

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU080618\
 Data File : VU025916.D
 Acq On : 07 Aug 2018 03:51
 Operator : MD/SY
 Sample : J4344-02
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 FL-0562

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	138	155	170	rBV	480947	533439	11.86%	2.015%
2	1.307	216	223	236	rVB	54388	58171	1.29%	0.220%
3	1.404	245	253	263	rBV	729497	749252	16.66%	2.831%
4	1.760	354	364	382	rBV	728572	947861	21.08%	3.581%
5	1.950	413	423	440	rVB	342575	458179	10.19%	1.731%
6	2.304	521	533	547	rVB3	29476	60123	1.34%	0.227%
7	2.712	647	660	664	rBV2	32928	62851	1.40%	0.237%
8	2.748	665	671	677	rVV	62295	112713	2.51%	0.426%
9	2.796	677	686	707	rVB	188353	359332	7.99%	1.358%
10	3.002	732	750	767	rBV	69577	148363	3.30%	0.561%
11	4.101	1072	1092	1115	rVB	206863	545619	12.14%	2.062%
12	4.889	1319	1337	1354	rBV	195806	444194	9.88%	1.678%
13	4.992	1354	1369	1388	rVV2	348930	849244	18.89%	3.209%
14	5.088	1389	1399	1415	rVB5	18888	46938	1.04%	0.177%
15	5.227	1426	1442	1450	rBV2	25146	63119	1.40%	0.238%
16	5.313	1458	1469	1489	rVB	191050	402579	8.95%	1.521%
17	5.487	1508	1523	1534	rBV4	25503	58823	1.31%	0.222%
18	5.558	1535	1545	1556	rVV4	21939	49987	1.11%	0.189%
19	5.628	1556	1567	1582	rVB5	20422	47201	1.05%	0.178%
20	5.892	1633	1649	1674	rBV	379378	751490	16.71%	2.839%
21	6.419	1794	1813	1826	rBV6	63006	157566	3.50%	0.595%
22	7.571	2148	2171	2189	rBV	708078	1276924	28.40%	4.825%
23	9.095	2632	2645	2666	rBV	573948	950983	21.15%	3.593%
24	9.252	2682	2694	2719	rBV	1140590	1818073	40.44%	6.869%
25	9.378	2719	2733	2763	rVB	2713298	4496036	100.00%	16.987%
26	10.169	2968	2979	2992	rBV	200020	310075	6.90%	1.172%
27	10.310	3011	3023	3039	rBV	596110	952656	21.19%	3.599%
28	10.590	3098	3110	3125	rBV	214847	331030	7.36%	1.251%
29	10.683	3127	3139	3157	rVV	609034	1297028	28.85%	4.901%
30	10.776	3157	3168	3188	rVB	474708	729721	16.23%	2.757%
31	10.976	3217	3230	3249	rVB	350559	555788	12.36%	2.100%
32	11.152	3273	3285	3299	rVB	1387374	2084374	46.36%	7.875%
33	11.432	3357	3372	3378	rBV	40534	60790	1.35%	0.230%
34	11.487	3378	3389	3404	rVV	797986	1250815	27.82%	4.726%

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35	11.574	3404	3416	3439	rVB	570897	869462	19.34%	3.285%
36	11.770	3465	3477	3486	rBV	160852	281255	6.26%	1.063%
37	11.812	3486	3490	3499	rVV2	49085	69766	1.55%	0.264%
38	11.876	3499	3510	3526	rVB6	55660	124111	2.76%	0.469%
39	12.149	3578	3595	3599	rBV	40940	60852	1.35%	0.230%
40	12.178	3600	3604	3614	rVV	49608	70601	1.57%	0.267%
41	12.255	3616	3628	3635	rVV	57591	92569	2.06%	0.350%
42	12.361	3650	3661	3689	rVB3	43641	130239	2.90%	0.492%
43	12.554	3711	3721	3731	rVB3	35605	64077	1.43%	0.242%
44	12.654	3731	3752	3760	rBV	50871	91686	2.04%	0.346%
45	12.705	3760	3768	3785	rVB	78656	122581	2.73%	0.463%
46	13.123	3889	3898	3913	rVB3	135635	217316	4.83%	0.821%
47	13.291	3941	3950	3961	rVB	70029	108227	2.41%	0.409%
48	13.744	4079	4091	4118	rVB	473309	784767	17.45%	2.965%
49	14.352	4266	4280	4292	rBV7	20397	47099	1.05%	0.178%
50	14.882	4434	4445	4456	rBV	97177	171349	3.81%	0.647%
51	15.065	4492	4502	4517	rBV	81605	139744	3.11%	0.528%

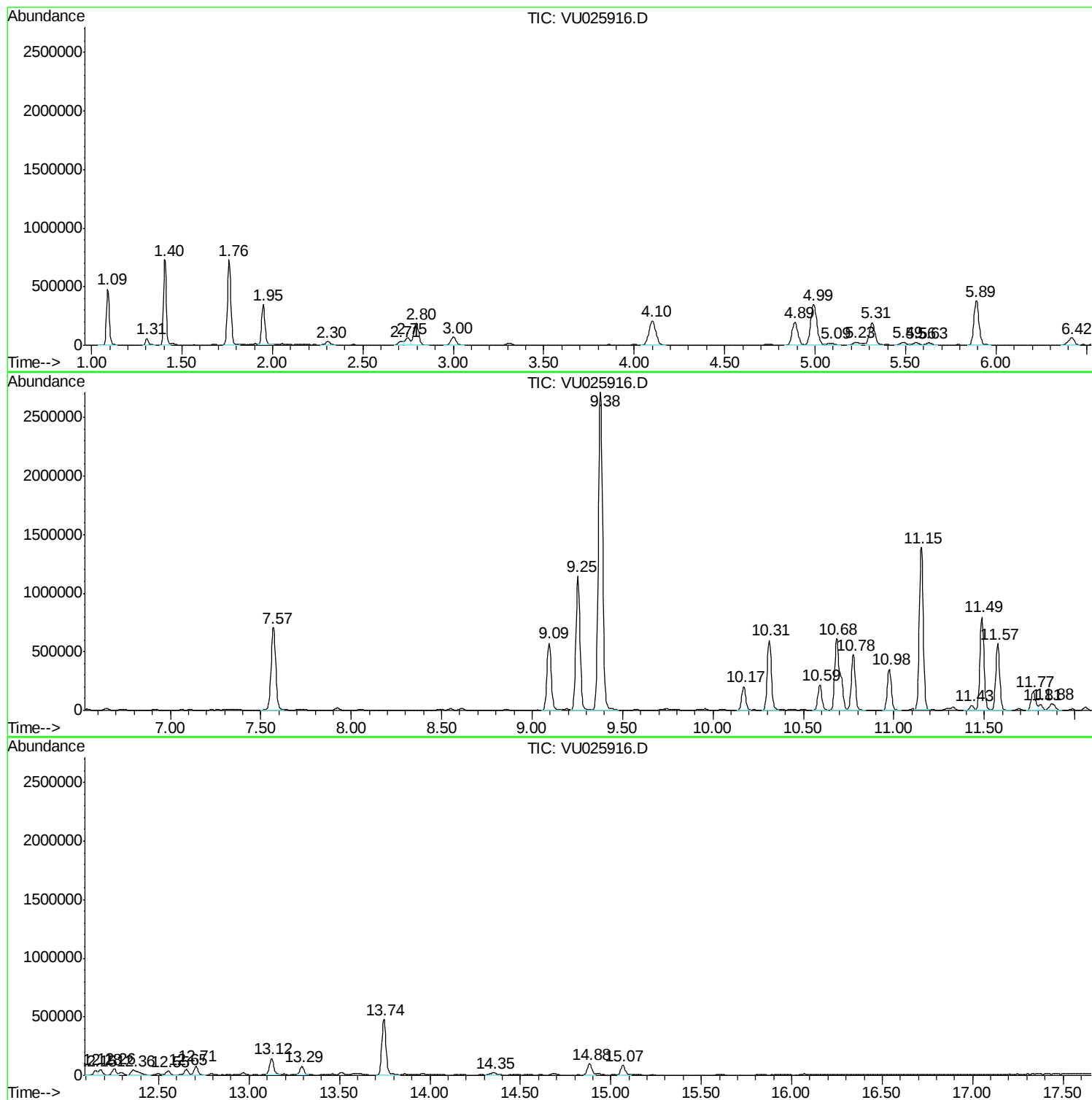
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.M
Quant Title : SW846 8260

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



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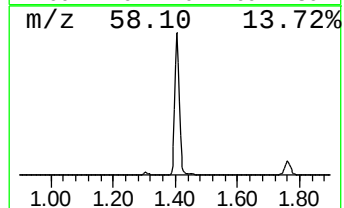
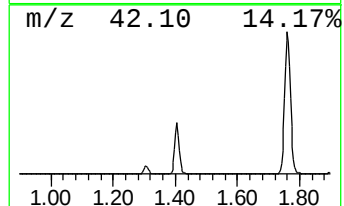
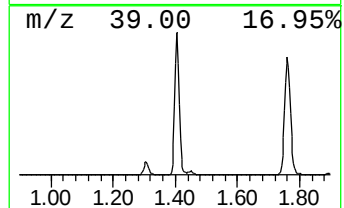
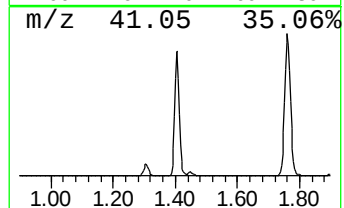
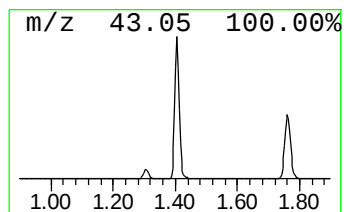
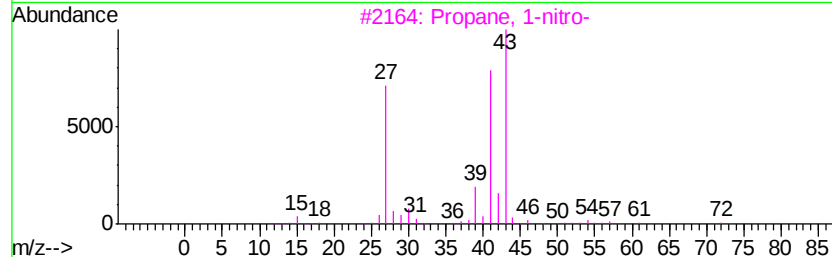
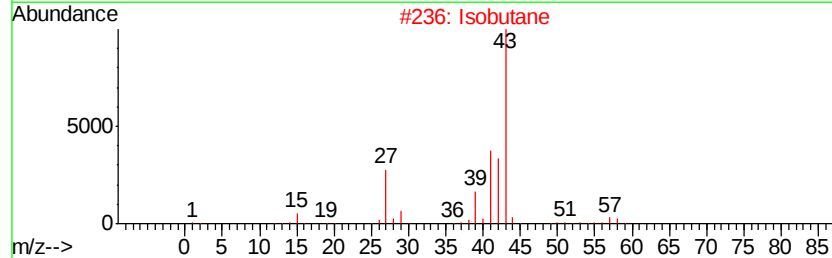
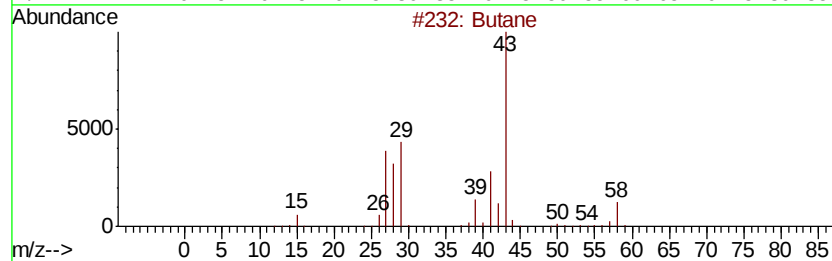
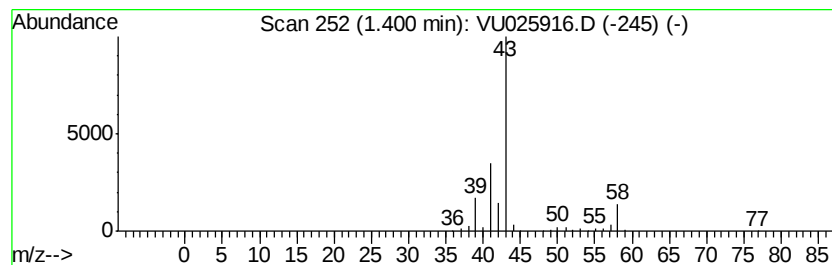
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.M
Quant Title : SW846 8260

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

Peak Number 2 Butane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.40	44.11 ug/l	749252	Pentafluorobenzene	4.99

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane		58	C4H10	000106-97-8	80
2	Isobutane		58	C4H10	000075-28-5	9
3	Propane, 1-nitro-		89	C3H7NO2	000108-03-2	9
4	Isopropylsulfonfyl chloride		142	C3H7ClO2S	010147-37-2	9
5	Diazene, dimethyl-		58	C2H6N2	000503-28-6	4



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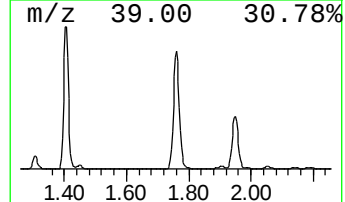
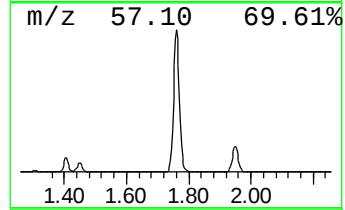
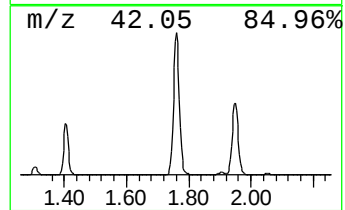
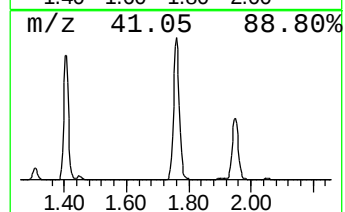
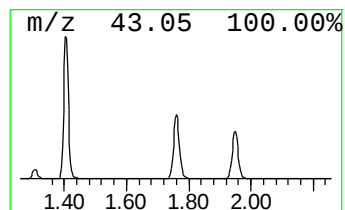
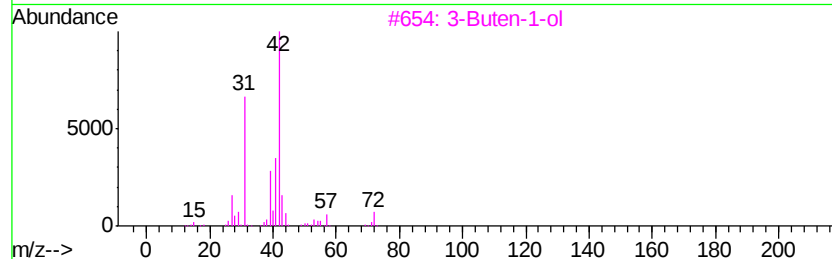
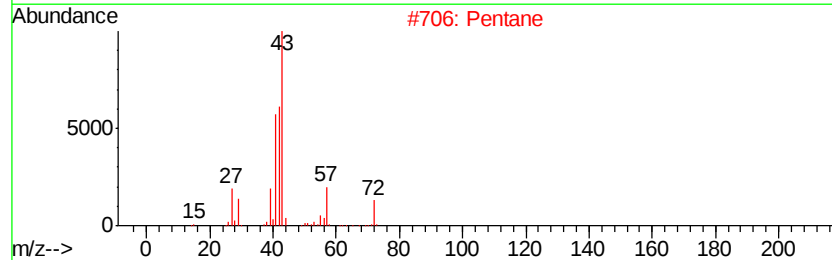
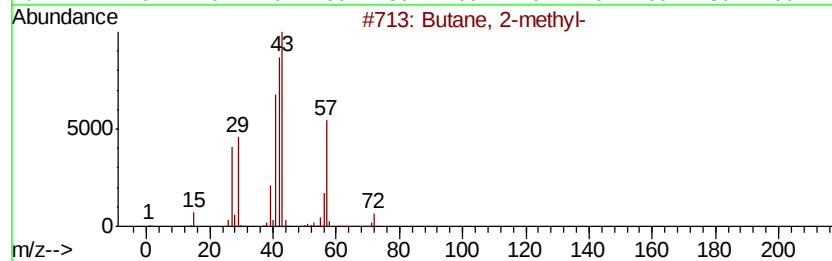
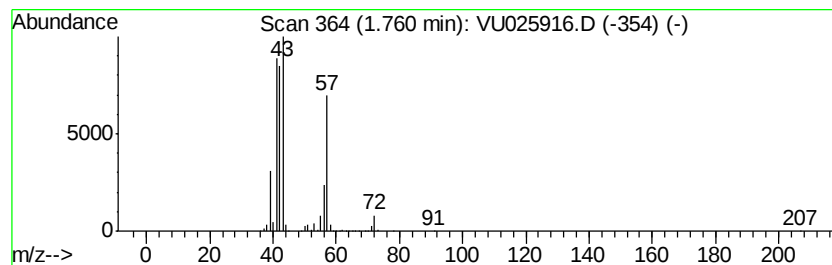
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.76	55.81 ug/l	947861	Pentafluorobenzene	4.99

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methyl-	72	C5H12	000078-78-4	91
2		Pentane	72	C5H12	000109-66-0	36
3		3-Buten-1-ol	72	C4H8O	000627-27-0	28
4		Butane, 1-chloro-2-methyl-	106	C5H11Cl	000616-13-7	16
5		1-Butene	56	C4H8	000106-98-9	9



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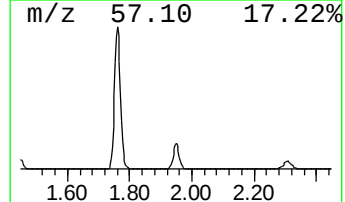
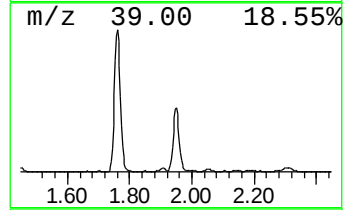
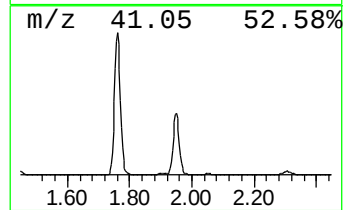
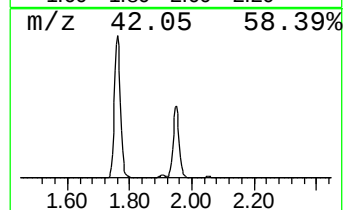
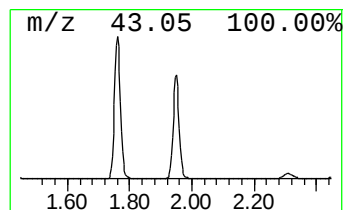
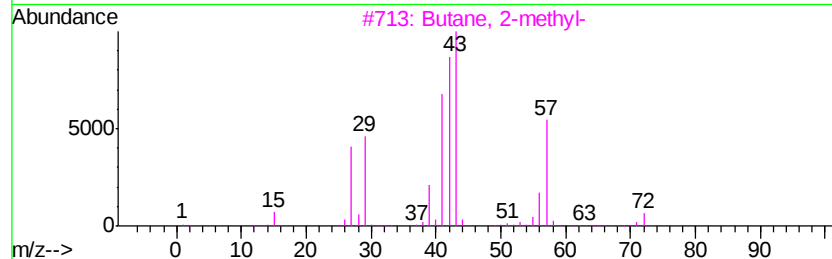
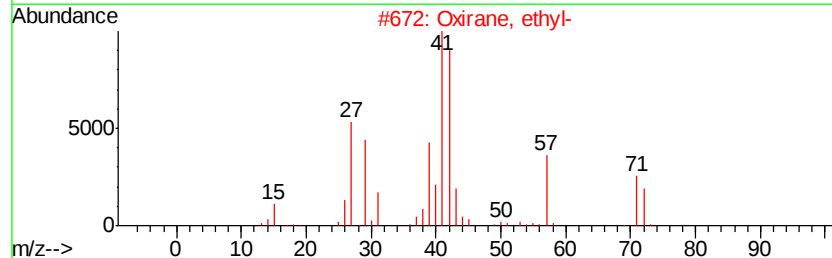
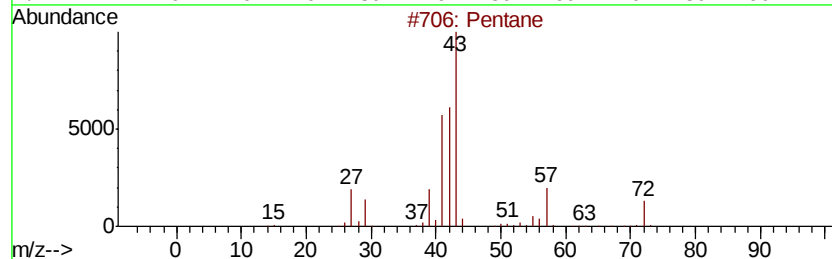
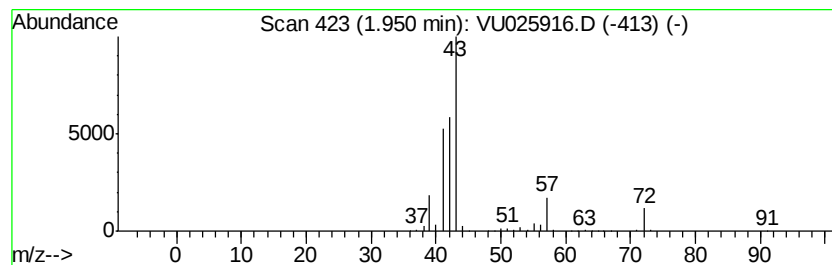
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Pentane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.95	26.98 ug/l	458179	Pentafluorobenzene	4.99

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane		72	C5H12	000109-66-0	91
2	Oxirane, ethyl-		72	C4H8O	000106-88-7	50
3	Butane, 2-methyl-		72	C5H12	000078-78-4	38
4	Diaziridine,3,3-dimethyl-		72	C3H8N2	004901-76-2	25
5	Isobutyl nitrite		103	C4H9NO2	000542-56-3	9



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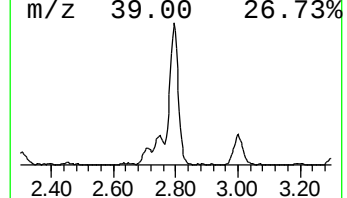
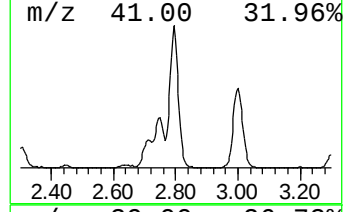
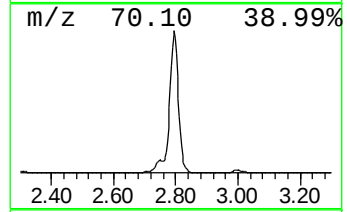
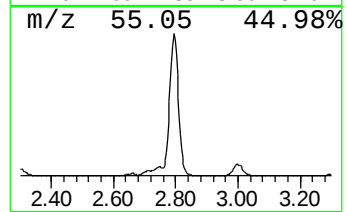
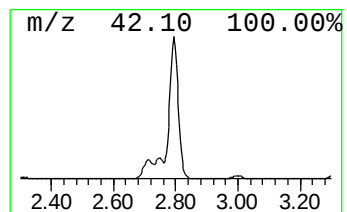
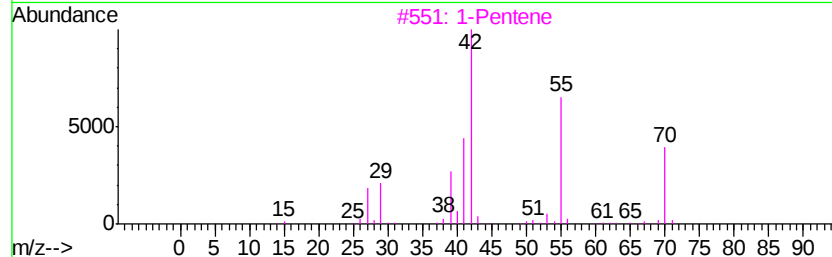
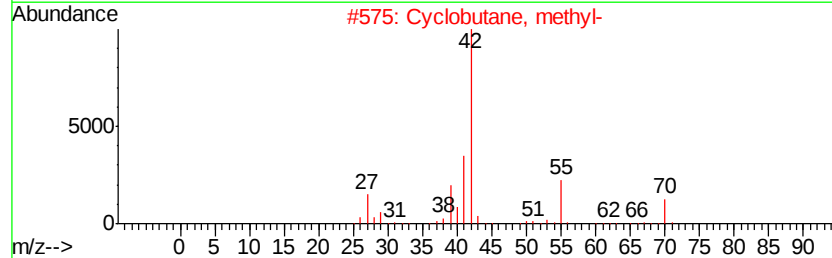
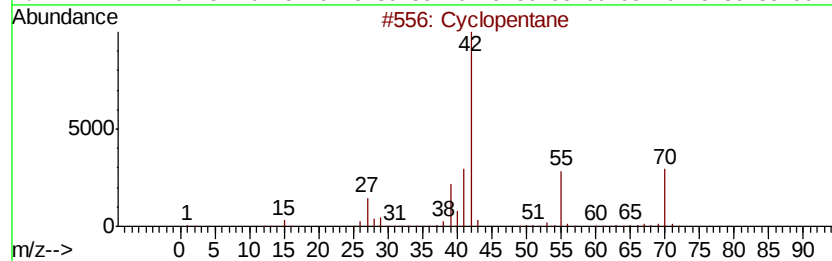
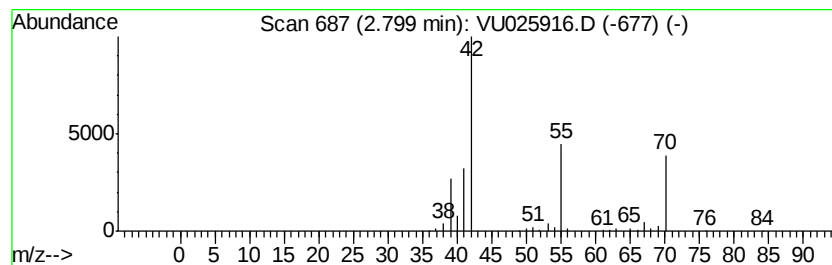
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Cyclopentane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.80	21.16 ug/l	359332	Pentafluorobenzene	4.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane	70	C5H10	000287-92-3	86
2		Cyclobutane, methyl-	70	C5H10	000598-61-8	78
3		1-Pentene	70	C5H10	000109-67-1	72
4		Cyclopropane, ethyl-	70	C5H10	001191-96-4	9
5		Methyl propargyl ether	70	C4H6O	000627-41-8	7



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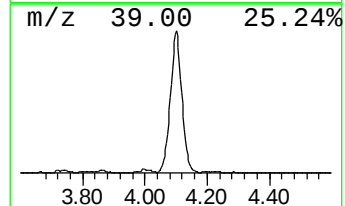
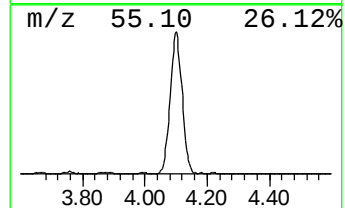
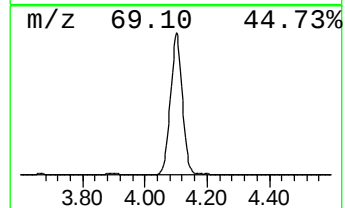
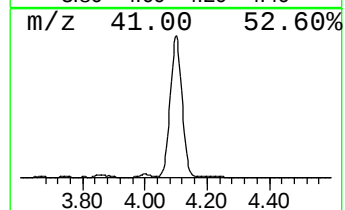
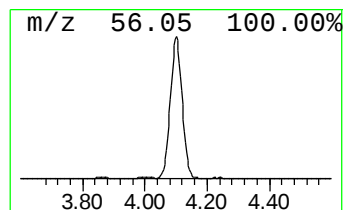
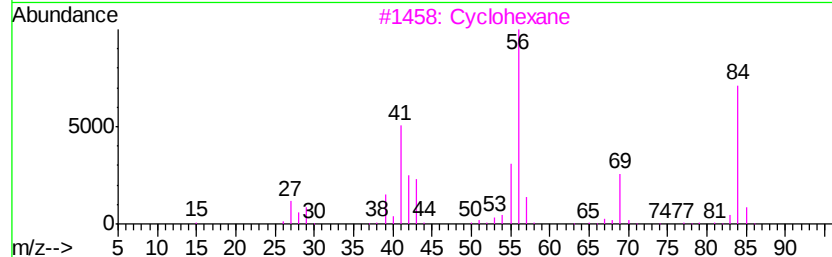
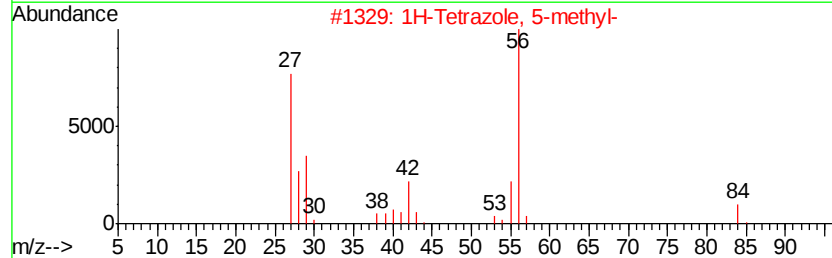
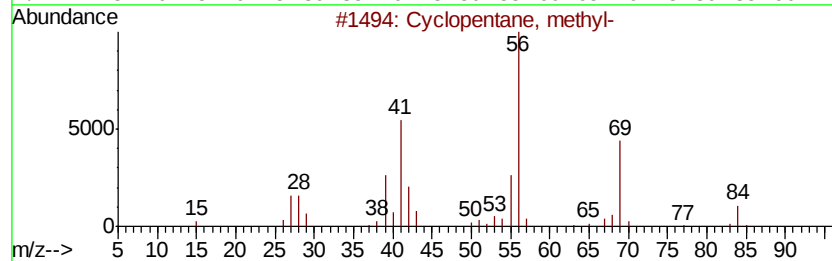
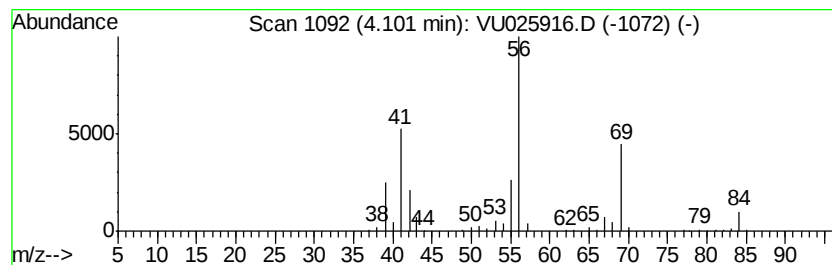
Instrument :
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ClientSampleId :
FL-0562

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 Cyclopentane, methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.		
4.10	32.12 ug/l	545619	Pentafluorobenzene	4.99		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cvclopentane, methyl-	84	C6H12	000096-37-7	91
2		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	80
3		Cvclohexane	84	C6H12	000110-82-7	64
4		Cvclobutane, ethyl-	84	C6H12	004806-61-5	64
5		Cyclobutane	56	C4H8	000287-23-0	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU080618\
Data File : VU025916.D
Acq On : 07 Aug 2018 03:51
Operator : MD/SY
Sample : J4344-02
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
FL-0562

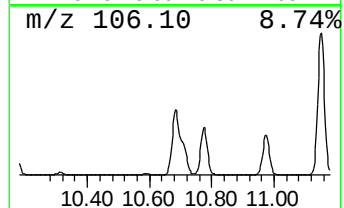
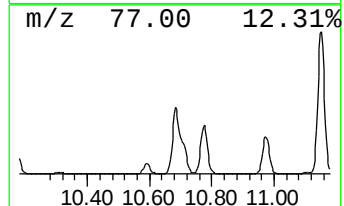
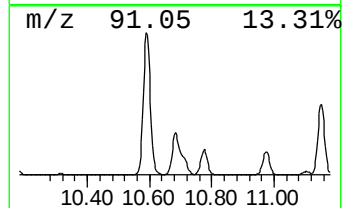
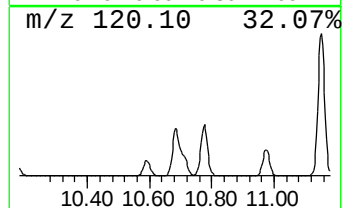
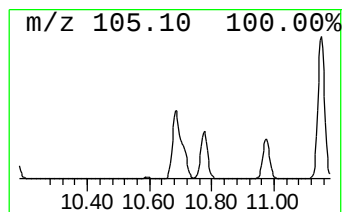
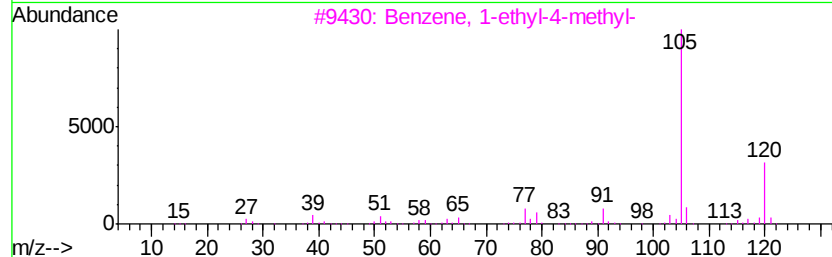
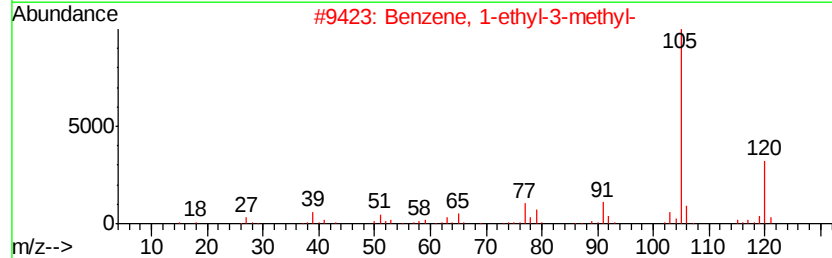
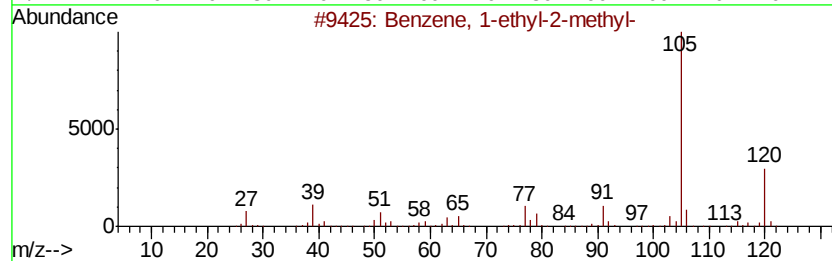
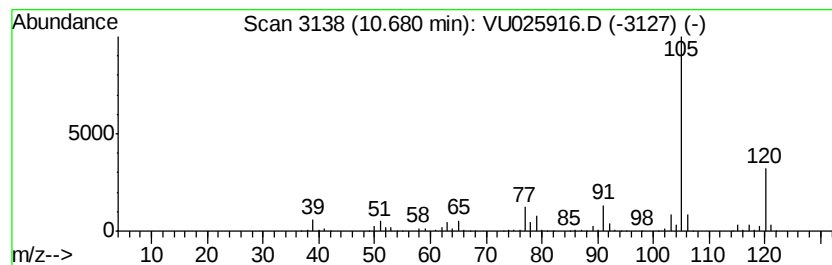
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.68	51.85 ug/l	1297030	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-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU080618\
Data File : VU025916.D
Acq On : 07 Aug 2018 03:51
Operator : MD/SY
Sample : J4344-02
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
FL-0562

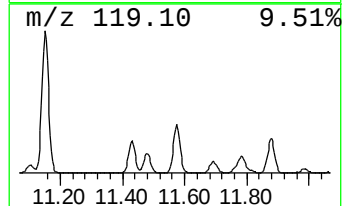
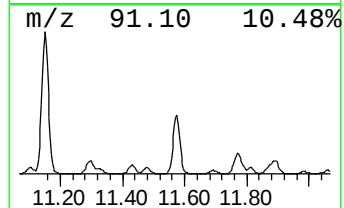
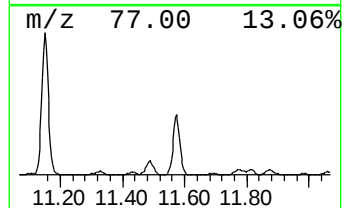
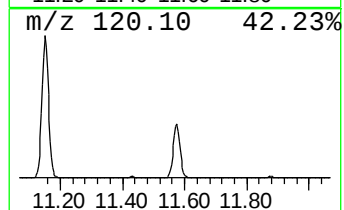
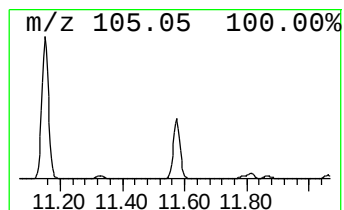
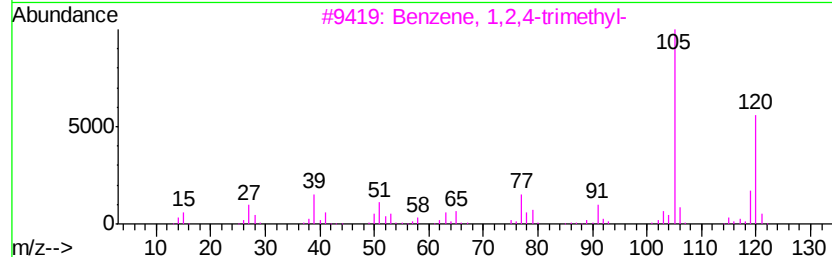
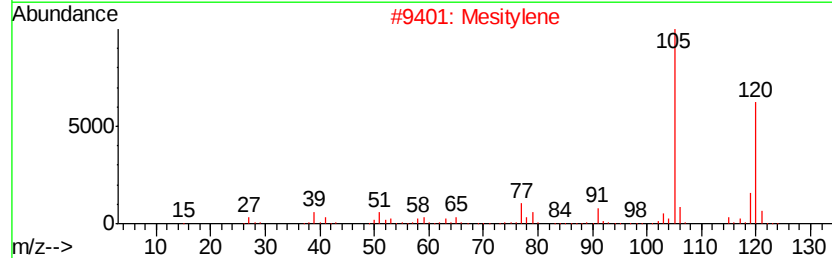
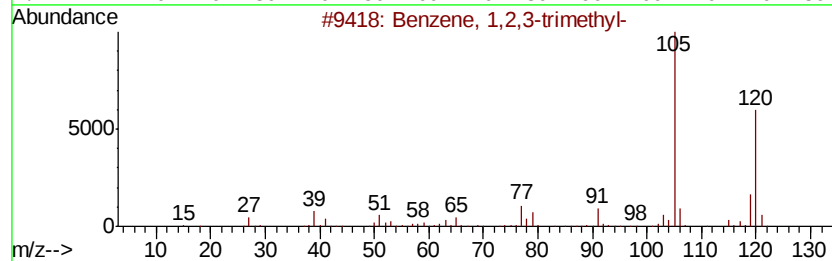
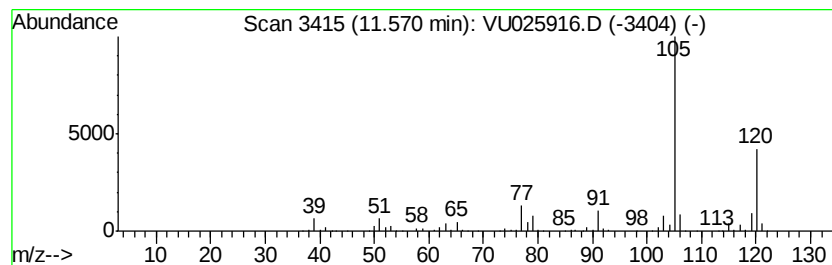
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,2,3-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	34.76 ug/l	869462	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	91
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU080618\
Data File : VU025916.D
Acq On : 07 Aug 2018 03:51
Operator : MD/SY
Sample : J4344-02
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
FL-0562

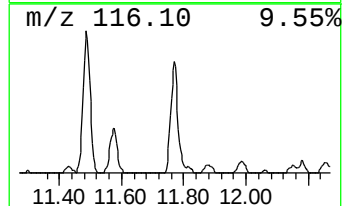
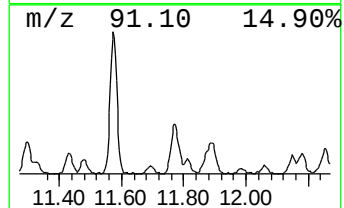
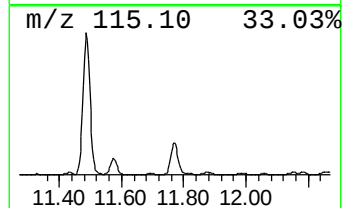
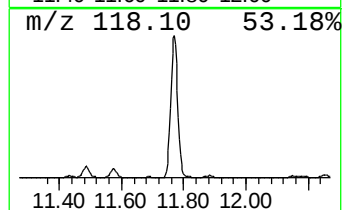
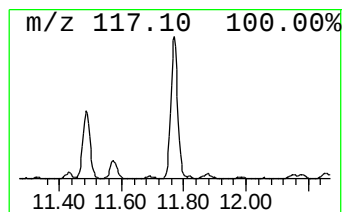
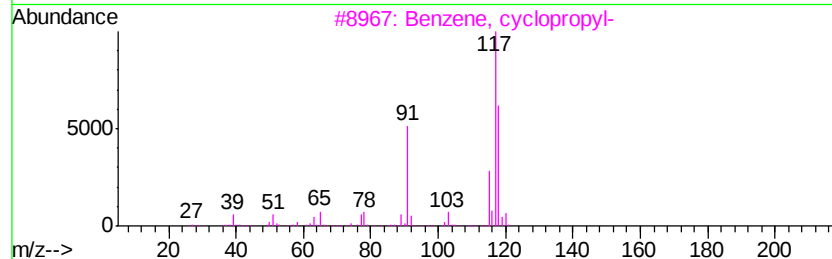
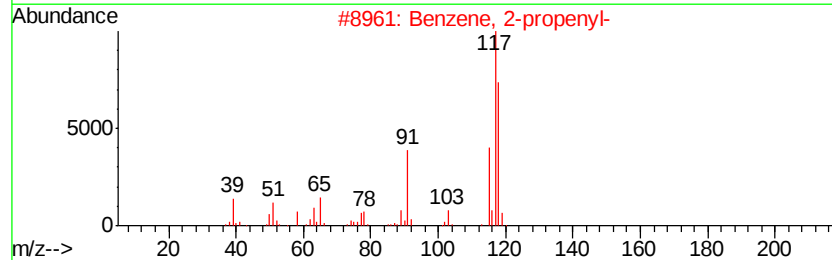
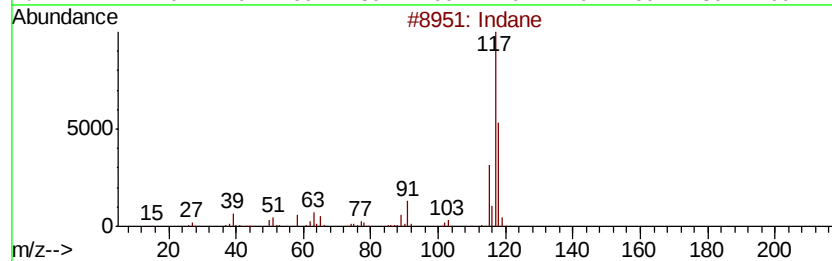
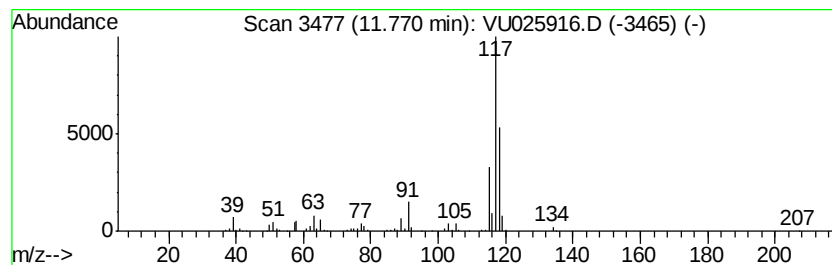
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.M
Quant Title : SW846 8260

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

Peak Number 10 Indane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	11.24 ug/l	281255	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	95
2		Benzene, 2-propenyl-	118	C9H10	000300-57-2	83
3		Benzene, cyclopropyl-	118	C9H10	000873-49-4	83
4		Deltacyclene	118	C9H10	007785-10-6	64
5		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU080618\
Data File : VU025916.D
Acq On : 07 Aug 2018 03:51
Operator : MD/SY
Sample : J4344-02
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
FL-0562

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
Butane	1.40	44.1	ug/l	749252	1	4.99	849244	50.0
Butane, 2-methyl-	1.76	55.8	ug/l	947861	1	4.99	849244	50.0
Pentane	1.95	27.0	ug/l	458179	1	4.99	849244	50.0
Cyclopentane	2.80	21.2	ug/l	359332	1	4.99	849244	50.0
Cyclopentane, met...	4.10	32.1	ug/l	545619	1	4.99	849244	50.0
Benzene, 1-ethyl-...	10.68	51.9	ug/l	1297030	4	11.49	1250820	50.0
Benzene, 1,2,3-tr...	11.57	34.8	ug/l	869462	4	11.49	1250820	50.0
Indane	11.77	11.2	ug/l	281255	4	11.49	1250820	50.0