

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU071218\
 Data File : VU025346.D
 Acq On : 12 Jul 2018 17:56
 Operator : MD/SY
 Sample : J3928-03
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 GROUNDWATER

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\82U070918W.M

Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.088	137	155	180	rBV	3478593	4000980	100.00%	36.747%
2	4.387	1168	1181	1200	rBV2	18960	47753	1.19%	0.439%
3	4.889	1320	1337	1353	rBV	176668	398325	9.96%	3.658%
4	4.985	1353	1367	1391	rVB	256372	567999	14.20%	5.217%
5	5.313	1454	1469	1498	rBV	187704	398920	9.97%	3.664%
6	5.889	1635	1648	1672	rBV	359378	716079	17.90%	6.577%
7	7.567	2156	2170	2187	rBV	650422	1132665	28.31%	10.403%
8	9.091	2632	2644	2669	rBV	502531	843179	21.07%	7.744%
9	10.310	3011	3023	3049	rBV	491285	794312	19.85%	7.295%
10	11.487	3377	3389	3406	rBV	560786	885211	22.12%	8.130%
11	11.574	3406	3416	3431	rVB	24676	40764	1.02%	0.374%
12	11.770	3464	3477	3497	rVB3	77327	168003	4.20%	1.543%
13	12.358	3649	3660	3680	rBV2	32209	59683	1.49%	0.548%
14	13.127	3887	3899	3910	rVB3	80780	126727	3.17%	1.164%
15	13.371	3965	3975	3983	rBV3	31069	48138	1.20%	0.442%
16	13.519	4010	4021	4033	rBV4	126249	235055	5.87%	2.159%
17	13.612	4045	4050	4067	rVB2	46466	78747	1.97%	0.723%
18	14.155	4208	4219	4234	rBV2	23498	43158	1.08%	0.396%
19	14.336	4264	4275	4285	rBV	52431	90241	2.26%	0.829%
20	14.480	4311	4320	4329	rVB2	27641	42251	1.06%	0.388%
21	14.545	4329	4340	4352	rVB5	40110	76626	1.92%	0.704%
22	14.692	4370	4386	4398	rBV6	22396	51273	1.28%	0.471%
23	15.593	4651	4666	4682	rVB3	22420	41964	1.05%	0.385%

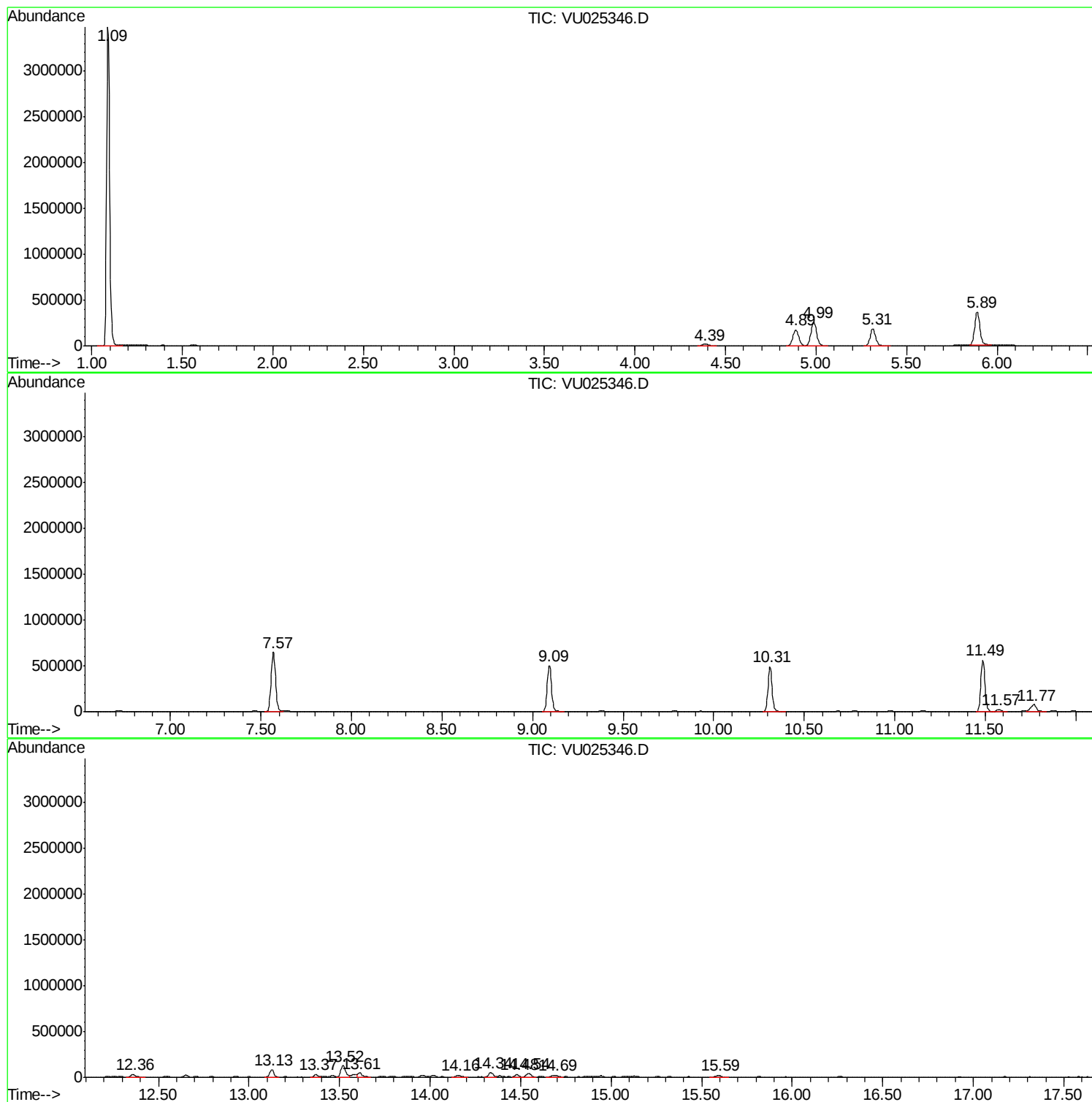
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Quant Title : SW846 8260

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



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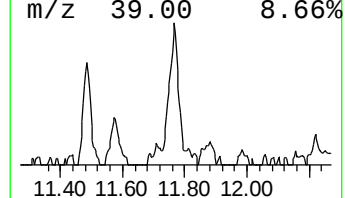
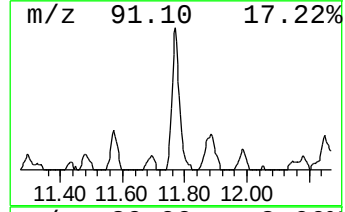
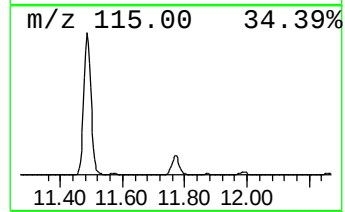
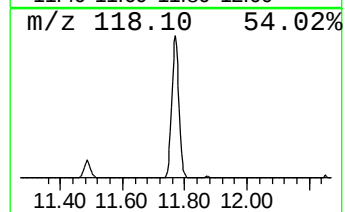
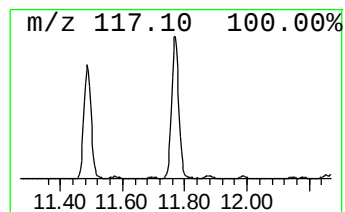
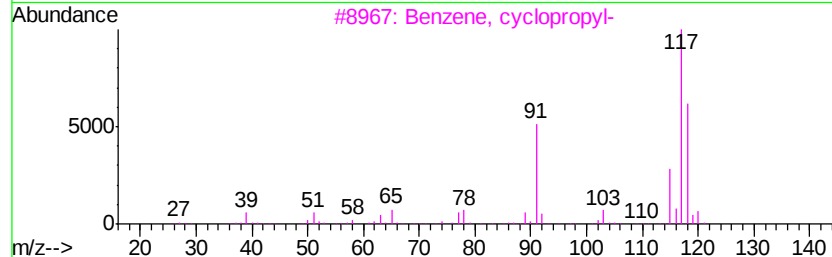
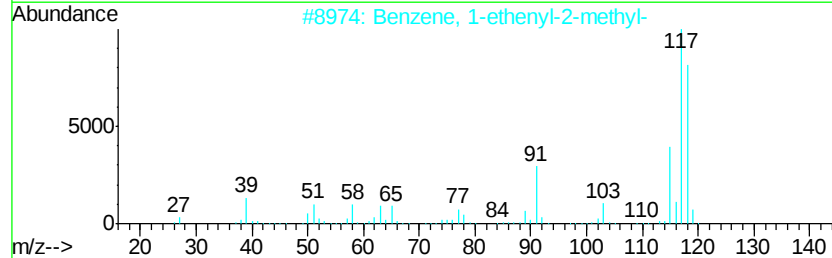
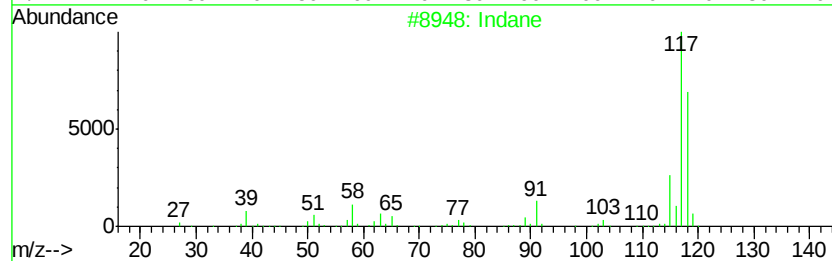
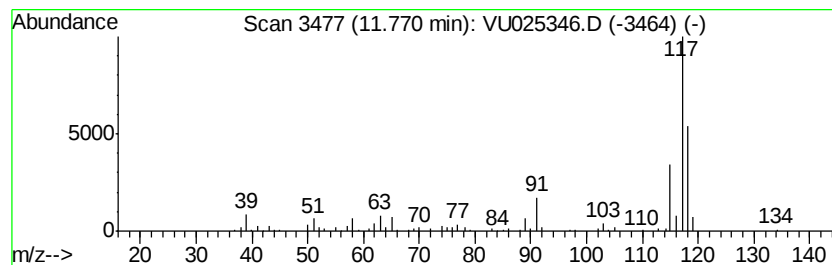
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260

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

Peak Number 2 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	9.49 ug/l	168003	1,4-Dichlorobenzene-d4	11.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	94
2	Benzene, 1-ethenyl-2-methyl-		118	C9H10	000611-15-4	78
3	Benzene, cyclopropyl-		118	C9H10	000873-49-4	74
4	Tetracyclo[3.3.1.0(2,8).0(4,6)]-	...	118	C9H10	1000191-13-7	74
5	Benzene, 1-propenyl-		118	C9H10	000637-50-3	74



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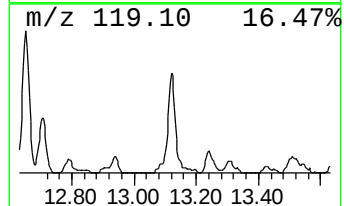
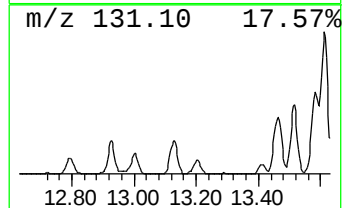
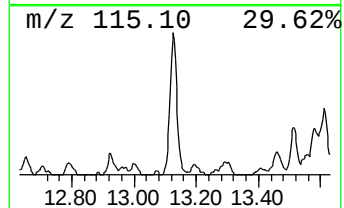
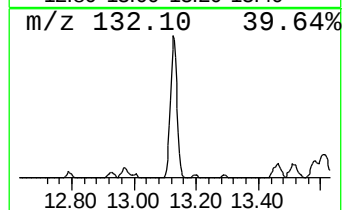
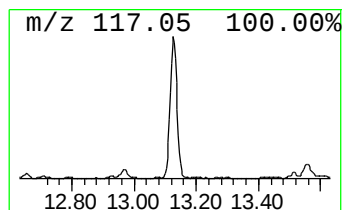
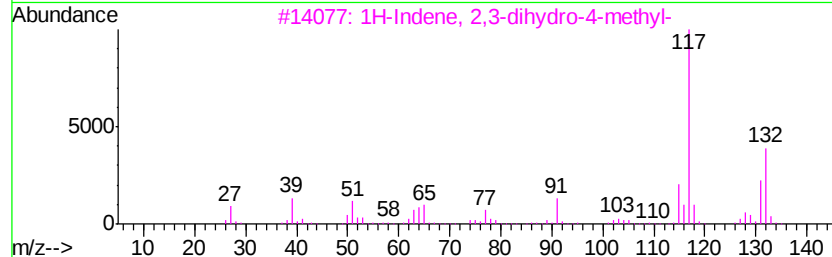
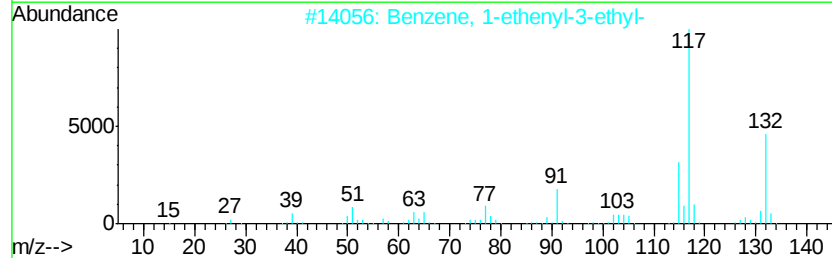
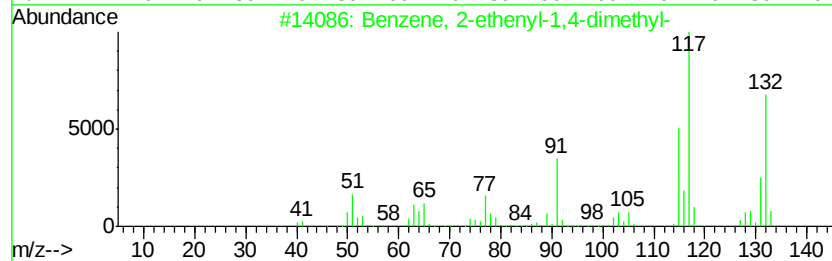
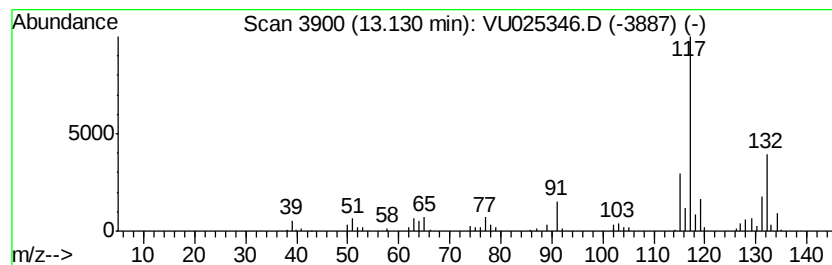
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TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.13	7.16 ug/l	126727	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	90
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	76
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76
4		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	74
5		Indan, 1-methyl-	132	C10H12	000767-58-8	74



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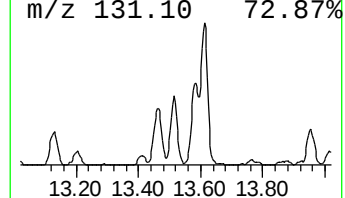
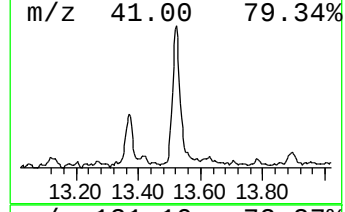
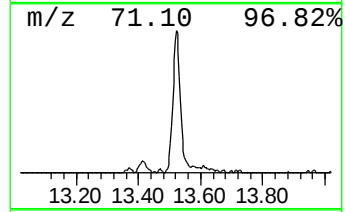
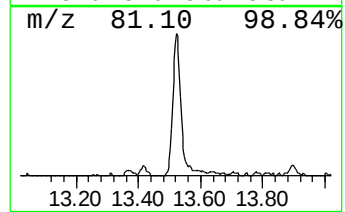
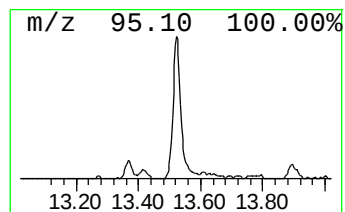
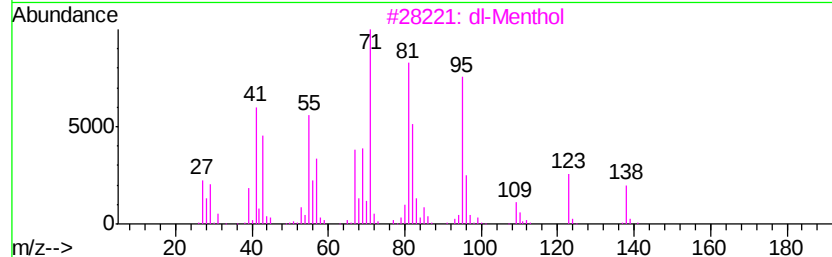
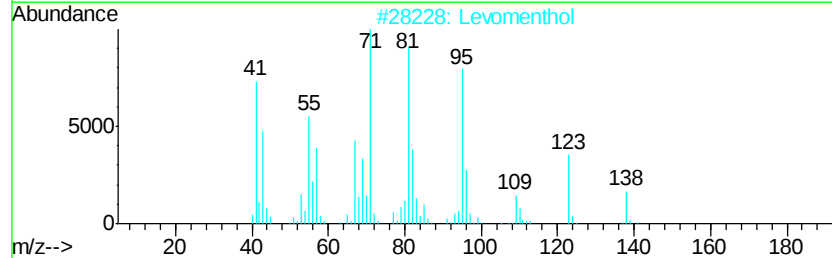
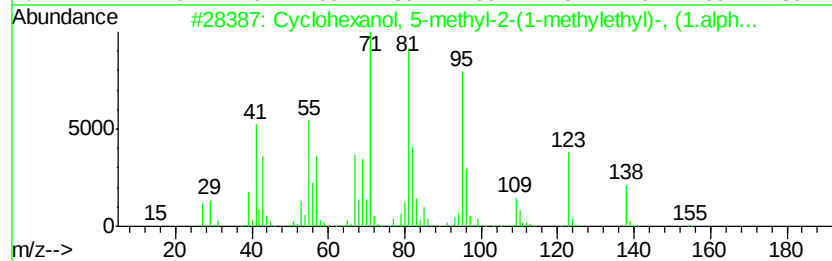
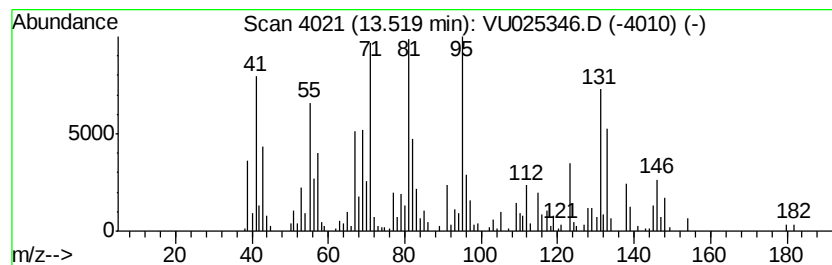
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Peak Number 4 Cyclohexanol, 5-methyl-2-(1... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.52	13.28 ug/l	235055	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanol, 5-methyl-2-(1-meth...	156	C10H20O	015356-70-4	70
2		Levomenthol	156	C10H20O	002216-51-5	70
3		dl-Menthol	156	C10H20O	000089-78-1	62
4		Cyclohexanol, 5-methyl-2-(1-meth...	156	C10H20O	002216-52-6	62
5		Cyclohexanol, 5-methyl-2-(1-meth...	156	C10H20O	001490-04-6	62



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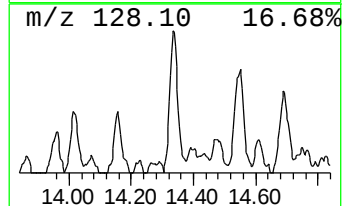
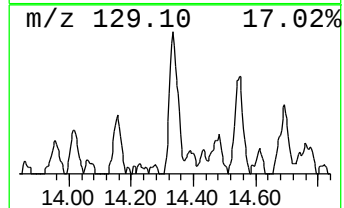
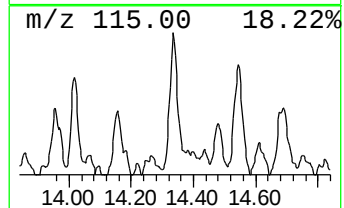
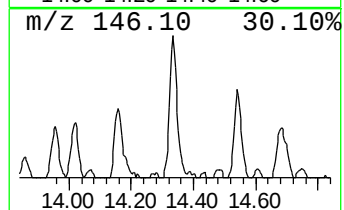
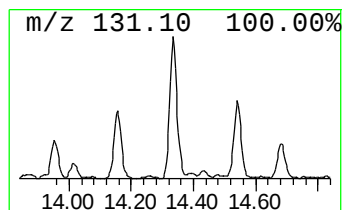
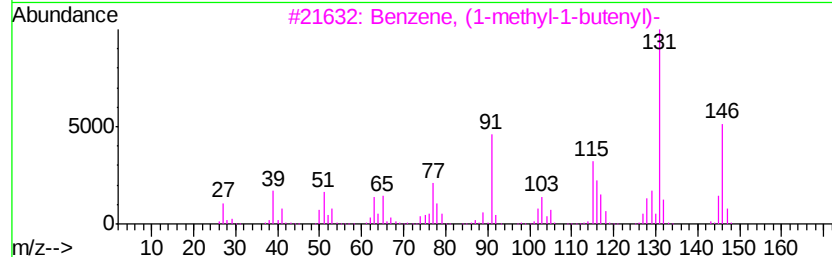
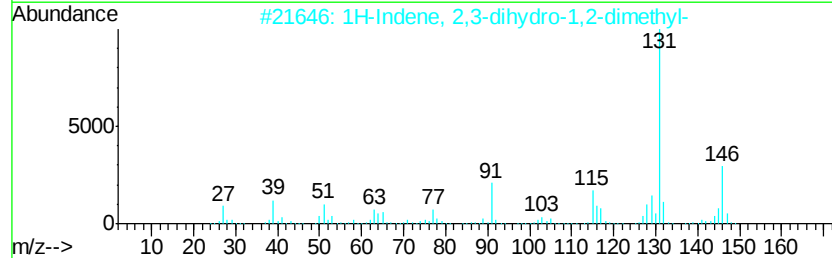
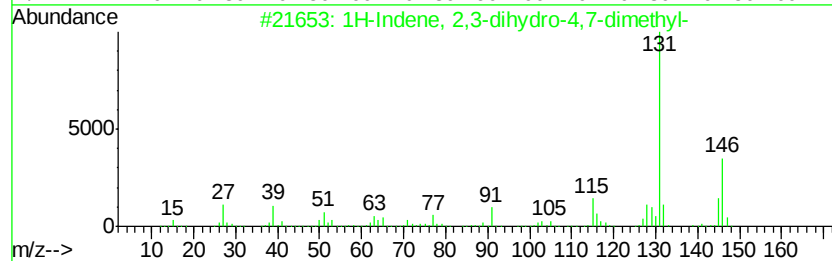
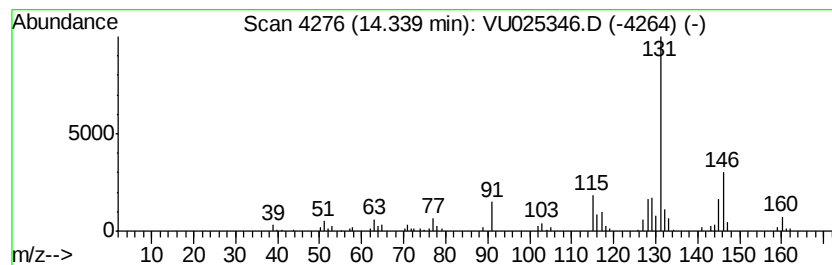
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Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 5 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.34	5.10 ug/l	90241	1,4-Dichlorobenzene-d4	11.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93
2		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	90
3		Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	90
4		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	90
5		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	90



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Indane	11.77	9.5	ug/l	168003	4	11.49	885211	50.0
Benzene, 2-etheny...	13.13	7.2	ug/l	126727	4	11.49	885211	50.0
Cyclohexanol, 5-m...	13.52	13.3	ug/l	235055	4	11.49	885211	50.0
1H-Indene, 2,3-di...	14.34	5.1	ug/l	90241	4	11.49	885211	50.0