

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU061818\
 Data File : VU024691.D
 Acq On : 18 Jun 2018 15:00
 Operator : MD/SY
 Sample : J3463-09
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 T11-MW-13-8.0-061218

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\82U061318W.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	147	155	181	rBV	15457734	19024035	100.00%	64.880%
2	3.001	729	750	770	rBV	138479	302075	1.59%	1.030%
3	4.889	1314	1337	1353	rBV	187712	426170	2.24%	1.453%
4	4.988	1353	1368	1389	rVB	253890	562501	2.96%	1.918%
5	5.313	1451	1469	1485	rBV	184877	390617	2.05%	1.332%
6	5.892	1634	1649	1672	rBV	365722	721007	3.79%	2.459%
7	7.570	2147	2171	2205	rBV	675337	1215239	6.39%	4.144%
8	9.094	2632	2645	2666	rBV	519699	873634	4.59%	2.979%
9	10.310	3012	3023	3041	rBV	500790	815000	4.28%	2.779%
10	11.487	3378	3389	3404	rBV	585239	928891	4.88%	3.168%
11	11.770	3462	3477	3497	rBV3	284221	564495	2.97%	1.925%
12	12.255	3616	3628	3634	rBV	189720	289758	1.52%	0.988%
13	12.358	3650	3660	3681	rBV3	189379	401573	2.11%	1.370%
14	12.651	3732	3751	3764	rBV	168224	296595	1.56%	1.012%
15	13.123	3885	3898	3909	rBV2	513805	824005	4.33%	2.810%
16	13.290	3940	3950	3967	rVB	155593	262277	1.38%	0.894%
17	13.744	4079	4091	4101	rBV	208091	350050	1.84%	1.194%
18	14.879	4432	4444	4455	rBV	229188	364957	1.92%	1.245%
19	15.065	4491	4502	4517	rBV	264470	440209	2.31%	1.501%
20	16.062	4800	4812	4819	rBV	165440	268830	1.41%	0.917%

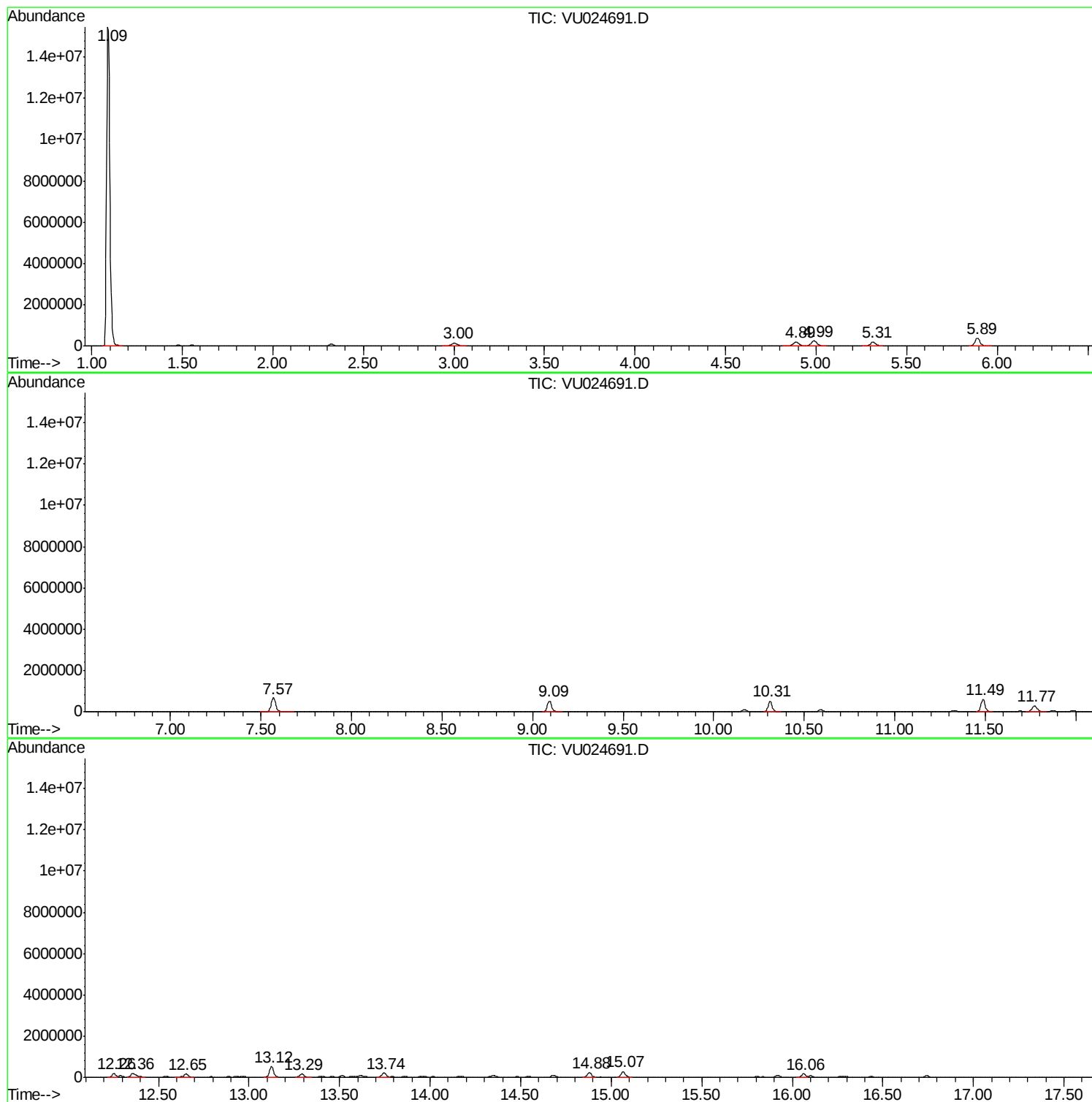
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

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



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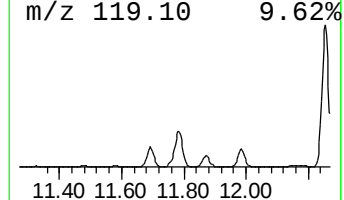
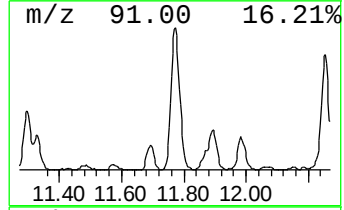
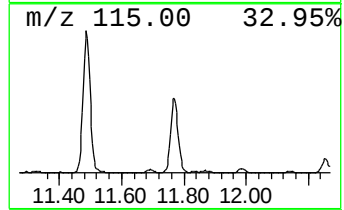
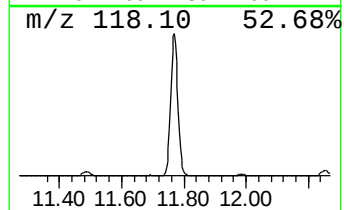
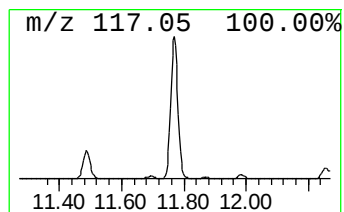
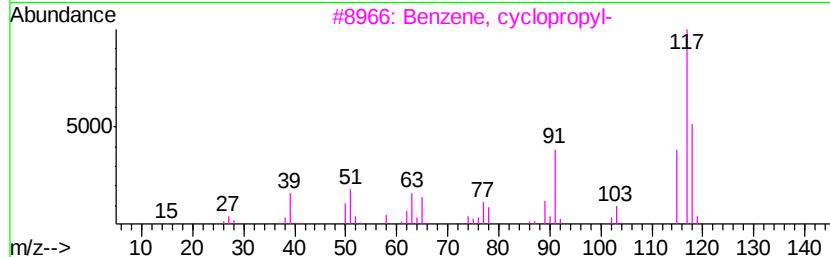
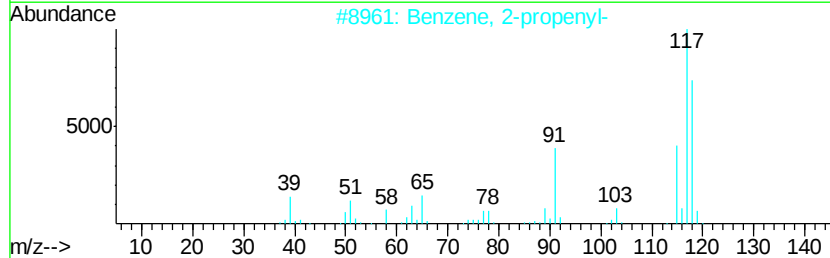
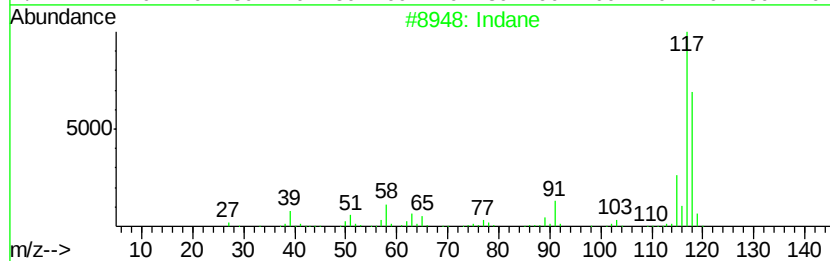
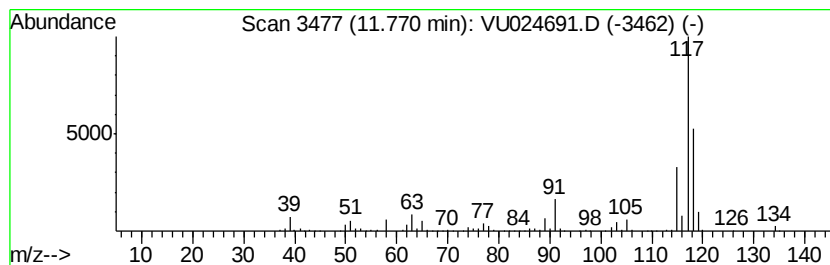
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Quant Title : SW846 8260

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

Peak Number 2 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	30.39 ug/l	564495	1,4-Dichlorobenzene-d4	11.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indane		118 C9H10	000496-11-7 83
2	Benzene, 2-propenyl-		118 C9H10	000300-57-2 83
3	Benzene, cyclopropyl-		118 C9H10	000873-49-4 81
4	Benzene, 1-ethenyl-2-methyl-		118 C9H10	000611-15-4 53
5	Benzene, 1-propenyl-		118 C9H10	000637-50-3 47



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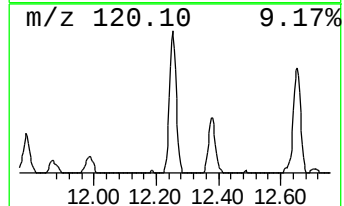
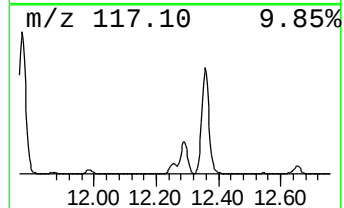
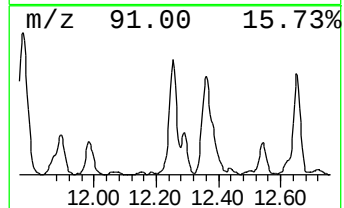
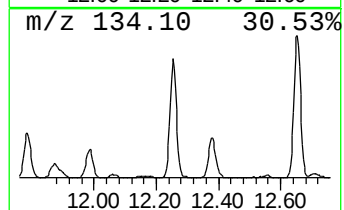
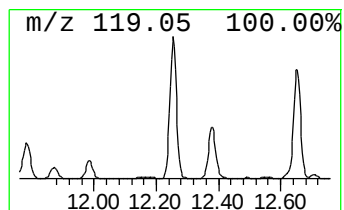
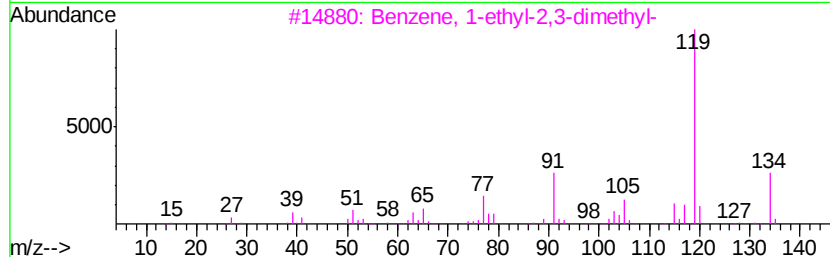
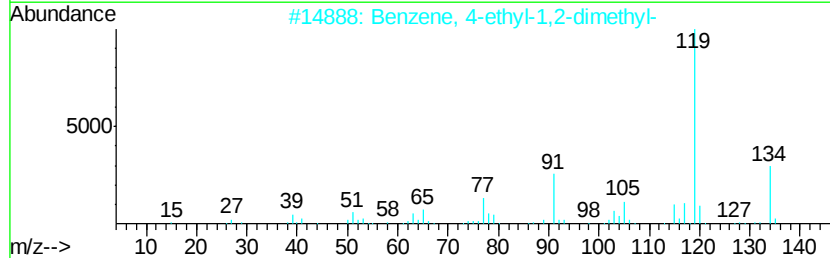
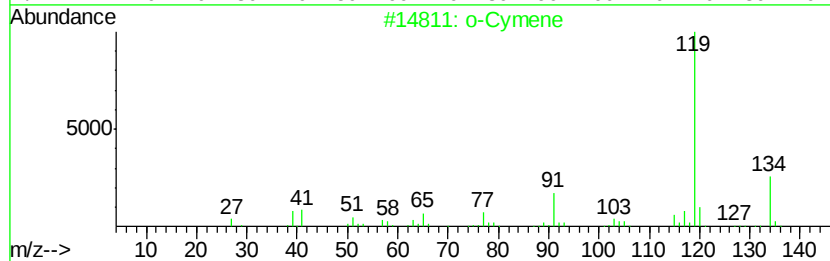
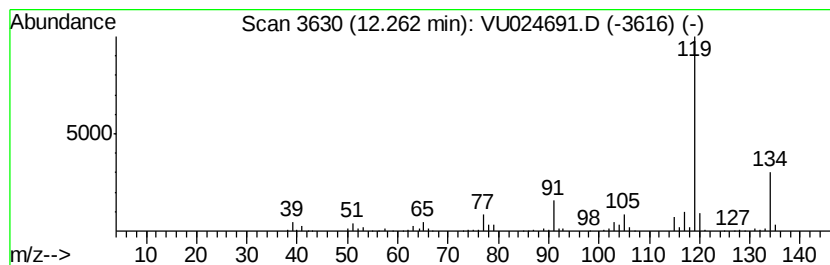
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Quant Title : SW846 8260

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

Peak Number 3 o-Cymene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.26	15.60 ug/l	289758	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	97
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



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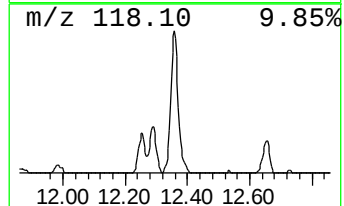
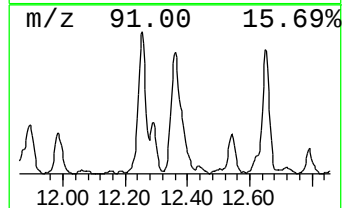
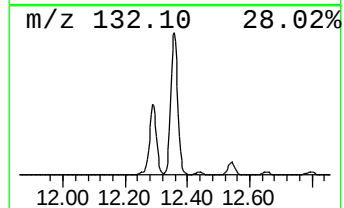
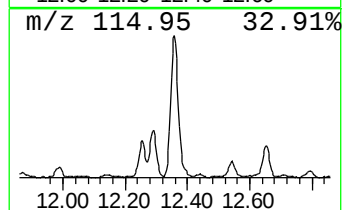
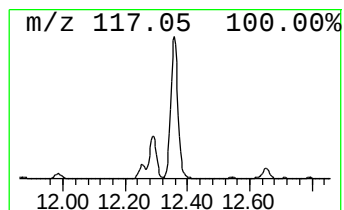
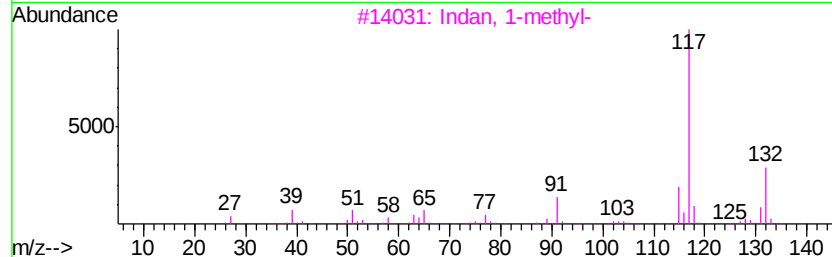
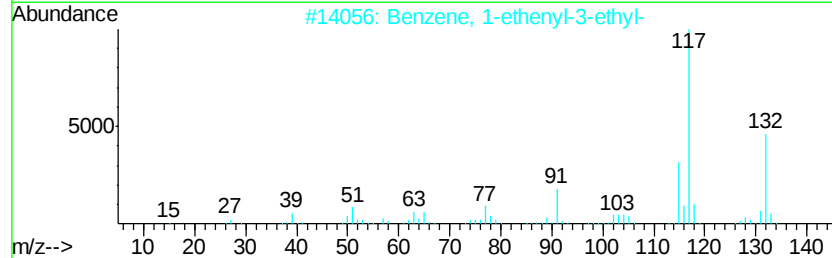
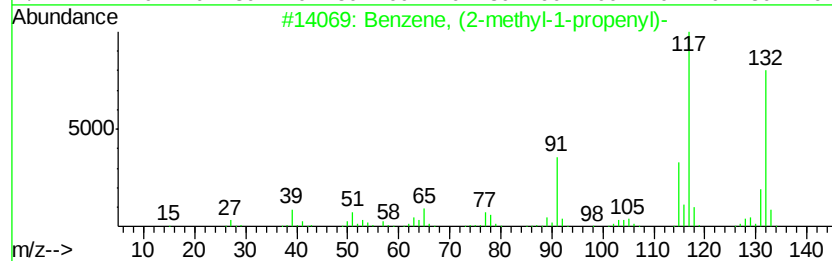
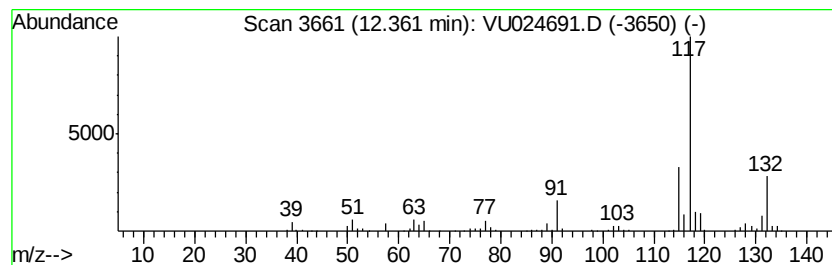
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Peak Number 4 Benzene, (2-methyl-1-propen... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	21.62 ug/l	401573	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
3		Indan, 1-methyl-	132	C10H12	000767-58-8	87
4		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83
5		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	83



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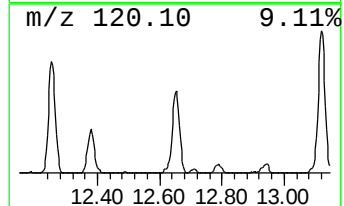
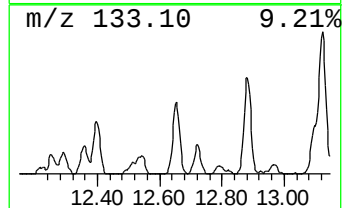
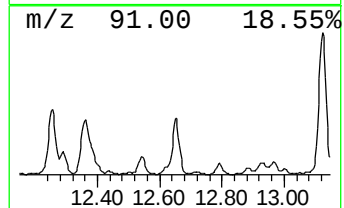
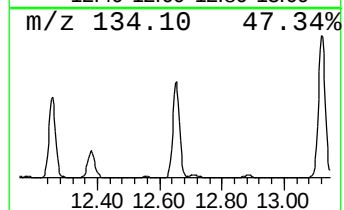
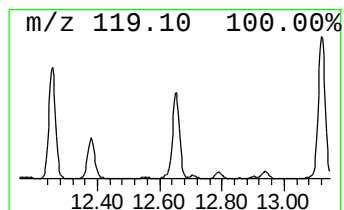
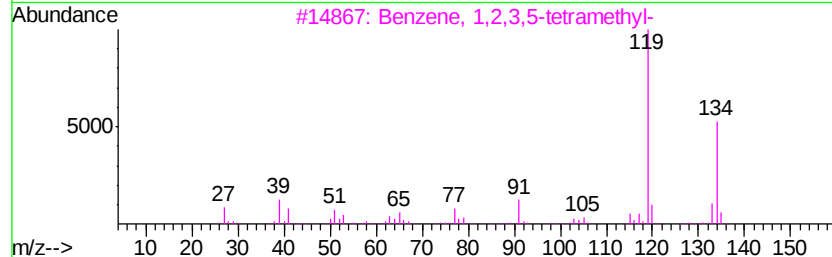
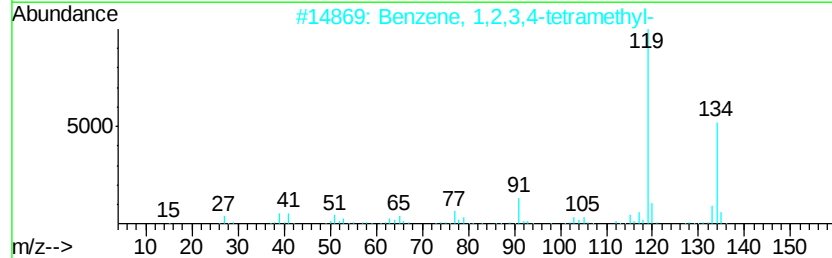
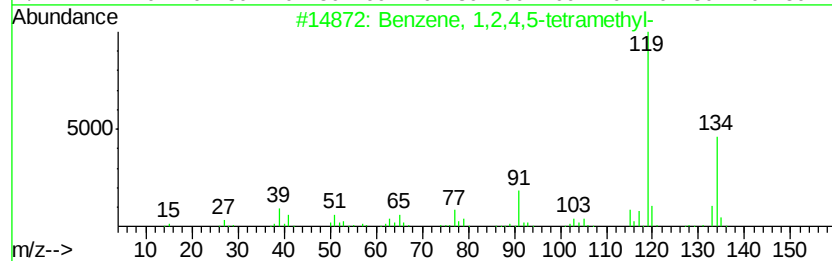
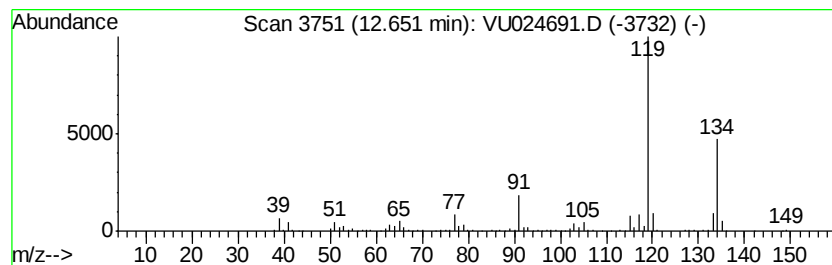
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	15.97 ug/l	296595	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91



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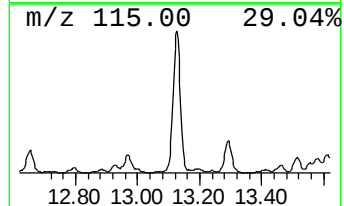
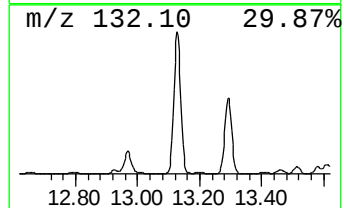
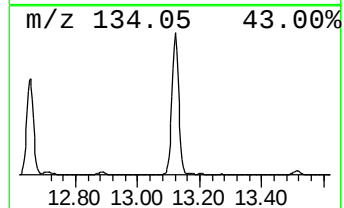
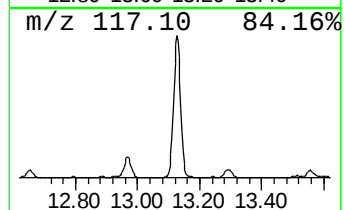
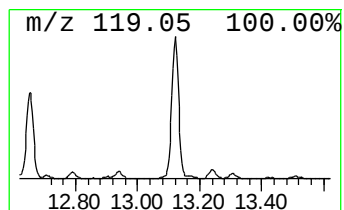
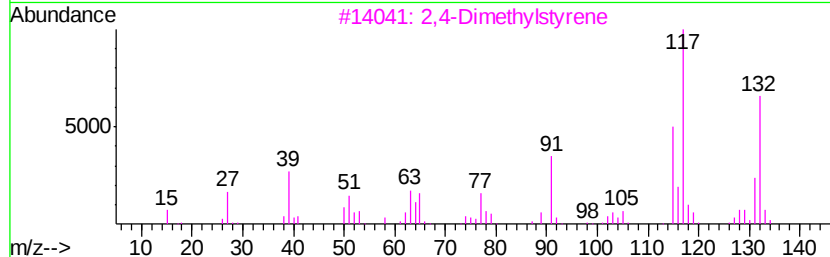
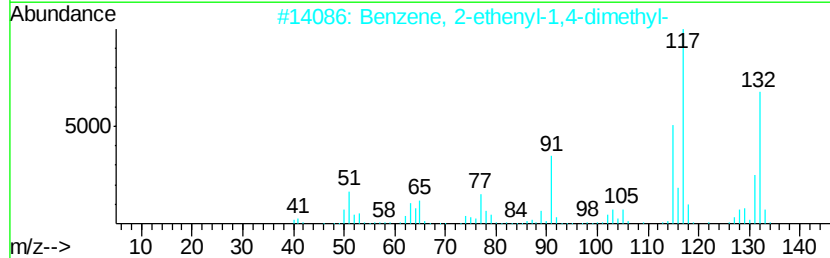
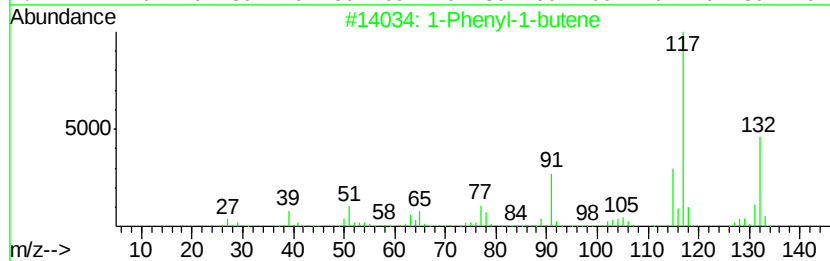
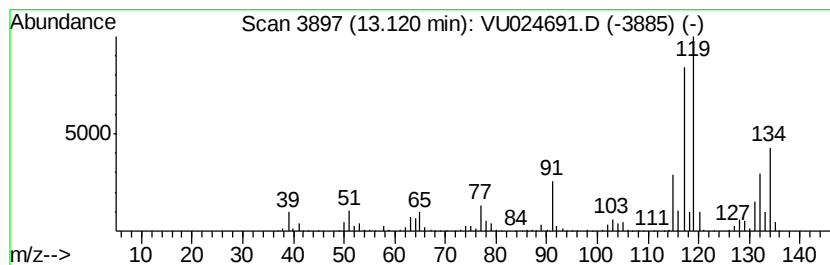
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TIC Integration Parameters: LSCINT.P

Peak Number 6 1-Phenyl-1-butene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	44.35 ug/l	824005	1,4-Dichlorobenzene-d4	11.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1-Phenyl-1-butene	132 C10H12	000824-90-8 86
2		Benzene, 2-ethenyl-1,4-dimethyl-	132 C10H12	002039-89-6 86
3		2,4-Dimethylstyrene	132 C10H12	002234-20-0 86
4		Benzene, 2-butenyl-	132 C10H12	001560-06-1 80
5		Benzene, 2-ethenyl-1,3-dimethyl-	132 C10H12	002039-90-9 64



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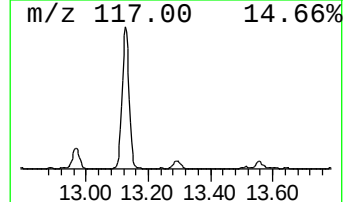
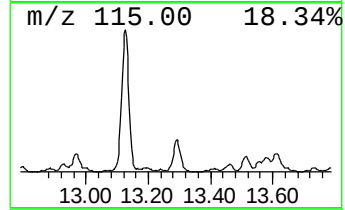
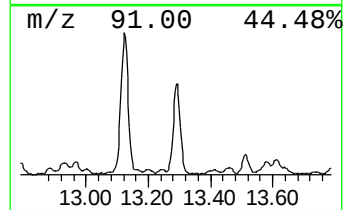
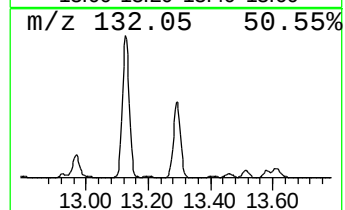
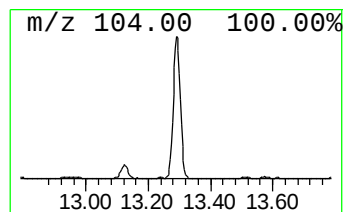
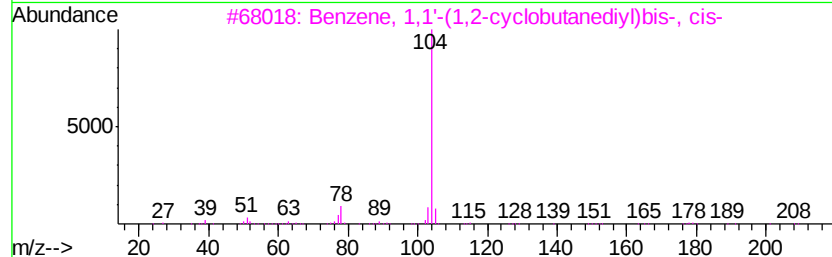
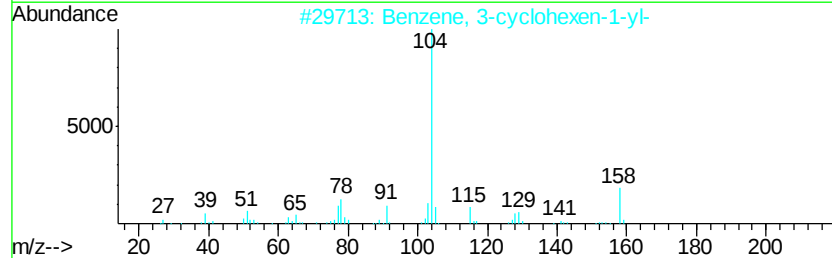
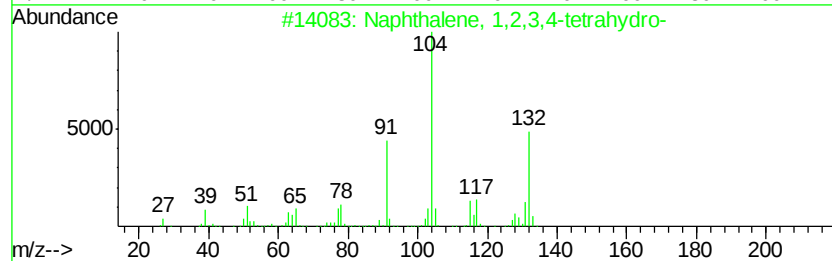
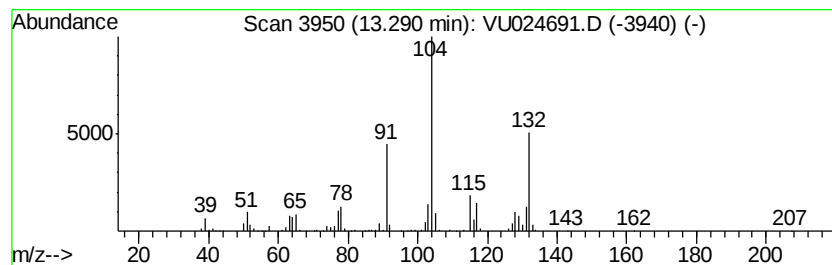
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Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1,2,3,4-tetra... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.29	14.12 ug/l	262277	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	91
2		Benzene, 3-cyclohexen-1-yl-	158	C12H14	004994-16-5	50
3		Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	007694-30-6	38
4		4H-Pyran-4-one 2,3-dihydro-2-phe...	174	C11H10O2	1000163-21-2	38
5		Acetic acid, 2-phenylethyl ester	164	C10H12O2	000103-45-7	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU061818\
Data File : VU024691.D
Acq On : 18 Jun 2018 15:00
Operator : MD/SY
Sample : J3463-09
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 13 Sample Multiplier: 1

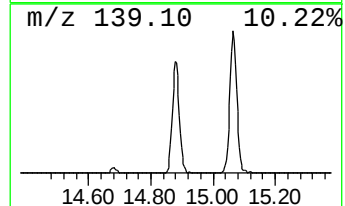
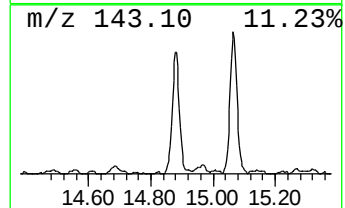
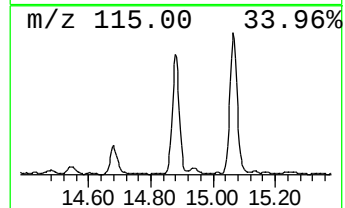
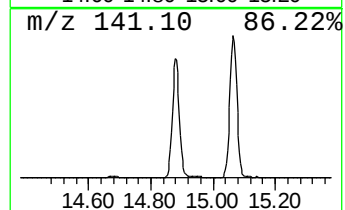
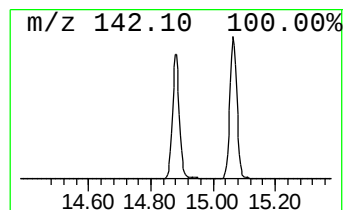
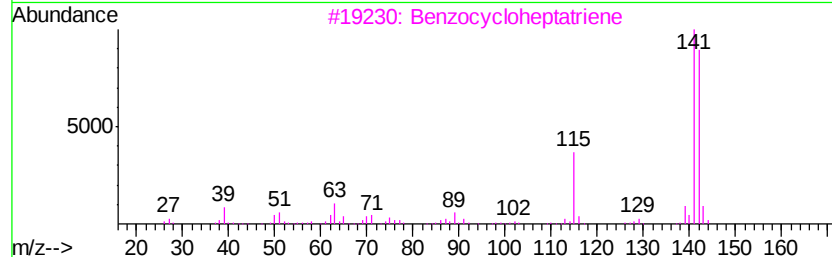
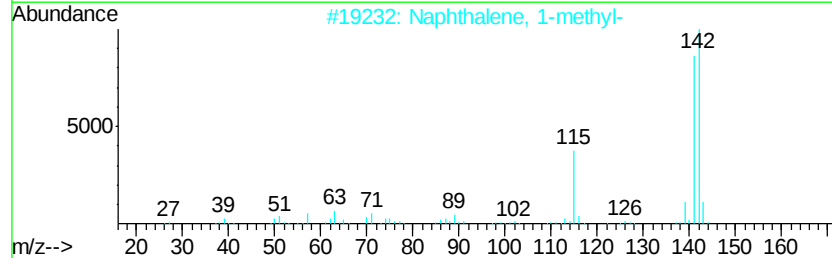
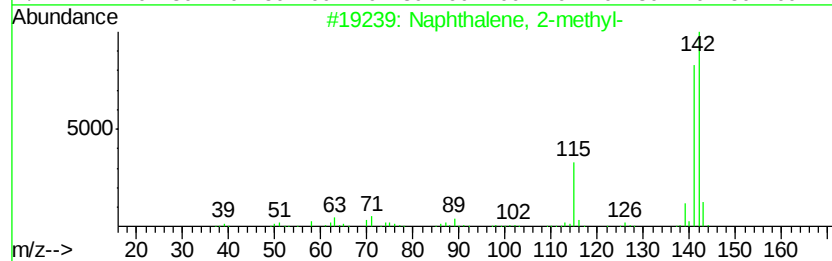
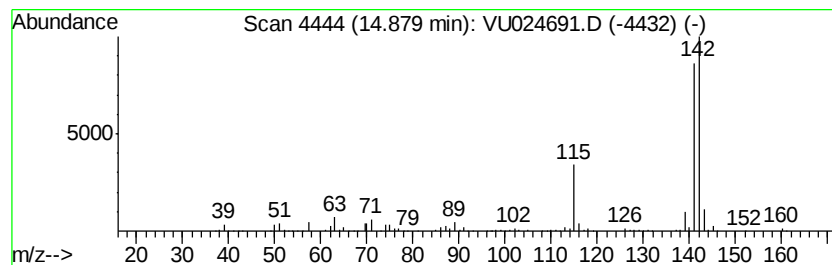
Instrument :
MSVOA_U
ClientSampleId :
T11-MW-13-8.0-061218

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.88	19.64 ug/l	364957	1,4-Dichlorobenzene-d4	11.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2-methyl-	142 C11H10	000091-57-6 96
2		Naphthalene, 1-methyl-	142 C11H10	000090-12-0 96
3		Benzocycloheptatriene	142 C11H10	000264-09-5 91
4		1,4-Methanonaphthalene, 1,4-dihy...	142 C11H10	004453-90-1 91
5		1H-Indene, 1-ethylidene-	142 C11H10	002471-83-2 59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU061818\
Data File : VU024691.D
Acq On : 18 Jun 2018 15:00
Operator : MD/SY
Sample : J3463-09
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
T11-MW-13-8.0-061218

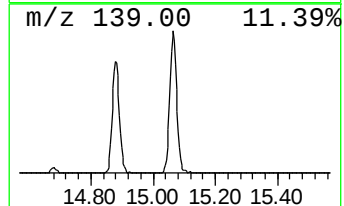
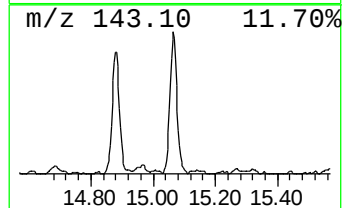
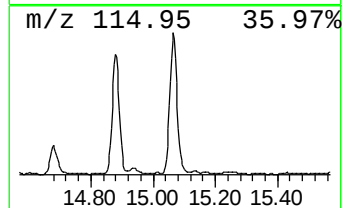
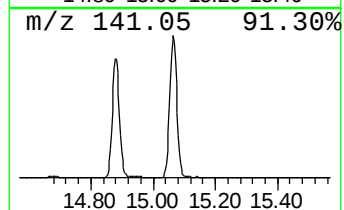
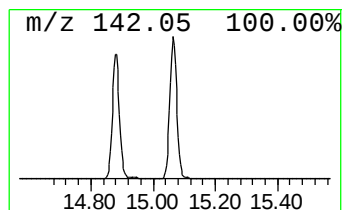
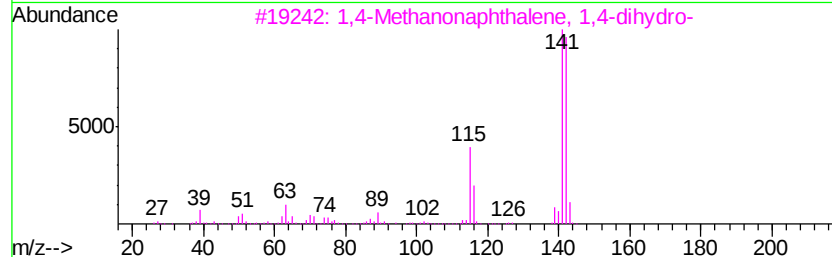
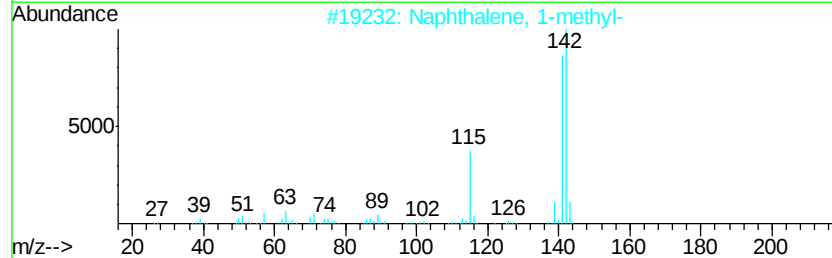
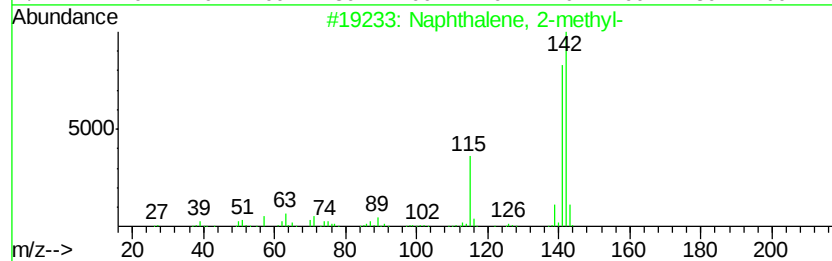
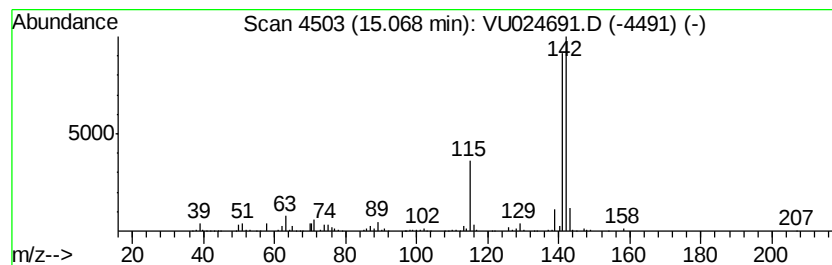
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Quant Title : SW846 8260

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

Peak Number 9 1,4-Methanonaphthalene, 1,4... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.07	23.70 ug/l	440209	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3		1,4-Methanonaphthalene, 1,4-dihv...	142	C11H10	004453-90-1	94
4		Benzocycloheptatriene	142	C11H10	000264-09-5	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU061818\
Data File : VU024691.D
Acq On : 18 Jun 2018 15:00
Operator : MD/SY
Sample : J3463-09
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
T11-MW-13-8.0-061218

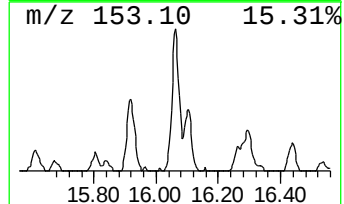
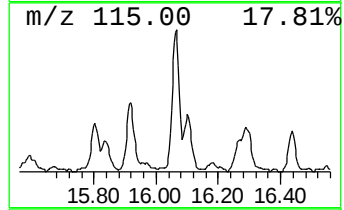
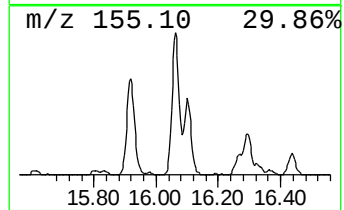
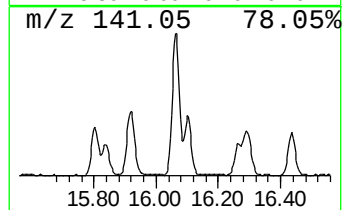
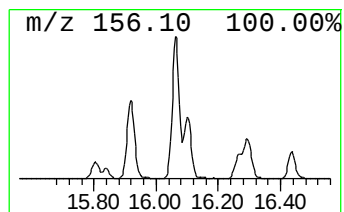
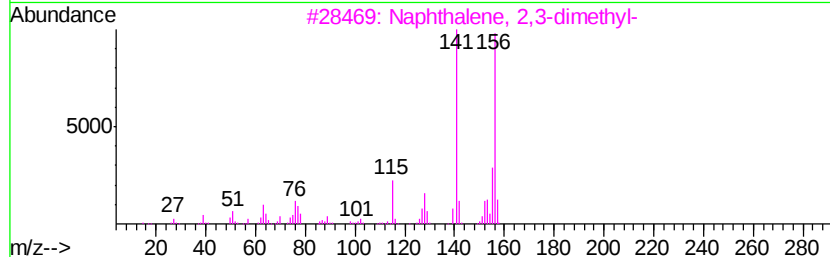
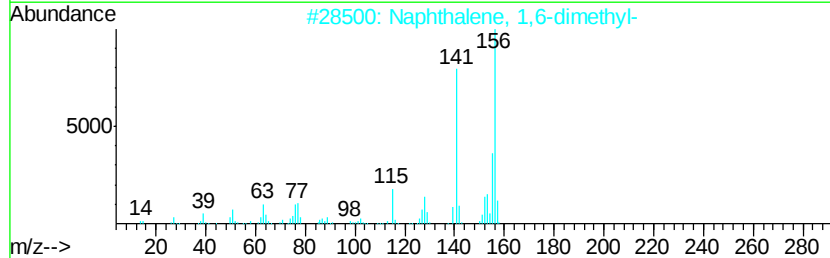
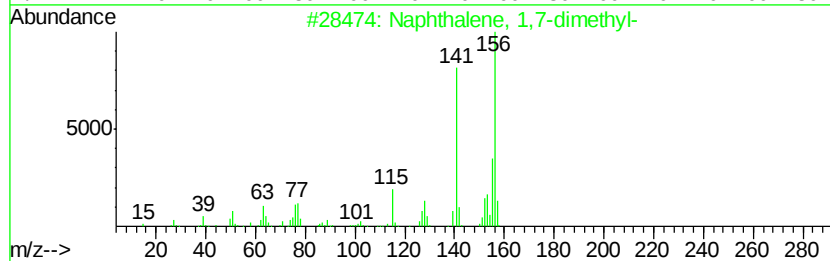
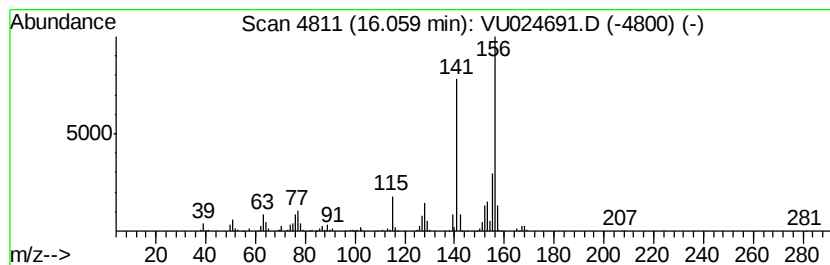
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,7-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.06	14.47 ug/l	268830	1,4-Dichlorobenzene-d4	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	98
2		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
3		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
5		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU061818\
Data File : VU024691.D
Acq On : 18 Jun 2018 15:00
Operator : MD/SY
Sample : J3463-09
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
T11-MW-13-8.0-061218

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.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
Indane	11.77	30.4	ug/l	564495	4	11.49	928891	50.0
o-Cymene	12.26	15.6	ug/l	289758	4	11.49	928891	50.0
Benzene, (2-methy...	12.36	21.6	ug/l	401573	4	11.49	928891	50.0
Benzene, 1,2,4,5-...	12.65	16.0	ug/l	296595	4	11.49	928891	50.0
1-Phenyl-1-butene	13.12	44.4	ug/l	824005	4	11.49	928891	50.0
Naphthalene, 1,2,...	13.29	14.1	ug/l	262277	4	11.49	928891	50.0
Naphthalene, 1-me...	14.88	19.6	ug/l	364957	4	11.49	928891	50.0
1,4-Methanonaphth...	15.07	23.7	ug/l	440209	4	11.49	928891	50.0
Naphthalene, 1,7-...	16.06	14.5	ug/l	268830	4	11.49	928891	50.0