

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091818\
 Data File : VX004654.D
 Acq On : 18 Sep 2018 13:35
 Operator : JC/MD
 Sample : J4898-01 10X
 Misc : 5.26g/5mL/100uL/5.0mL/MSVOA_X/MEOH
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 ROLLOFF01-COMP-FILL

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\82X091118W.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	5.659	727	739	755	rVB	176445	493753	1.34%	0.537%
2	6.141	811	818	831	rVB	488064	1243656	3.37%	1.352%
3	6.860	925	936	948	rBV	299776	682488	1.85%	0.742%
4	8.720	1233	1241	1246	rBV	634934	1020109	2.76%	1.109%
5	8.787	1246	1252	1265	rVB	1705659	2568062	6.96%	2.791%
6	10.116	1464	1470	1479	rBV	577685	778794	2.11%	0.846%
7	10.250	1487	1492	1502	rBV	958531	1287140	3.49%	1.399%
8	10.360	1502	1510	1516	rBV	1882548	2500672	6.77%	2.718%
9	10.701	1561	1566	1580	rVB	689085	892670	2.42%	0.970%
10	11.140	1632	1638	1649	rBV	609444	772143	2.09%	0.839%
11	11.433	1681	1686	1693	rBV	256670	434033	1.18%	0.472%
12	11.506	1693	1698	1710	rVB	402701	617939	1.67%	0.672%
13	11.811	1742	1748	1752	rVB	936348	1115455	3.02%	1.212%
14	11.865	1752	1757	1760	rBV	418877	521003	1.41%	0.566%
15	12.018	1775	1782	1787	rBV	687731	908744	2.46%	0.988%
16	12.079	1787	1792	1798	rVB	810722	997730	2.70%	1.084%
17	12.469	1847	1856	1862	rBV	4986725	6361978	17.23%	6.914%
18	13.024	1942	1947	1951	rBV	850642	1181403	3.20%	1.284%
19	13.402	2004	2009	2013	rVV	913218	1083043	2.93%	1.177%
20	13.481	2017	2022	2028	rVB	945142	1227474	3.32%	1.334%
21	13.841	2072	2081	2086	rBV	27218382	36923229	100.00%	40.128%
22	13.932	2092	2096	2102	rVB	840648	1168715	3.17%	1.270%
23	14.700	2217	2222	2232	rVB	8107200	9905101	26.83%	10.765%
24	14.840	2239	2245	2257	rVB	4046604	5165286	13.99%	5.614%
25	15.273	2306	2316	2321	rBV	986543	1338845	3.63%	1.455%
26	15.548	2355	2361	2372	rBV	1246762	2136743	5.79%	2.322%
27	15.688	2372	2384	2388	rVV	1409574	2384868	6.46%	2.592%
28	15.724	2388	2390	2397	rVB	850165	1185230	3.21%	1.288%
29	15.920	2412	2422	2433	rVB	453705	1113898	3.02%	1.211%
30	16.072	2441	2447	2456	rVB	233404	407002	1.10%	0.442%
31	16.176	2456	2464	2476	rBV2	1436695	2497045	6.76%	2.714%
32	16.419	2494	2504	2514	rBV2	535278	1100451	2.98%	1.196%

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Integration Parameters: RTEINT.P

Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Title : SW846 8260

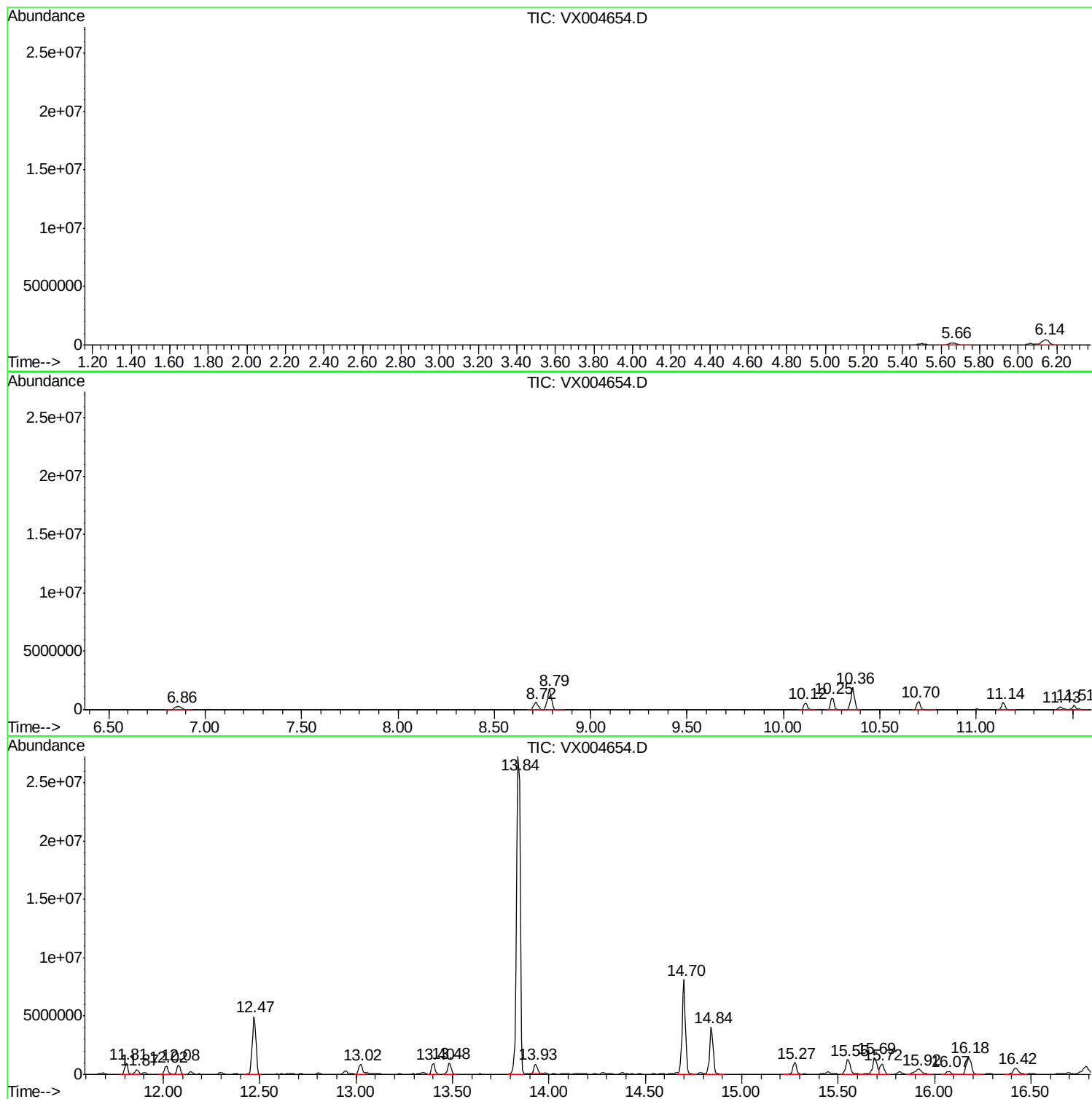
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Quant Title : SW846 8260

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



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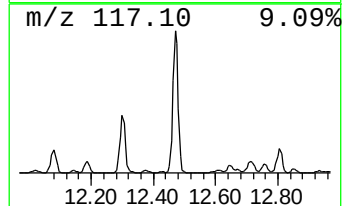
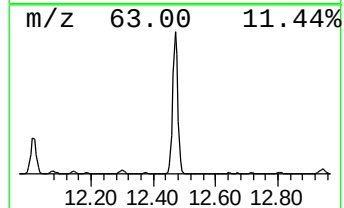
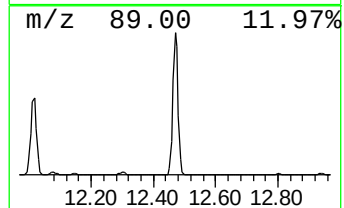
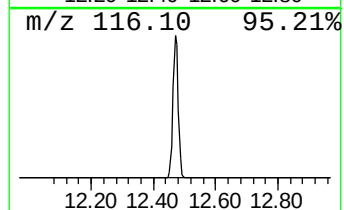
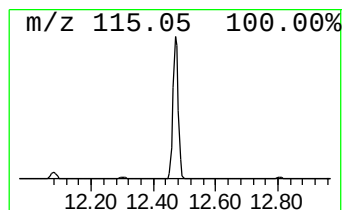
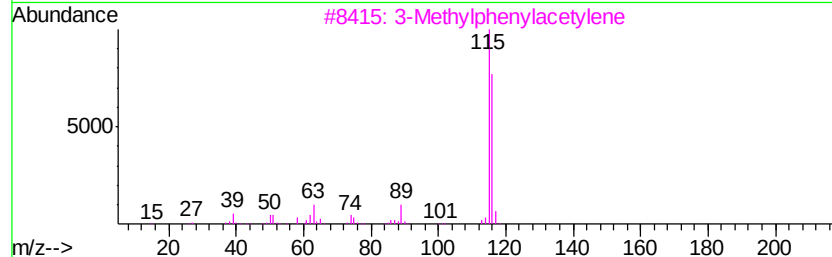
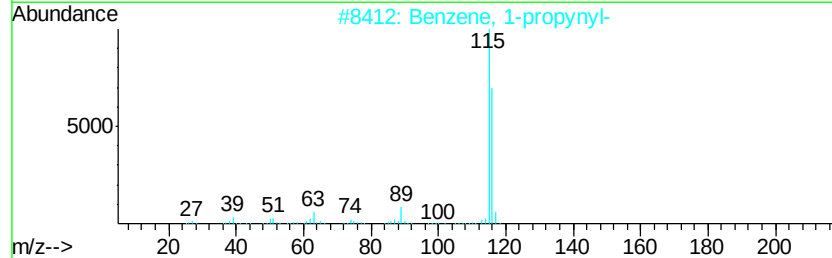
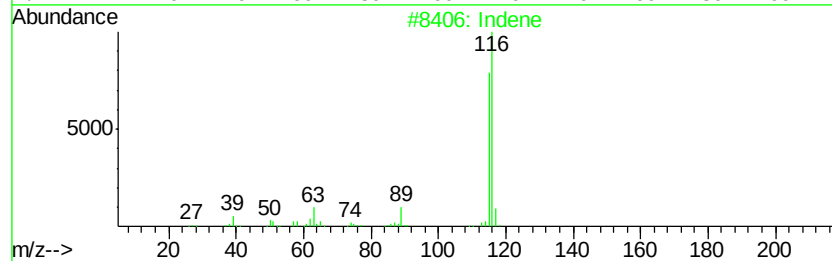
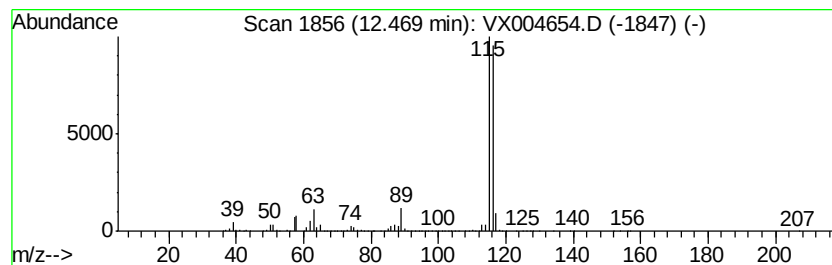
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 1 Indene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.47	318.82 ug/l	6361980	1,4-Dichlorobenzene-d4	12.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene		116	C9H8	000095-13-6	97
2	Benzene, 1-propynyl-		116	C9H8	000673-32-5	94
3	3-Methylphenylacetylene		116	C9H8	000766-82-5	91
4	Benzene, 1-ethynyl-4-methyl-		116	C9H8	000766-97-2	91
5	2-Methylphenylacetylene		116	C9H8	000766-47-2	83



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MSVOA_X
ClientSampleId :
ROLLOFF01-COMP-FILL

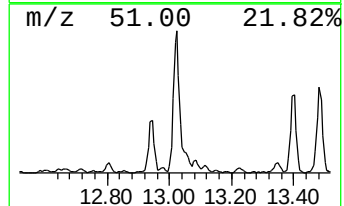
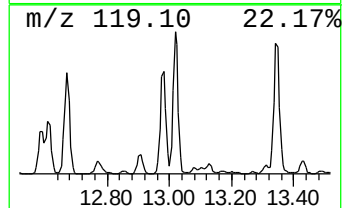
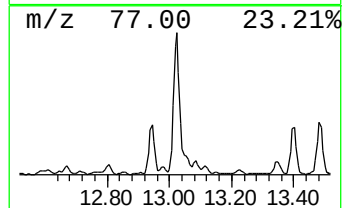
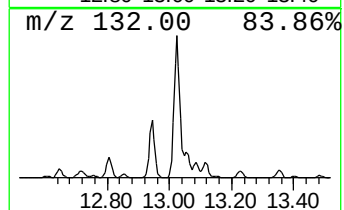
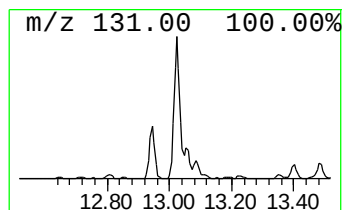
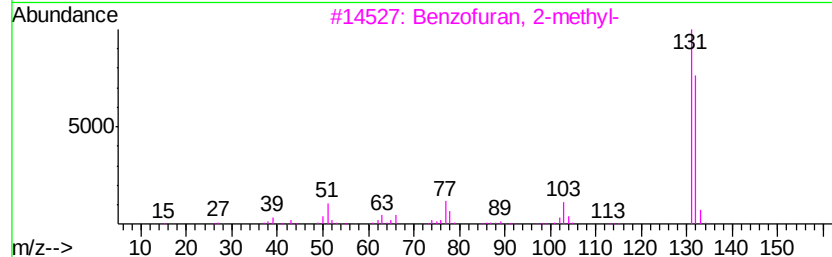
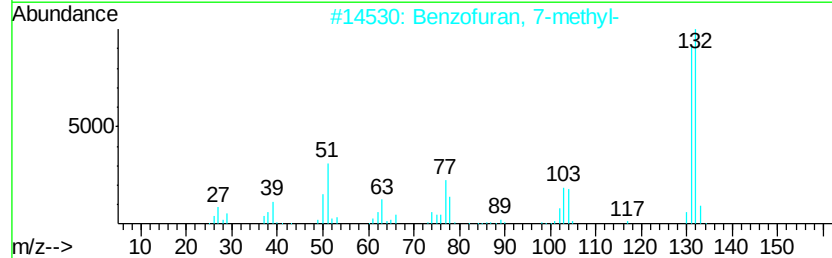
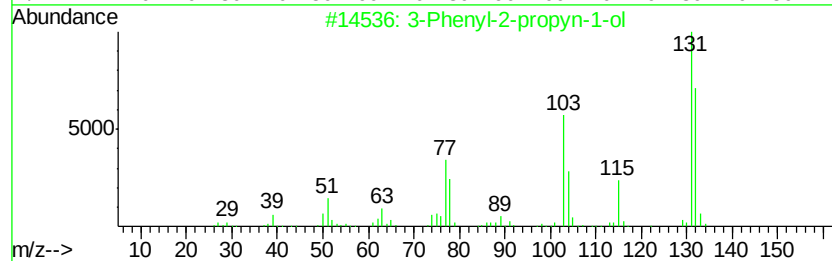
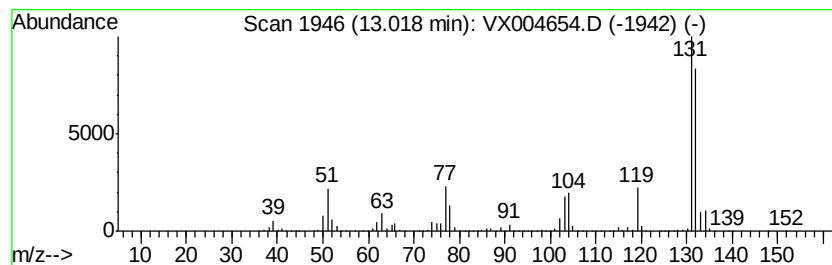
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 2 3-Phenyl-2-propyn-1-ol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	59.20 ug/l	1181400	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	93
2		Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	93
3		Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90
4		Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	86
5		2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	83



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ALS Vial : 11 Sample Multiplier: 1

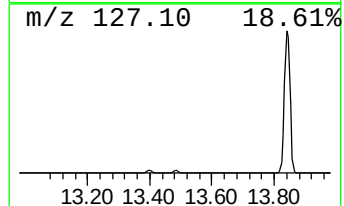
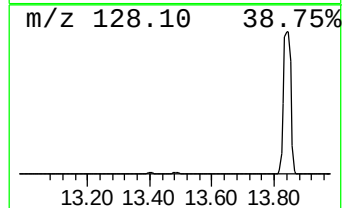
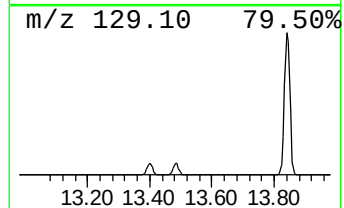
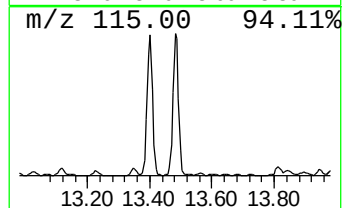
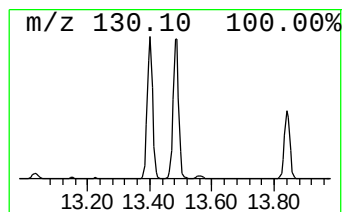
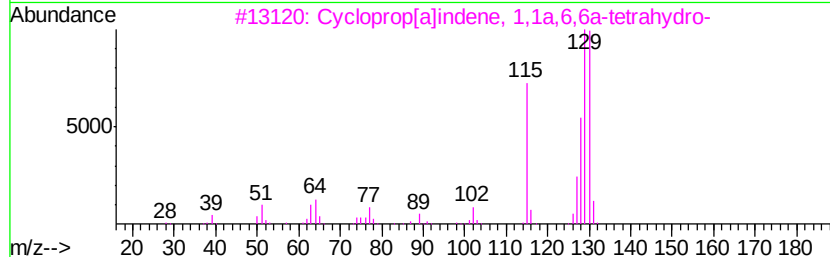
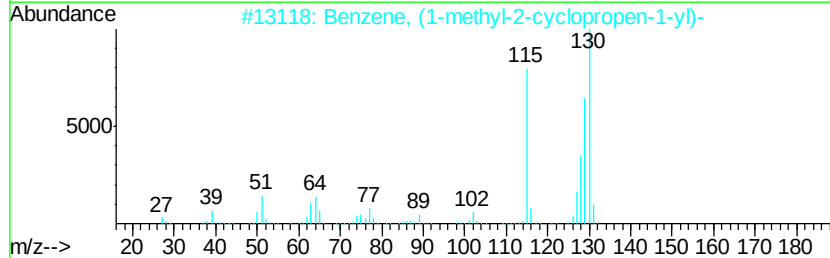
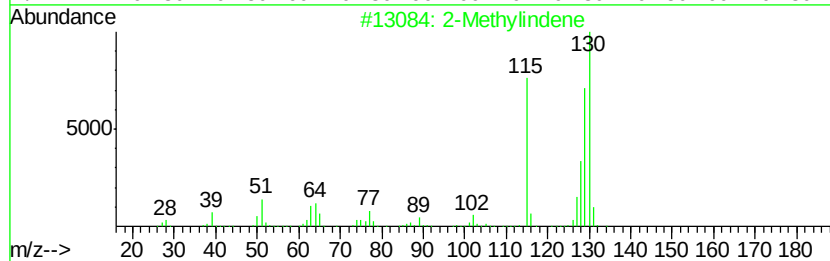
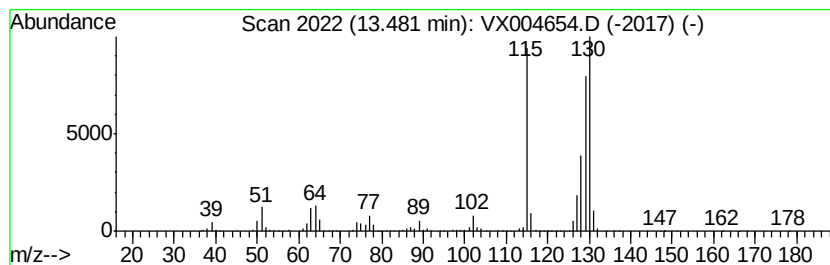
Instrument :
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ClientSampleId :
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Methylindene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.48	61.51 ug/l	1227470	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2-Methylindene	130 C10H10	002177-47-1	97
2	Benzene, (1-methyl-2-cyclopropen...	130 C10H10	065051-83-4	97
3	Cycloprop[alindene, 1,1a,6,6a-te...	130 C10H10	015677-15-3	95
4	1,4-Dihydronaphthalene	130 C10H10	000612-17-9	95
5	1H-Indene, 1-methyl-	130 C10H10	000767-59-9	94



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Misc : 5.26g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
ROLLOFF01-COMP-FILL

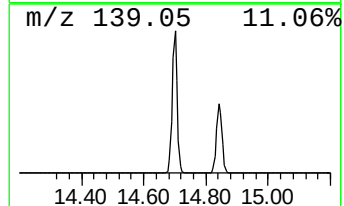
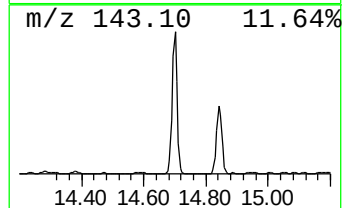
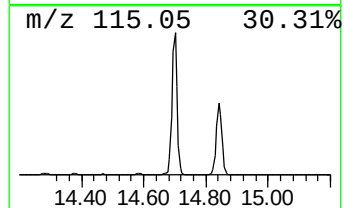
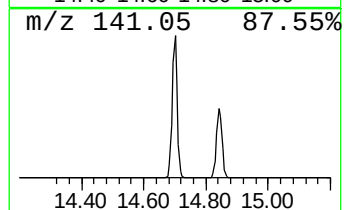
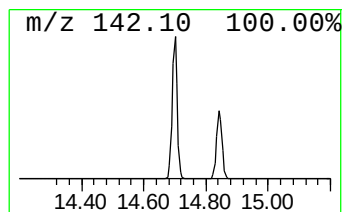
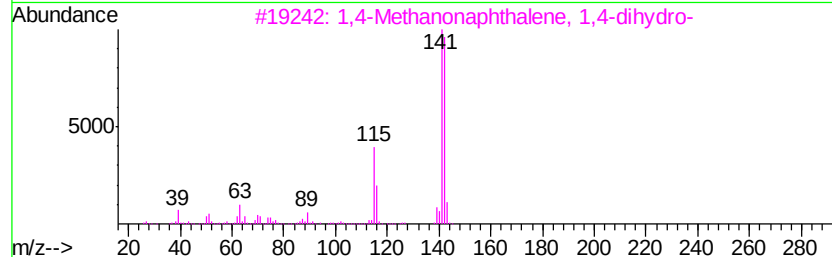
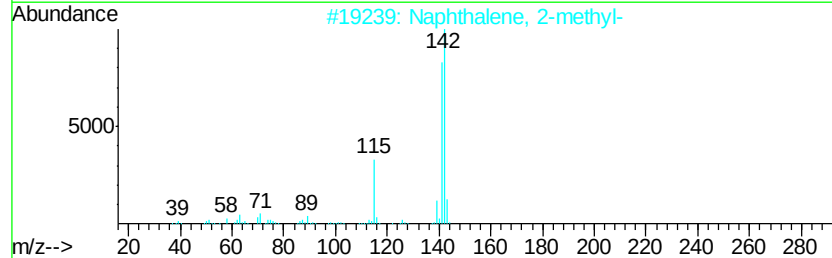
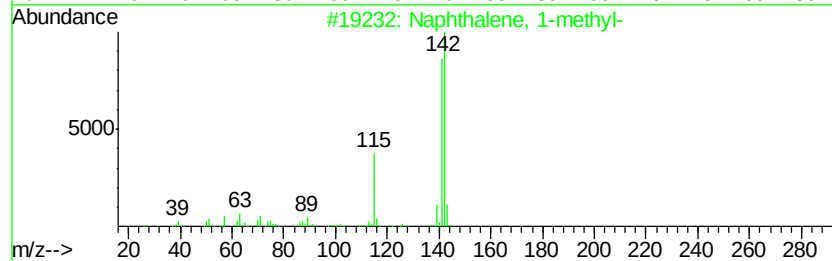
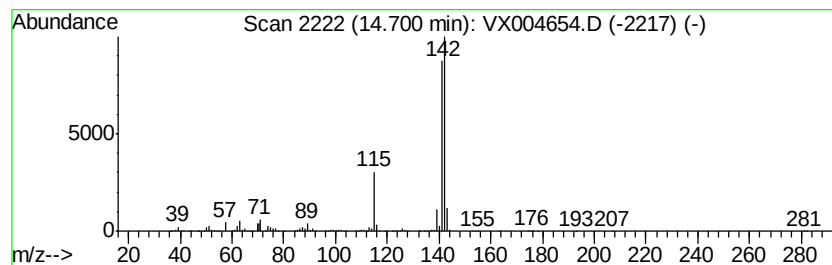
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.70	496.38 ug/l	9905100	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	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	64



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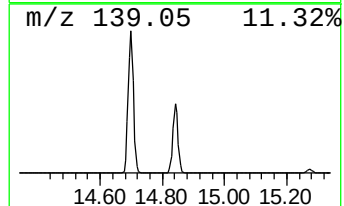
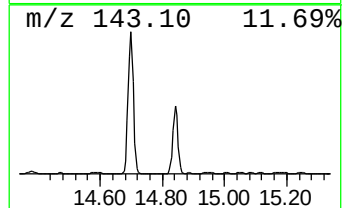
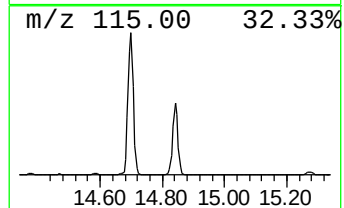
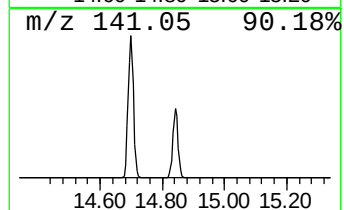
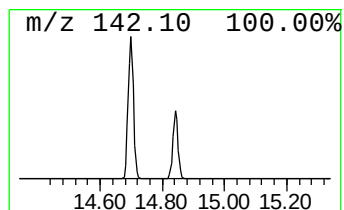
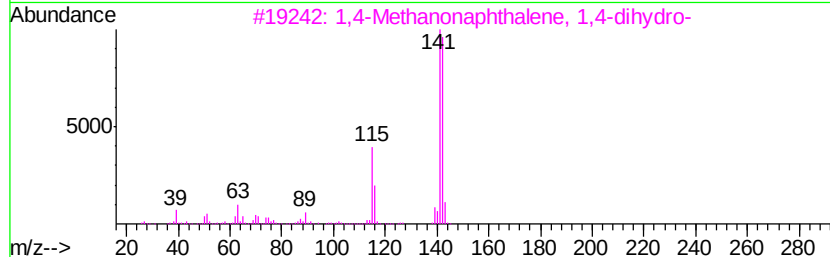
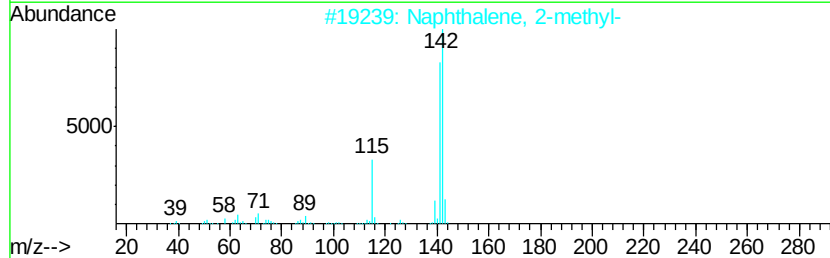
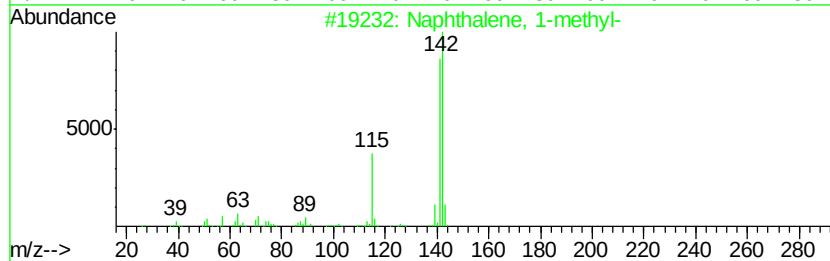
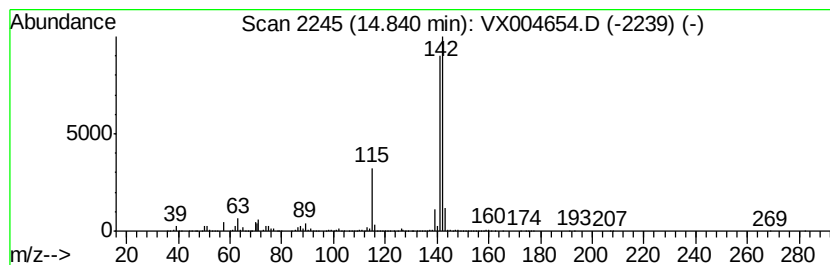
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Naphthalene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.84	258.85 ug/l	5165290	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	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	49



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091818\
Data File : VX004654.D
Acq On : 18 Sep 2018 13:35
Operator : JC/MD
Sample : J4898-01 10X
Misc : 5.26g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 11 Sample Multiplier: 1

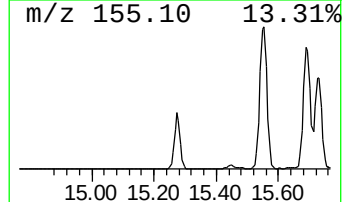
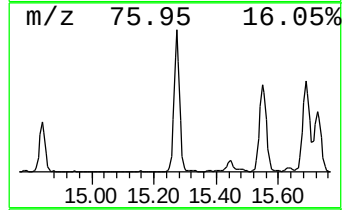
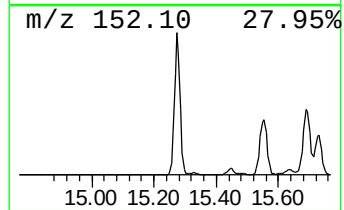
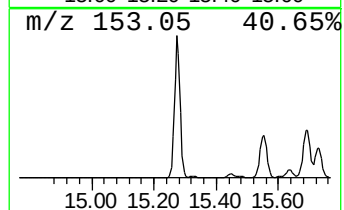
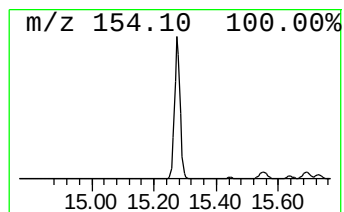
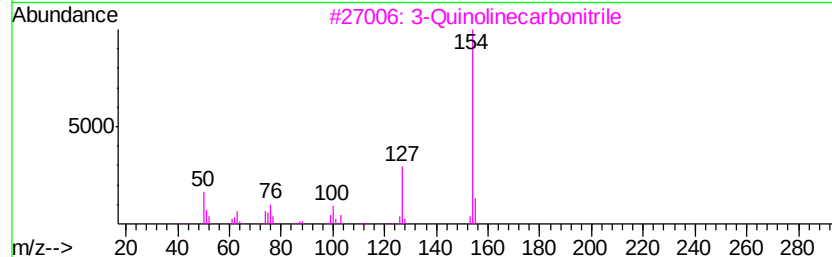
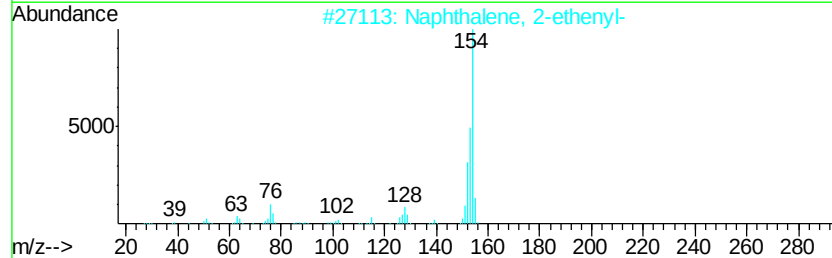
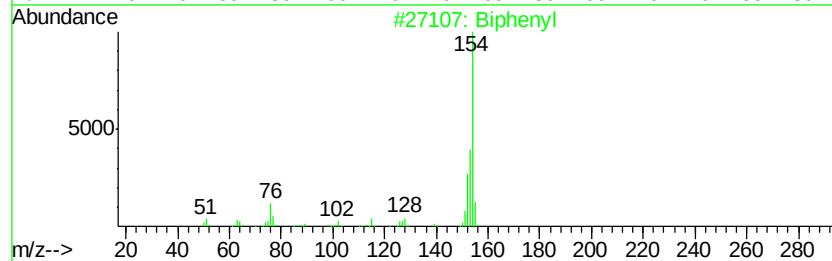
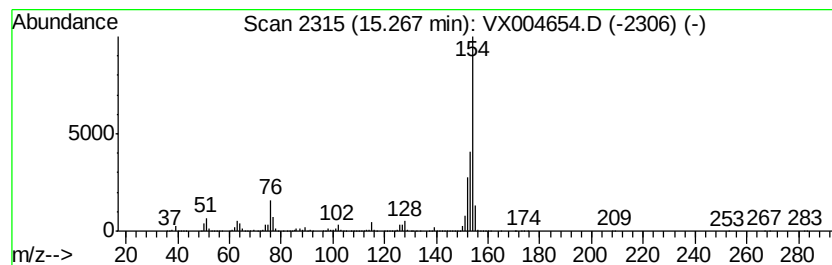
Instrument :
MSVOA_X
ClientSampleId :
ROLLOFF01-COMP-FILL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 6 Biphenyl Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.27	67.09 ug/l	1338850	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Biphenyl	154 C12H10	000092-52-4	95
2	Naphthalene, 2-ethenyl-	154 C12H10	000827-54-3	87
3	3-Quinolinecarbonitrile	154 C10H6N2	034846-64-5	47
4	Acenaphthene	154 C12H10	000083-32-9	47
5	1-(2,2-Difluoroethenyl)-4-methyl...	154 C9H8F2	1000305-64-2	45



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091818\
Data File : VX004654.D
Acq On : 18 Sep 2018 13:35
Operator : JC/MD
Sample : J4898-01 10X
Misc : 5.26µ/5mL/100µL/5.0mL/MSVOA_X/MEOH
ALS Vial : 11 Sample Multiplier: 1

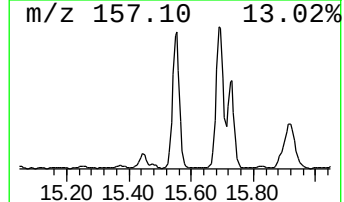
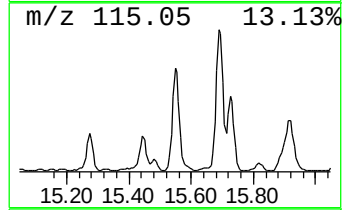
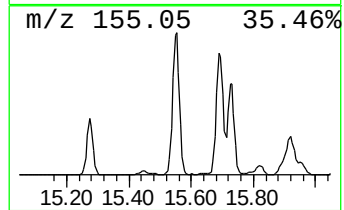
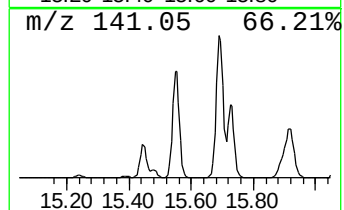
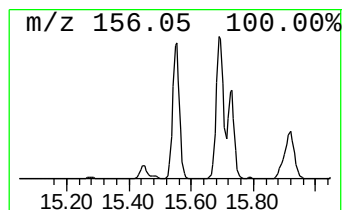
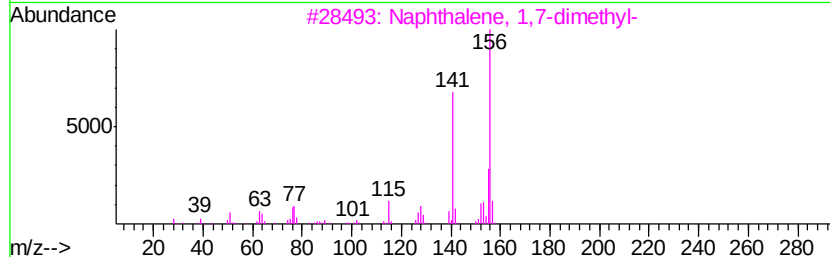
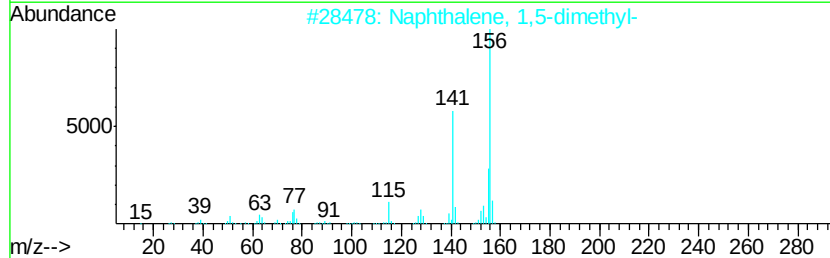
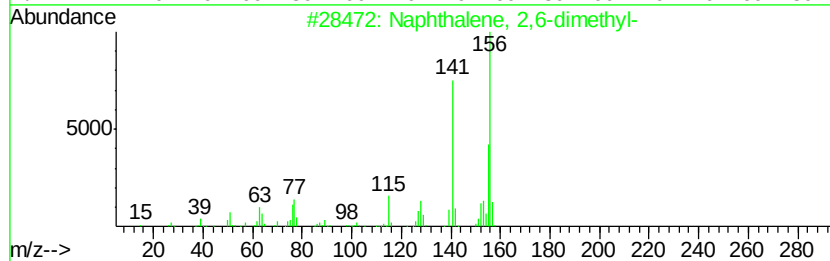
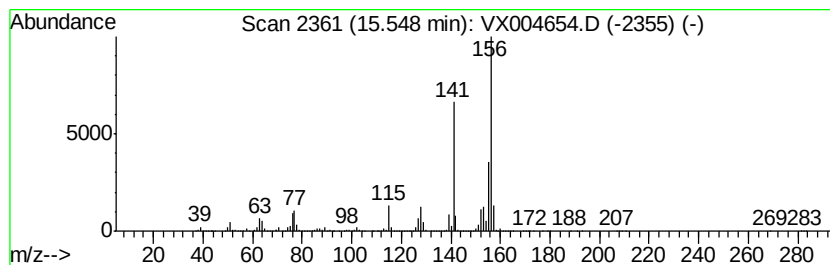
Instrument :
MSVOA_X
ClientSampleId :
ROLLOFF01-COMP-FILL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 2,6-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.55	107.08 ug/l	2136740	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 98
2		Naphthalene, 1,5-dimethyl-	156 C12H12	000571-61-9 97
3		Naphthalene, 1,7-dimethyl-	156 C12H12	000575-37-1 97
4		Naphthalene, 2,7-dimethyl-	156 C12H12	000582-16-1 97
5		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 97



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091818\
Data File : VX004654.D
Acq On : 18 Sep 2018 13:35
Operator : JC/MD
Sample : J4898-01 10X
Misc : 5.26g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 11 Sample Multiplier: 1

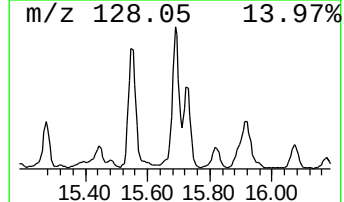
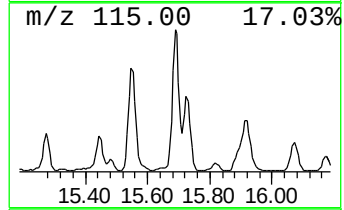
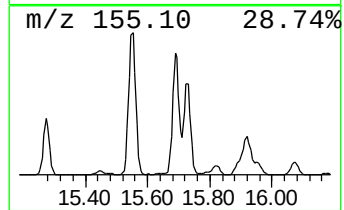
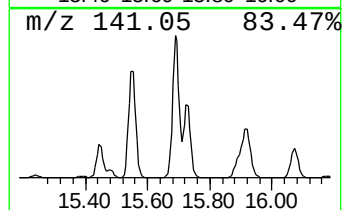
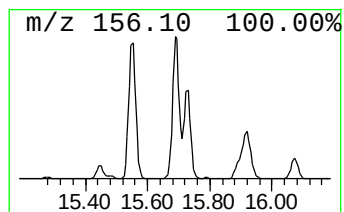
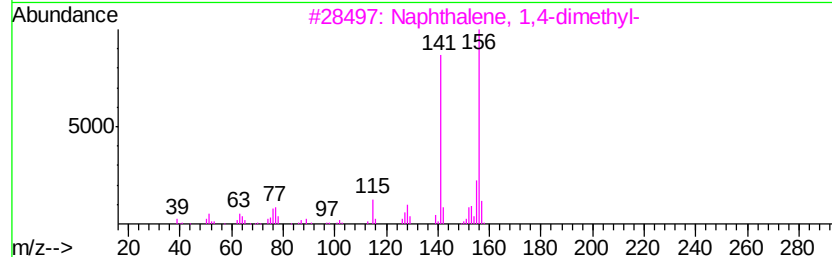
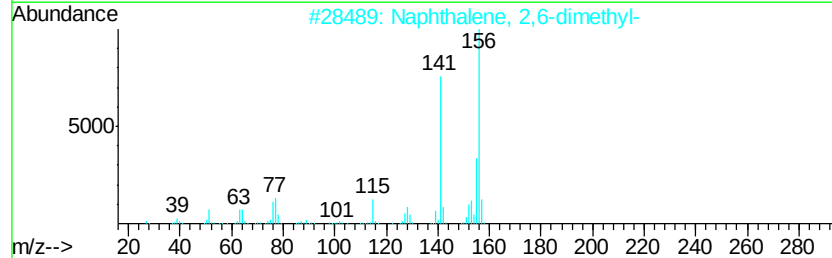
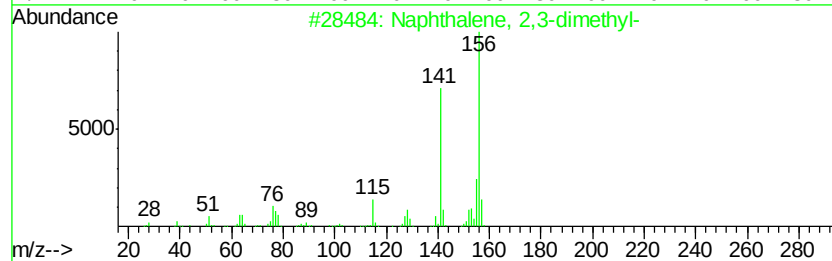
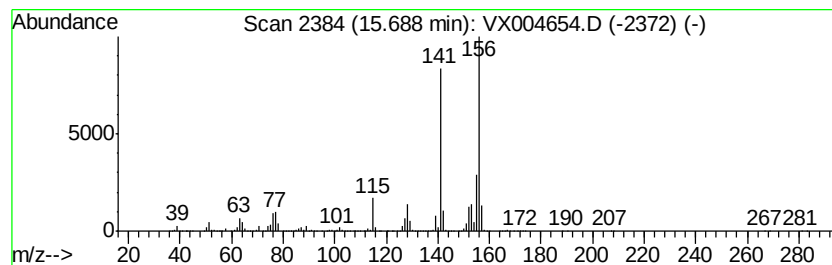
Instrument :
MSVOA_X
ClientSampleId :
ROLLOFF01-COMP-FILL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 2,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.69	119.52 ug/l	2384870	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 97
2		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 97
3		Naphthalene, 1,4-dimethyl-	156 C12H12	000571-58-4 97
4		Naphthalene, 1,3-dimethyl-	156 C12H12	000575-41-7 97
5		Naphthalene, 1,7-dimethyl-	156 C12H12	000575-37-1 96



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091818\
Data File : VX004654.D
Acq On : 18 Sep 2018 13:35
Operator : JC/MD
Sample : J4898-01 10X
Misc : 5.26g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
ROLLOFF01-COMP-FILL

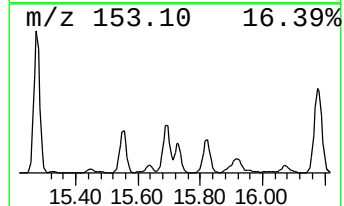
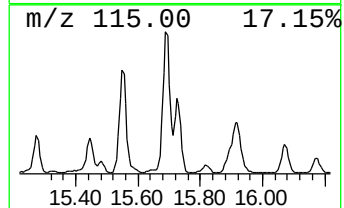
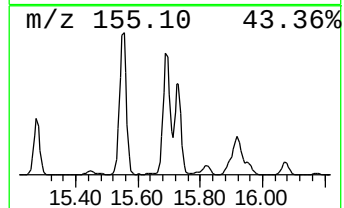
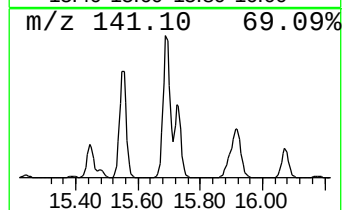
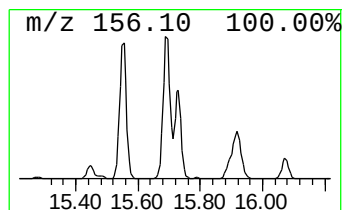
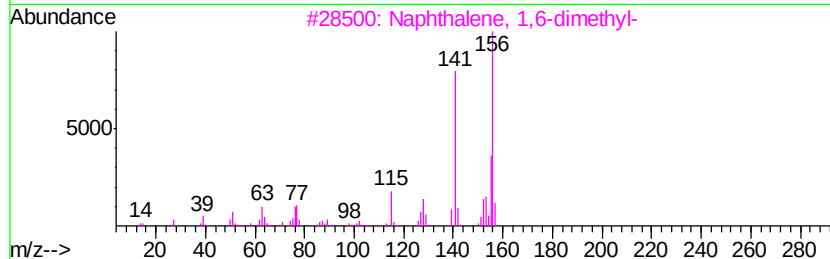
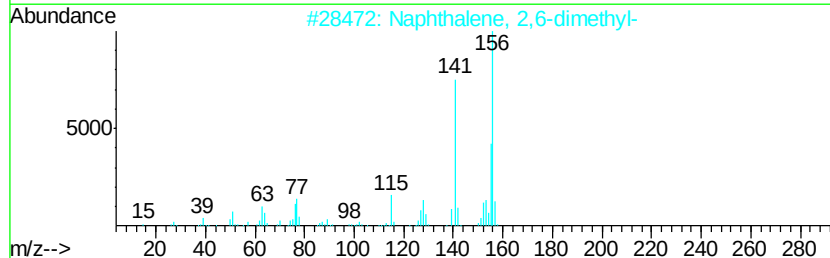
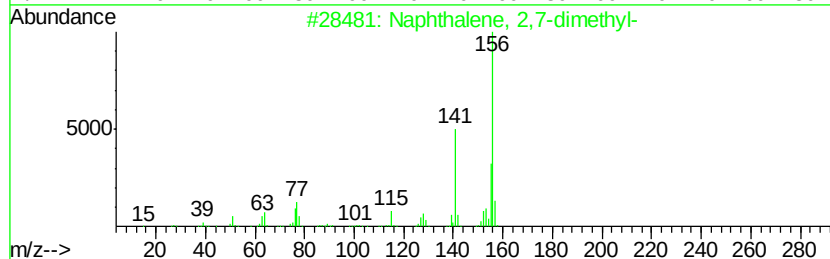
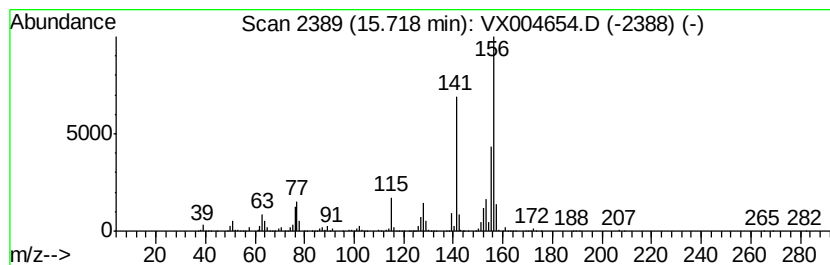
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 2,7-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.72	59.40 ug/l	1185230	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
2		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
5		Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	96



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091818\
Data File : VX004654.D
Acq On : 18 Sep 2018 13:35
Operator : JC/MD
Sample : J4898-01 10X
Misc : 5.26g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
ROLLOFF01-COMP-FILL

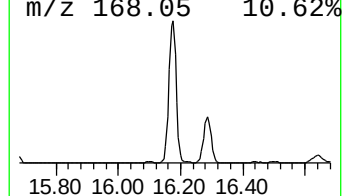
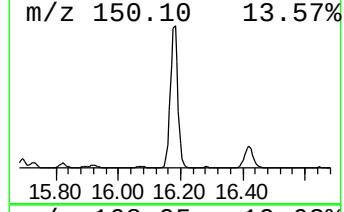
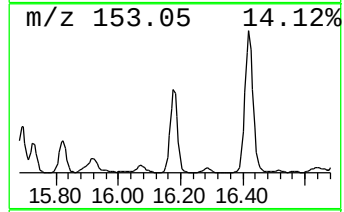
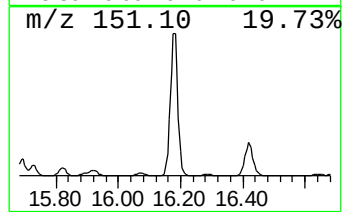
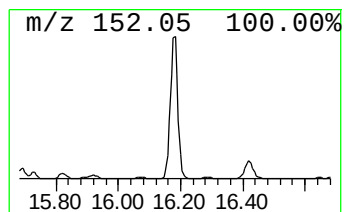
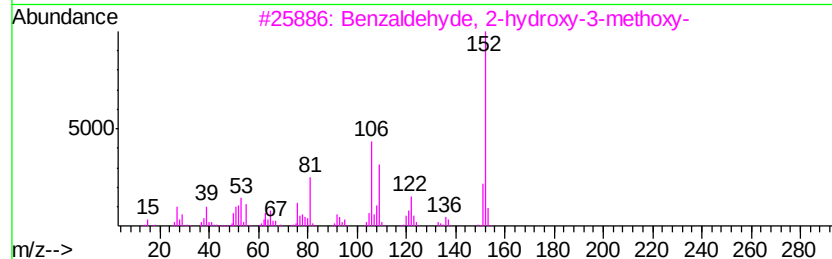
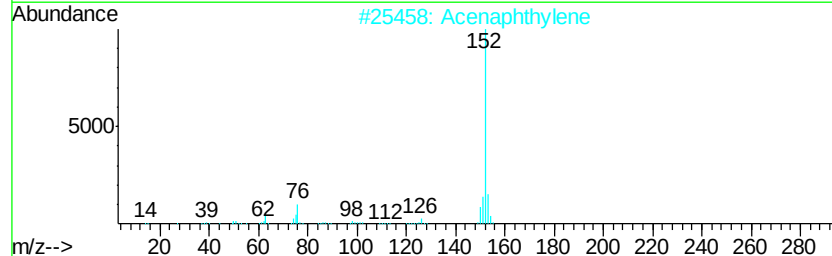
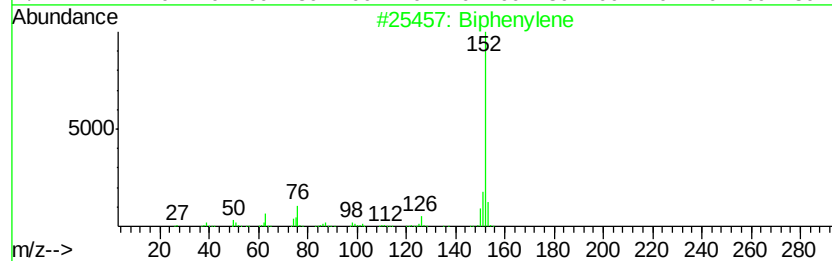
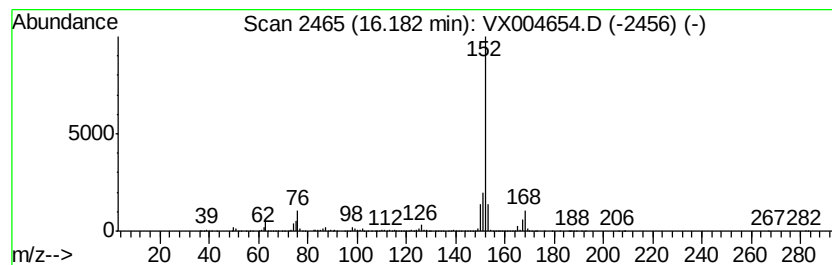
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 10 Biphenylene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.18	125.14 ug/l	2497050	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Biphenylene	152	C12H8	000259-79-0	93
2		Acenaphthylene	152	C12H8	000208-96-8	91
3		Benzaldehyde, 2-hydroxy-3-methoxy-	152	C8H8O3	000148-53-8	59
4		5,6-Dihydro-4-methylthieno(2,3-d...	152	C7H8N2S	092204-06-3	53
5		1-acenaphthenol, trifluoroacetat...	266	C14H9F3O2	1000376-45-2	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX091818\
Data File : VX004654.D
Acq On : 18 Sep 2018 13:35
Operator : JC/MD
Sample : J4898-01 10X
Misc : 5.26g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
ROLLOFF01-COMP-FILL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.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
Indene	12.47	318.8	ug/l	6361980	4	12.08	997730	50.0
3-Phenyl-2-propyn...	13.02	59.2	ug/l	1181400	4	12.08	997730	50.0
2-Methylindene	13.48	61.5	ug/l	1227470	4	12.08	997730	50.0
Naphthalene, 1-me...	14.70	496.4	ug/l	9905100	4	12.08	997730	50.0
Naphthalene, 2-me...	14.84	258.9	ug/l	5165290	4	12.08	997730	50.0
Biphenyl	15.27	67.1	ug/l	1338850	4	12.08	997730	50.0
Naphthalene, 2,6-...	15.55	107.1	ug/l	2136740	4	12.08	997730	50.0
Naphthalene, 2,3-...	15.69	119.5	ug/l	2384870	4	12.08	997730	50.0
Naphthalene, 2,7-...	15.72	59.4	ug/l	1185230	4	12.08	997730	50.0
Biphenylene	16.18	125.1	ug/l	2497050	4	12.08	997730	50.0