

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN050820\
 Data File : VN061360.D
 Acq On : 8 May 2020 14:36
 Operator : JC/MD
 Sample : L2544-11
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SS038MW009-200505

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_N\METHODS\82N042720W.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	1.635	3	14	133	rBV	4184860	15813867	100.00%	59.426%
2	7.548	1835	1853	1864	rBV2	70404	175289	1.11%	0.659%
3	7.625	1864	1877	1894	rVB2	133935	313339	1.98%	1.177%
4	7.989	1972	1990	2013	rBV2	106928	265815	1.68%	0.999%
5	8.551	2148	2165	2184	rBV	218239	452703	2.86%	1.701%
6	10.059	2621	2634	2652	rBV	353051	633264	4.00%	2.380%
7	11.406	3032	3053	3068	rBV2	2044894	4060257	25.68%	15.258%
8	12.377	3342	3355	3367	rVB	272022	462629	2.93%	1.738%
9	13.320	3637	3648	3663	rVB	334156	565136	3.57%	2.124%
10	13.419	3669	3679	3691	rBV	176112	285030	1.80%	1.071%
11	13.487	3692	3700	3705	rVV	129292	199662	1.26%	0.750%
12	13.641	3737	3748	3758	rBV5	96420	187062	1.18%	0.703%
13	13.863	3808	3817	3826	rVB	217781	322886	2.04%	1.213%
14	13.921	3826	3835	3841	rBV	182978	289441	1.83%	1.088%
15	13.969	3841	3850	3860	rVB2	772901	1241407	7.85%	4.665%
16	14.107	3882	3893	3901	rBV	155743	259552	1.64%	0.975%
17	14.184	3901	3917	3932	rVB3	243057	531426	3.36%	1.997%
18	14.570	4024	4037	4048	rBV3	314660	552254	3.49%	2.075%

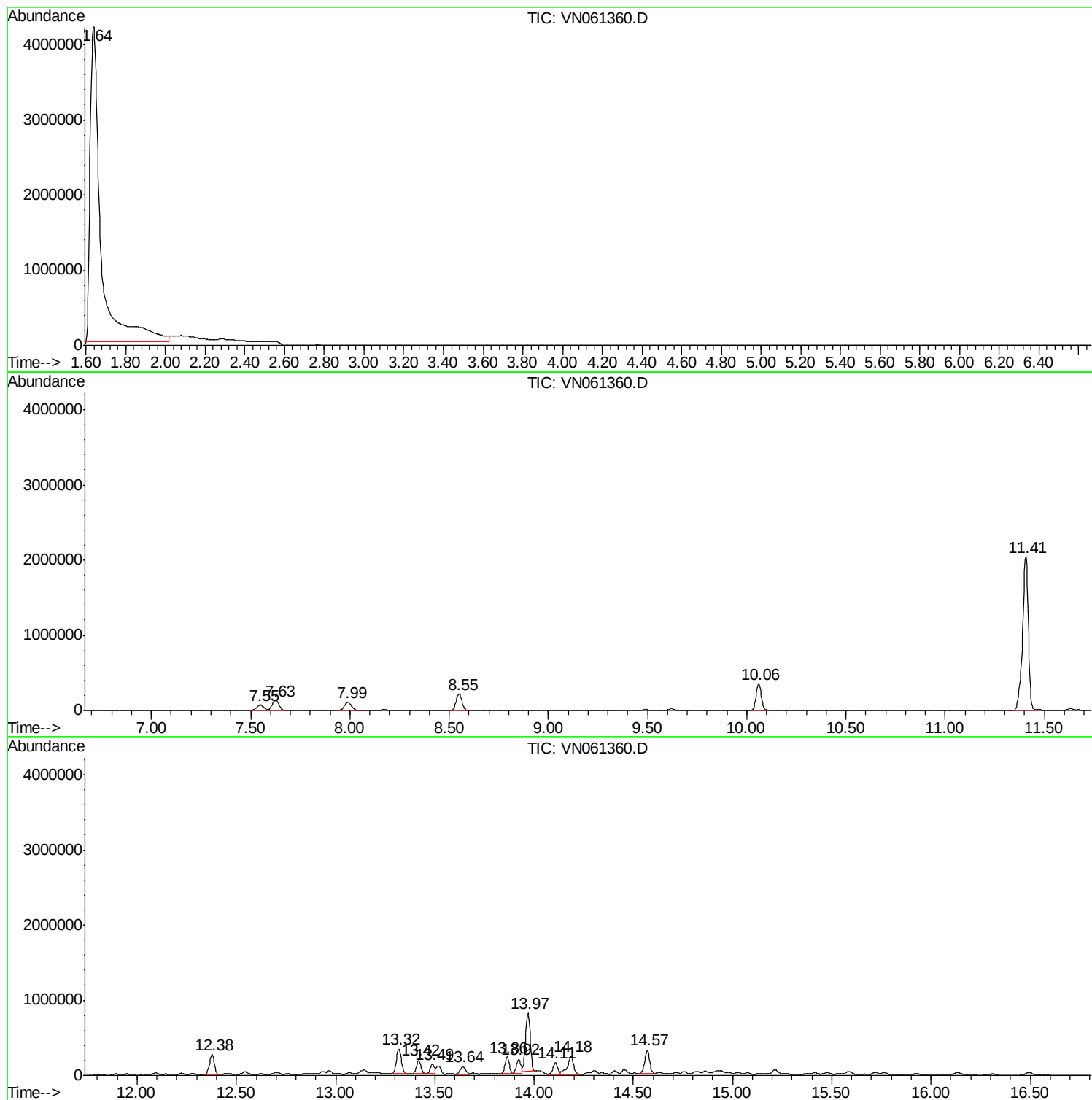
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N042720W.M
Quant Title : SW846 8260

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



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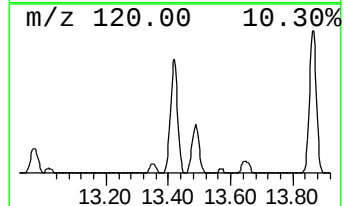
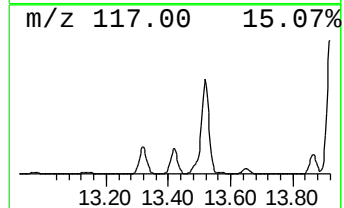
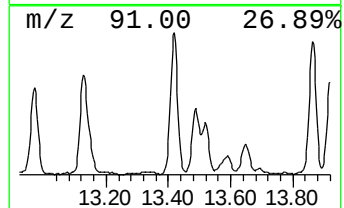
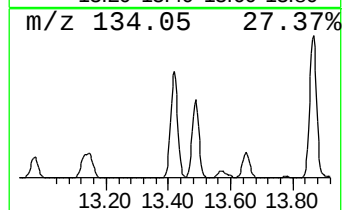
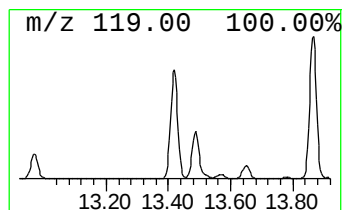
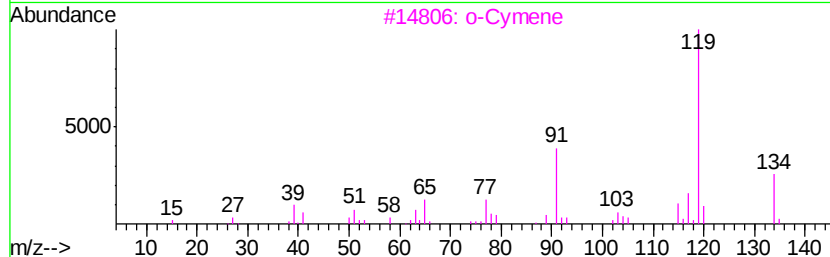
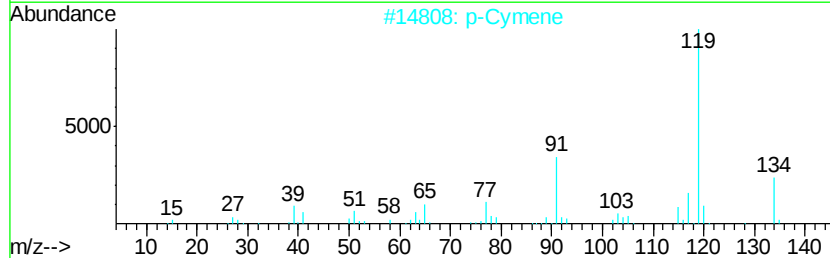
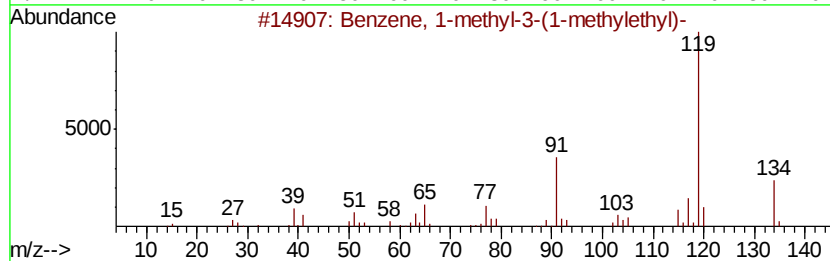
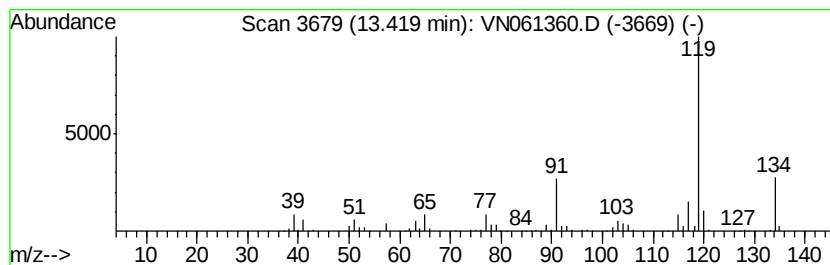
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Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1-methyl-3-(1-meth... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.42	25.22 ug/l	285030	1,4-Dichlorobenzene-d4	13.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1-methyl-3-(1-methylethyl-)	134 C10H14	000535-77-3	95
2	p-Cymene	134 C10H14	000099-87-6	95
3	o-Cymene	134 C10H14	000527-84-4	95
4	Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4	91
5	Benzene, 4-ethyl-1,2-dimethyl-	134 C10H14	000934-80-5	91



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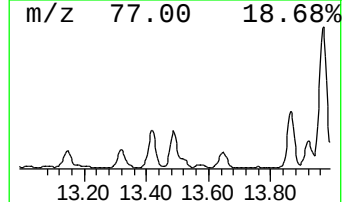
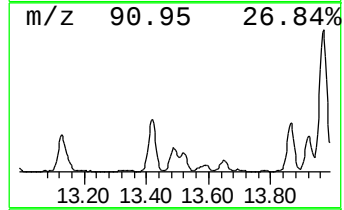
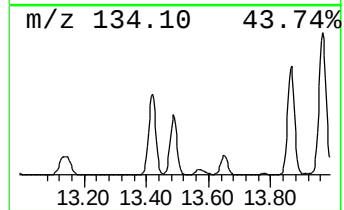
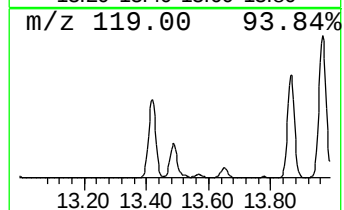
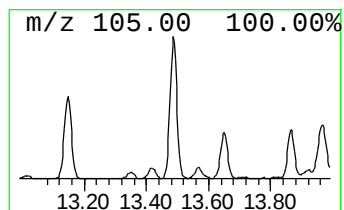
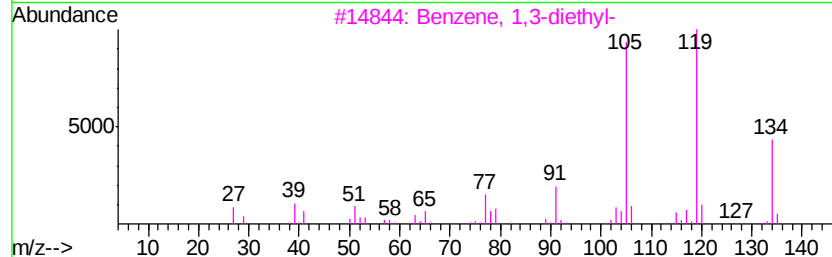
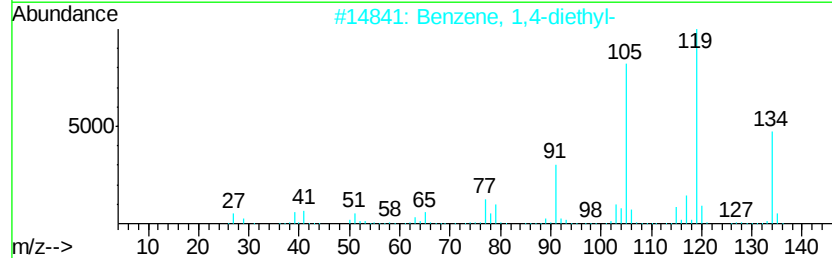
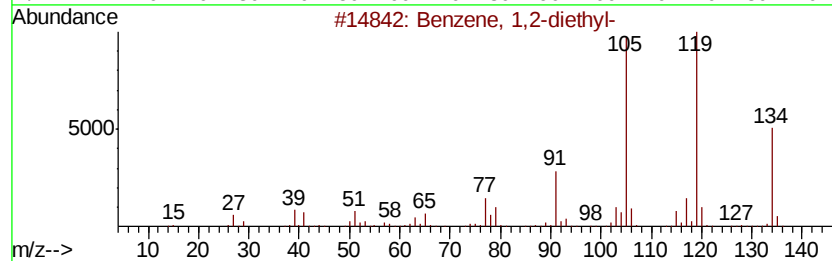
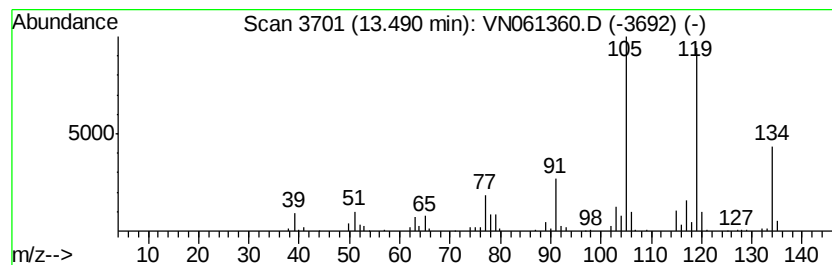
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TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1,2-diethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.49	17.67 ug/l	199662	1,4-Dichlorobenzene-d4	13.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2-diethyl-	134 C10H14	000135-01-3 96
2		Benzene, 1,4-diethyl-	134 C10H14	000105-05-5 96
3		Benzene, 1,3-diethyl-	134 C10H14	000141-93-5 95
4		Benzene, 4-ethyl-1,2-dimethyl-	134 C10H14	000934-80-5 91
5		Benzene, 1-ethyl-3,5-dimethyl-	134 C10H14	000934-74-7 90



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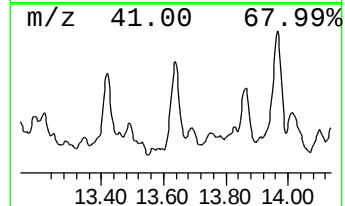
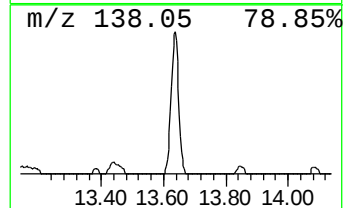
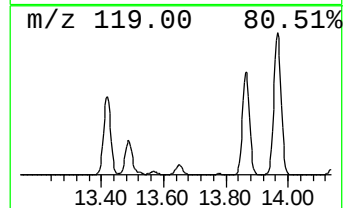
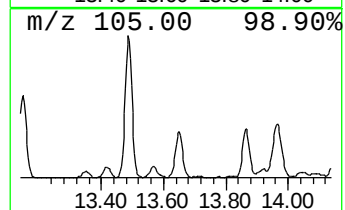
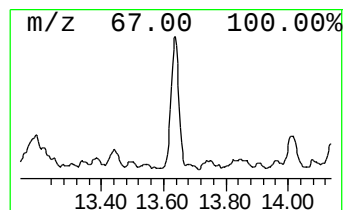
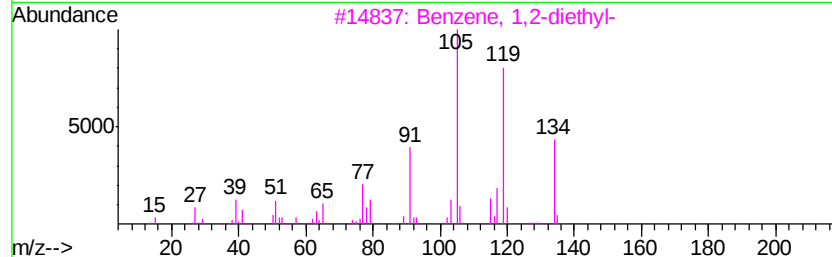
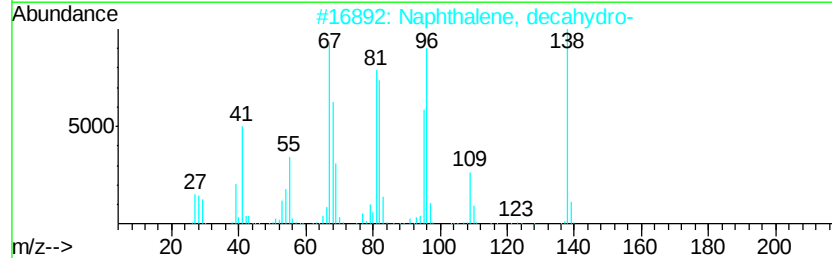
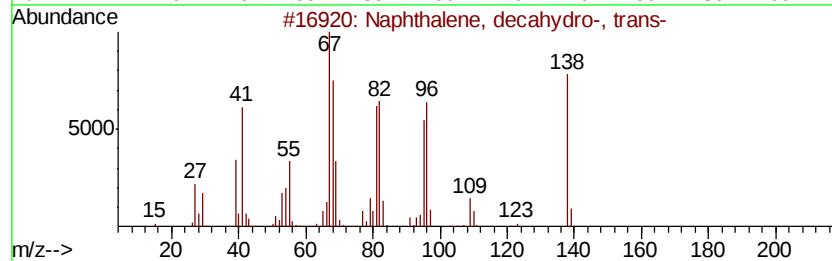
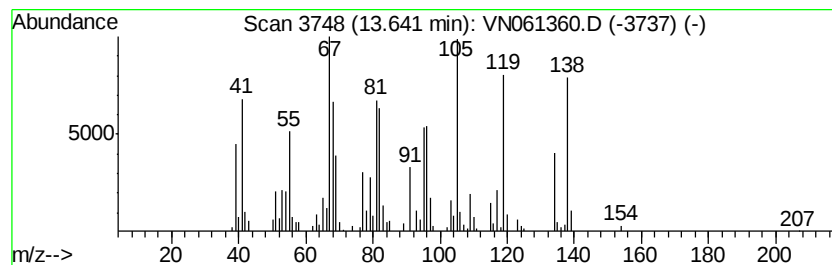
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TIC Integration Parameters: LSCINT.P

Peak Number 4 Naphthalene, decahydro-, tr... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.64	16.55 ug/l	187062	1,4-Dichlorobenzene-d4	13.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, decahydro-, trans-	138 C10H18	000493-02-7 96
2		Naphthalene, decahydro-	138 C10H18	000091-17-8 93
3		Benzene, 1,2-diethyl-	134 C10H14	000135-01-3 60
4		Spiro[4.5]decane	138 C10H18	000176-63-6 60
5		Cyclohexane, 1-methyl-3-(1-methy...	138 C10H18	024399-15-3 60



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Misc : 5.00mL/MSVOA N/WATER
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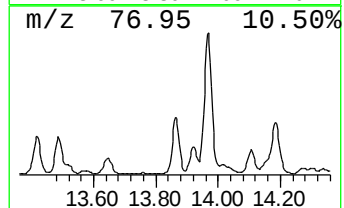
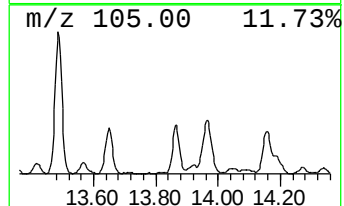
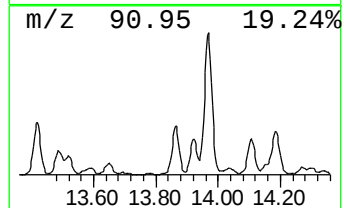
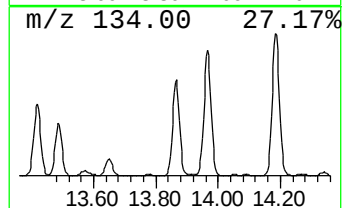
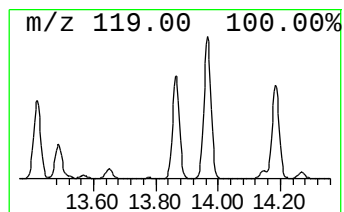
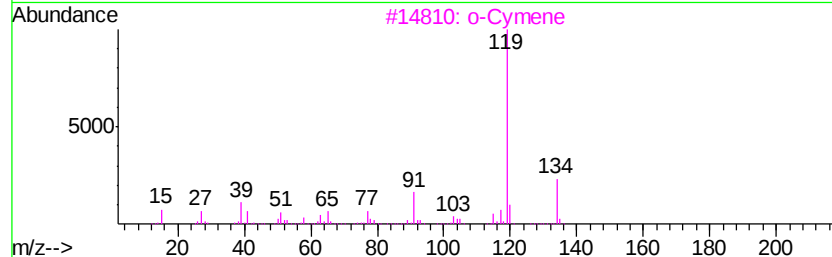
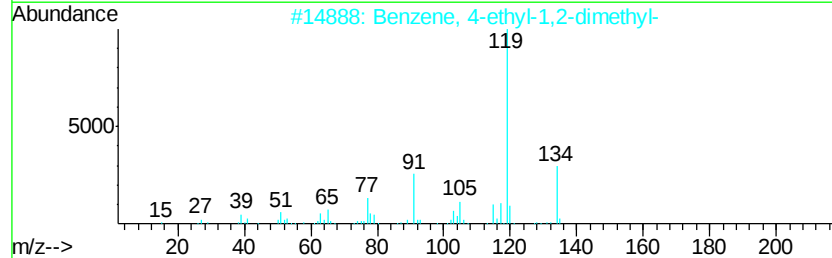
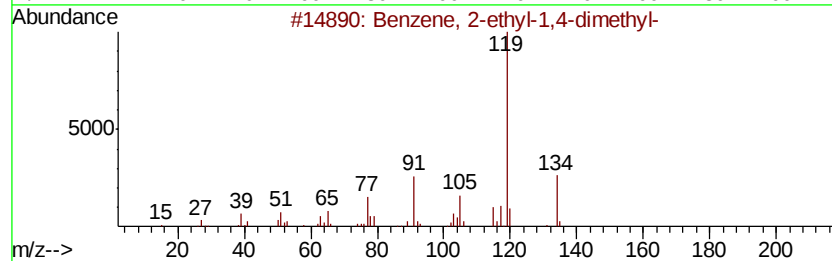
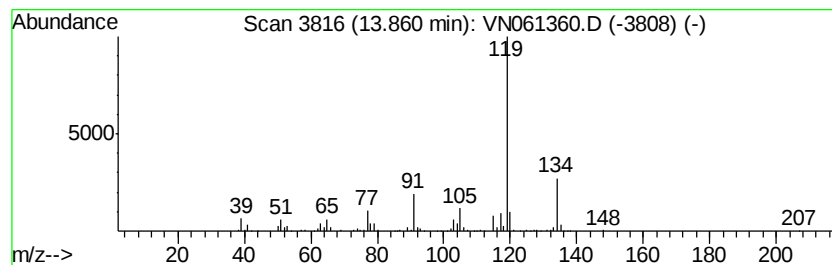
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	28.57 ug/l	322886	1,4-Dichlorobenzene-d4	13.32

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3	o-Cymene	134	C10H14	000527-84-4	95
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5	p-Cymene	134	C10H14	000099-87-6	95



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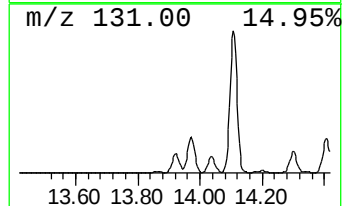
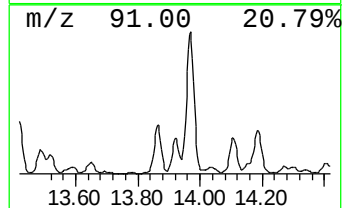
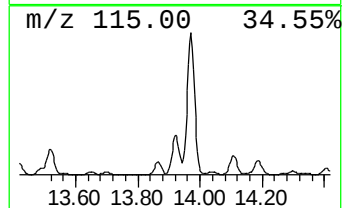
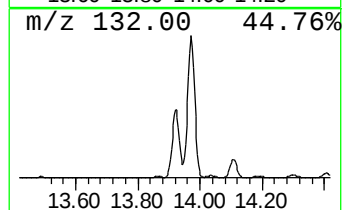
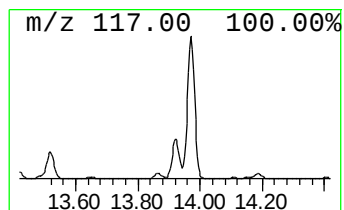
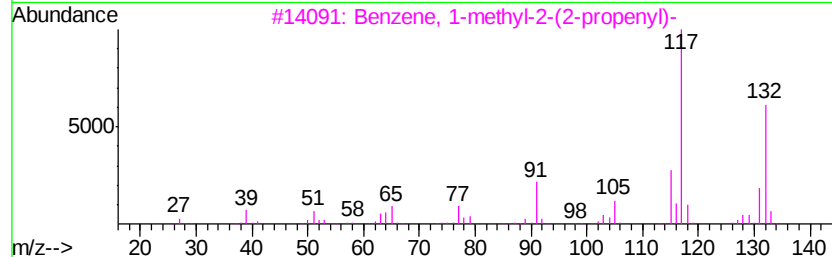
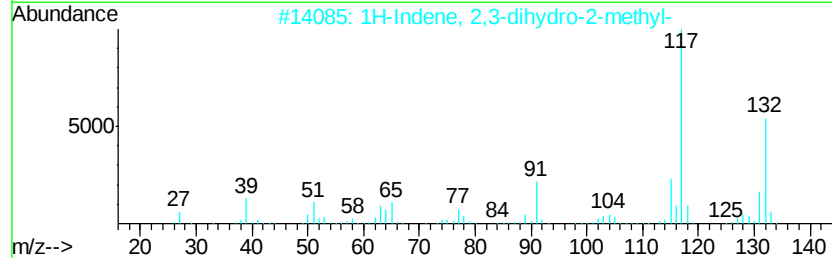
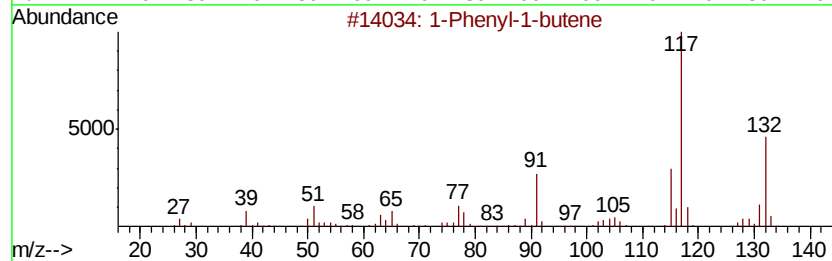
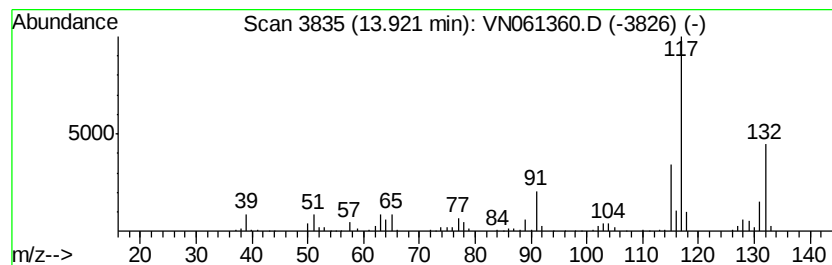
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.92	25.61 ug/l	289441	1,4-Dichlorobenzene-d4	13.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Phenyl-1-butene	132 C10H12	000824-90-8	94
2	1H-Indene, 2,3-dihydro-2-methyl-	132 C10H12	000824-63-5	93
3	Benzene, 1-methyl-2-(2-propenyl)-	132 C10H12	001587-04-8	91
4	Benzene, (2-methyl-2-propenyl)-	132 C10H12	003290-53-7	91
5	Benzene, (2-methyl-1-propenyl)-	132 C10H12	000768-49-0	91



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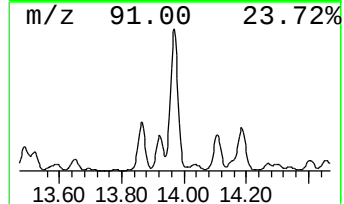
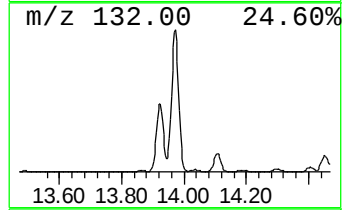
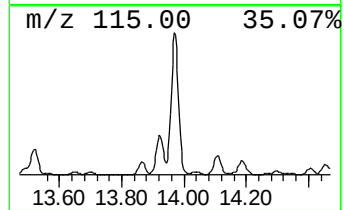
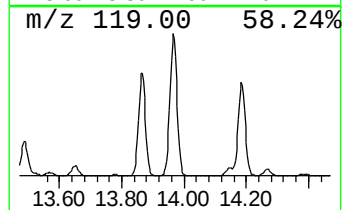
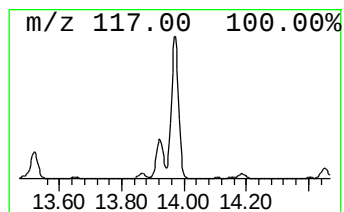
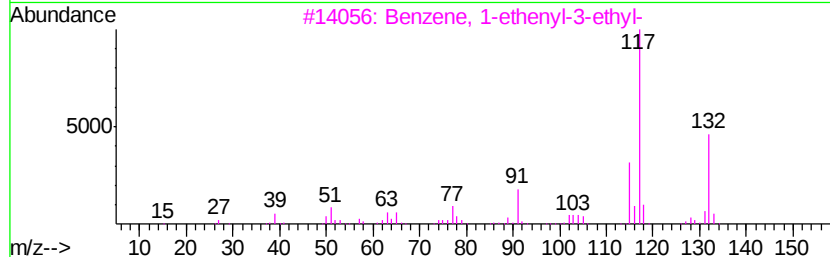
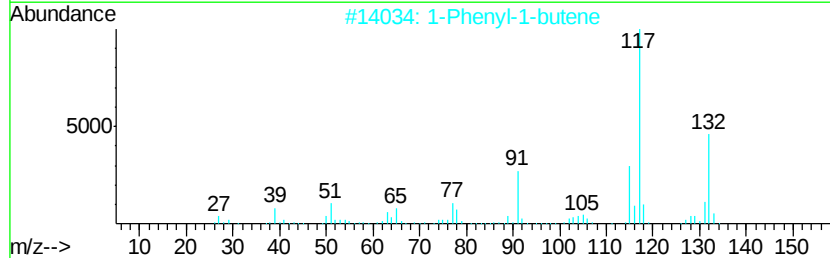
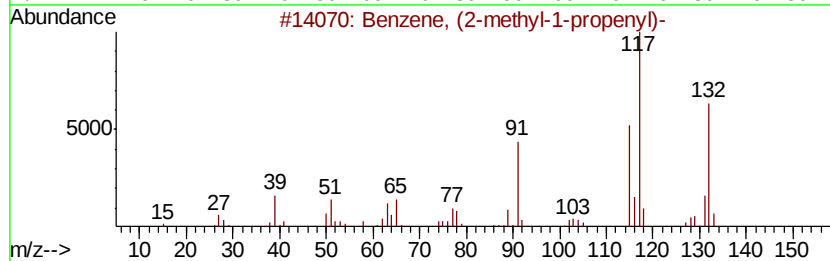
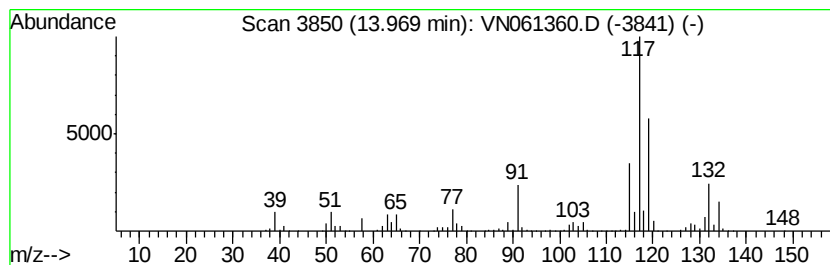
Instrument :
MSVOA_N
ClientSampleId :
SS038MW009-200505

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N042720W.M
Quant Title : SW846 8260

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

Peak Number 7 Benzene, (2-methyl-1-propen... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	109.83 ug/l	1241410	1,4-Dichlorobenzene-d4	13.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (2-methyl-1-propenyl)-	132 C10H12	000768-49-0 86
2		1-Phenyl-1-butene	132 C10H12	000824-90-8 86
3		Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4 70
4		Indan, 1-methyl-	132 C10H12	000767-58-8 64
5		3-Phenylbut-1-ene	132 C10H12	000934-10-1 64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN050820\
Data File : VN061360.D
Acq On : 8 May 2020 14:36
Operator : JC/MD
Sample : L2544-11
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 12 Sample Multiplier: 1

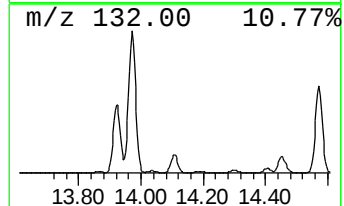
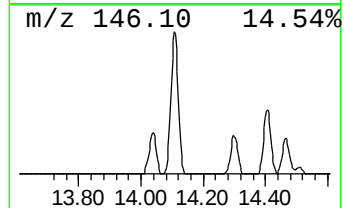
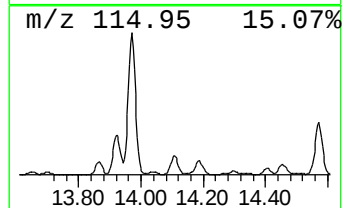
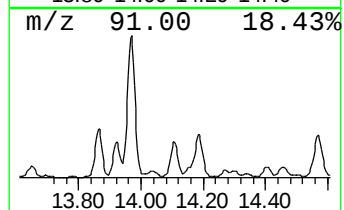
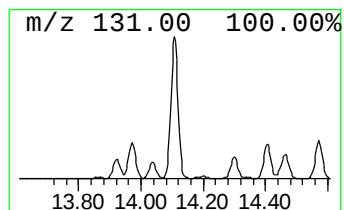
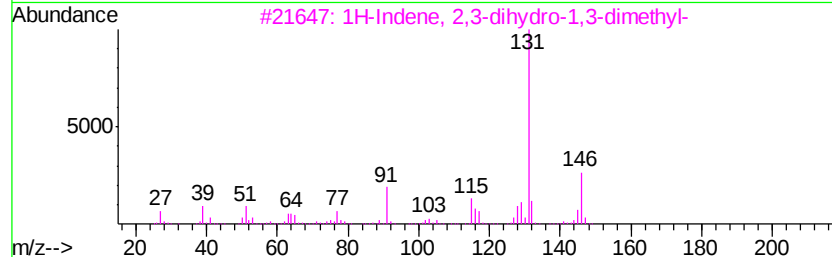
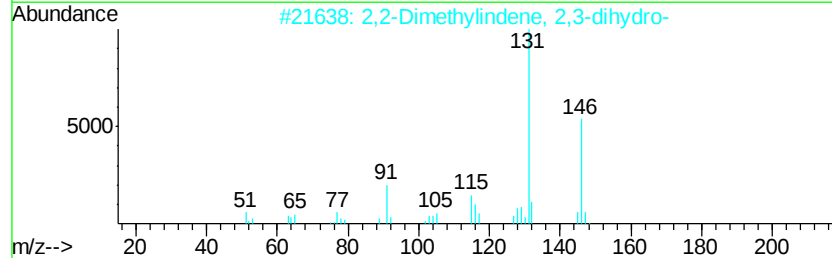
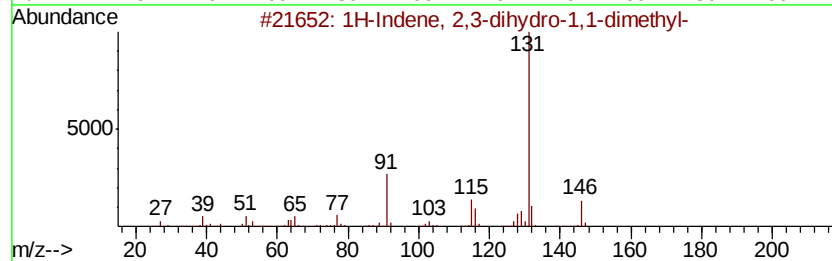
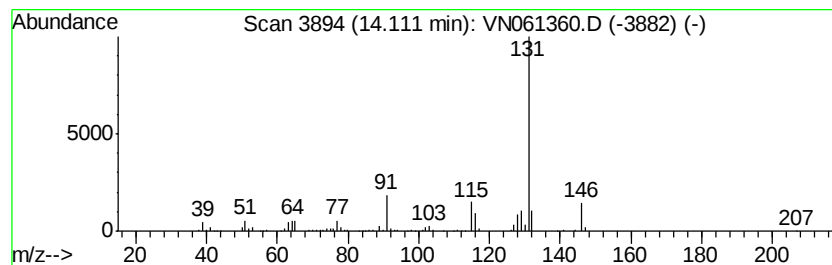
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ClientSampleId :
SS038MW009-200505

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N042720W.M
Quant Title : SW846 8260

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

Peak Number 8 1H-Indene, 2,3-dihydro-1,1-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.11	22.96 ug/l	259552	1,4-Dichlorobenzene-d4	13.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-1,1-dimet...	146 C11H14	004912-92-9	97
2	2,2-Dimethylindene, 2,3-dihydro-	146 C11H14	020836-11-7	91
3	1H-Indene, 2,3-dihydro-1,3-dimet...	146 C11H14	004175-53-5	91
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9	91
5	Cyclopropane, 2-bromo-1-methyl-1...	210 C10H11Br	055091-64-0	78



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN050820\
Data File : VN061360.D
Acq On : 8 May 2020 14:36
Operator : JC/MD
Sample : L2544-11
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SS038MW009-200505

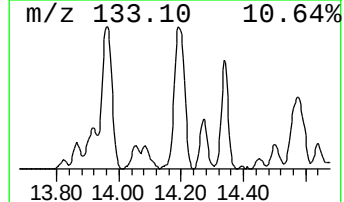
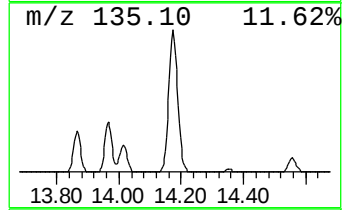
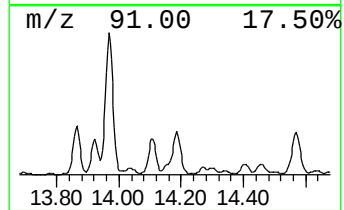
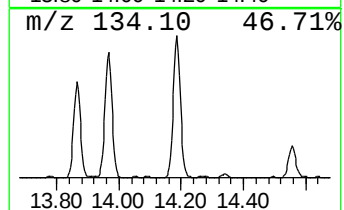
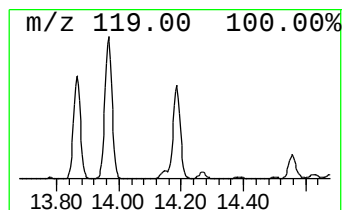
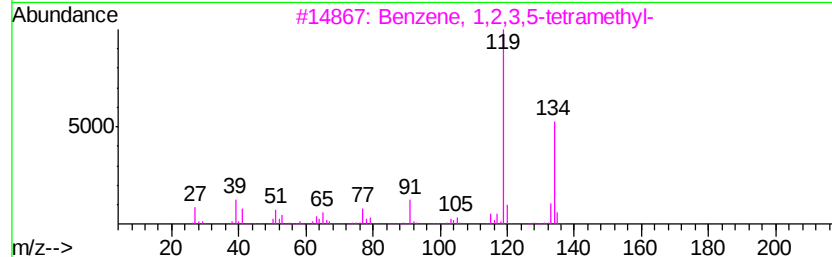
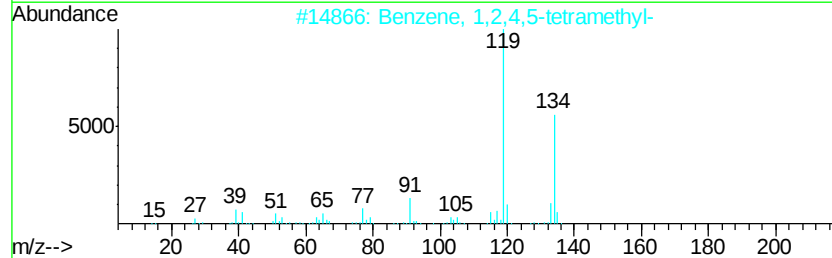
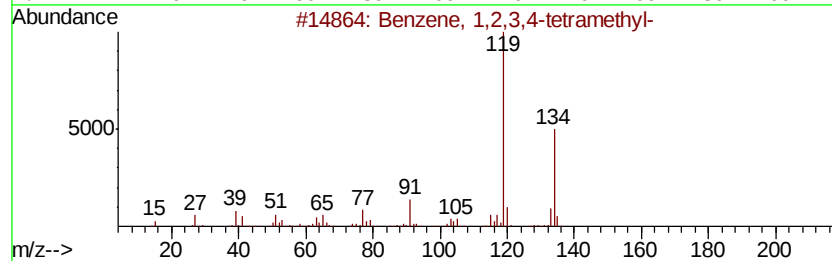
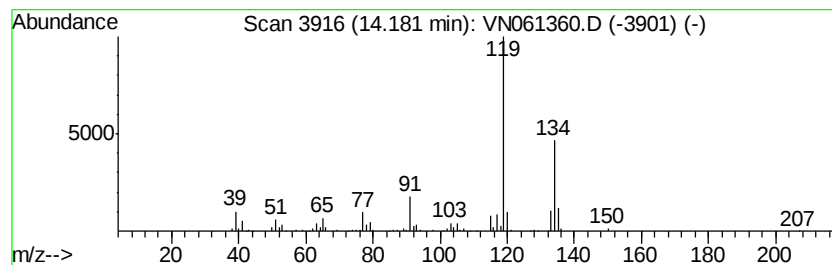
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.18	47.02 ug/l	531426	1,4-Dichlorobenzene-d4	13.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4			o-Cymene	134	C10H14	000527-84-4	94
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN050820\
Data File : VN061360.D
Acq On : 8 May 2020 14:36
Operator : JC/MD
Sample : L2544-11
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SS038MW009-200505

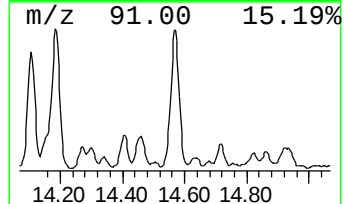
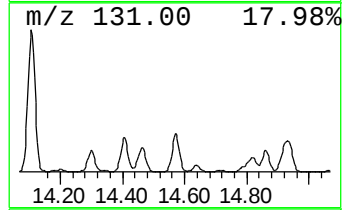
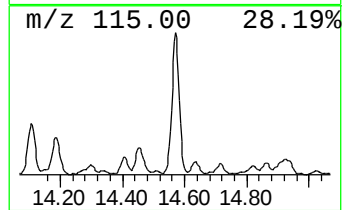
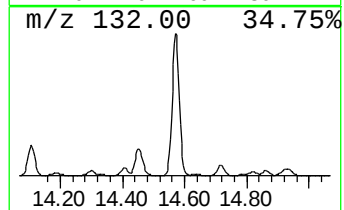
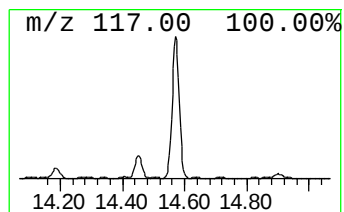
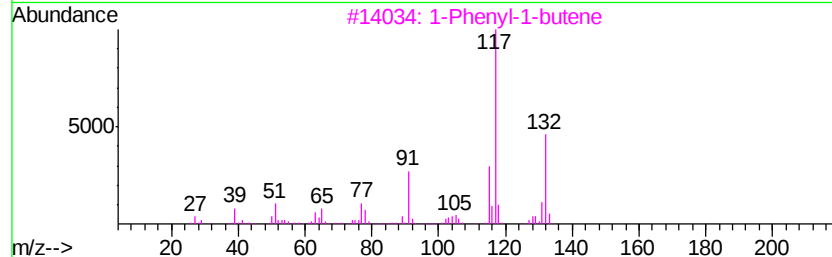
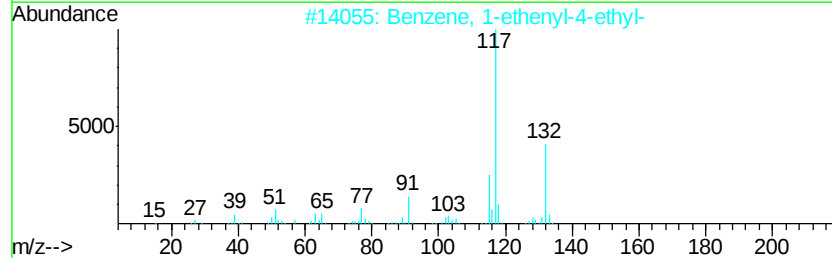
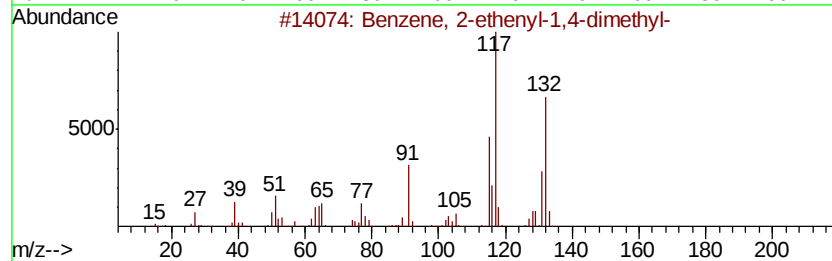
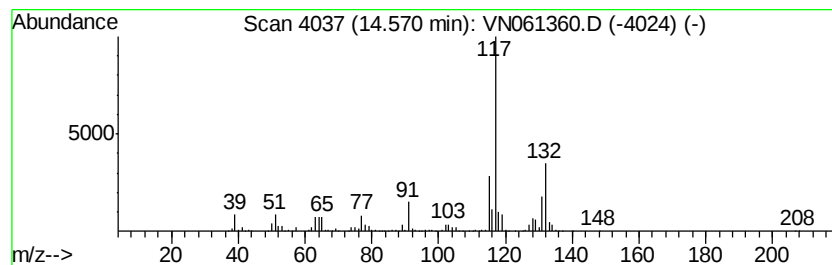
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	48.86 ug/l	552254	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3		1-Phenyl-1-butene	132	C10H12	000824-90-8	94
4		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
5		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN050820\
Data File : VN061360.D
Acq On : 8 May 2020 14:36
Operator : JC/MD
Sample : L2544-11
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SS038MW009-200505

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N042720W.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
Benzene, 1-methyl...	13.42	25.2	ug/l	285030	4	13.32	565136	50.0
Benzene, 1,2-diet...	13.49	17.7	ug/l	199662	4	13.32	565136	50.0
Naphthalene, deca...	13.64	16.6	ug/l	187062	4	13.32	565136	50.0
Benzene, 2-ethyl-...	13.86	28.6	ug/l	322886	4	13.32	565136	50.0
1-Phenyl-1-butene	13.92	25.6	ug/l	289441	4	13.32	565136	50.0
Benzene, (2-methy...	13.97	109.8	ug/l	1241410	4	13.32	565136	50.0
1H-Indene, 2,3-di...	14.11	23.0	ug/l	259552	4	13.32	565136	50.0
Benzene, 1,2,3,4-...	14.18	47.0	ug/l	531426	4	13.32	565136	50.0
Benzene, 2-etheny...	14.57	48.9	ug/l	552254	4	13.32	565136	50.0