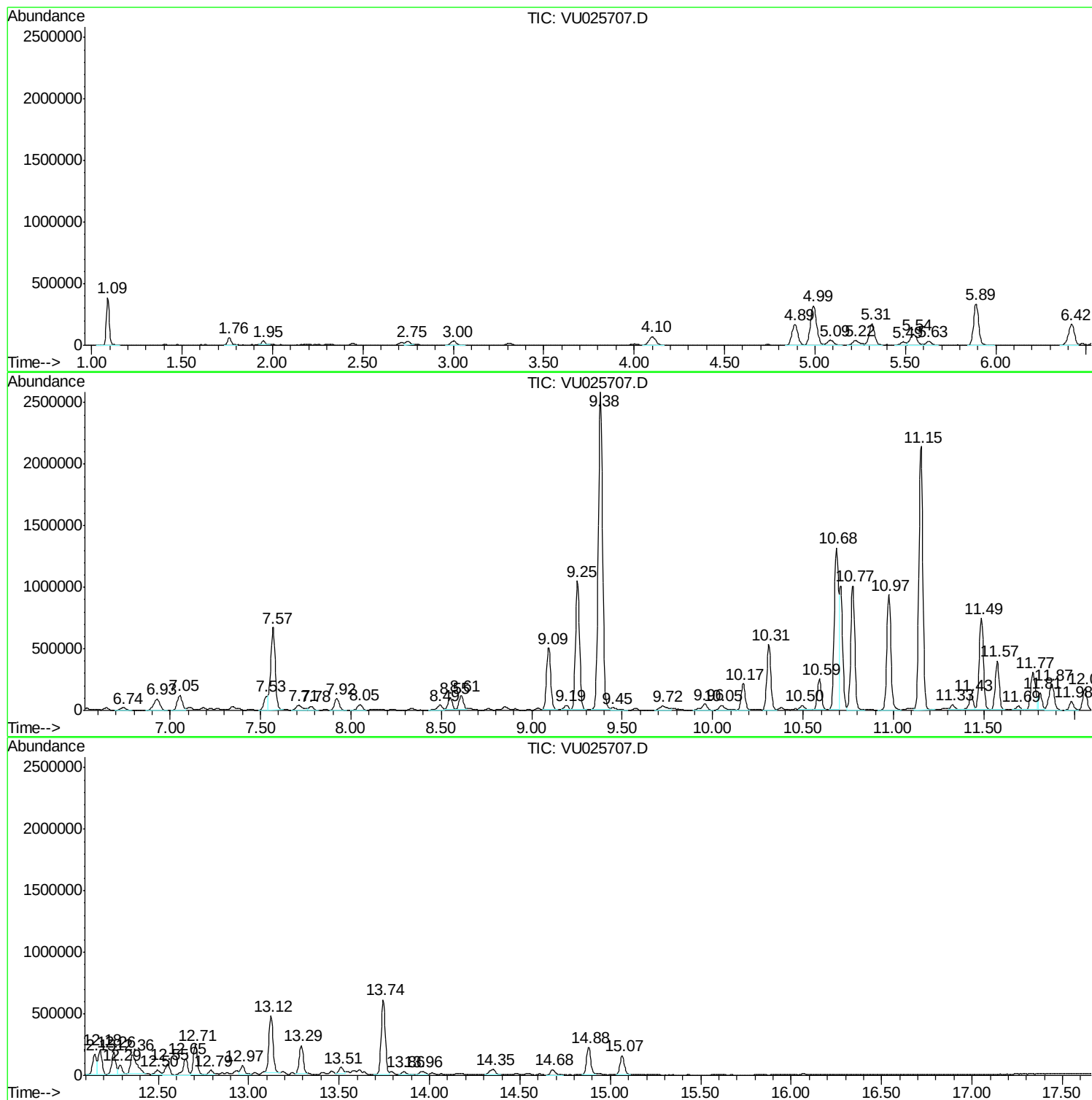
Sum of corrected areas: 32937032

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Instrument :
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ClientSampleId :
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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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU072718\
Data File : VU025707.D
Acq On : 28 Jul 2018 06:55
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Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 53 Sample Multiplier: 1

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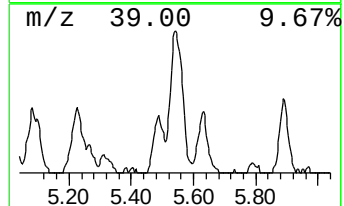
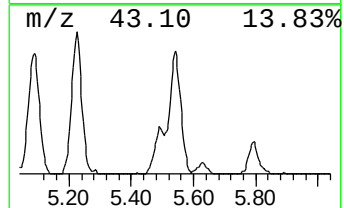
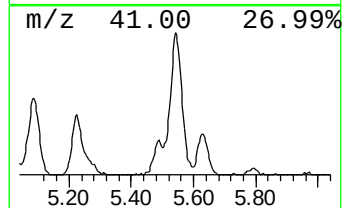
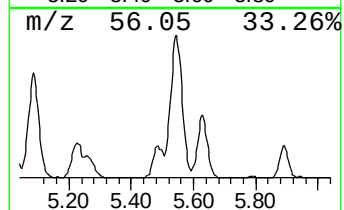
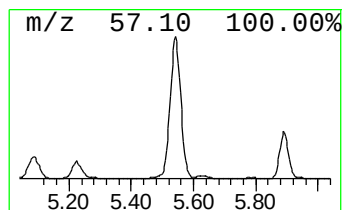
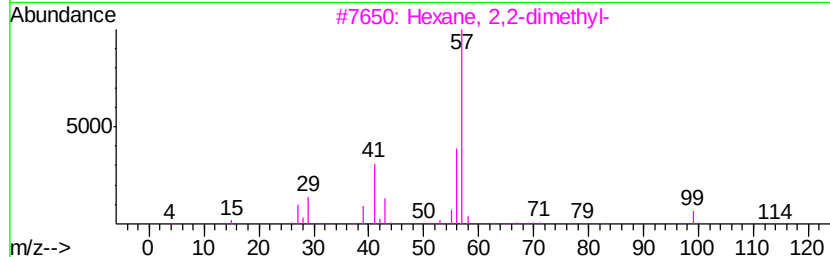
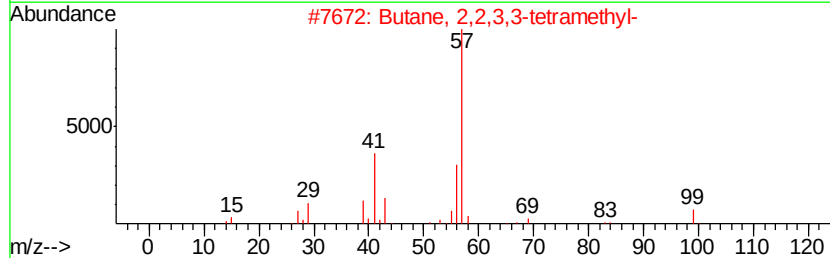
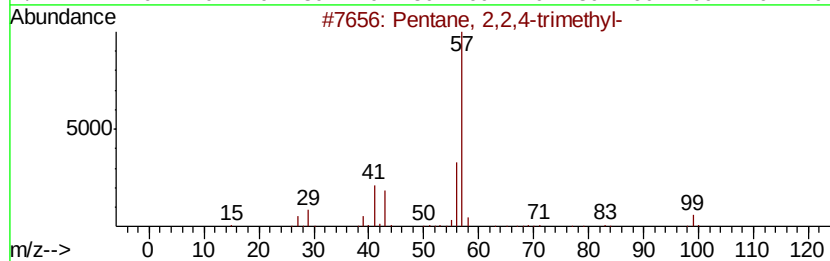
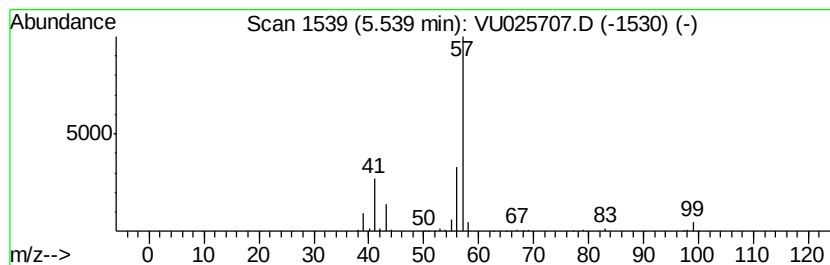
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.M
Quant Title : SW846 8260

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

Peak Number 2 Pentane, 2,2,4-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.54	17.46 ug/l	231415	1,4-Difluorobenzene	5.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
2		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
3		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
4		Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	64
5		Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59



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Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 53 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampled :
FL-0514

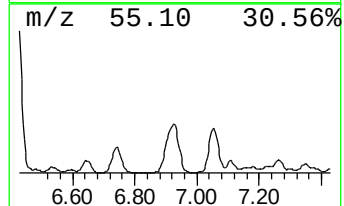
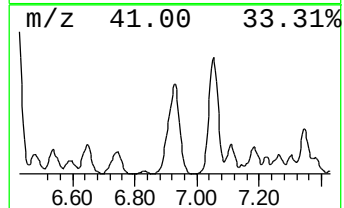
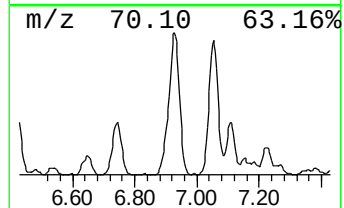
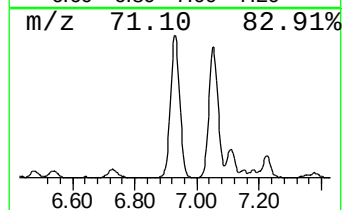
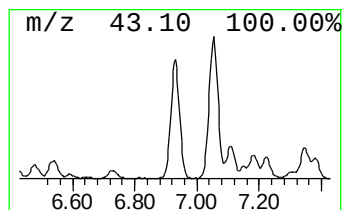
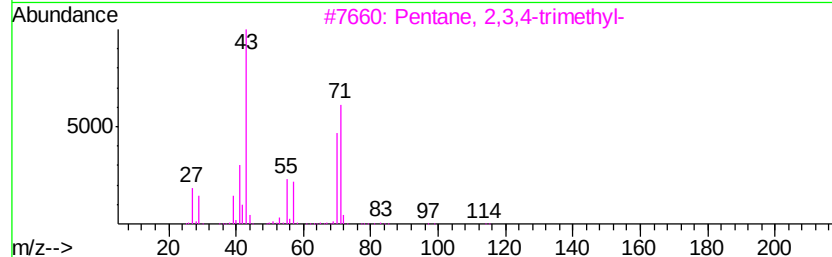
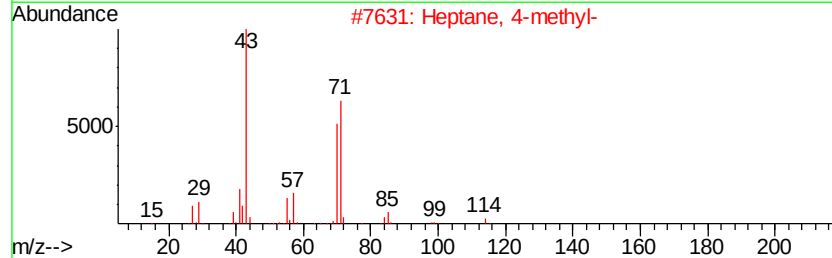
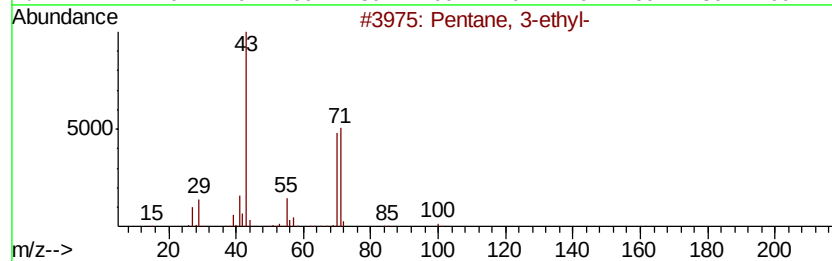
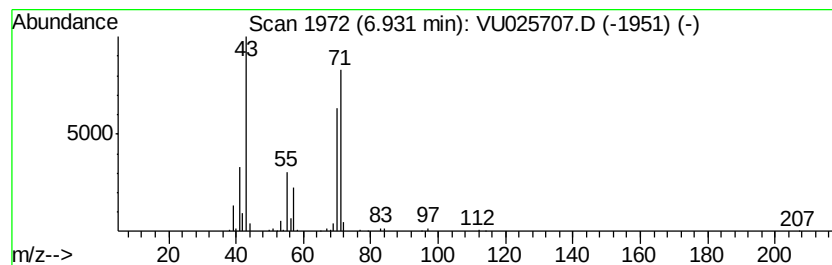
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.M
Quant Title : SW846 8260

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

Peak Number 3 Pentane, 3-ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.93	16.80 ug/l	222607	1,4-Difluorobenzene	5.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 3-ethyl-	100	C7H16	000617-78-7	83
2		Heptane, 4-methyl-	114	C8H18	000589-53-7	78
3		Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	78
4		Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	74
5		Hexane, 3-methyl-	100	C7H16	000589-34-4	72



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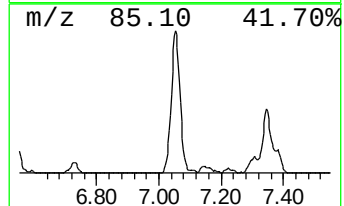
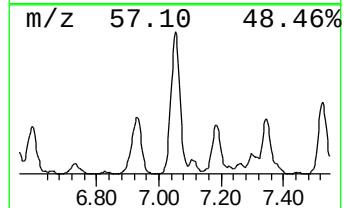
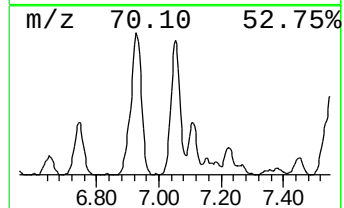
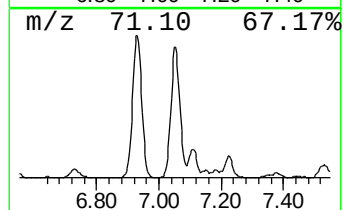
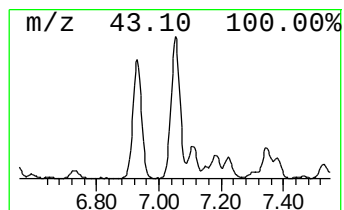
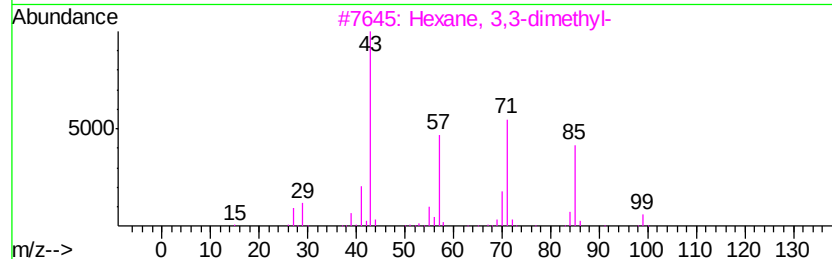
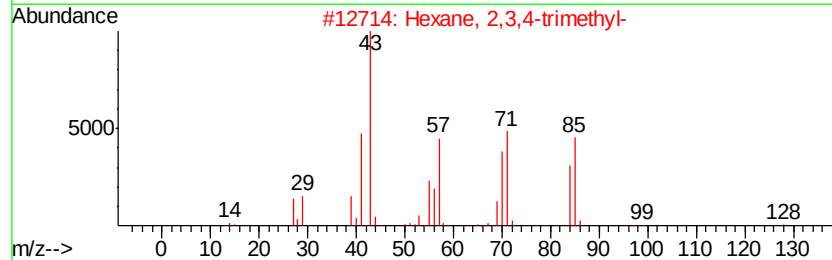
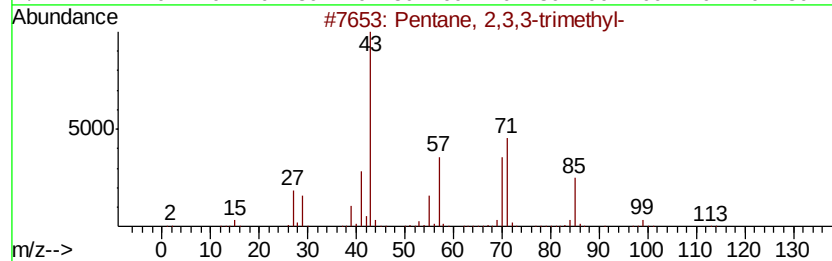
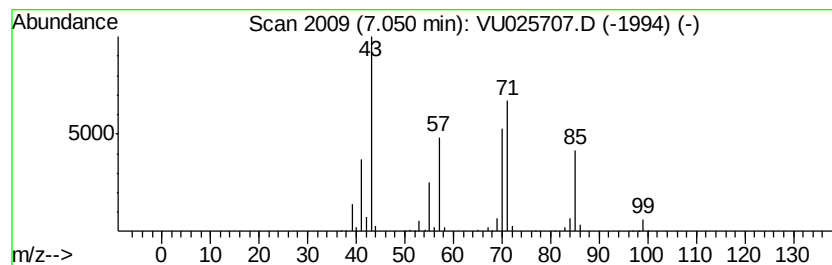
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Quant Title : SW846 8260

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

Peak Number 4 Pentane, 2,3,3-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.05	18.57 ug/l	246080	1,4-Difluorobenzene	5.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2		Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	78
3		Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
4		Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	50
5		Undecane, 2,4-dimethyl-	184	C13H28	017312-80-0	47



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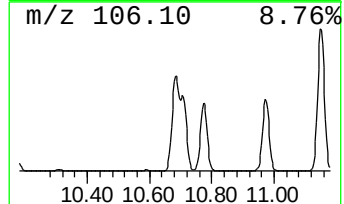
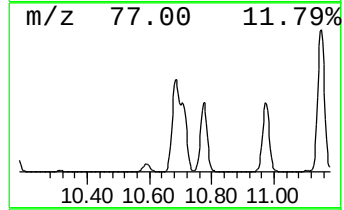
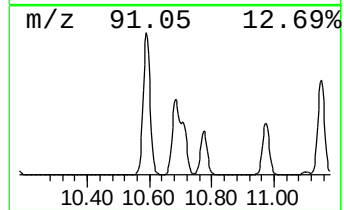
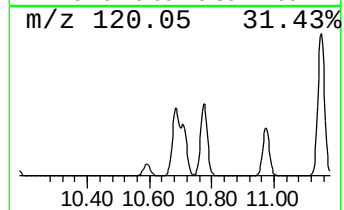
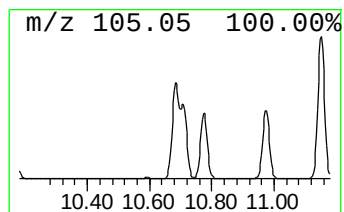
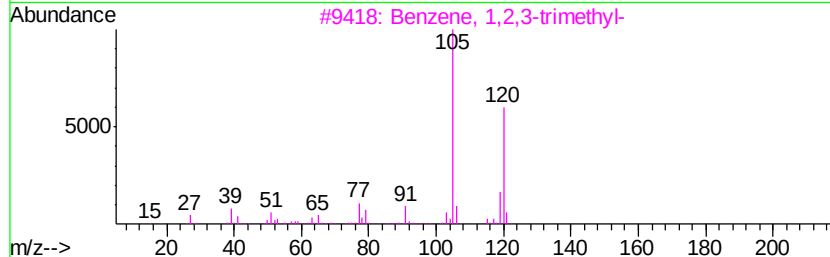
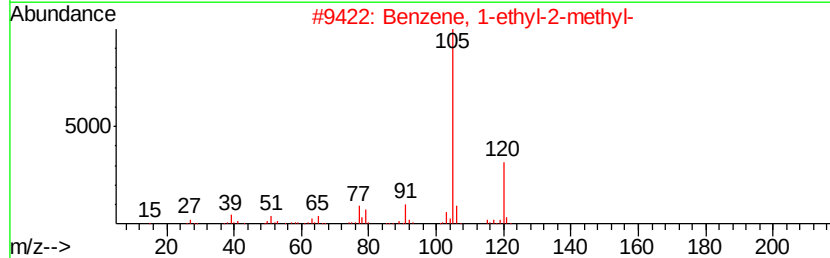
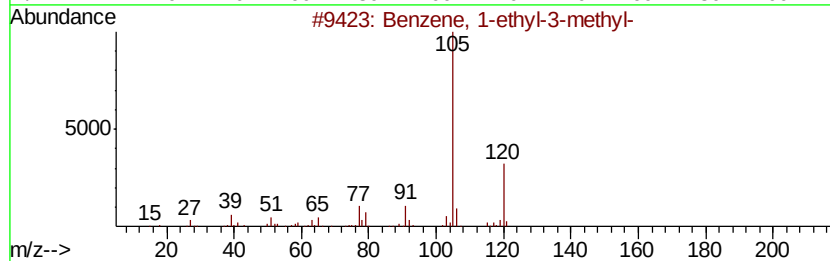
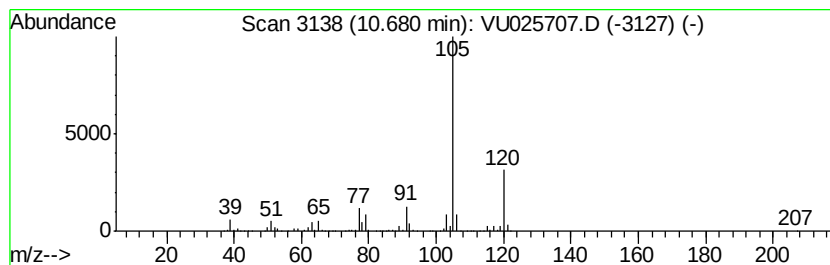
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethyl-3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.68	89.70 ug/l	2148480	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		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU072718\
Data File : VU025707.D
Acq On : 28 Jul 2018 06:55
Operator : MD/SY
Sample : J4208-04
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 53 Sample Multiplier: 1

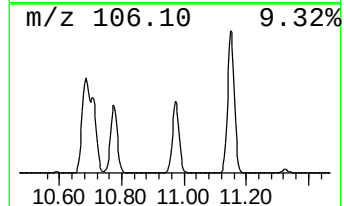
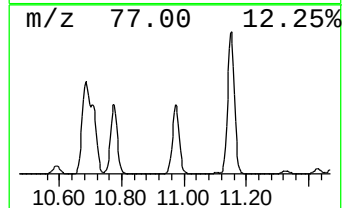
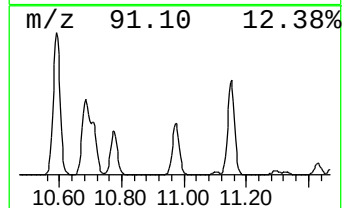
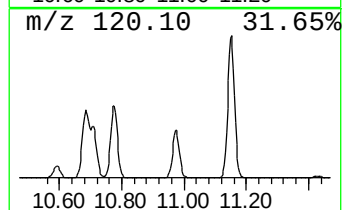
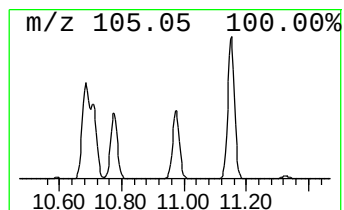
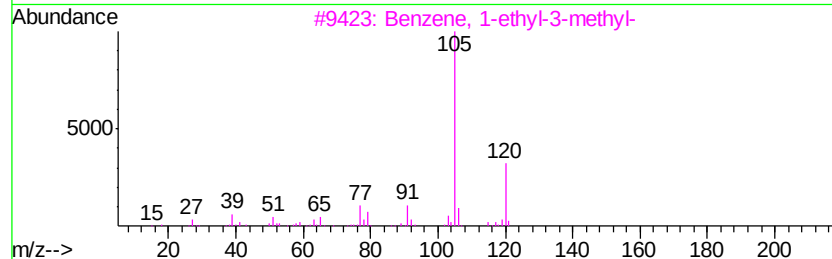
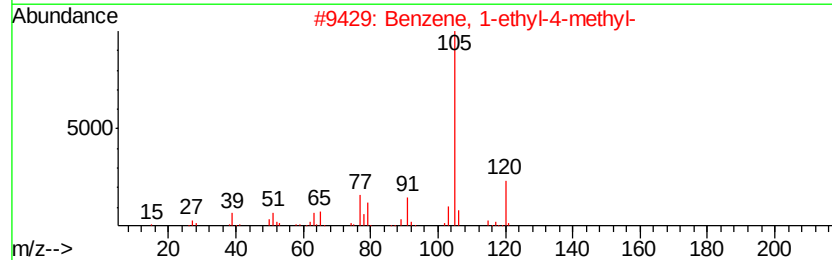
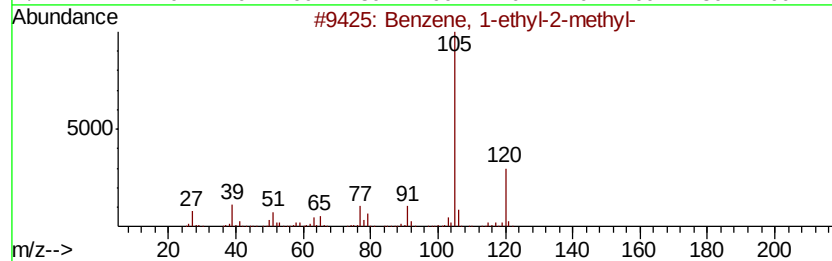
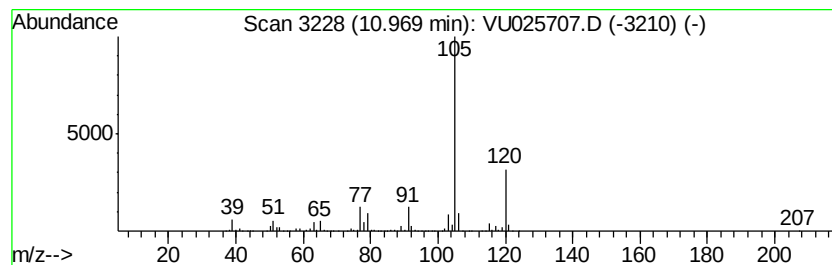
Instrument :
MSVOA_U
ClientSampleId :
FL-0514

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 6 Benzene, 1-ethyl-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.97	62.18 ug/l	1489390	1,4-Dichlorobenzene-d4	11.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethyl-2-methyl-	120 C9H12	000611-14-3 95
2		Benzene, 1-ethyl-4-methyl-	120 C9H12	000622-96-8 94
3		Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4 91
4		Mesitylene	120 C9H12	000108-67-8 91
5		Benzene, 1,2,3-trimethyl-	120 C9H12	000526-73-8 91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU072718\
Data File : VU025707.D
Acq On : 28 Jul 2018 06:55
Operator : MD/SY
Sample : J4208-04
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 53 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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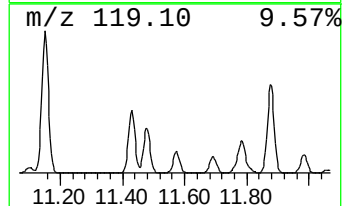
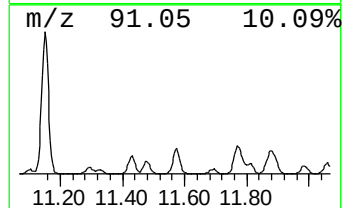
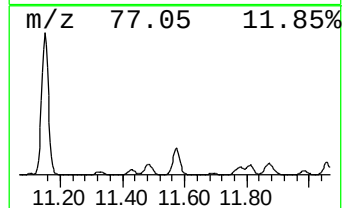
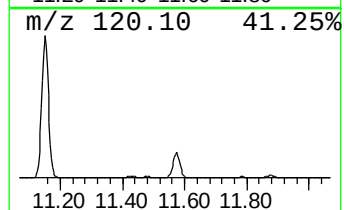
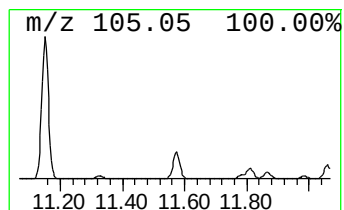
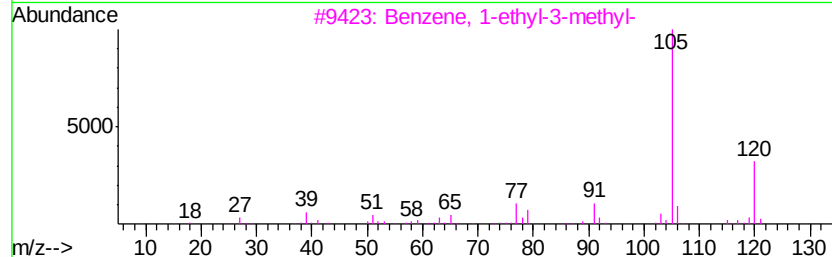
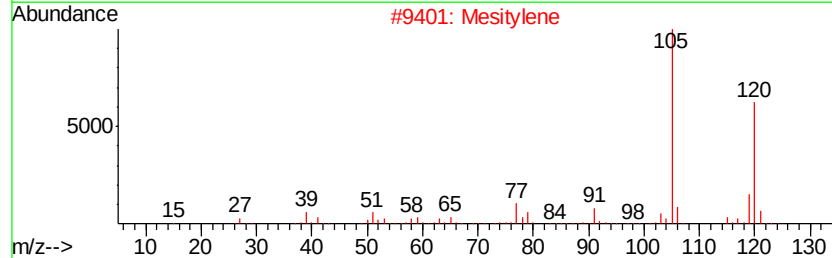
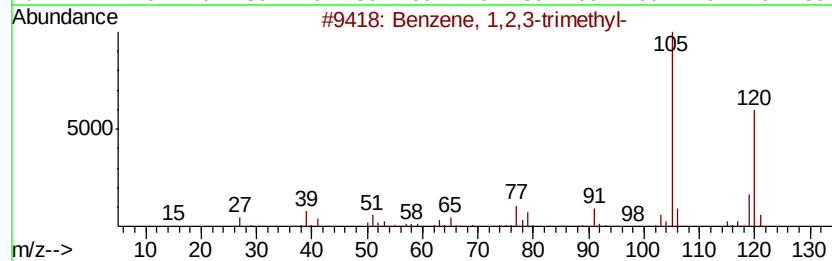
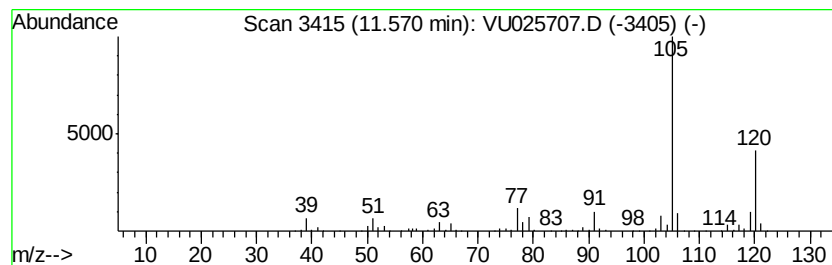
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	24.79 ug/l	593814	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	94
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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU072718\
Data File : VU025707.D
Acq On : 28 Jul 2018 06:55
Operator : MD/SY
Sample : J4208-04
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 53 Sample Multiplier: 1

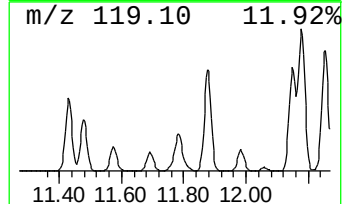
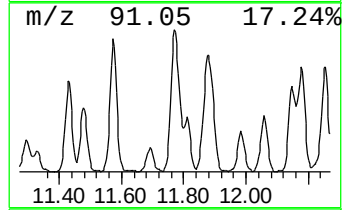
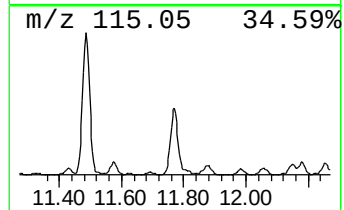
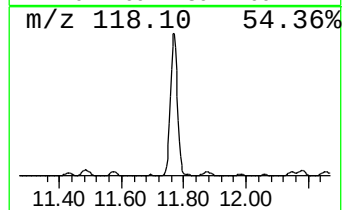
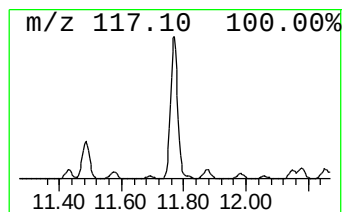
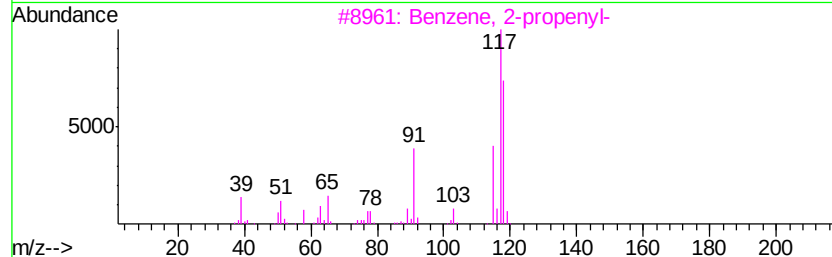
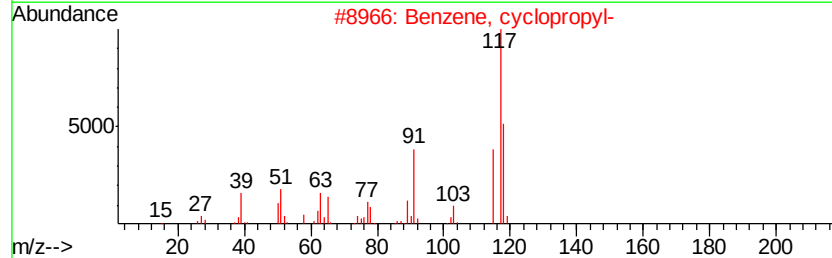
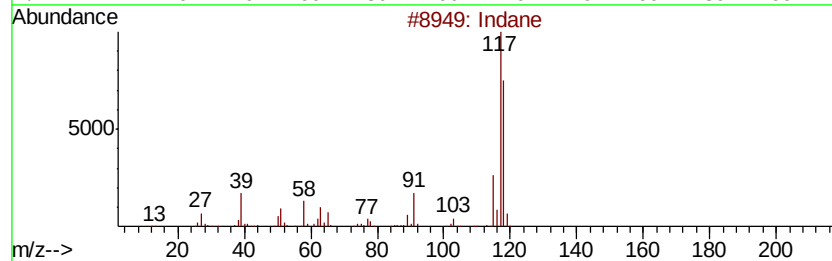
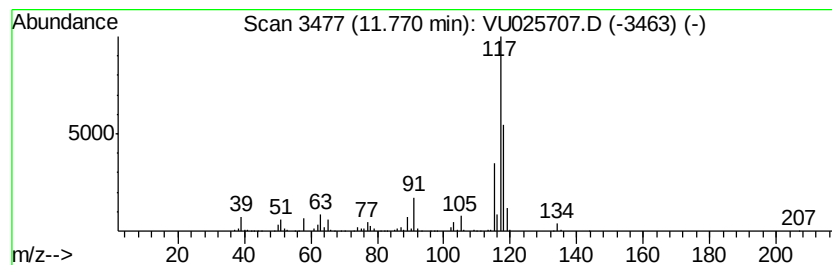
Instrument :
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ClientSampleId :
FL-0514

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 8 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	23.95 ug/l	573647	1,4-Dichlorobenzene-d4	11.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indane		118 C9H10	000496-11-7 81
2	Benzene, cyclopropyl-		118 C9H10	000873-49-4 76
3	Benzene, 2-propenyl-		118 C9H10	000300-57-2 72
4	Benzene, 1-ethenyl-2-methyl-		118 C9H10	000611-15-4 68
5	Deltacyclene		118 C9H10	007785-10-6 64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU072718\
Data File : VU025707.D
Acq On : 28 Jul 2018 06:55
Operator : MD/SY
Sample : J4208-04
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 53 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
FL-0514

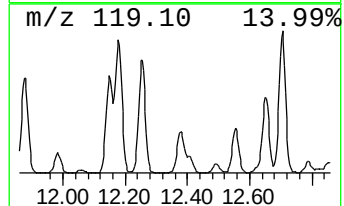
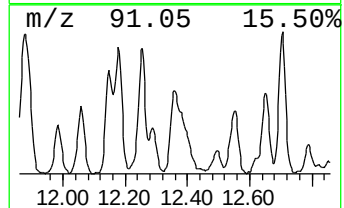
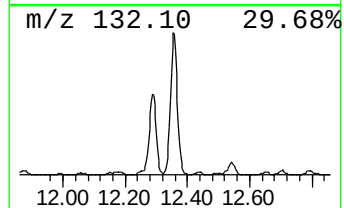
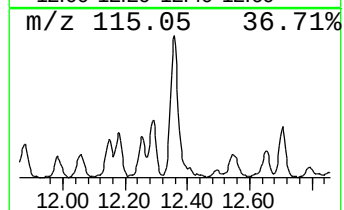
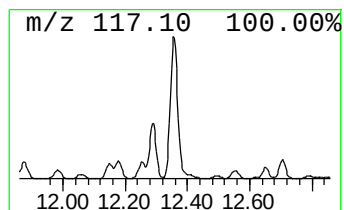
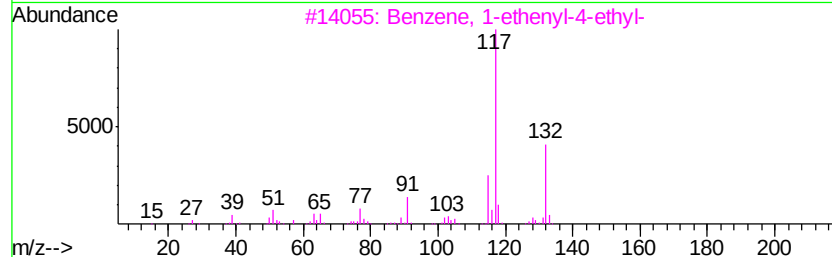
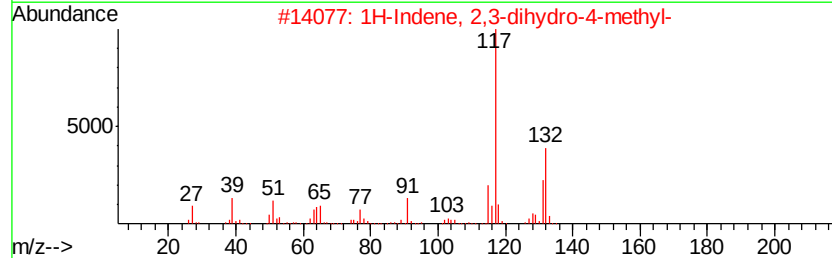
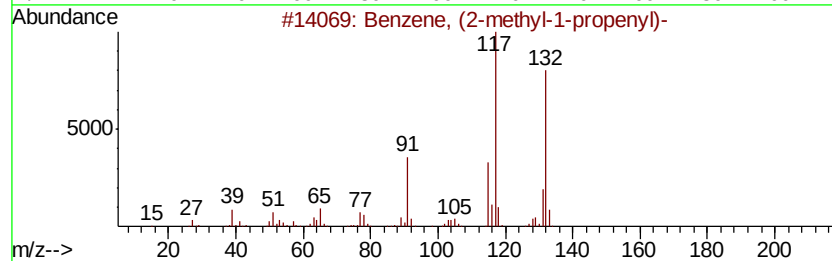
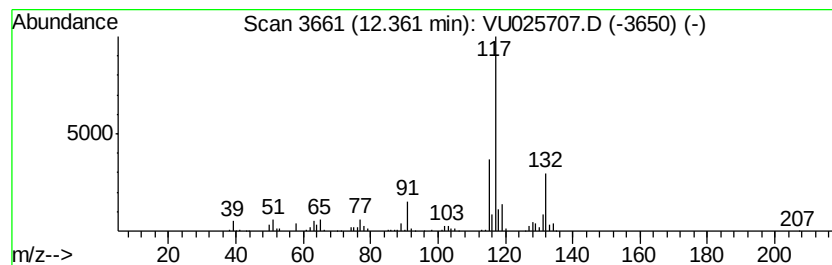
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, (2-methyl-1-propen... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	16.50 ug/l	395141	1,4-Dichlorobenzene-d4	11.49

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	91
2	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
5	Indan, 1-methyl-	132	C10H12	000767-58-8	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU072718\
Data File : VU025707.D
Acq On : 28 Jul 2018 06:55
Operator : MD/SY
Sample : J4208-04
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 53 Sample Multiplier: 1

Instrument :
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ClientSampleId :
FL-0514

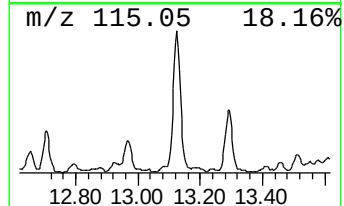
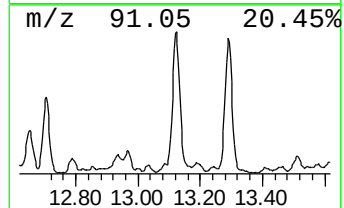
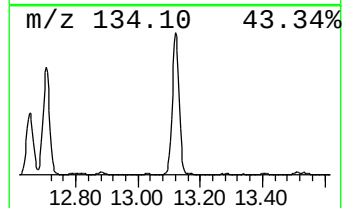
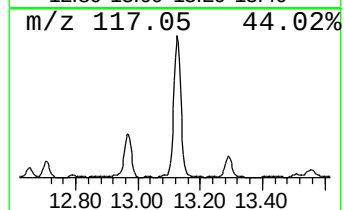
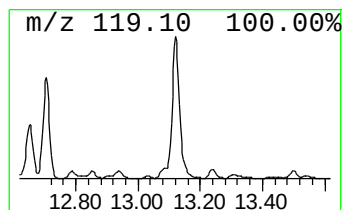
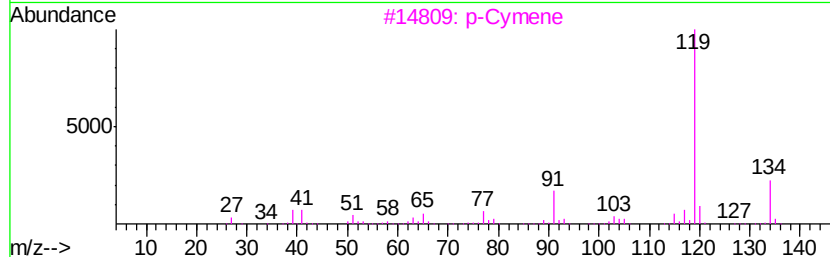
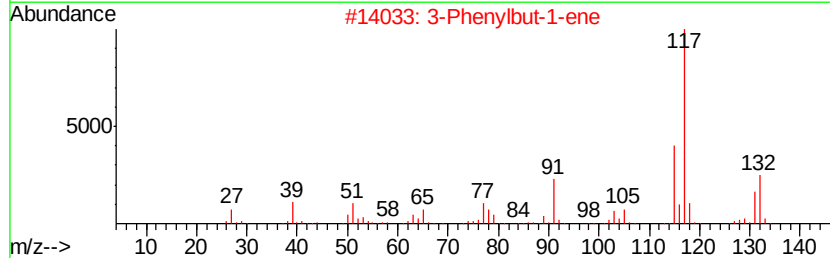
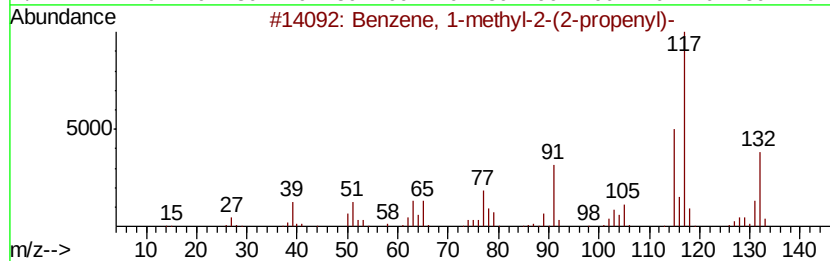
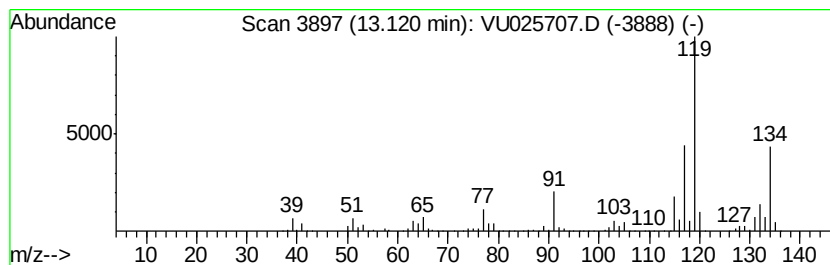
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1-methyl-2-(2-prop... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	30.44 ug/l	728968	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
3		p-Cymene	134	C10H14	000099-87-6	64
4		o-Cymene	134	C10H14	000527-84-4	64
5		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU072718\
Data File : VU025707.D
Acq On : 28 Jul 2018 06:55
Operator : MD/SY
Sample : J4208-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 53 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
FL-0514

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
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Pentane, 3-ethyl-	6.93	16.8	ug/l	222607	2	5.89	662522	50.0
Pentane, 2,3,3-tr...	7.05	18.6	ug/l	246080	2	5.89	662522	50.0
Benzene, 1-ethyl-...	10.68	89.7	ug/l	2148480	4	11.49	1197570	50.0
Benzene, 1-ethyl-...	10.97	62.2	ug/l	1489390	4	11.49	1197570	50.0
Benzene, 1,2,3-tr...	11.57	24.8	ug/l	593814	4	11.49	1197570	50.0
Indane	11.77	23.9	ug/l	573647	4	11.49	1197570	50.0
Benzene, (2-methy...	12.36	16.5	ug/l	395141	4	11.49	1197570	50.0
Benzene, 1-methyl...	13.12	30.4	ug/l	728968	4	11.49	1197570	50.0