

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080122\
 Data File : VX030322.D
 Acq On : 01 Aug 2022 15:07
 Operator : JC/MD
 Sample : N3964-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 HMGP-WATER-1

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\82X072522W.M
 Title : SW846 8260

Signal : TIC: VX030322.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.069	153	161	180	rBV	1861709	2960075	35.92%	11.620%
2	2.703	258	265	277	rBV	1380859	2413010	29.28%	9.472%
3	3.117	320	333	350	rBV	524252	1208298	14.66%	4.743%
4	4.721	583	596	615	rBV	2922179	8239955	100.00%	32.347%
5	5.391	695	706	722	rBV2	130393	385990	4.68%	1.515%
6	5.556	722	733	746	rVB	126873	352559	4.28%	1.384%
7	5.958	788	799	809	rBV2	166364	440098	5.34%	1.728%
8	6.769	921	932	950	rBV	230622	585612	7.11%	2.299%
9	8.653	1234	1241	1249	rBV	672952	1034950	12.56%	4.063%
10	10.055	1465	1471	1481	rBV	506951	677323	8.22%	2.659%
11	10.305	1504	1512	1519	rBV	115077	160600	1.95%	0.630%
12	11.085	1634	1640	1648	rBV	507602	652155	7.91%	2.560%
13	12.024	1788	1794	1801	rBV	641177	781489	9.48%	3.068%
14	12.244	1822	1830	1838	rBV3	59888	109433	1.33%	0.430%
15	12.414	1853	1858	1863	rBV	199987	242332	2.94%	0.951%
16	13.426	2019	2024	2029	rBV2	79523	100850	1.22%	0.396%
17	13.774	2074	2081	2092	rBV	2733741	3558827	43.19%	13.970%
18	14.640	2218	2223	2234	rVB	435528	529575	6.43%	2.079%
19	14.780	2240	2246	2252	rBV	473068	598283	7.26%	2.349%
20	15.207	2310	2316	2322	rBV	91688	120058	1.46%	0.471%
21	15.615	2375	2383	2386	rBV	74481	120971	1.47%	0.475%
22	16.328	2492	2500	2509	rBV	109082	201563	2.45%	0.791%

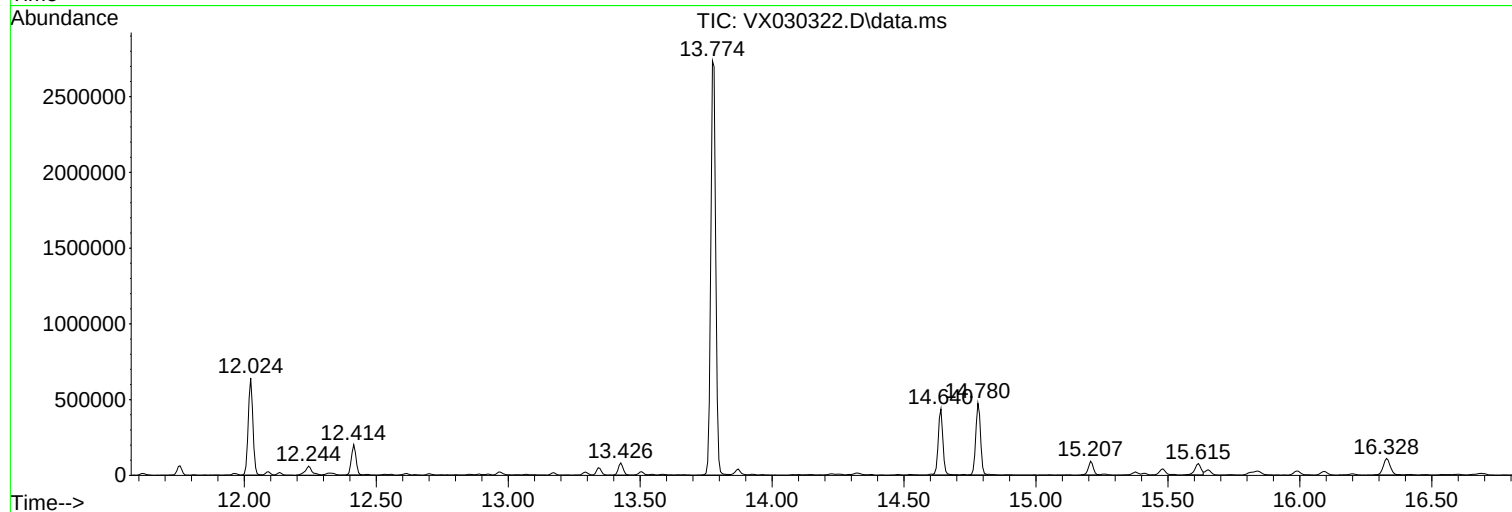
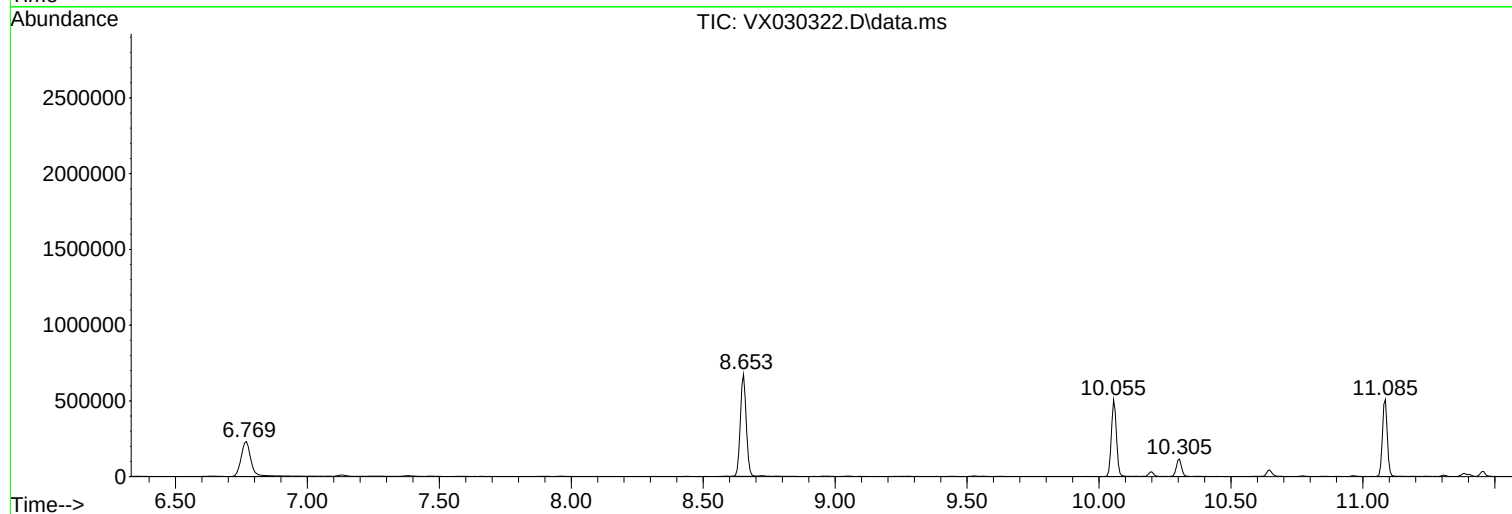
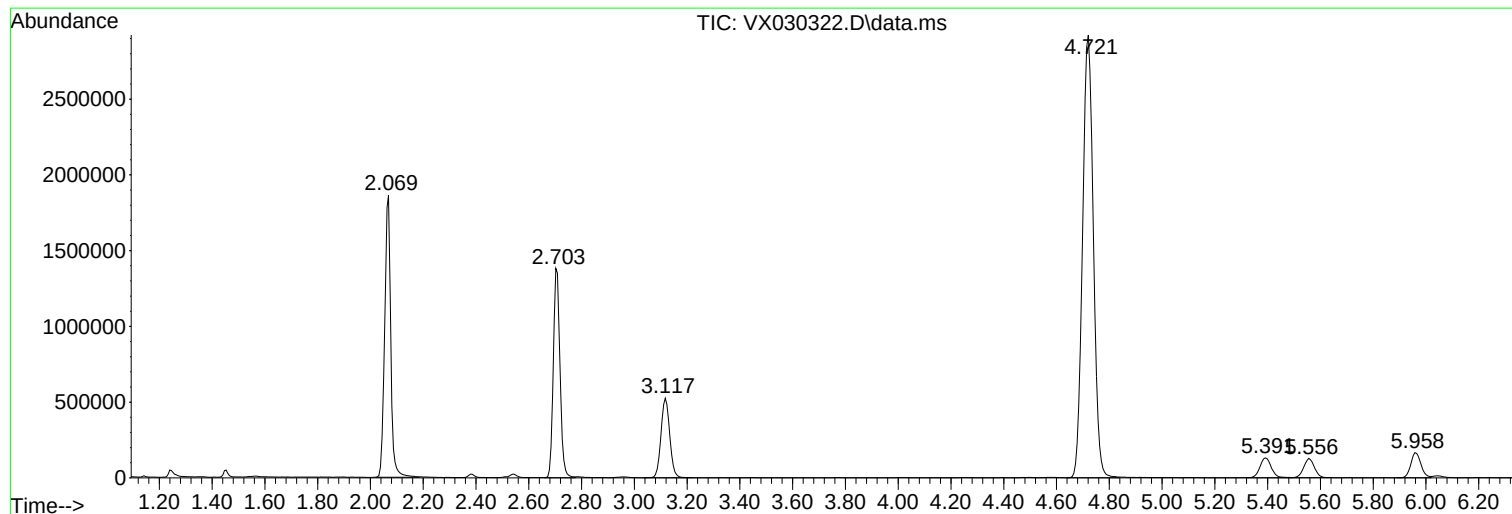
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ClientSampleId :
HMGP-WATER-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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ClientSampleId :
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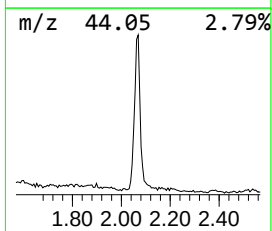
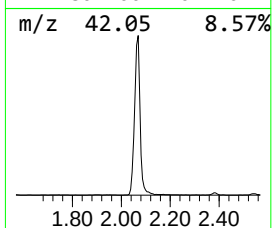
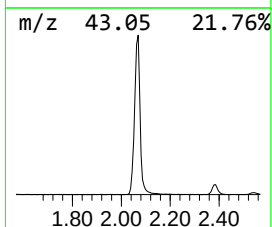
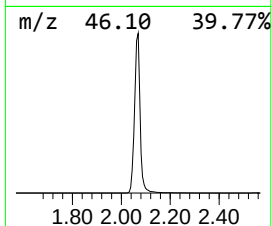
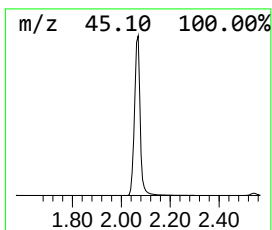
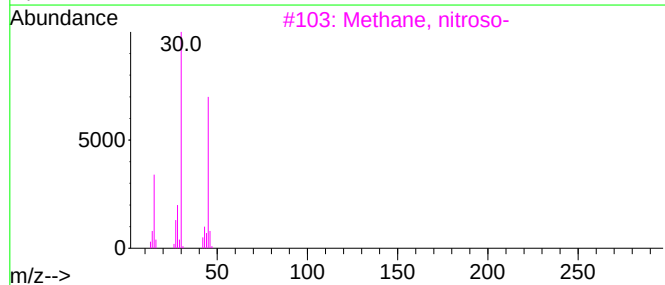
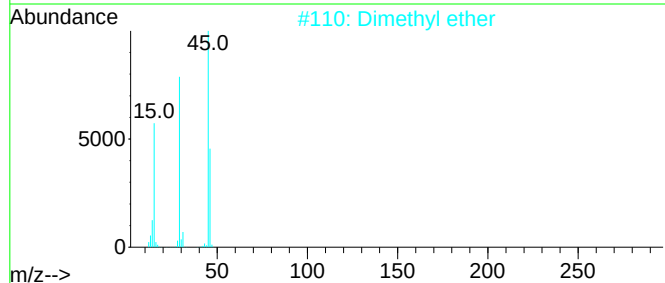
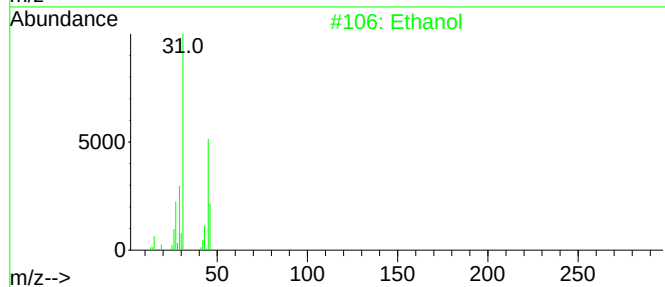
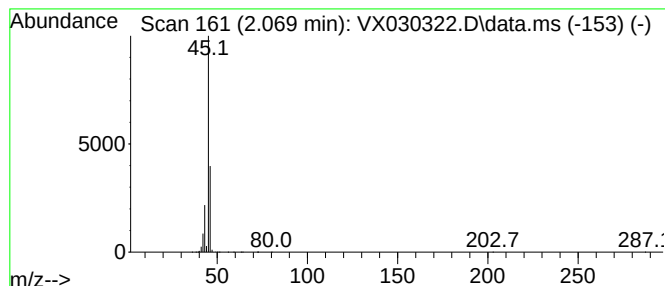
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.069	419.80 ug/l	2960080	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol	46	C2H6O	000064-17-5	64
2			Dimethyl ether	46	C2H6O	000115-10-6	9
3			Methane, nitroso-	45	CH3NO	000865-40-7	4
4			Formic acid	46	CH2O2	000064-18-6	4
5			Oxalic acid	90	C2H2O4	000144-62-7	4



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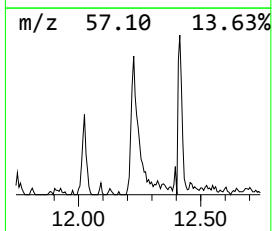
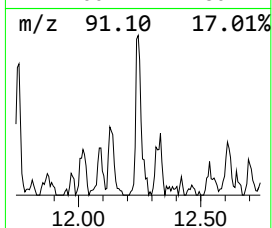
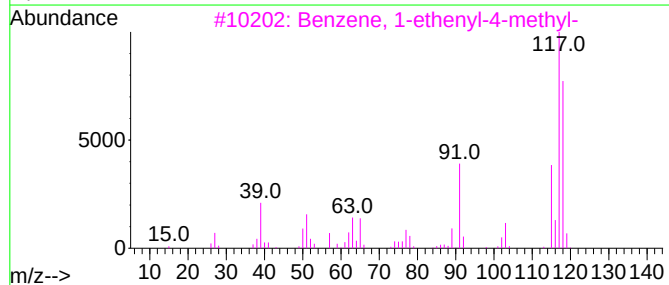
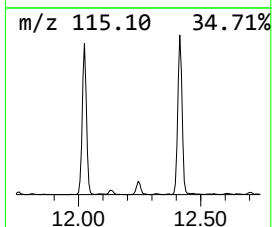
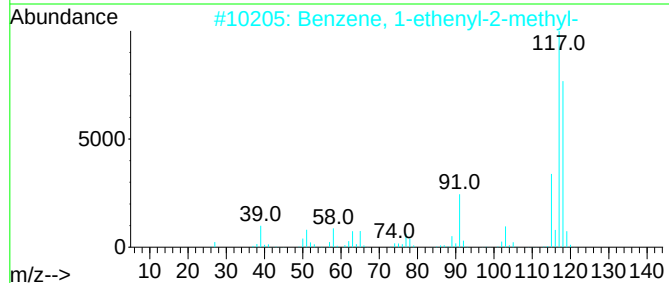
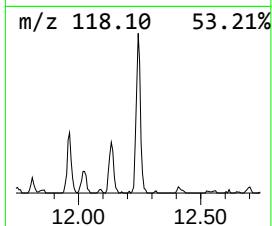
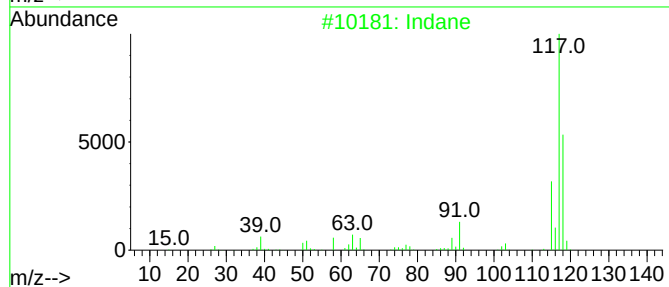
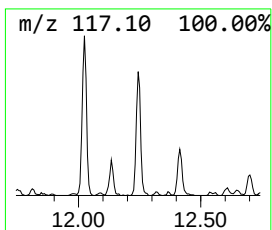
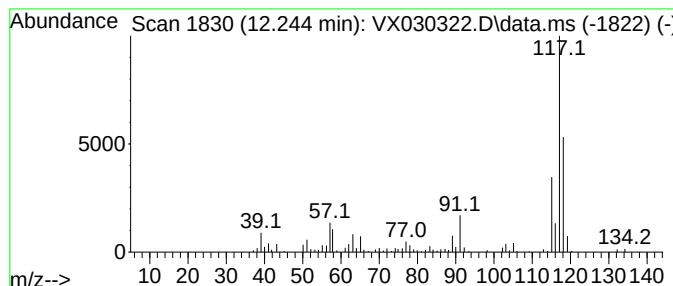
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Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 2 Indane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	7.00 ug/l	109433	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	95
2			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	76
3			Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	74
4			Delta-cyclene	118	C9H10	007785-10-6	72
5			1,3-Methanopentalene, 1,2,3,5-te...	118	C9H10	128600-88-4	64



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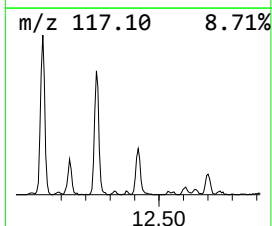
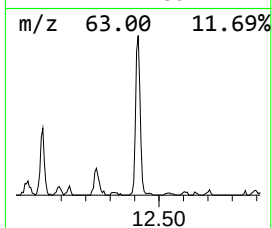
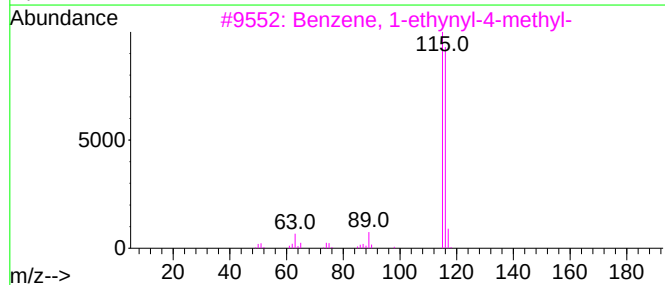
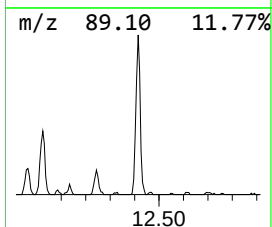
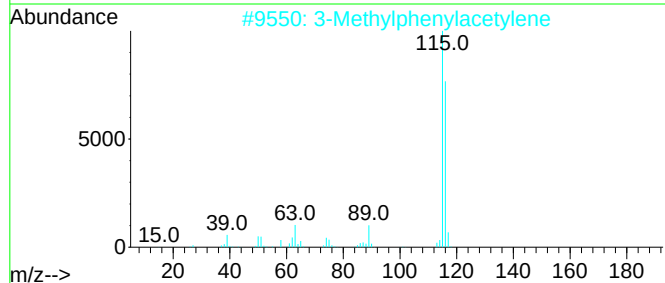
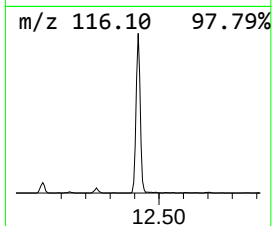
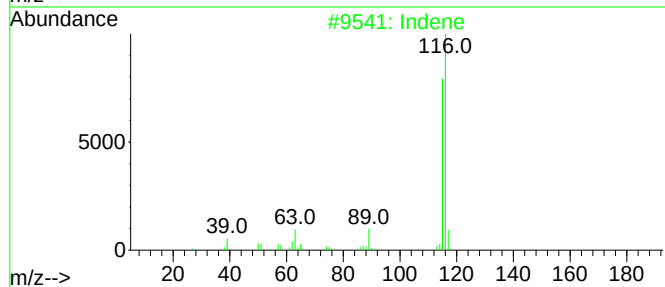
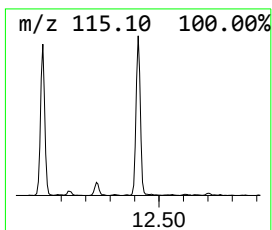
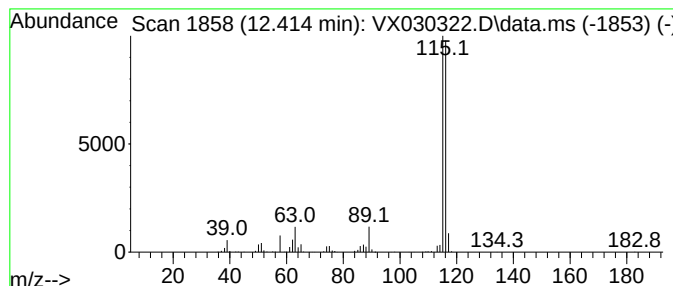
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Peak Number 3 Indene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.414	15.50 ug/l	242332	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indene	116	C9H8	000095-13-6	97
2			3-Methylphenylacetylene	116	C9H8	000766-82-5	91
3			Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	91
4			Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
5			Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	91



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ALS Vial : 12 Sample Multiplier: 1

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ClientSampleId :
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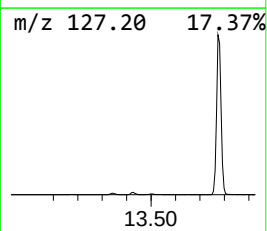
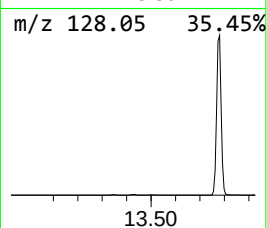
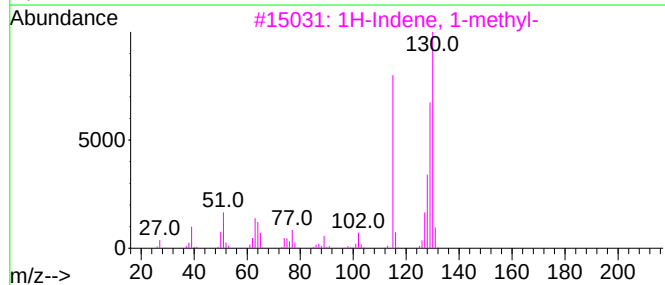
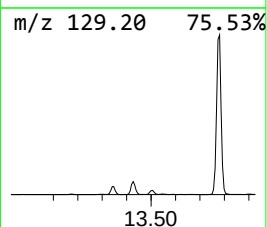
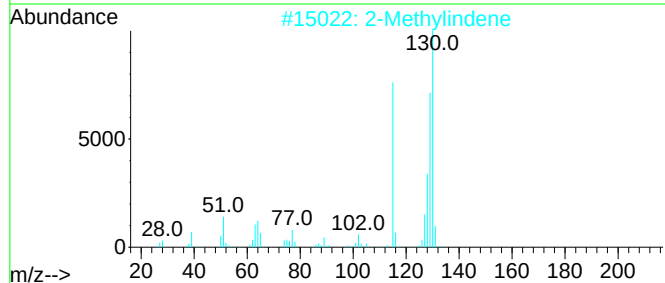
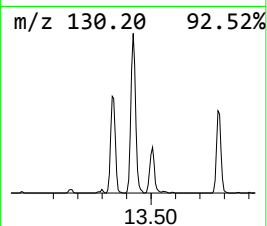
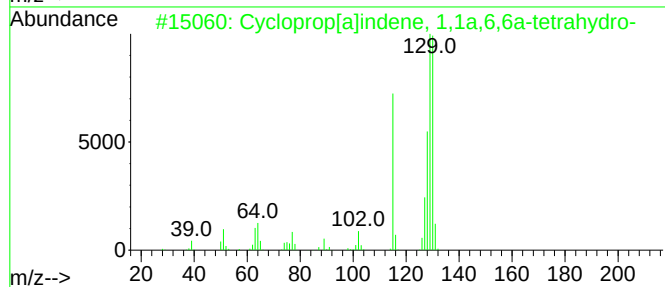
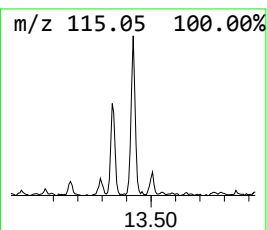
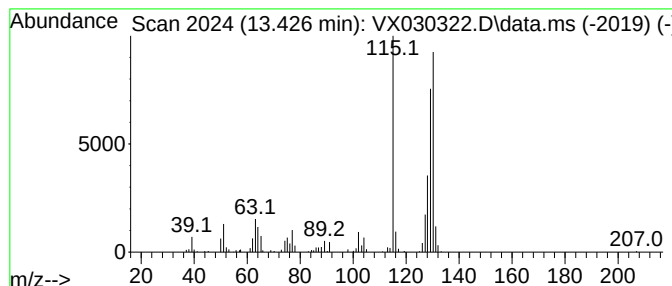
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Cycloprop[a]indene, 1,1a,6,... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.426	6.45 ug/l	100850	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	95
2			2-Methylindene	130	C10H10	002177-47-1	95
3			1H-Indene, 1-methyl-	130	C10H10	000767-59-9	94
4			1,4-Dihydronaphthalene	130	C10H10	000612-17-9	93
5			1H-Indene, 3-methyl-	130	C10H10	000767-60-2	93



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ClientSampleId :
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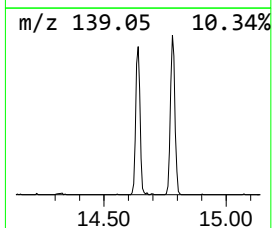
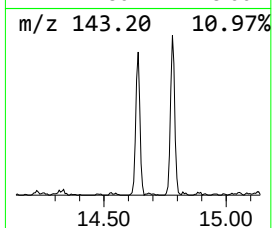
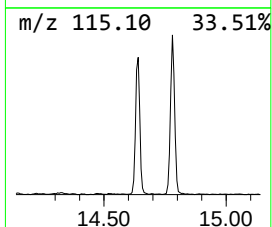
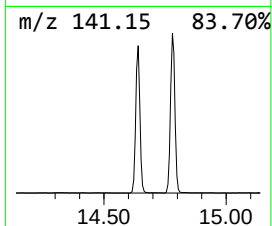
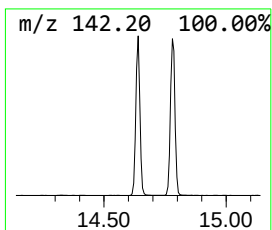
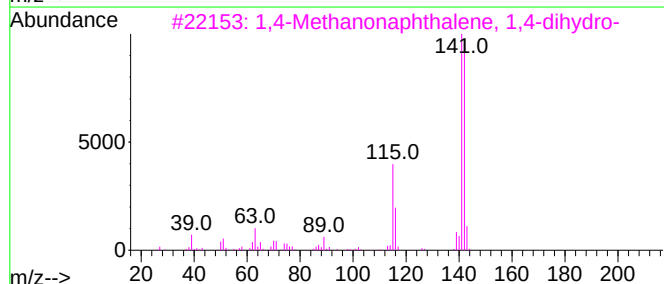
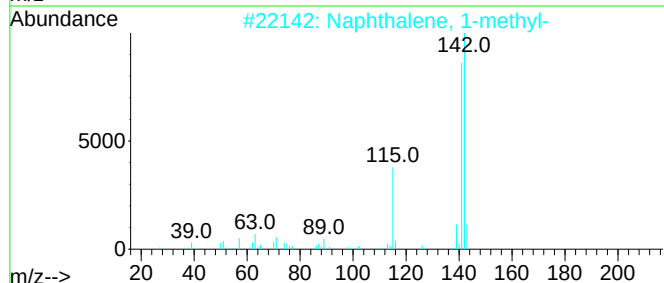
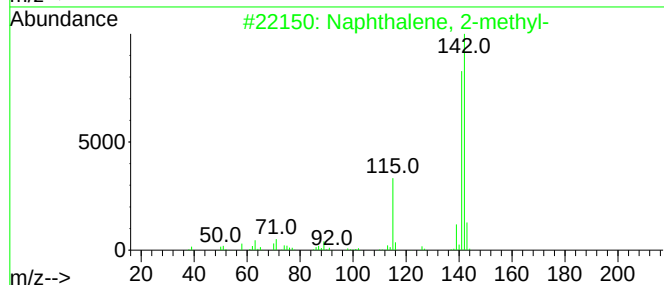
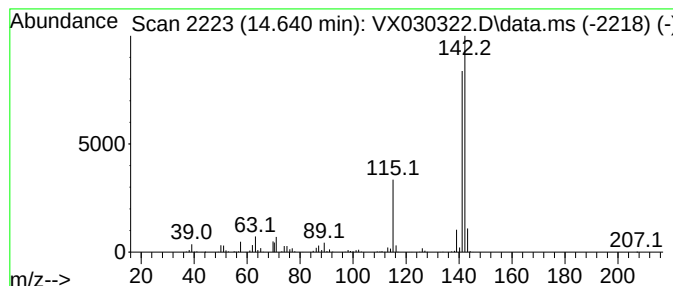
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Naphthalene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.640	33.88 ug/l	529575	1,4-Dichlorobenzene-d4	12.024

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-dihy...	142	C11H10	004453-90-1	93
4			Benzocycloheptatriene	142	C11H10	000264-09-5	90
5			Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	59



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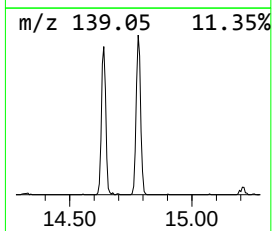
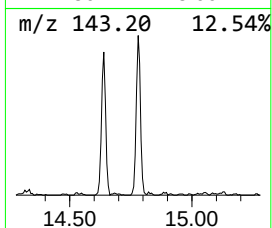
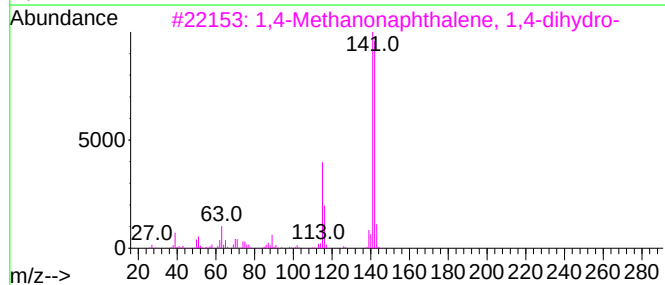
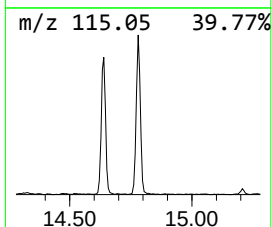
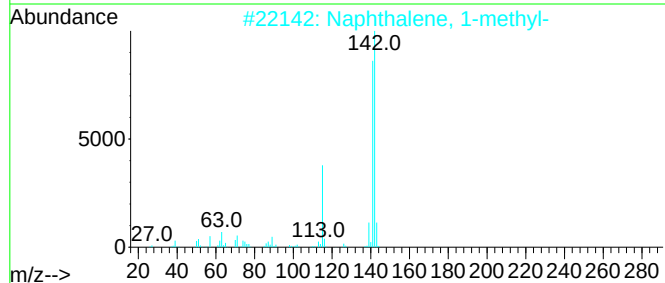
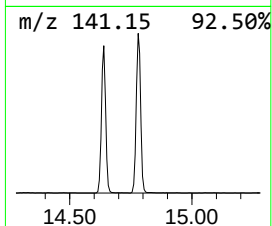
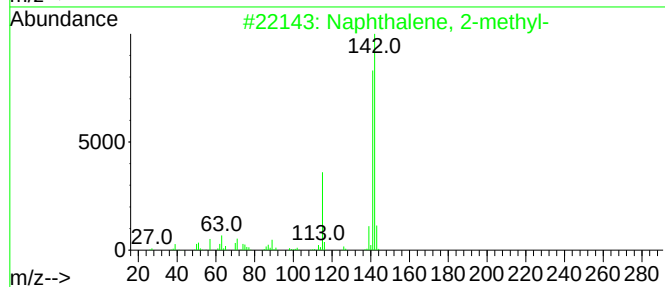
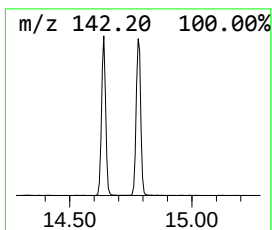
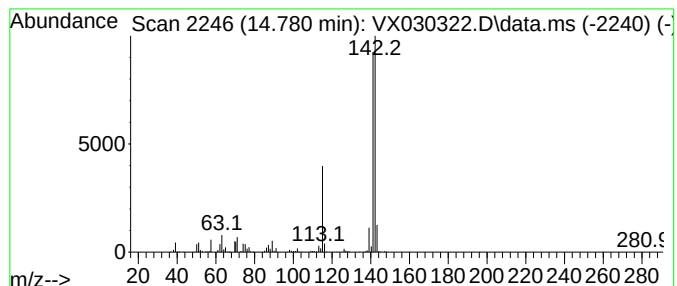
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 1,4-Methanonaphthalene, 1,4... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.780	38.28 ug/l	598283	1,4-Dichlorobenzene-d4	12.024

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-dihy...	142	C11H10	004453-90-1	93
4			Benzocycloheptatriene	142	C11H10	000264-09-5	91
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080122\
Data File : VX030322.D
Acq On : 01 Aug 2022 15:07
Operator : JC/MD
Sample : N3964-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
HMGP-WATER-1

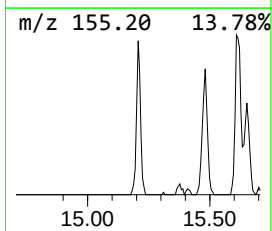
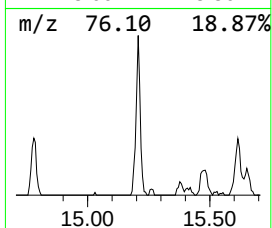
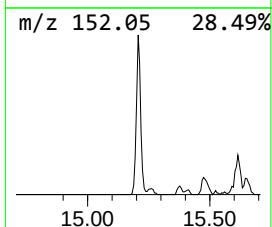
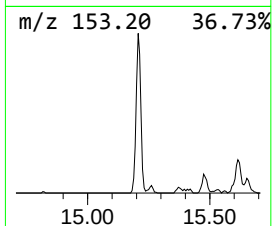
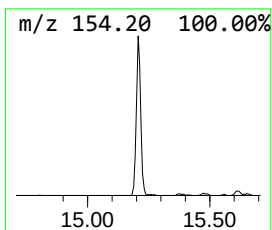
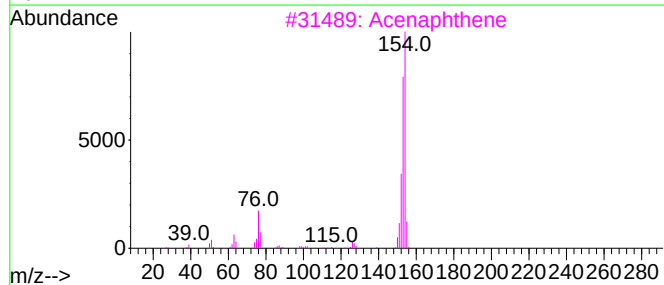
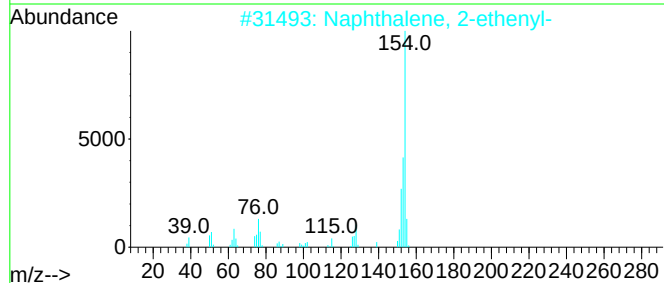
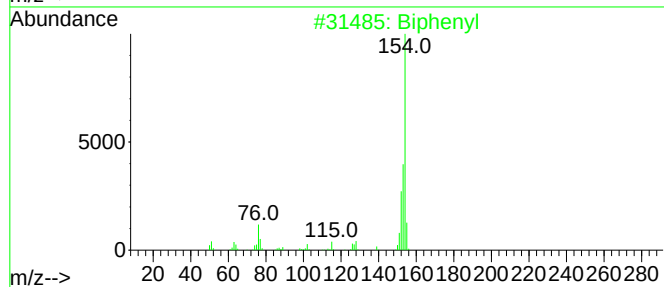
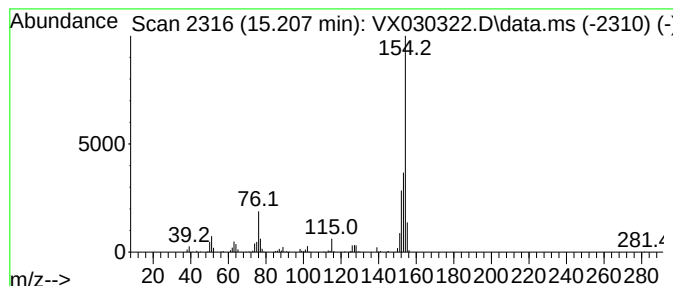
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Biphenyl Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.207	7.68 ug/l	120058	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Biphenyl	154	C12H10	000092-52-4	94
2		Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	74
3		Acenaphthene	154	C12H10	000083-32-9	62
4		3-Quinolincarbonitrile	154	C10H6N2	034846-64-5	47
5		Sebacic acid, 2,6-dimethoxypheny...	450	C26H42O6	1000354-75-6	28



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080122\
Data File : VX030322.D
Acq On : 01 Aug 2022 15:07
Operator : JC/MD
Sample : N3964-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
HMGP-WATER-1

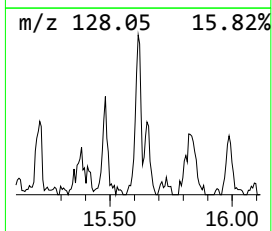
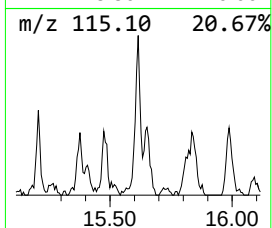
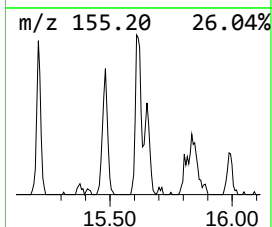
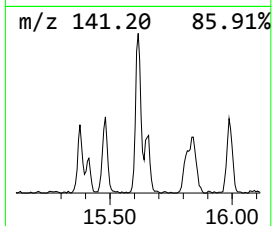
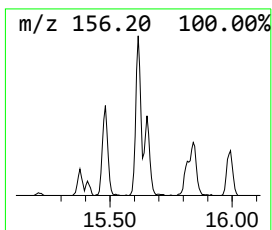
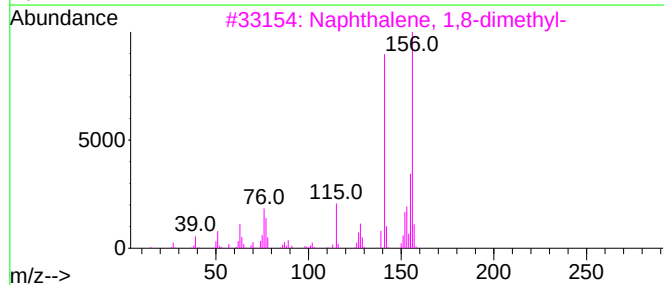
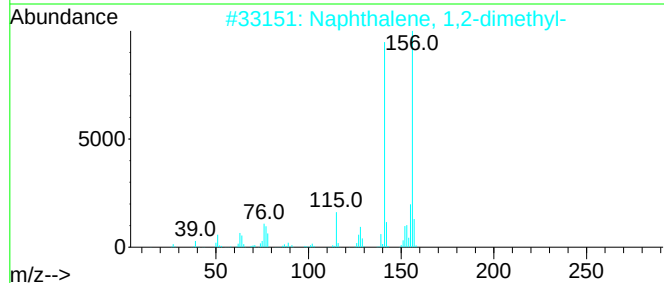
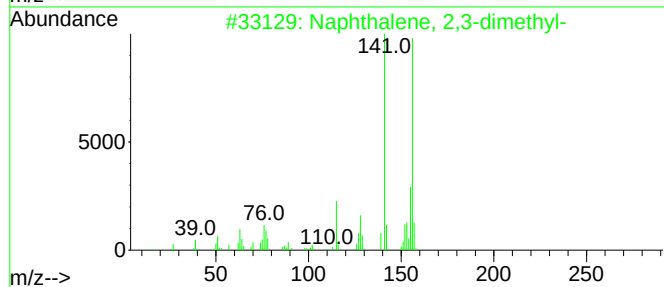
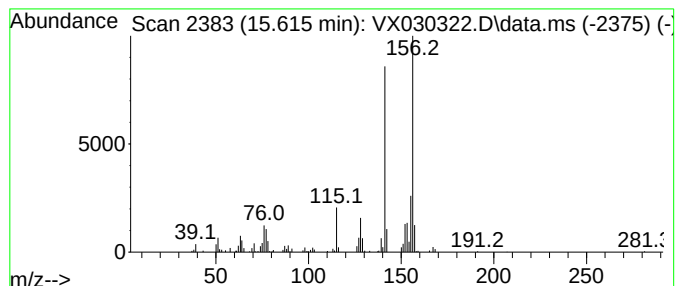
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Naphthalene, 2,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.615	7.74 ug/l	120971	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2			Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	98
3			Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	98
4			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
5			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080122\
Data File : VX030322.D
Acq On : 01 Aug 2022 15:07
Operator : JC/MD
Sample : N3964-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
HMGP-WATER-1

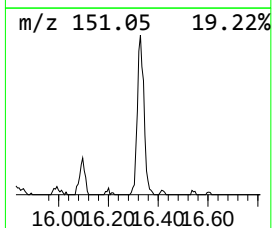
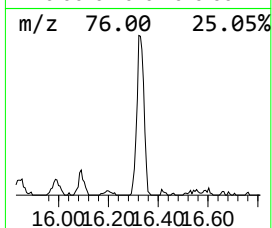
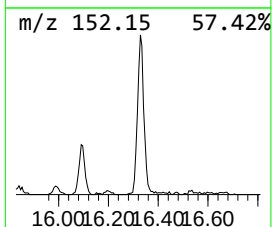
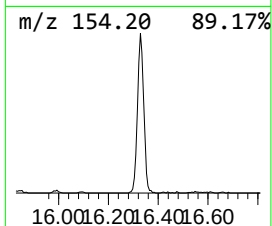
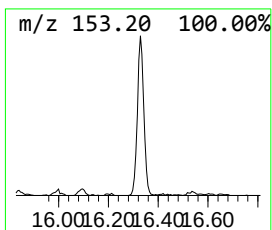
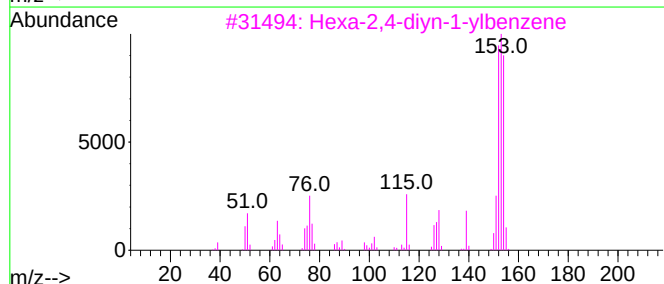
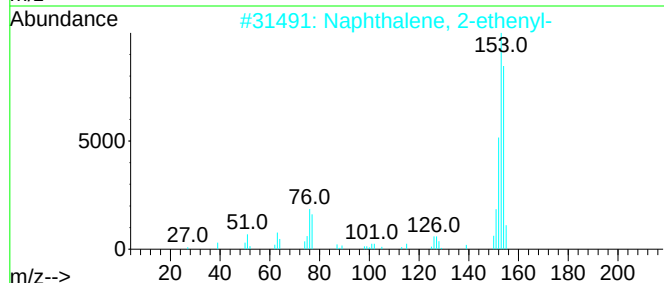
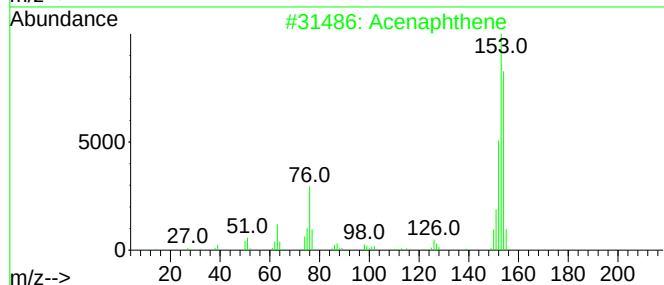
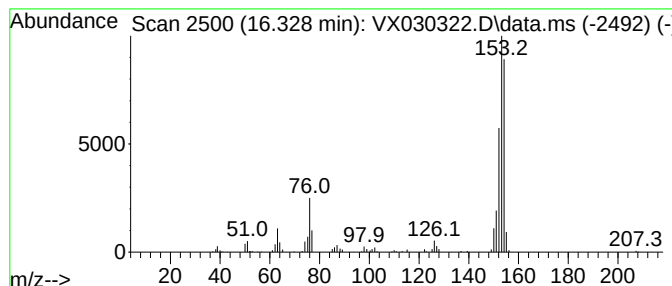
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Acenaphthene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.328	12.90 ug/l	201563	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acenaphthene	154	C12H10	000083-32-9	93
2			Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	62
3			Hexa-2,4-diyne-1-ylbenzene	154	C12H10	000520-74-1	58
4			1,4-Ethenonaphthalene, 1,4-dihydro-	154	C12H10	007322-47-6	58
5			3-Fluoro-para-anisaldehyde	154	C8H7FO2	000351-54-2	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080122\
Data File : VX030322.D
Acq On : 01 Aug 2022 15:07
Operator : JC/MD
Sample : N3964-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
HMGP-WATER-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol	2.069	419.8	ug/l	2960080	1	5.556	352559	50.0
Indane	12.244	7.0	ug/l	109433	4	12.024	781489	50.0
Indene	12.414	15.5	ug/l	242332	4	12.024	781489	50.0
Cycloprop[a]ind...	13.426	6.5	ug/l	100850	4	12.024	781489	50.0
Naphthalene, 1-...	14.640	33.9	ug/l	529575	4	12.024	781489	50.0
1,4-Methanonaph...	14.780	38.3	ug/l	598283	4	12.024	781489	50.0
Biphenyl	15.207	7.7	ug/l	120058	4	12.024	781489	50.0
Naphthalene, 2,...	15.615	7.7	ug/l	120971	4	12.024	781489	50.0
Acenaphthene	16.328	12.9	ug/l	201563	4	12.024	781489	50.0