

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100418\
 Data File : VU027425.D
 Acq On : 04 Oct 2018 21:57
 Operator : MD/SY
 Sample : J5232-03
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MW-3

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	1.305	34	41	55	rVB	37746	37386	1.29%	0.151%
2	1.402	61	71	82	rBV	142365	149744	5.16%	0.605%
3	1.759	172	182	198	rBV	355902	467619	16.11%	1.889%
4	1.948	233	241	255	rVB3	96813	164165	5.65%	0.663%
5	2.051	264	273	284	rVB	78527	105818	3.64%	0.427%
6	2.138	291	300	307	rBV	44865	64333	2.22%	0.260%
7	2.186	307	315	329	rVB	103816	155770	5.36%	0.629%
8	2.305	341	352	372	rVB4	17173	39220	1.35%	0.158%
9	2.447	384	396	416	rBV	45997	83202	2.87%	0.336%
10	2.707	460	477	482	rBV4	44296	133365	4.59%	0.539%
11	2.746	483	489	495	rVV	72775	130307	4.49%	0.526%
12	2.794	495	504	535	rVB	199319	424245	14.61%	1.714%
13	3.000	551	568	586	rBV2	114105	247990	8.54%	1.002%
14	3.556	731	741	758	rVB3	36228	81670	2.81%	0.330%
15	3.659	758	773	785	rBV2	27226	63423	2.18%	0.256%
16	3.733	785	796	811	rVV5	17474	45029	1.55%	0.182%
17	3.894	833	846	863	rVB3	24159	60645	2.09%	0.245%
18	4.096	891	909	926	rVV2	149685	402116	13.85%	1.624%
19	4.189	927	938	966	rVB5	19944	60040	2.07%	0.243%
20	4.784	1114	1123	1139	rVV2	19468	44840	1.54%	0.181%
21	4.884	1139	1154	1172	rVV	110372	256383	8.83%	1.036%
22	4.990	1172	1187	1204	rVV2	214187	534159	18.40%	2.158%
23	5.083	1205	1216	1241	rVB3	41530	121460	4.18%	0.491%
24	5.222	1246	1259	1274	rBV3	27534	72036	2.48%	0.291%
25	5.312	1275	1287	1299	rBV	130656	258893	8.92%	1.046%
26	5.392	1299	1312	1326	rBV	618882	1293203	44.54%	5.224%
27	5.623	1374	1384	1399	rVB4	30262	70275	2.42%	0.284%
28	5.787	1421	1435	1452	rBV5	17213	41917	1.44%	0.169%
29	5.887	1453	1466	1478	rVV	233719	472243	16.26%	1.908%
30	5.951	1479	1486	1508	rVV5	49910	144188	4.97%	0.582%
31	6.222	1556	1570	1588	rVV4	23209	52558	1.81%	0.212%
32	6.414	1609	1630	1645	rBV4	84504	234384	8.07%	0.947%
33	6.643	1690	1701	1714	rVB3	18158	34405	1.18%	0.139%
34	6.829	1744	1759	1775	rBV5	22552	65430	2.25%	0.264%

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Integrator: RTE
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35	6.926	1775	1789	1804	rVB4	22390	55003	1.89%	0.222%
36	7.067	1811	1833	1844	rBV5	55793	147367	5.08%	0.595%
37	7.299	1895	1905	1916	rBV	56246	100479	3.46%	0.406%
38	7.366	1918	1926	1937	rVV	38806	69022	2.38%	0.279%
39	7.431	1938	1946	1968	rVB4	24860	57355	1.98%	0.232%
40	7.566	1968	1988	2002	rBV	456046	849204	29.25%	3.431%
41	7.640	2003	2011	2023	rVV	96636	169992	5.85%	0.687%
42	7.717	2023	2035	2062	rVB4	52783	120516	4.15%	0.487%
43	8.096	2145	2153	2167	rVB	23047	36705	1.26%	0.148%
44	8.173	2167	2177	2190	rBV4	15206	35598	1.23%	0.144%
45	8.398	2230	2247	2265	rBV	281609	533636	18.38%	2.156%
46	8.848	2378	2387	2409	rVB5	13644	30912	1.06%	0.125%
47	9.025	2429	2442	2451	rBV3	31548	65302	2.25%	0.264%
48	9.090	2452	2462	2480	rVV	349768	626142	21.56%	2.530%
49	9.170	2480	2487	2500	rVB3	44018	75935	2.62%	0.307%
50	9.250	2500	2512	2535	rBV	1654442	2693302	92.76%	10.880%
51	9.379	2543	2552	2563	rVB	72793	119192	4.11%	0.482%
52	9.781	2665	2677	2694	rVB	78984	134660	4.64%	0.544%
53	10.167	2788	2797	2808	rVB	85855	136341	4.70%	0.551%
54	10.308	2830	2841	2854	rBV	352174	565495	19.48%	2.284%
55	10.414	2866	2874	2891	rVB6	17896	47180	1.62%	0.191%
56	10.588	2919	2928	2944	rVB	107592	188705	6.50%	0.762%
57	10.726	2962	2971	2977	rBV4	48235	83544	2.88%	0.338%
58	10.771	2977	2985	2995	rVV	192463	294598	10.15%	1.190%
59	10.971	3035	3047	3072	rBV	660810	1033527	35.60%	4.175%
60	11.151	3091	3103	3117	rVB	1909902	2903532	100.00%	11.730%
61	11.350	3152	3165	3175	rVB6	16510	44210	1.52%	0.179%
62	11.485	3196	3207	3221	rVB	450879	710282	24.46%	2.869%
63	11.572	3223	3234	3244	rBV	110822	171288	5.90%	0.692%
64	11.768	3280	3295	3316	rBV	1254820	2108465	72.62%	8.518%
65	11.871	3316	3327	3342	rVB3	56008	110579	3.81%	0.447%
66	11.980	3342	3361	3373	rBV	112007	193835	6.68%	0.783%
67	12.057	3377	3385	3396	rVB2	32775	53829	1.85%	0.217%
68	12.147	3401	3413	3418	rBV	91267	141643	4.88%	0.572%
69	12.250	3434	3445	3452	rBV	186731	289144	9.96%	1.168%
70	12.286	3453	3456	3466	rVV2	55556	70364	2.42%	0.284%
71	12.353	3467	3477	3499	rVB	185658	326695	11.25%	1.320%

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Sampling : 1 Min Area: 3 % of largest Peak
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72	12.549	3530	3538	3549	rVB3	18733	33546	1.16%	0.136%
73	12.649	3550	3569	3577	rBV	104302	172502	5.94%	0.697%
74	12.704	3577	3586	3597	rVB	141862	224255	7.72%	0.906%
75	12.964	3658	3667	3677	rVB	117458	167096	5.75%	0.675%
76	13.125	3707	3717	3729	rVV2	348682	564304	19.44%	2.280%
77	13.186	3730	3736	3746	rVB3	28954	47636	1.64%	0.192%
78	13.289	3760	3768	3786	rVB3	53066	94916	3.27%	0.383%
79	13.456	3812	3820	3829	rBV3	24720	39911	1.37%	0.161%
80	13.511	3829	3837	3843	rBV3	37780	58957	2.03%	0.238%
81	13.742	3897	3909	3920	rBV	618040	964646	33.22%	3.897%
82	13.861	3937	3946	3965	rVB	49267	93114	3.21%	0.376%
83	14.150	4028	4036	4050	rVB2	21328	38648	1.33%	0.156%
84	14.334	4085	4093	4107	rVB3	21173	38109	1.31%	0.154%
85	14.478	4130	4138	4147	rVV2	21395	32670	1.13%	0.132%
86	14.546	4150	4159	4170	rVB5	22569	43505	1.50%	0.176%
87	14.877	4253	4262	4272	rBV2	24208	42220	1.45%	0.171%
88	15.064	4310	4320	4332	rBV	57886	91498	3.15%	0.370%

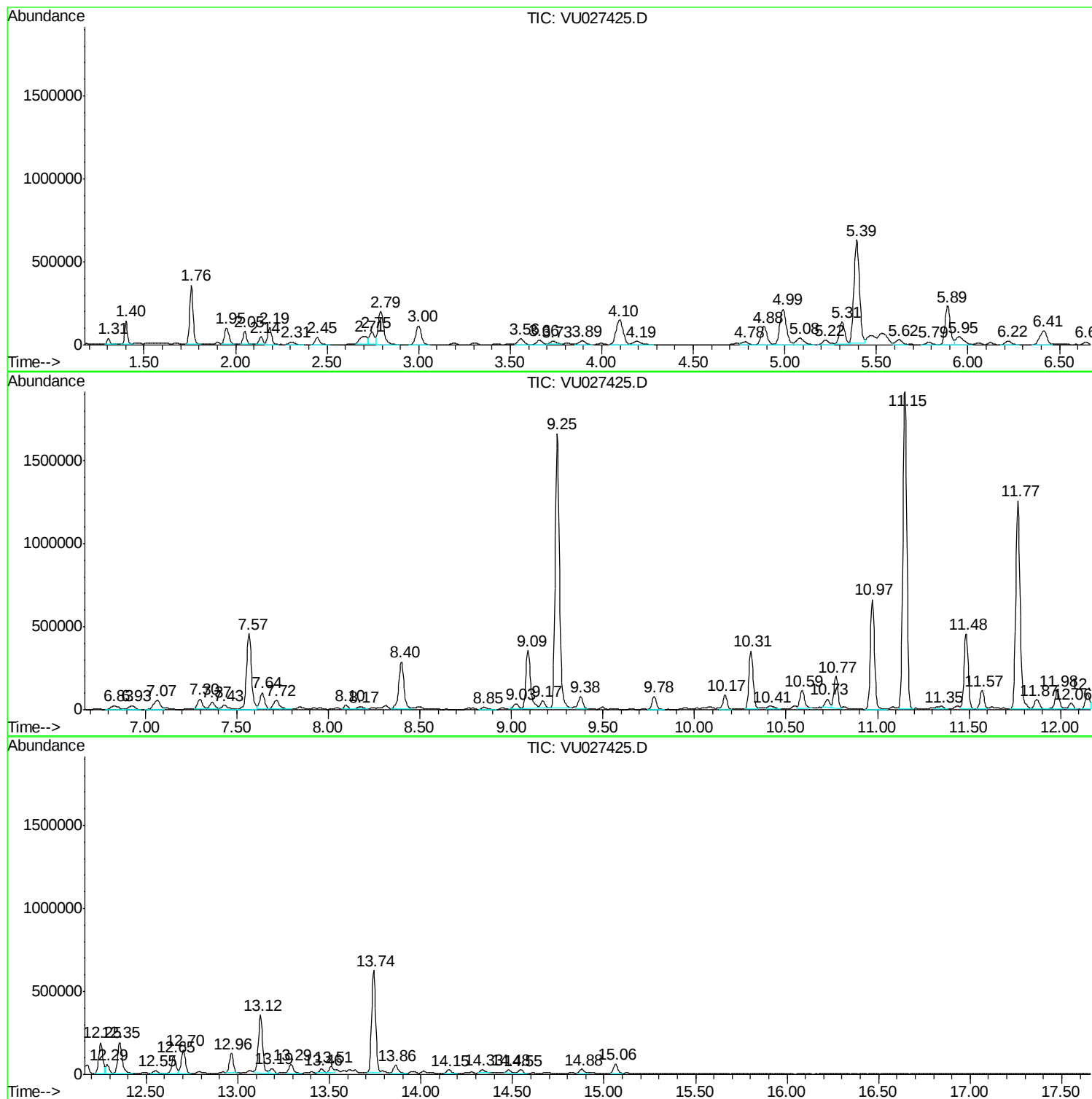
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

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



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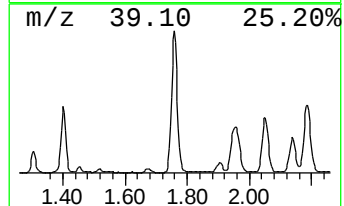
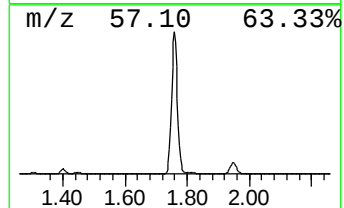
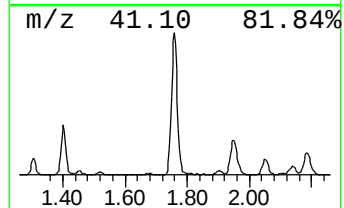
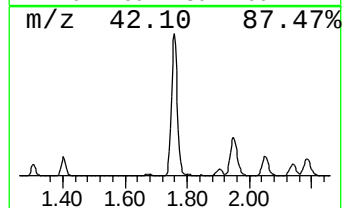
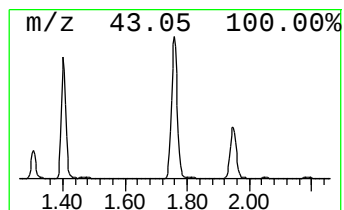
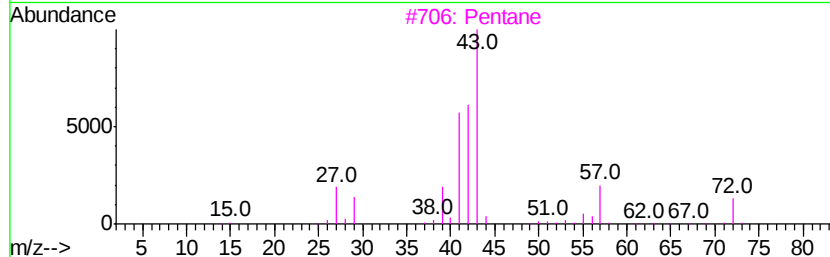
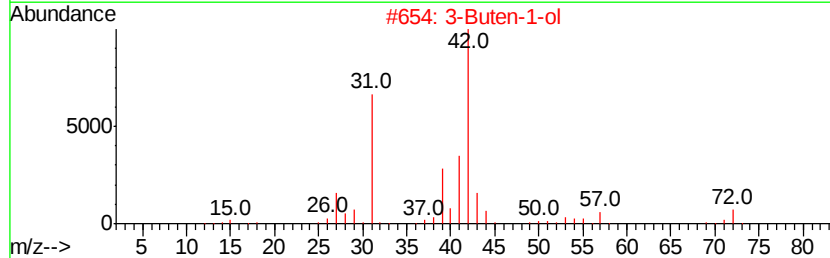
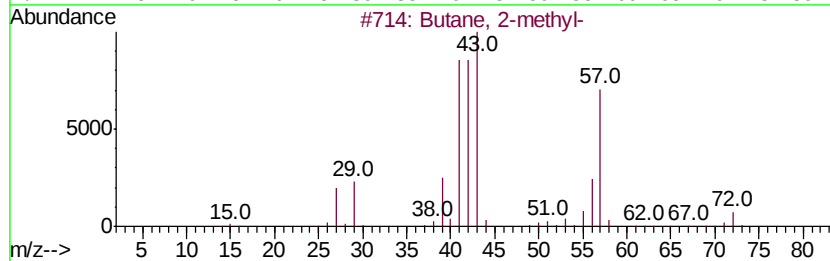
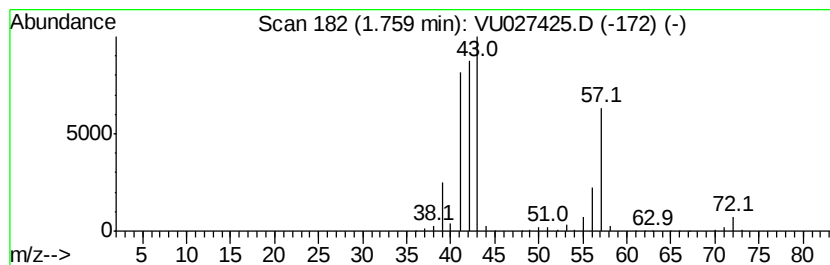
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.76	43.77 ug/l	467619	Pentafluorobenzene	4.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methyl-	72	C5H12	000078-78-4	94
2		3-Buten-1-ol	72	C4H8O	000627-27-0	47
3		Pentane	72	C5H12	000109-66-0	39
4		Methane, isocyanato-	57	C2H3NO	000624-83-9	9
5		1-Butene	56	C4H8	000106-98-9	9



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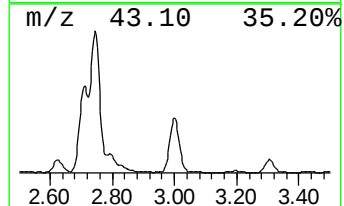
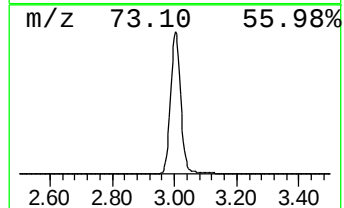
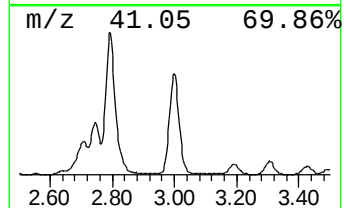
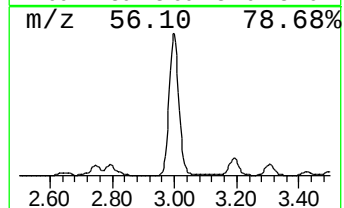
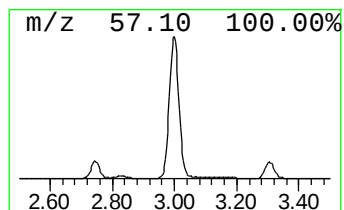
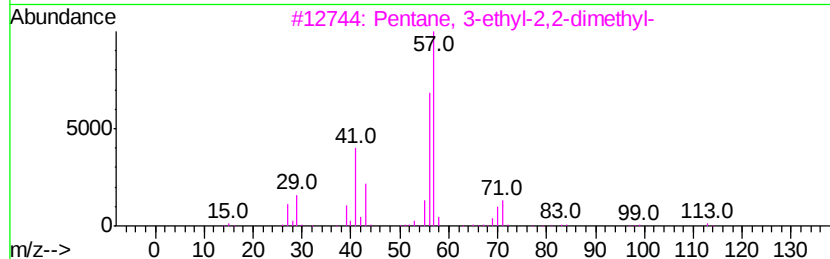
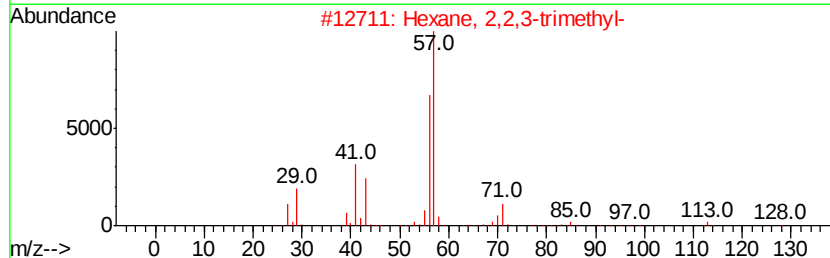
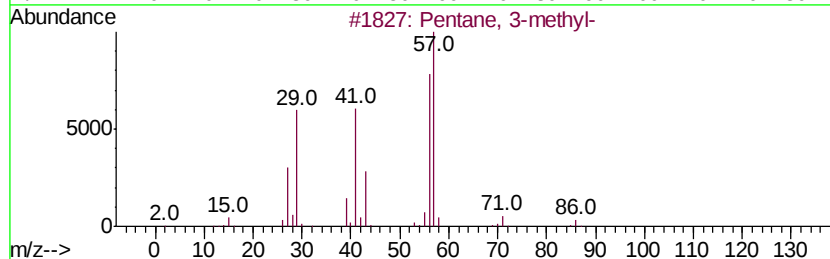
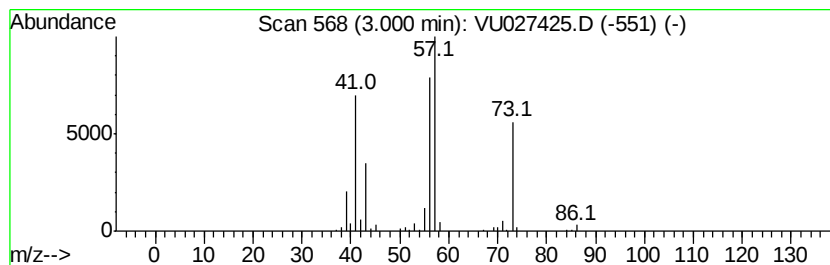
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Pentane, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.00	23.21 ug/l	247990	Pentafluorobenzene	4.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 3-methyl-	86	C6H14	000096-14-0	87
2		Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	72
3		Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	56
4		1-Cyclopentyl-2,2-dimethyl-1-pro...	156	C10H20O	337966-85-5	50
5		Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	39



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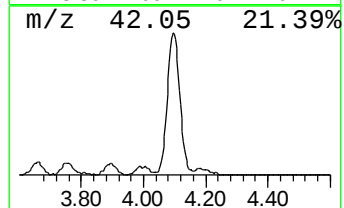
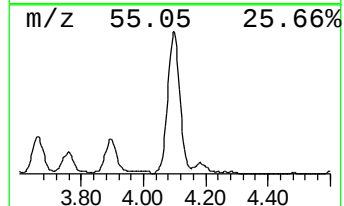
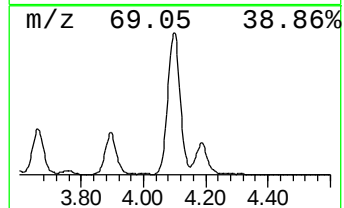
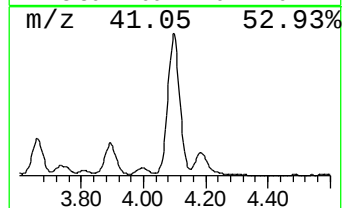
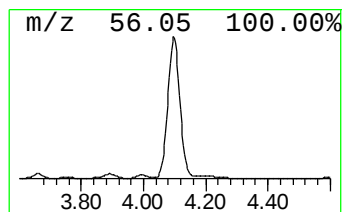
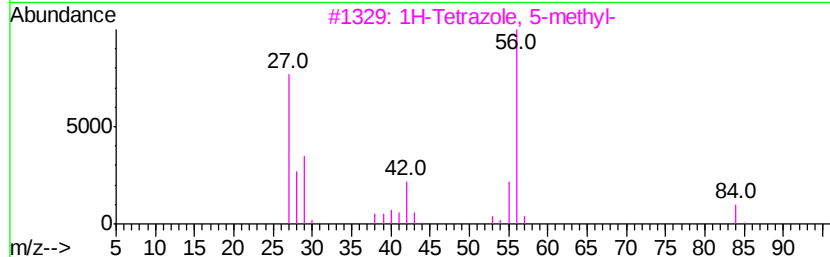
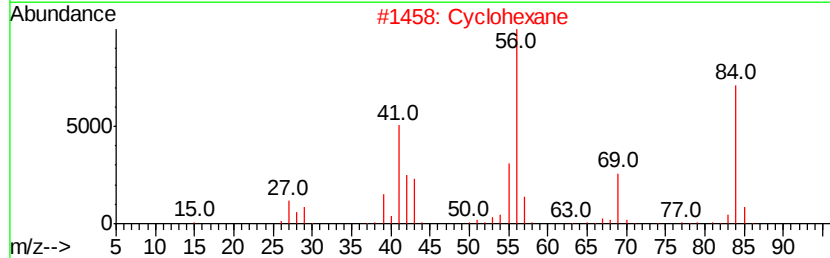
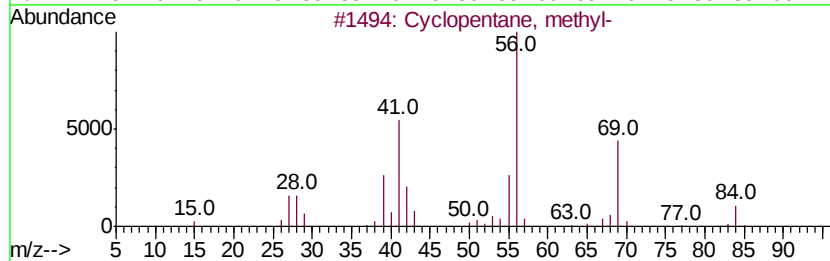
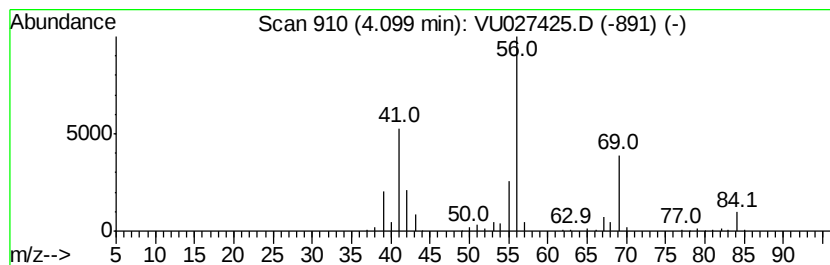
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Quant Title : SW846 8260

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

Peak Number 3 Cyclopentane, methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.10	37.64 ug/l	402116	Pentafluorobenzene	4.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2		Cyclohexane	84	C6H12	000110-82-7	78
3		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
4		Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
5		Aziridine, 1-(2-buten-2-yl)-	97	C6H11N	1000158-95-2	56



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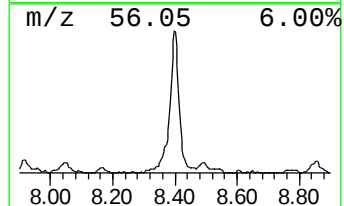
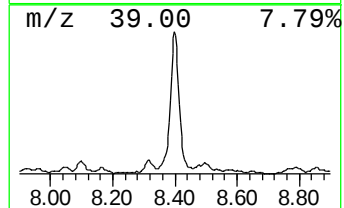
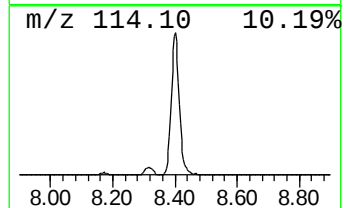
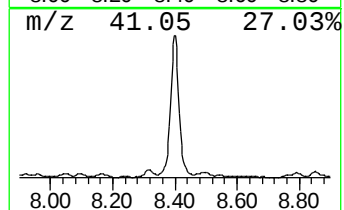
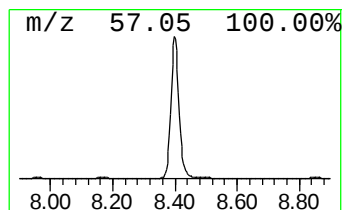
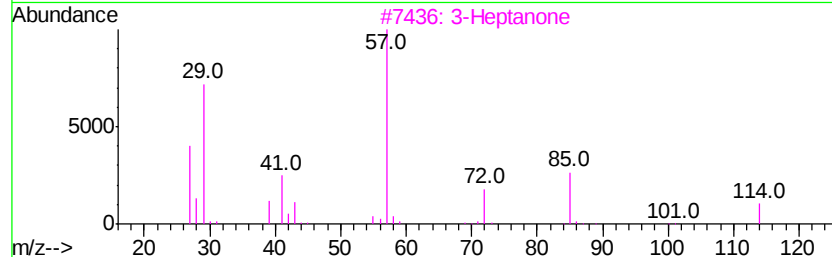
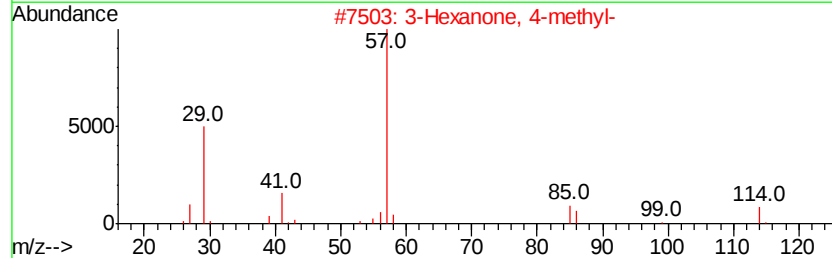
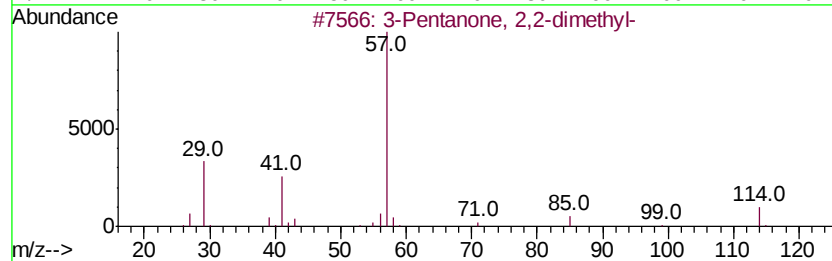
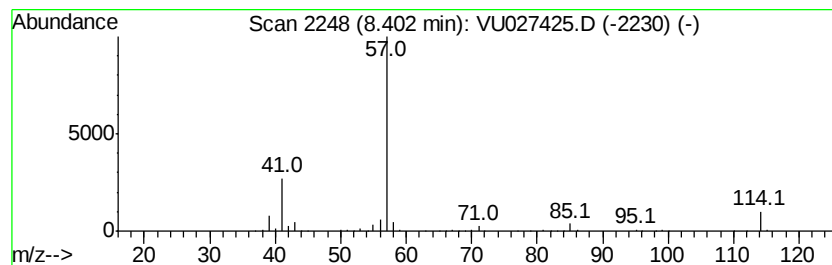
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Quant Title : SW846 8260

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

Peak Number 4 3-Pentanone, 2,2-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.40	42.61 ug/l	533636	Chlorobenzene-d5	9.09

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Pentanone, 2,2-dimethyl-	114	C7H14O	000564-04-5	87
2			3-Hexanone, 4-methyl-	114	C7H14O	017042-16-9	53
3			3-Heptanone	114	C7H14O	000106-35-4	53
4			3-Hexanone, 5-methyl-	114	C7H14O	000623-56-3	53
5			2,2,5,5-Tetramethyl-3-hexanone	156	C10H20O	000868-91-7	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100418\
Data File : VU027425.D
Acq On : 04 Oct 2018 21:57
Operator : MD/SY
Sample : J5232-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampled :
MW-3

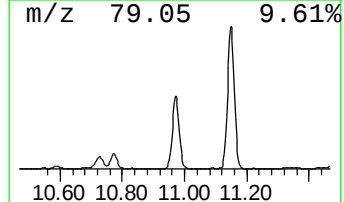
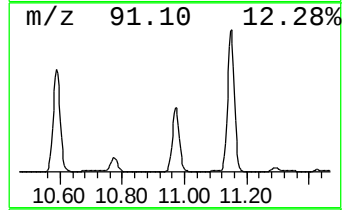
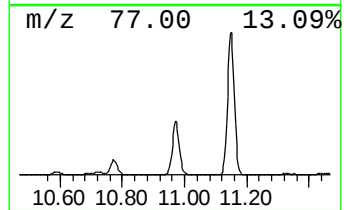
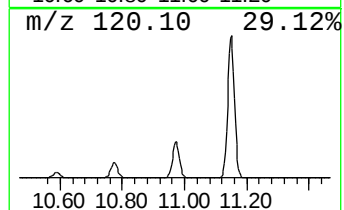
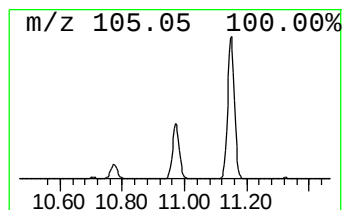
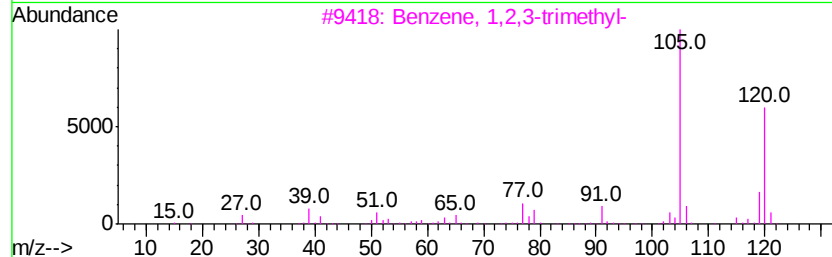
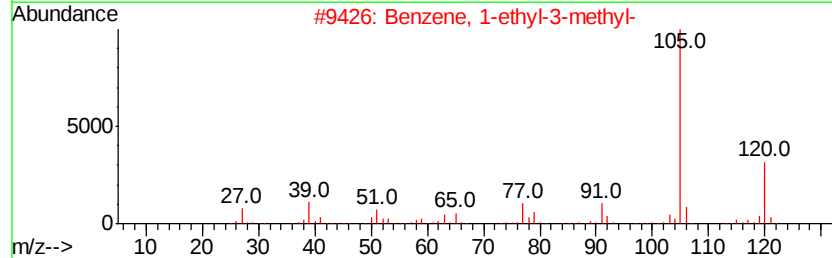
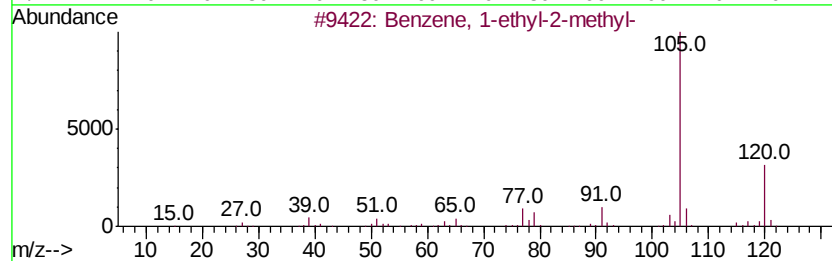
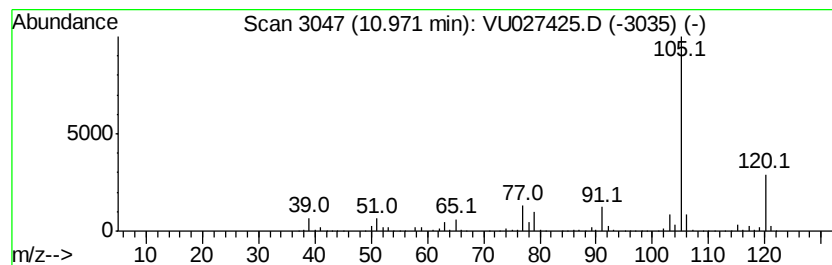
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.97	72.75 ug/l	1033530	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
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\VU100418\
Data File : VU027425.D
Acq On : 04 Oct 2018 21:57
Operator : MD/SY
Sample : J5232-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

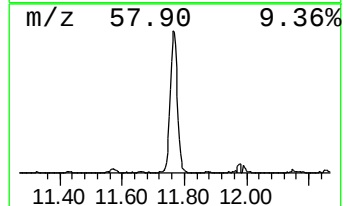
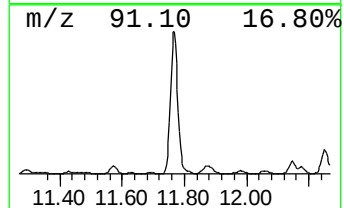
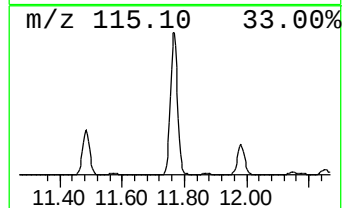
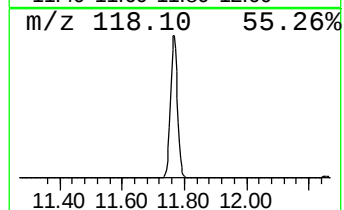
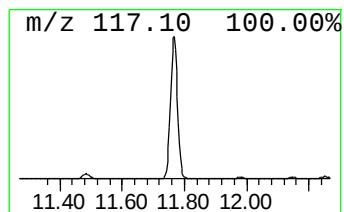
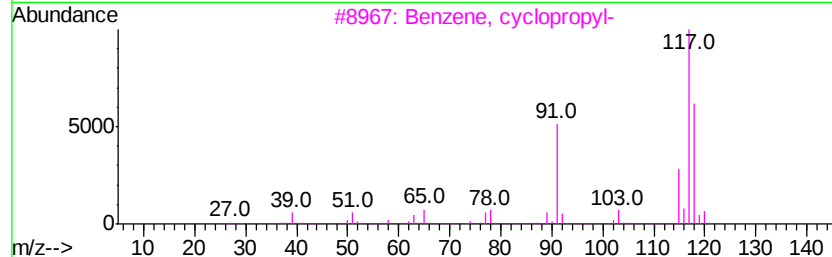
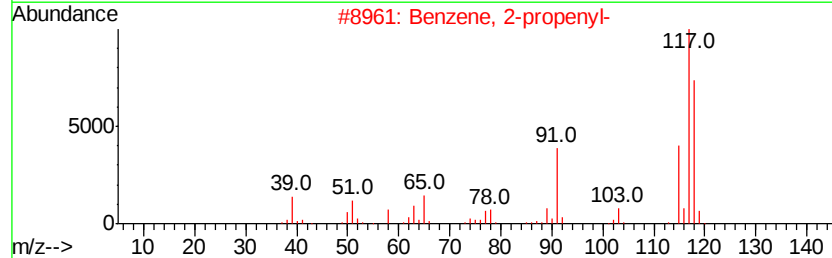
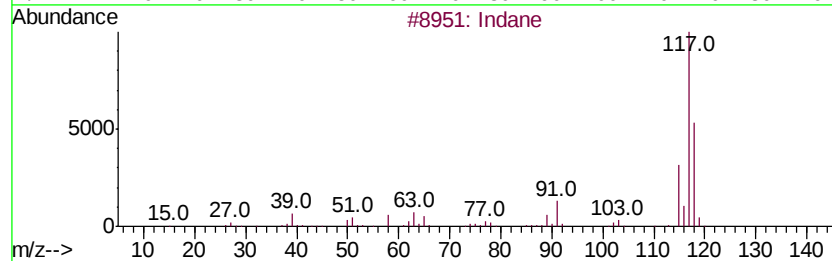
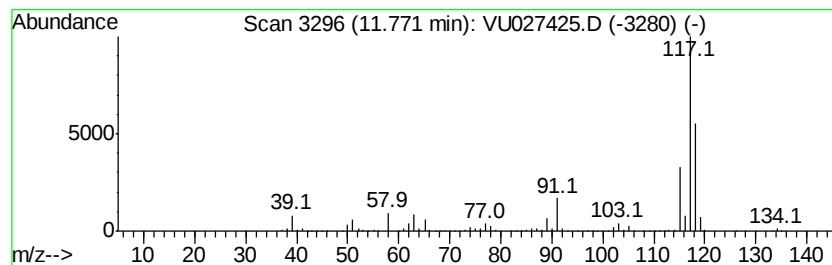
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Quant Title : SW846 8260

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

Peak Number 6 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	148.42 ug/l	2108470	1,4-Dichlorobenzene-d4	11.48

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	93
2	Benzene, 2-propenyl-		118	C9H10	000300-57-2	74
3	Benzene, cyclopropyl-		118	C9H10	000873-49-4	72
4	Deltacyclene		118	C9H10	007785-10-6	68
5	Benzene, 1-ethenyl-4-methyl-		118	C9H10	000622-97-9	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100418\
Data File : VU027425.D
Acq On : 04 Oct 2018 21:57
Operator : MD/SY
Sample : J5232-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

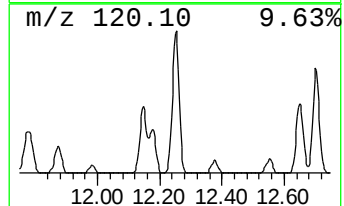
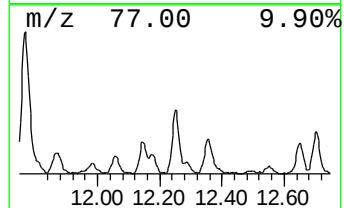
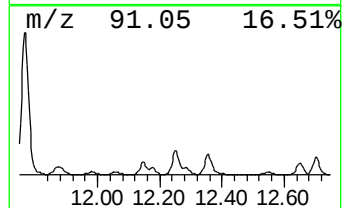
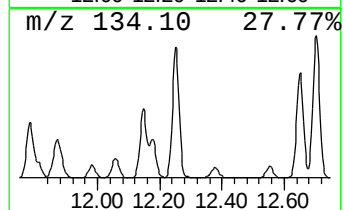
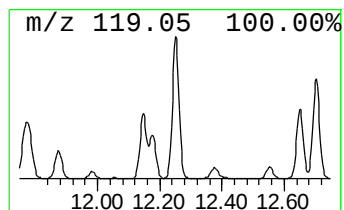
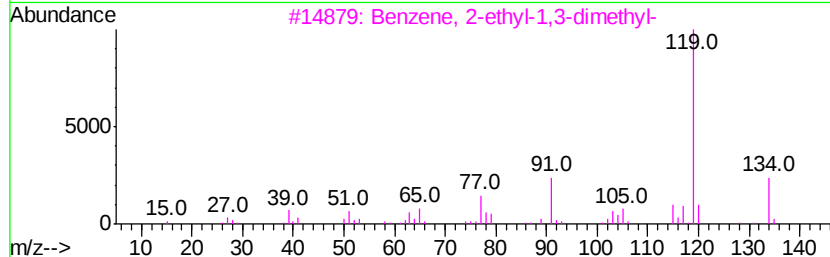
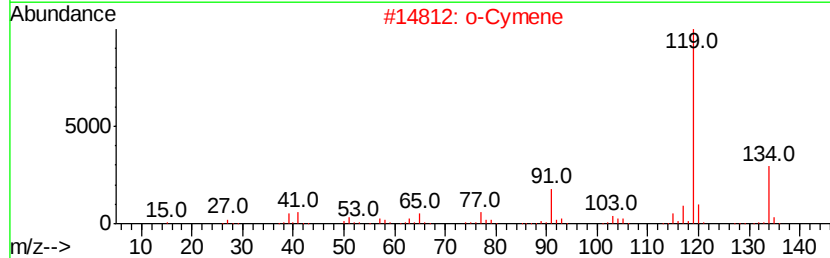
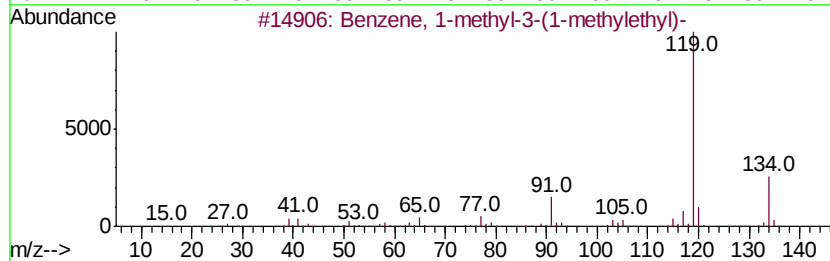
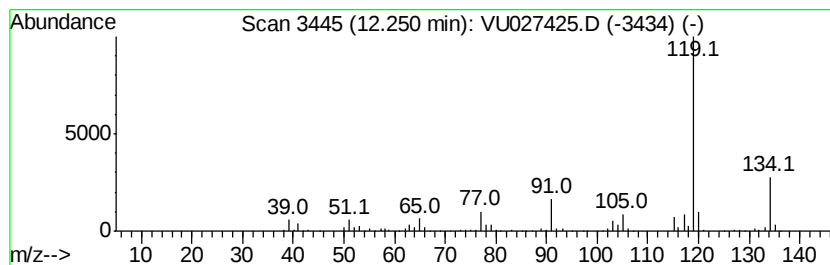
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-methyl-3-(1-meth... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.25	20.35 ug/l	289144	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
2		o-Cymene	134	C10H14	000527-84-4	95
3		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100418\
Data File : VU027425.D
Acq On : 04 Oct 2018 21:57
Operator : MD/SY
Sample : J5232-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

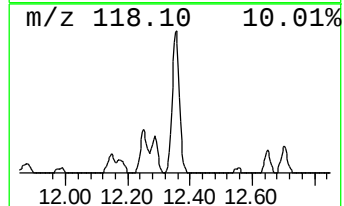
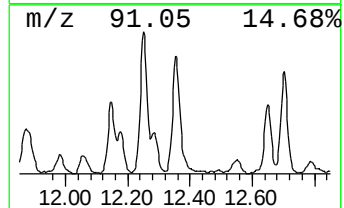
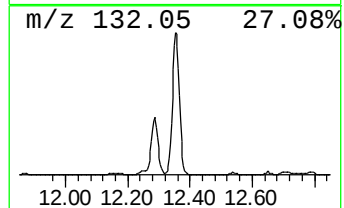
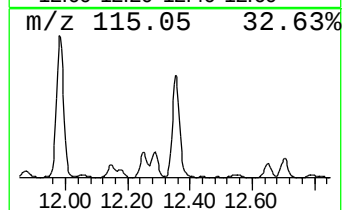
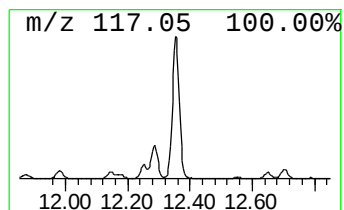
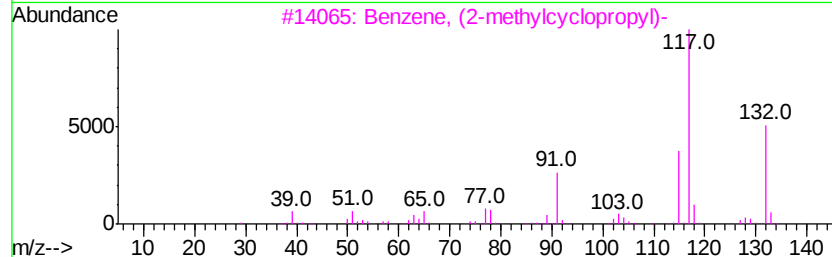
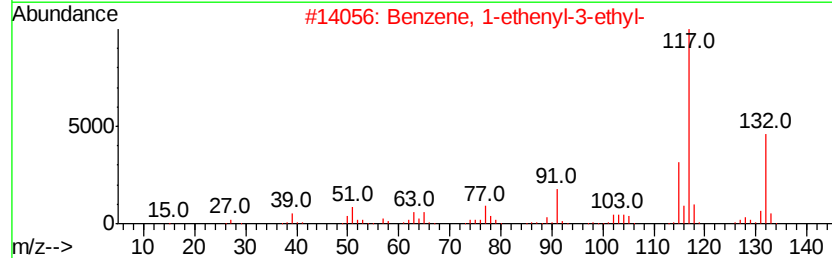
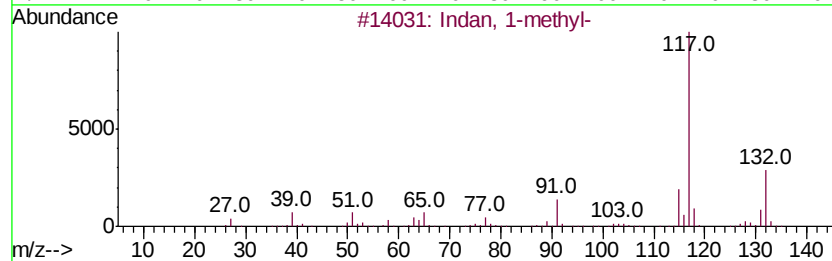
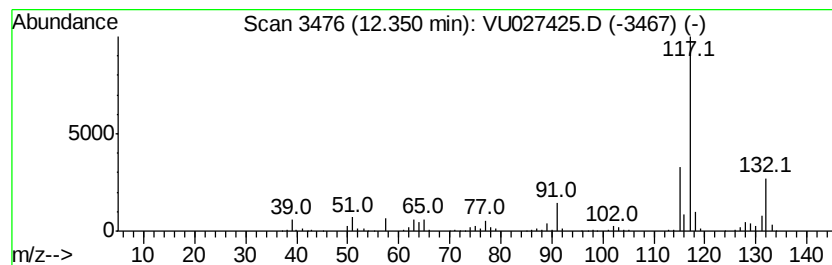
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.35	23.00 ug/l	326695	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indan, 1-methyl-	132	C10H12	000767-58-8	93
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
3		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	86
4		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	86
5		Benzene, (1-methyl-1-propenyl)-, ...	132	C10H12	000768-00-3	83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100418\
Data File : VU027425.D
Acq On : 04 Oct 2018 21:57
Operator : MD/SY
Sample : J5232-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

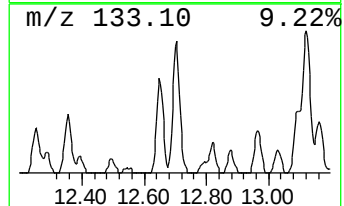
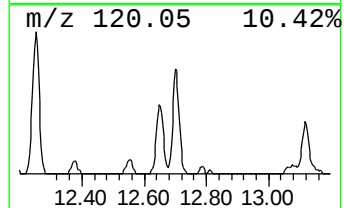
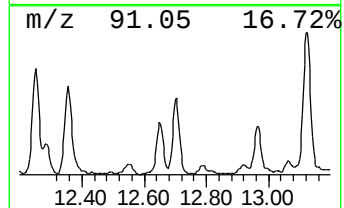
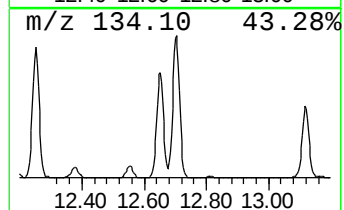
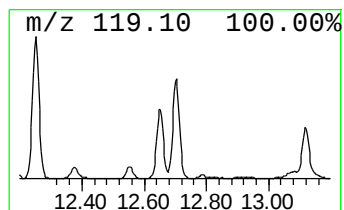
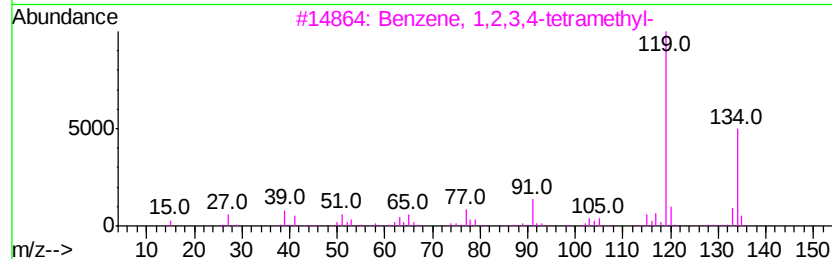
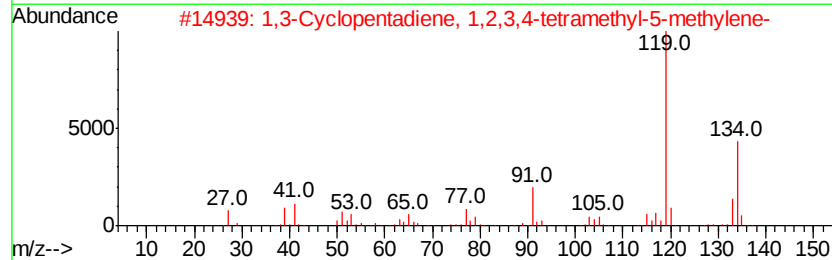
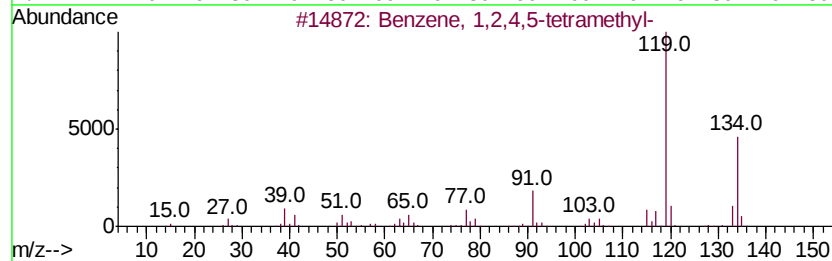
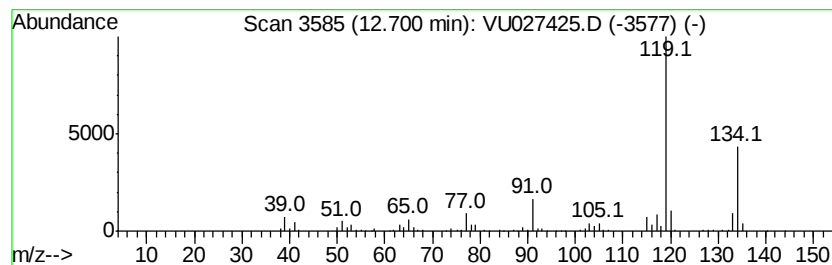
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	15.79 ug/l	224255	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		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	95
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100418\
Data File : VU027425.D
Acq On : 04 Oct 2018 21:57
Operator : MD/SY
Sample : J5232-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

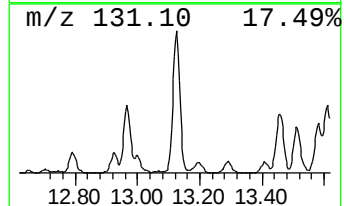
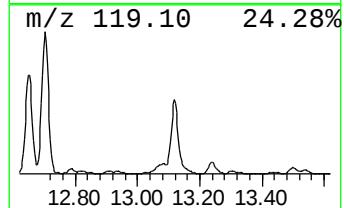
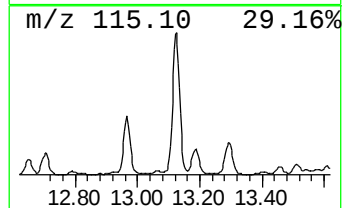
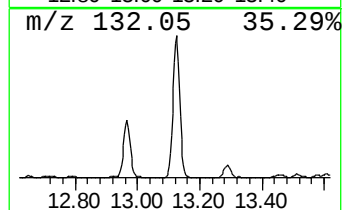
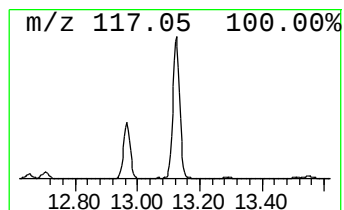
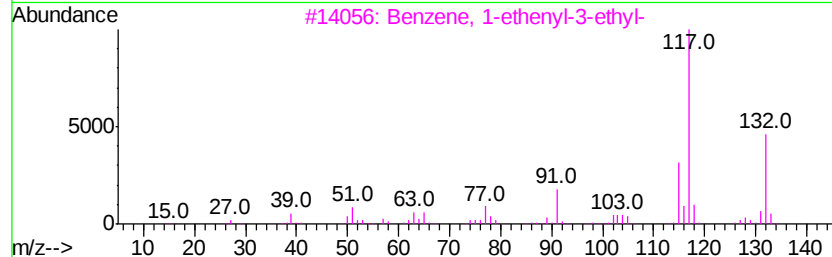
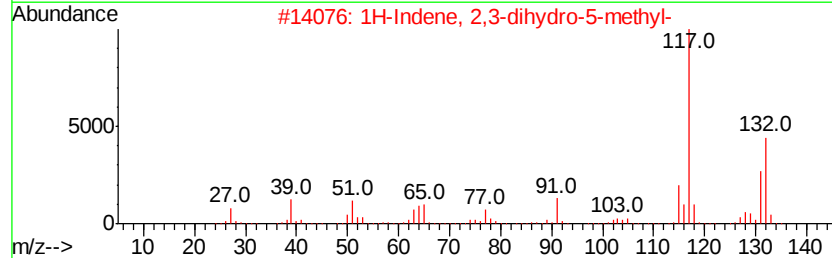
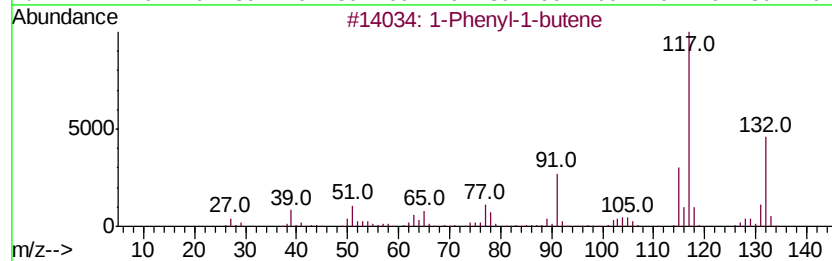
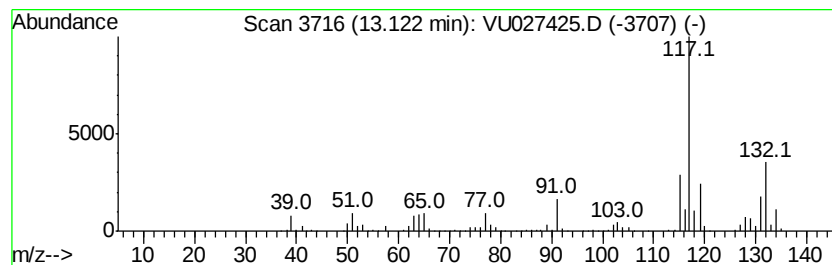
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Quant Title : SW846 8260

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

Peak Number 10 1-Phenyl-1-butene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	39.72 ug/l	564304	1,4-Dichlorobenzene-d4	11.48

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Phenyl-1-butene	132	C10H12	000824-90-8	90
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	87
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
4		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	81
5		2,4-Dimethylstyrene	132	C10H12	002234-20-0	78



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU100418\
Data File : VU027425.D
Acq On : 04 Oct 2018 21:57
Operator : MD/SY
Sample : J5232-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

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
Butane, 2-methyl-	1.76	43.8	ug/l	467619	1	4.99	534159	50.0
Pentane, 3-methyl-	3.00	23.2	ug/l	247990	1	4.99	534159	50.0
Cyclopentane, met...	4.10	37.6	ug/l	402116	1	4.99	534159	50.0
3-Pentanone, 2,2-...	8.40	42.6	ug/l	533636	3	9.09	626142	50.0
Benzene, 1-ethyl-...	10.97	72.8	ug/l	1033530	4	11.48	710282	50.0
Indane	11.77	148.4	ug/l	2108470	4	11.48	710282	50.0
Benzene, 1-methyl...	12.25	20.4	ug/l	289144	4	11.48	710282	50.0
Indan, 1-methyl-	12.35	23.0	ug/l	326695	4	11.48	710282	50.0
Benzene, 1,2,4,5-...	12.70	15.8	ug/l	224255	4	11.48	710282	50.0
1-Phenyl-1-butene	13.12	39.7	ug/l	564304	4	11.48	710282	50.0