

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX061518\
 Data File : VX002617.D
 Acq On : 16 Jun 2018 00:35
 Operator : JC/MD
 Sample : J3467-05
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PR-MW-05-061118

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_X\METHOD\82X061518W.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	1.075	76	80	93	rBV	1238703	1483646	2.48%	1.098%
2	1.788	192	197	208	rVB	418230	605924	1.01%	0.448%
3	4.398	616	625	638	rVB	210186	611055	1.02%	0.452%
4	5.672	826	834	864	rVB2	218359	865895	1.45%	0.641%
5	6.867	1021	1030	1037	rBV	307874	755094	1.26%	0.559%
6	8.720	1328	1334	1339	rBV	756202	1150500	1.93%	0.851%
7	8.787	1339	1345	1358	rVB	3583882	5637921	9.44%	4.171%
8	10.116	1553	1563	1570	rBV	656437	914871	1.53%	0.677%
9	10.256	1580	1586	1597	rBV	13449504	17954309	30.05%	13.282%
10	10.372	1597	1605	1610	rBV	42885767	59745807	100.00%	44.200%
11	11.024	1705	1712	1723	rBV	1523156	1983890	3.32%	1.468%
12	11.140	1723	1731	1737	rBV	738490	924695	1.55%	0.684%
13	11.366	1759	1768	1774	rBV	1270157	1610803	2.70%	1.192%
14	11.439	1774	1780	1787	rVV	5007392	8482908	14.20%	6.276%
15	11.512	1787	1792	1804	rVB	2608257	3148243	5.27%	2.329%
16	11.811	1836	1841	1847	rVB	10952074	12948775	21.67%	9.579%
17	12.079	1880	1885	1891	rVB	906688	1213153	2.03%	0.897%
18	12.146	1891	1896	1901	rBV	2320562	2763270	4.63%	2.044%
19	12.305	1915	1922	1929	rBV2	2403241	3361188	5.63%	2.487%
20	12.378	1929	1934	1941	rVB2	565940	841421	1.41%	0.622%
21	12.670	1977	1982	1986	rBV	589892	717266	1.20%	0.531%
22	13.024	2036	2040	2046	rVB	637120	765818	1.28%	0.567%
23	13.353	2089	2094	2099	rVB2	865199	1191734	1.99%	0.882%
24	13.841	2166	2174	2182	rBV	2912663	3724708	6.23%	2.756%
25	14.701	2307	2315	2320	rBV	854912	1022330	1.71%	0.756%
26	14.841	2333	2338	2344	rBV	578943	747806	1.25%	0.553%

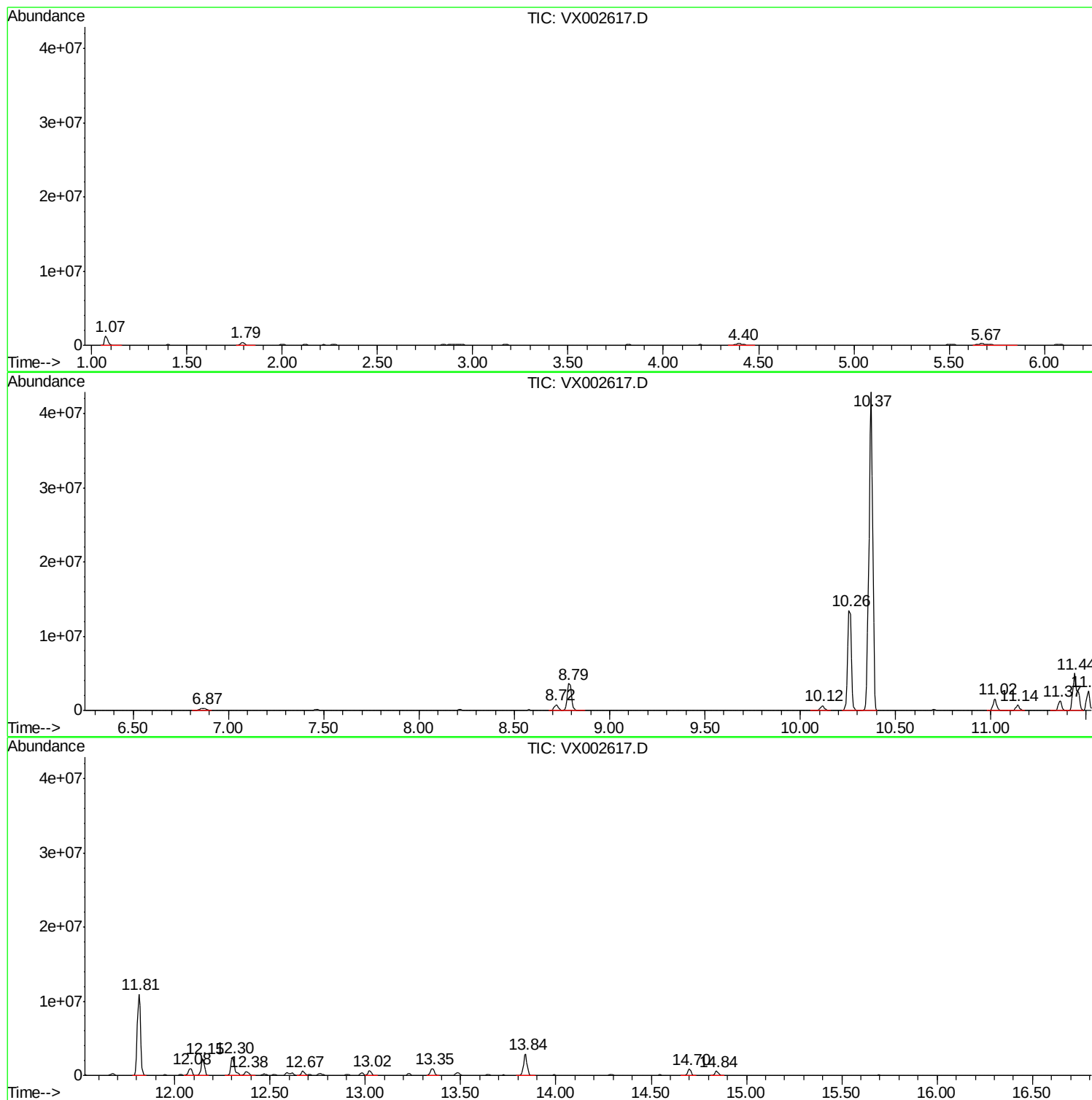
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061518W.M
Quant Title : SW846 8260

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



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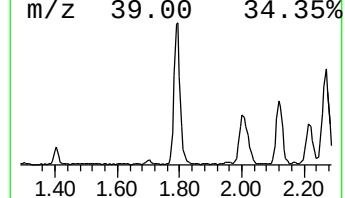
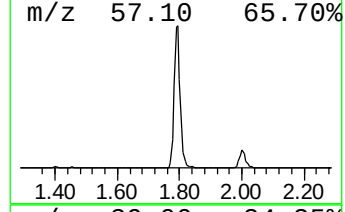
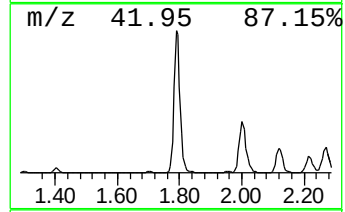
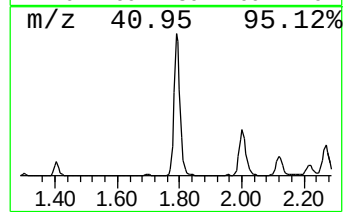
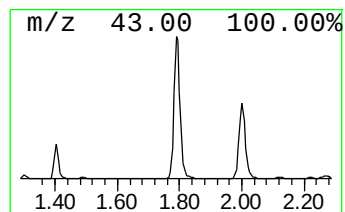
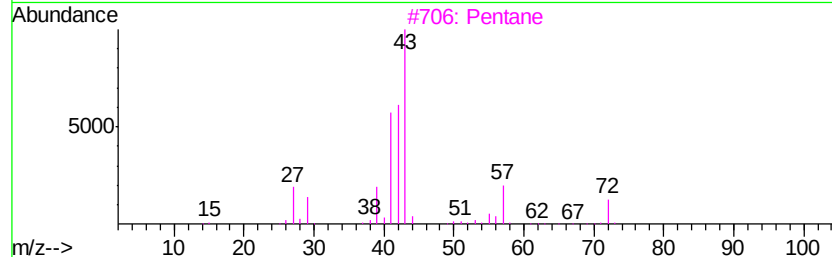
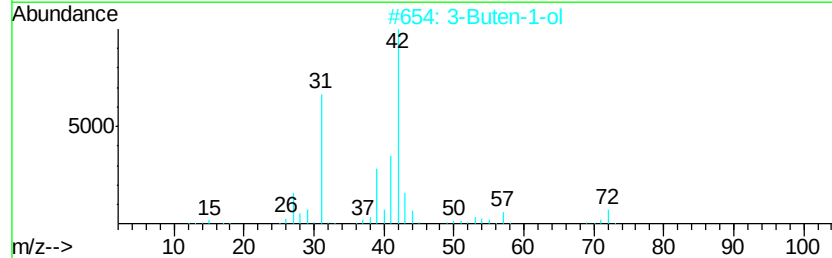
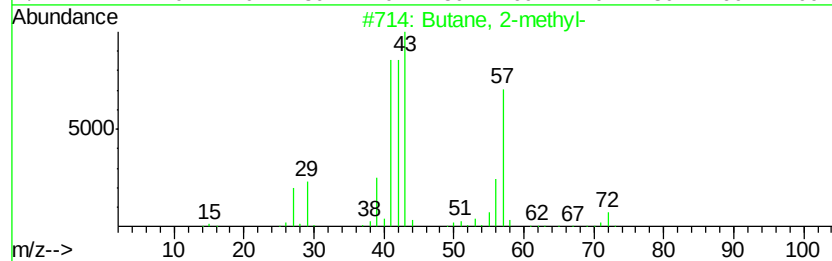
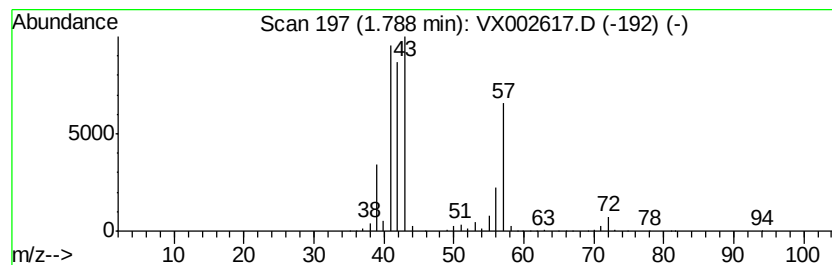
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061518W.M
Quant Title : SW846 8260

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

Peak Number 2 Butane, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.79	34.99 ug/l	605924	Pentafluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methyl-	72	C5H12	000078-78-4	91
2		3-Buten-1-ol	72	C4H8O	000627-27-0	47
3		Pentane	72	C5H12	000109-66-0	36
4		Butane, 1-chloro-2-methyl-	106	C5H11Cl	000616-13-7	16
5		1-Butene	56	C4H8	000106-98-9	14



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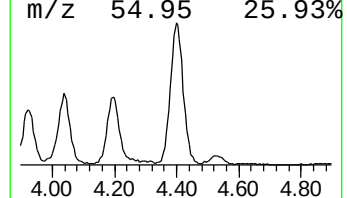
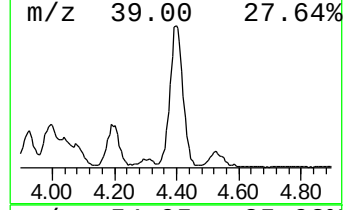
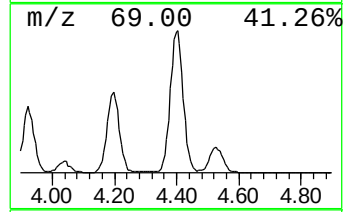
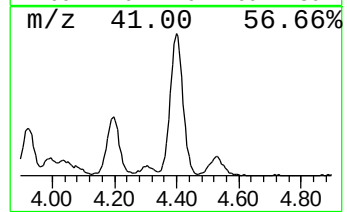
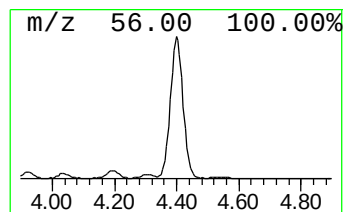
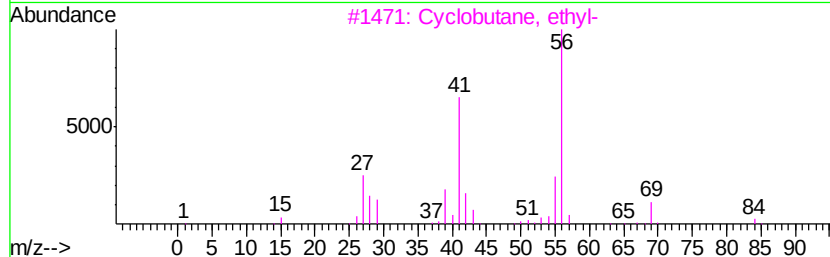
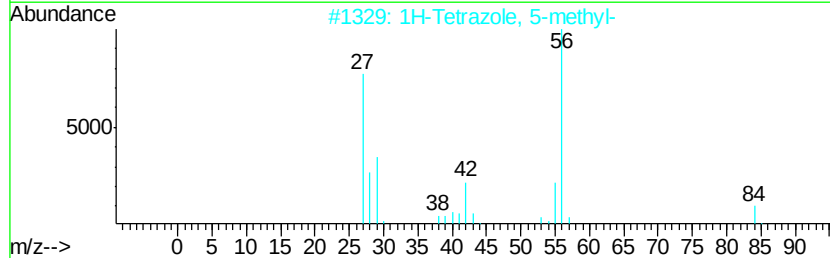
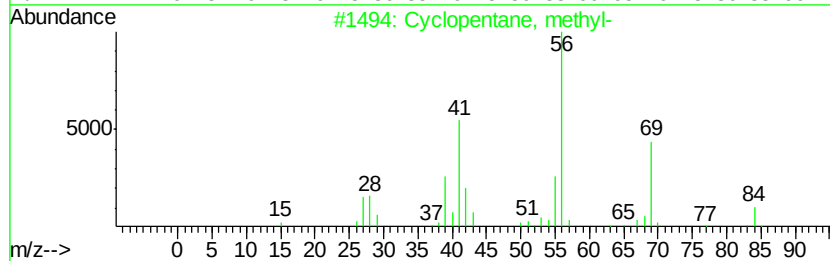
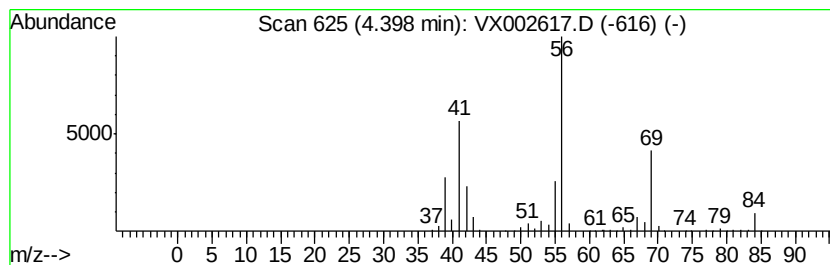
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Quant Title : SW846 8260

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

Peak Number 3 Cyclopentane, methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.40	35.28 ug/l	611055	Pentafluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	80
3		Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
4		Cyclohexane	84	C6H12	000110-82-7	64
5		Cyclobutane	56	C4H8	000287-23-0	50



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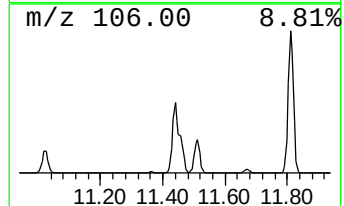
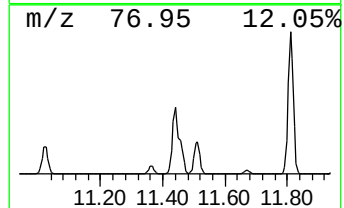
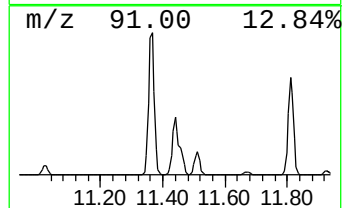
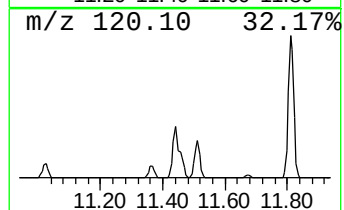
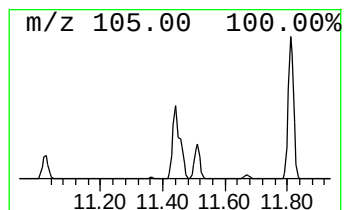
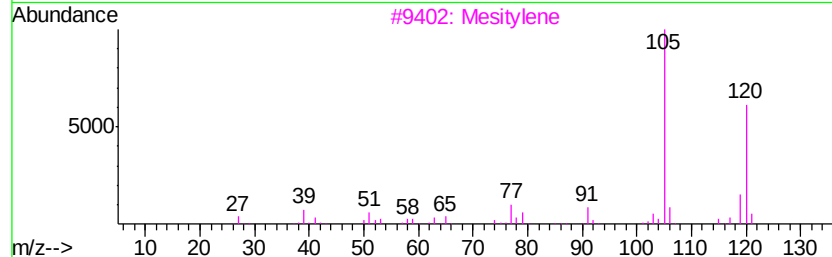
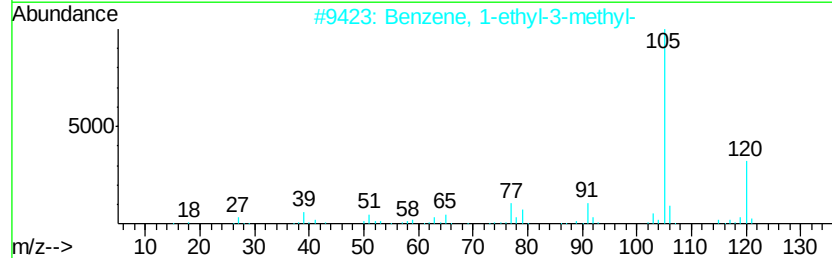
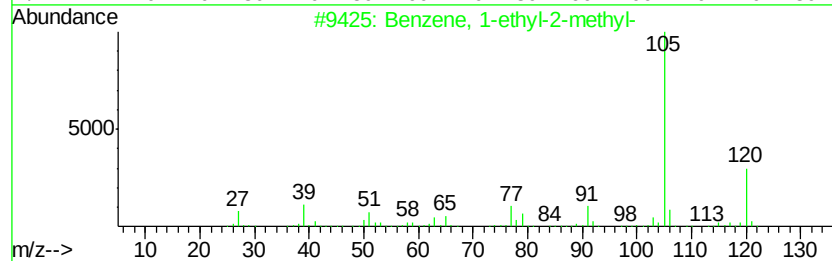
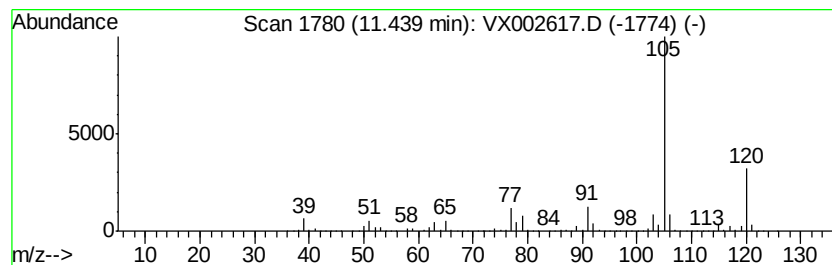
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TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.44	349.62 ug/l	8482910	1,4-Dichlorobenzene-d4	12.09

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	95
3		Mesitylene	120	C9H12	000108-67-8	91
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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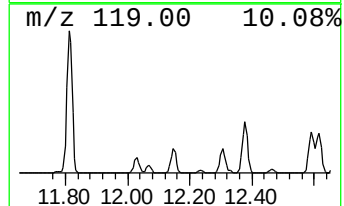
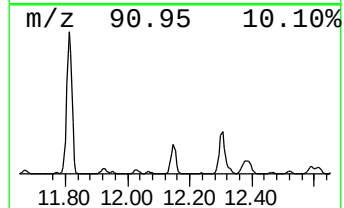
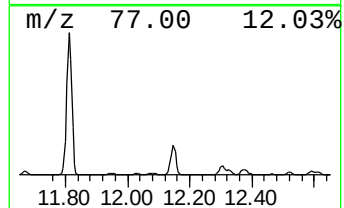
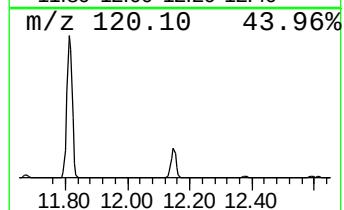
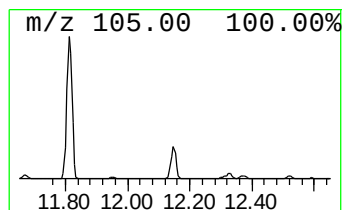
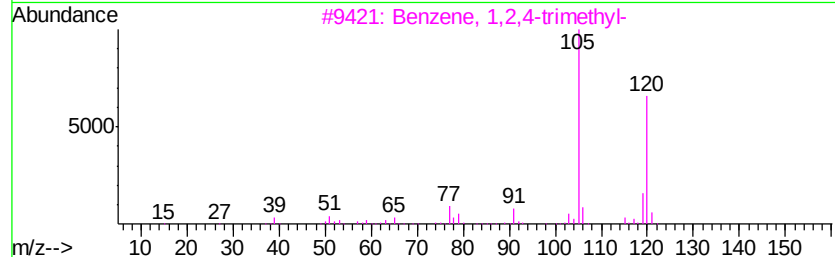
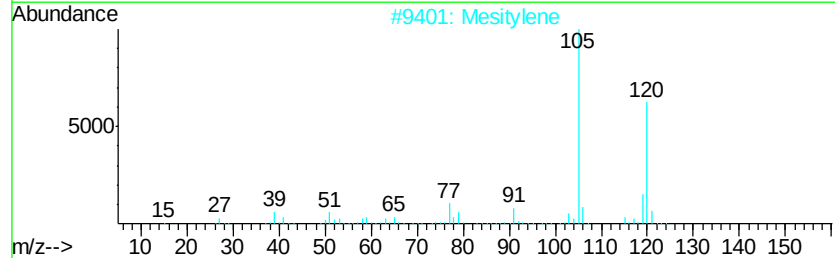
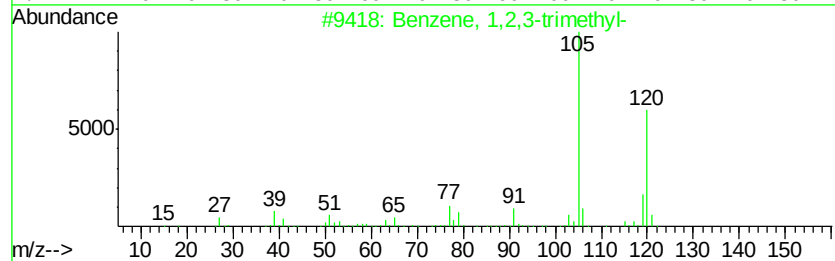
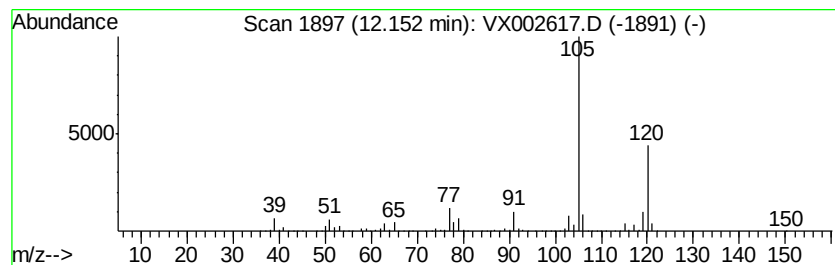
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	113.89 ug/l	2763270	1,4-Dichlorobenzene-d4	12.09

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, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



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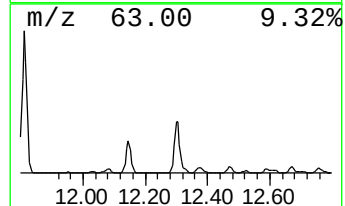
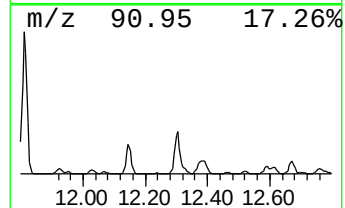
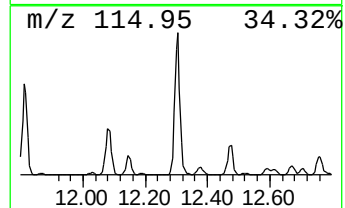
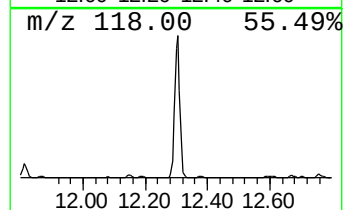
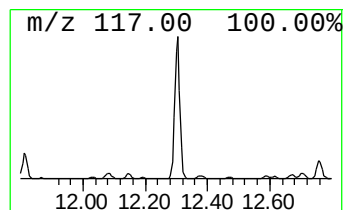
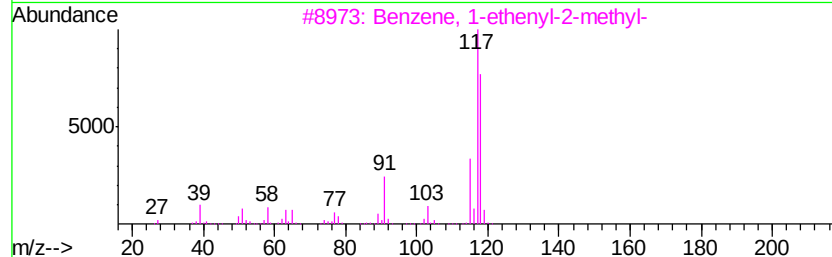
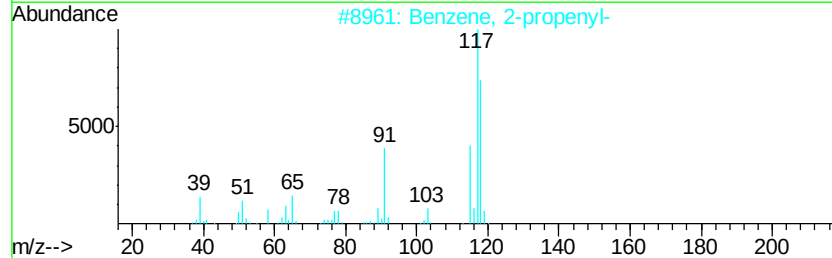
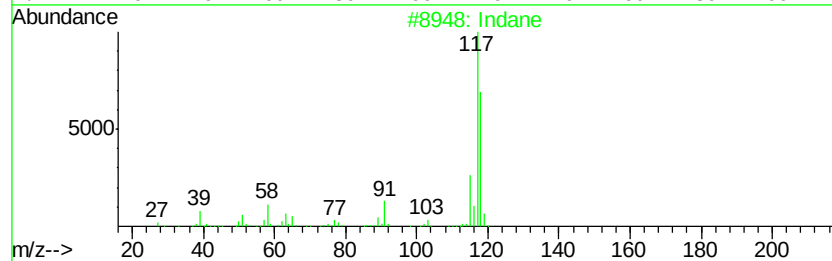
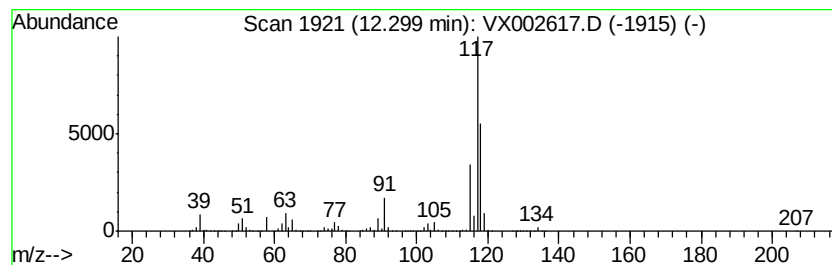
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Quant Title : SW846 8260

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

Peak Number 6 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	138.53 ug/l	3361190	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	87
2		Benzene, 2-propenyl-	118	C9H10	000300-57-2	87
3		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	81
4		Benzene, cyclopropyl-	118	C9H10	000873-49-4	81
5		Deltacyclene	118	C9H10	007785-10-6	64



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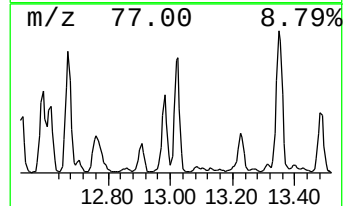
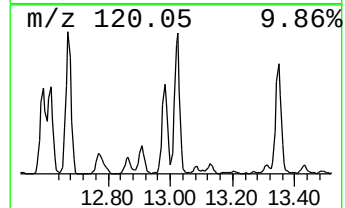
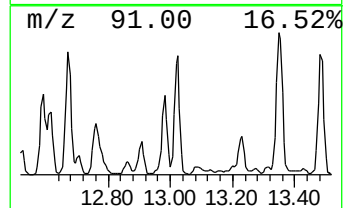
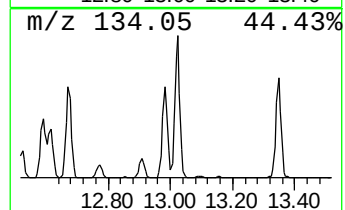
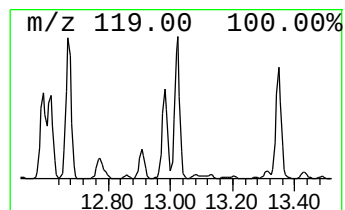
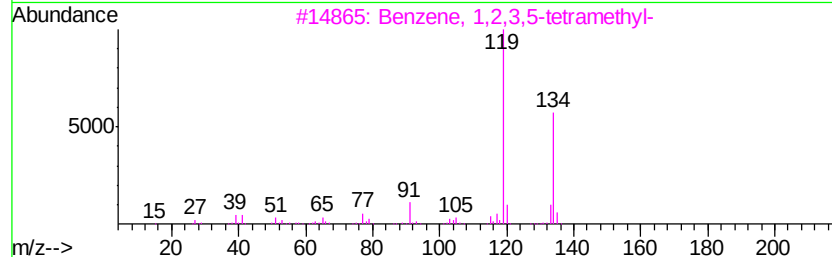
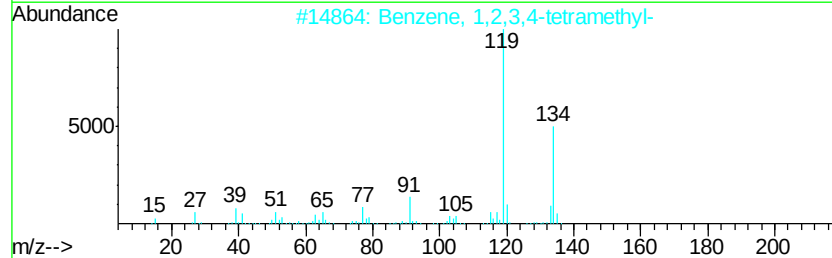
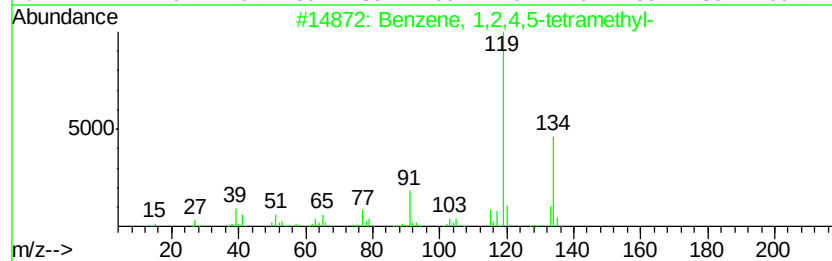
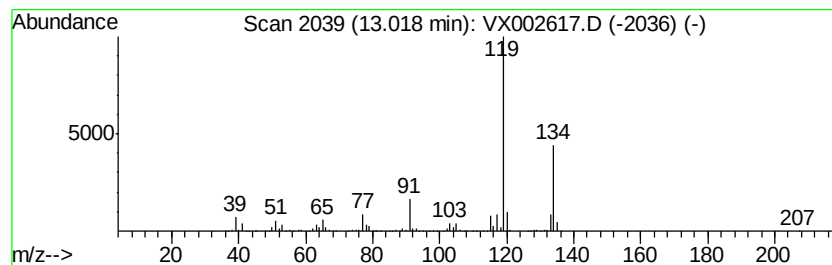
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	31.56 ug/l	765818	1,4-Dichlorobenzene-d4	12.09

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	97
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061518\
Data File : VX002617.D
Acq On : 16 Jun 2018 00:35
Operator : JC/MD
Sample : J3467-05
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PR-MW-05-061118

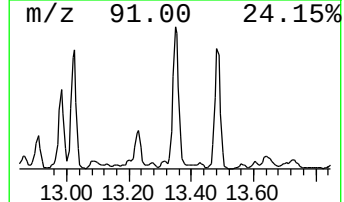
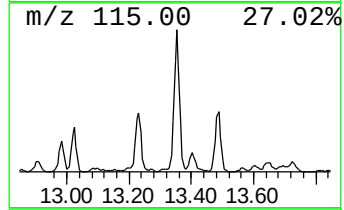
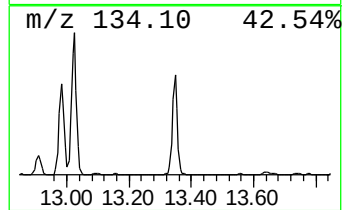
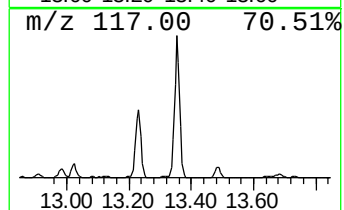
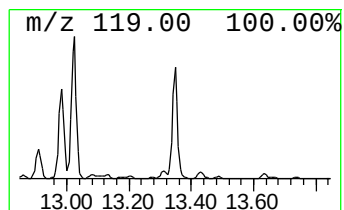
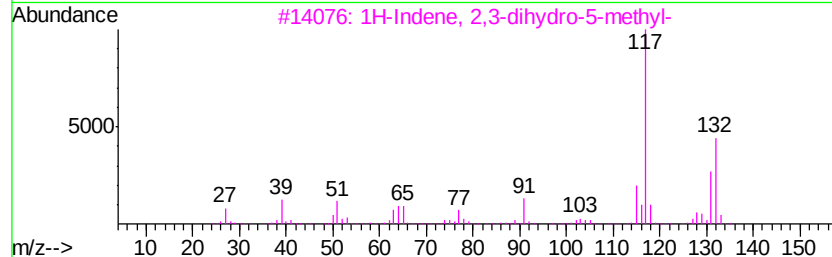
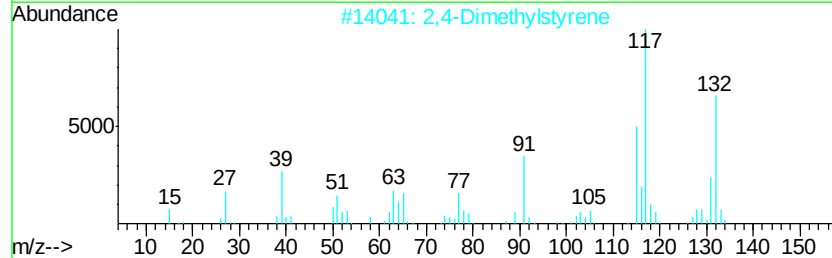
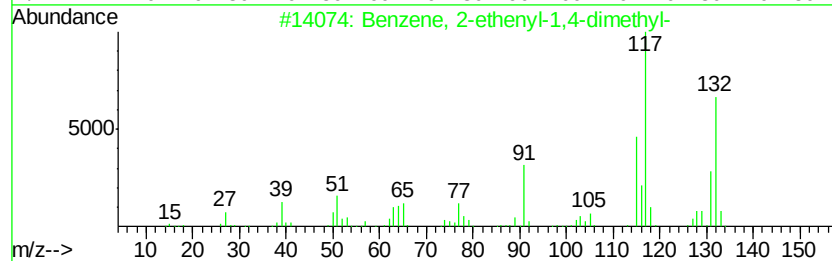
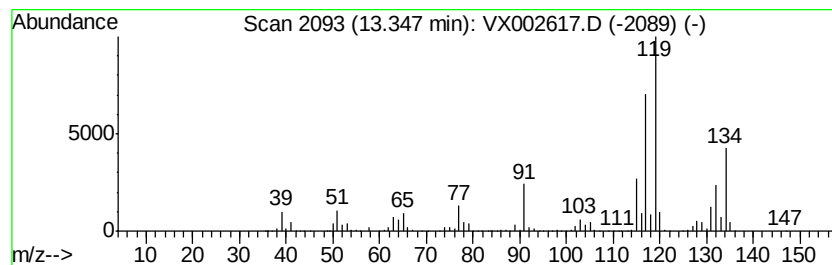
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061518W.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	49.12 ug/l	1191730	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2		2,4-Dimethylstyrene	132	C10H12	002234-20-0	95
3		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
4		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70
5		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061518\
Data File : VX002617.D
Acq On : 16 Jun 2018 00:35
Operator : JC/MD
Sample : J3467-05
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 26 Sample Multiplier: 1

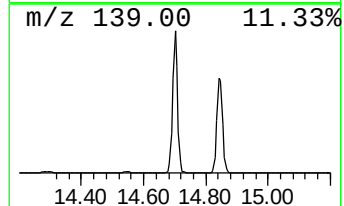
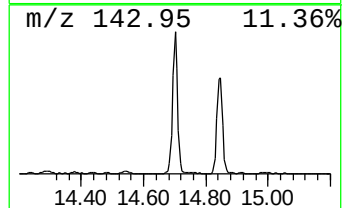
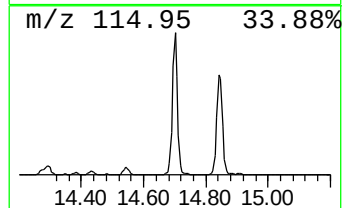
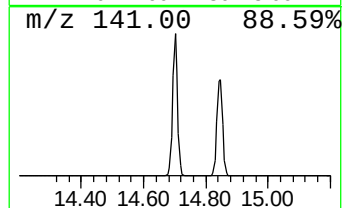
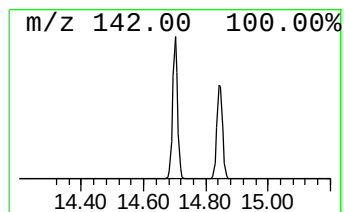
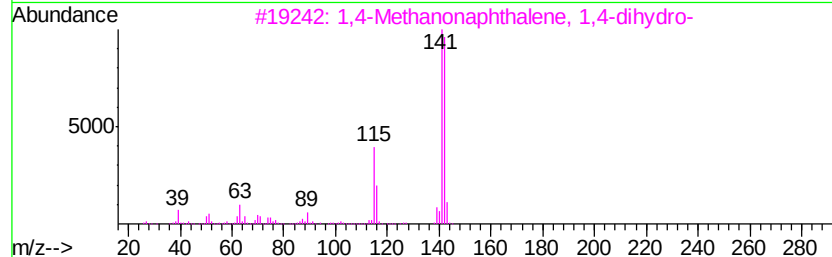
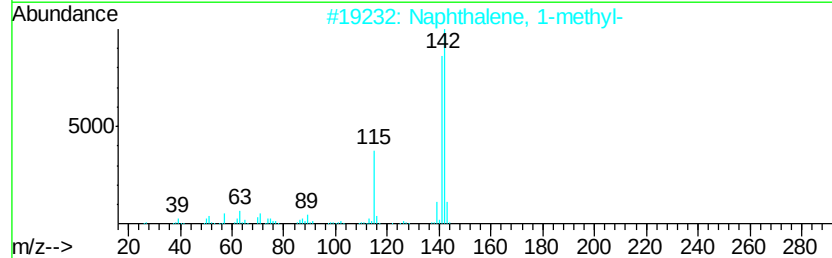
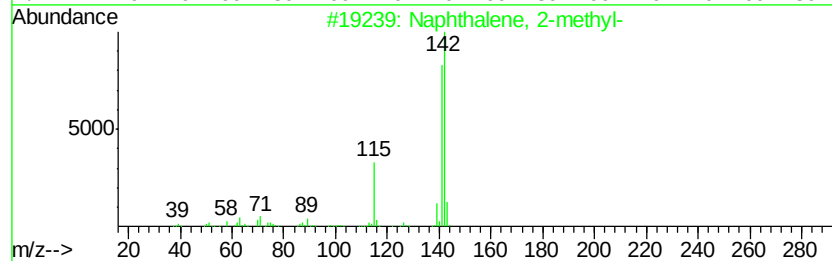
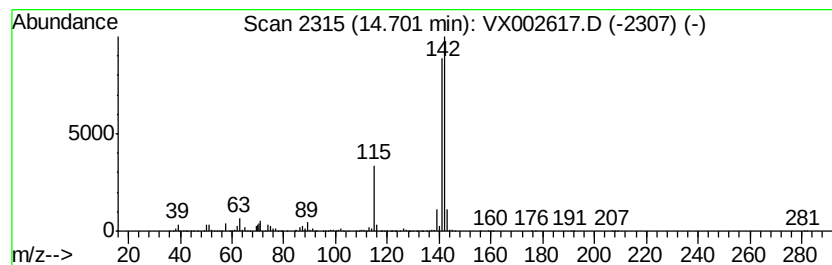
Instrument :
MSVOA_X
ClientSampleId :
PR-MW-05-061118

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061518W.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.70	42.14 ug/l	1022330	1,4-Dichlorobenzene-d4	12.09
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		1,4-Methanonaphthalene, 1,4-dihv...	142 C11H10	004453-90-1 91
4		Benzocycloheptatriene	142 C11H10	000264-09-5 91
5		1H-Indene, 1-ethylidene-	142 C11H10	002471-83-2 50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX061518\
Data File : VX002617.D
Acq On : 16 Jun 2018 00:35
Operator : JC/MD
Sample : J3467-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PR-MW-05-061118

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061518W.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-	1.79	35.0	ug/l	605924	1	5.67	865895	50.0
Cyclopentane, met...	4.40	35.3	ug/l	611055	1	5.67	865895	50.0
Benzene, 1-ethyl-...	11.44	349.6	ug/l	8482910	4	12.09	1213150	50.0
Benzene, 1,2,3-tr...	12.15	113.9	ug/l	2763270	4	12.09	1213150	50.0
Indane	12.30	138.5	ug/l	3361190	4	12.09	1213150	50.0
Benzene, 1,2,4,5-...	13.02	31.6	ug/l	765818	4	12.09	1213150	50.0
Benzene, 2-etheny...	13.35	49.1	ug/l	1191730	4	12.09	1213150	50.0
Naphthalene, 1-me...	14.70	42.1	ug/l	1022330	4	12.09	1213150	50.0