

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN050820\
 Data File : VN061357.D
 Acq On : 8 May 2020 13:26
 Operator : JC/MD
 Sample : L2544-05
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 ST008MW037-200504

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_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	1.638	3	15	74	rBV	3913158	13255620	100.00%	57.472%
2	7.548	1837	1853	1864	rBV	72082	176343	1.33%	0.765%
3	7.625	1865	1877	1891	rVV	133854	315167	2.38%	1.366%
4	7.989	1973	1990	2018	rBV	94055	221585	1.67%	0.961%
5	8.172	2035	2047	2067	rVB	59511	168504	1.27%	0.731%
6	8.551	2149	2165	2192	rBV	223661	458461	3.46%	1.988%
7	10.059	2621	2634	2659	rVB	358719	656165	4.95%	2.845%
8	11.381	3032	3045	3065	rBV2	340818	717118	5.41%	3.109%
9	12.378	3343	3355	3367	rBV	269210	444454	3.35%	1.927%
10	12.969	3533	3539	3549	rVB	87850	135079	1.02%	0.586%
11	13.140	3578	3592	3619	rVB3	282813	732846	5.53%	3.177%
12	13.320	3635	3648	3663	rBV	335184	582492	4.39%	2.525%
13	13.419	3663	3679	3691	rVV	324035	524313	3.96%	2.273%
14	13.487	3691	3700	3717	rVV2	109389	187611	1.42%	0.813%
15	13.866	3801	3818	3828	rBV	504097	784260	5.92%	3.400%
16	13.969	3840	3850	3861	rVV2	760189	1221203	9.21%	5.295%
17	14.107	3881	3893	3901	rBV	234234	379799	2.87%	1.647%
18	14.185	3901	3917	3935	rVB	501948	916435	6.91%	3.973%
19	14.570	4024	4037	4048	rBV	575536	978573	7.38%	4.243%
20	14.934	4138	4150	4168	rVB2	85202	208457	1.57%	0.904%

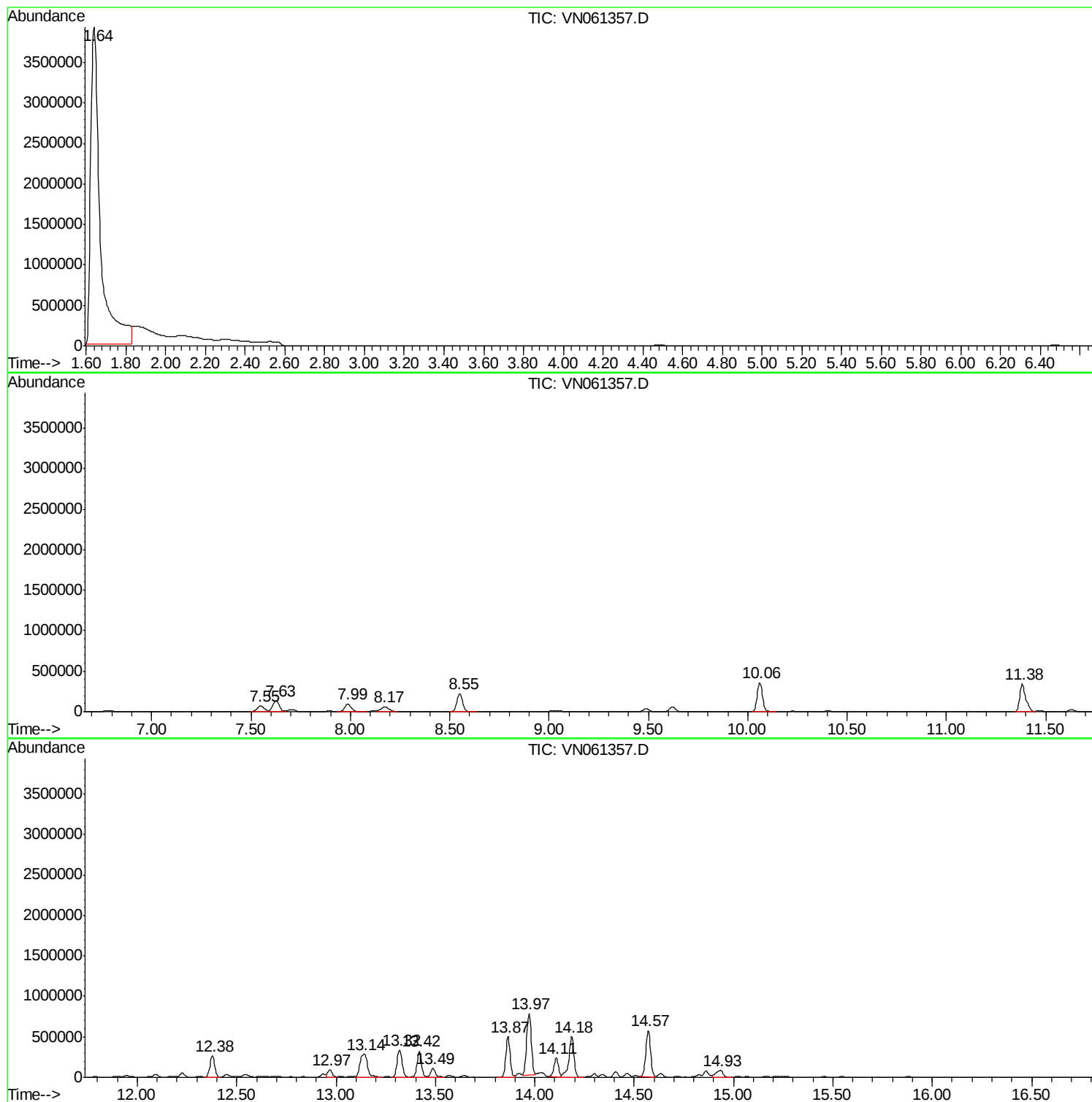
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N042720W.M
Quant Title : SW846 8260

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



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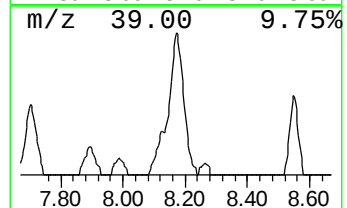
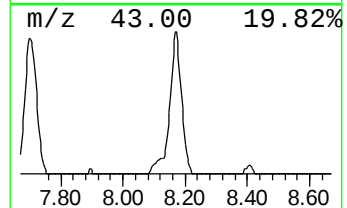
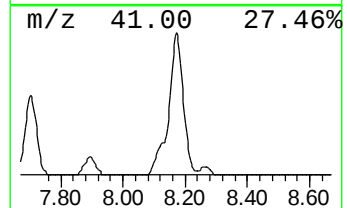
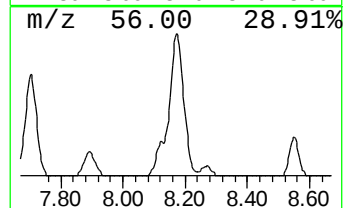
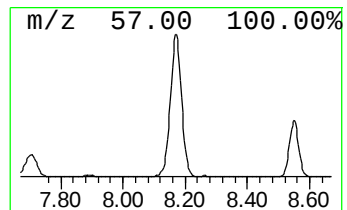
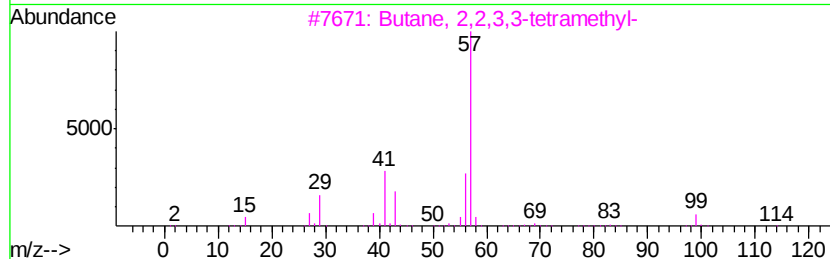
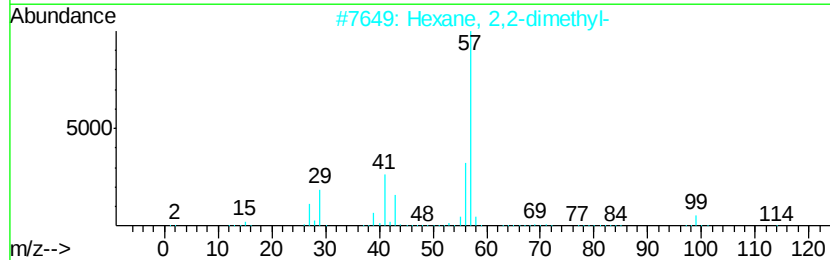
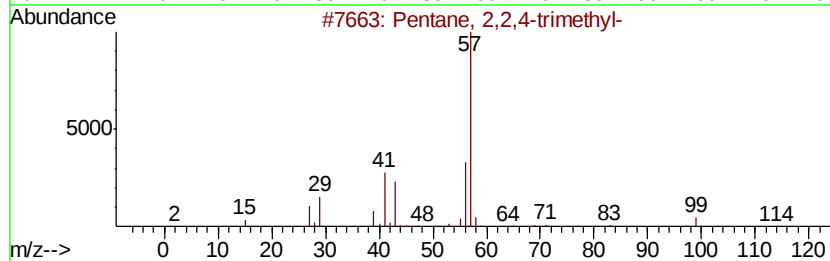
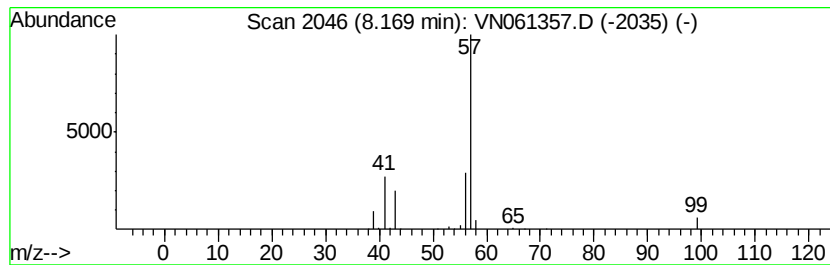
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Quant Title : SW846 8260

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

Peak Number 2 Pentane, 2,2,4-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.17	18.38 ug/l	168504	1,4-Difluorobenzene	8.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	64
2		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	64
3		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	43
4		Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	39
5		Acetamide, N-2-propenyl-	99	C5H9NO	000692-33-1	23



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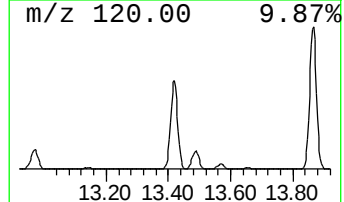
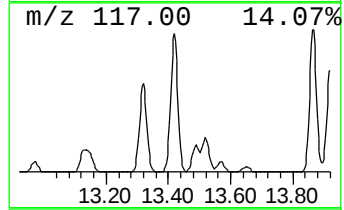
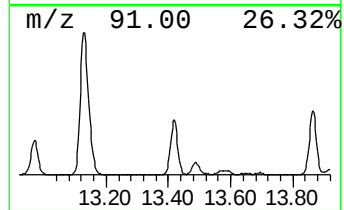
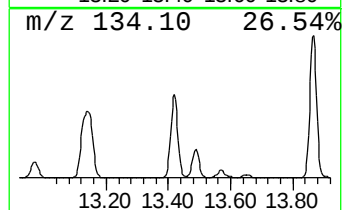
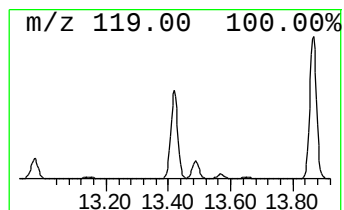
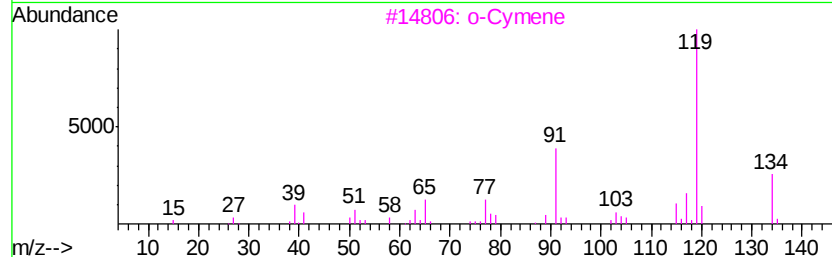
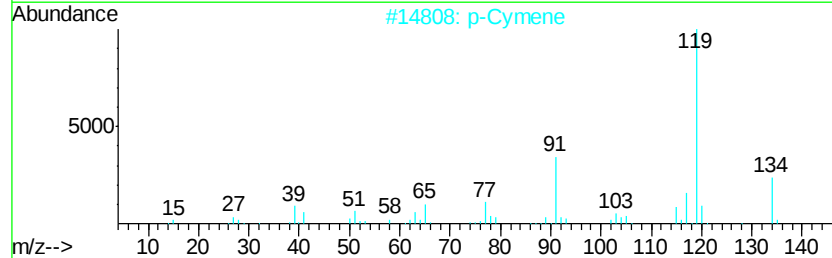
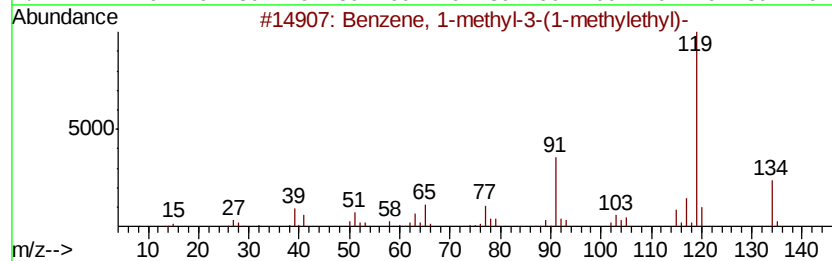
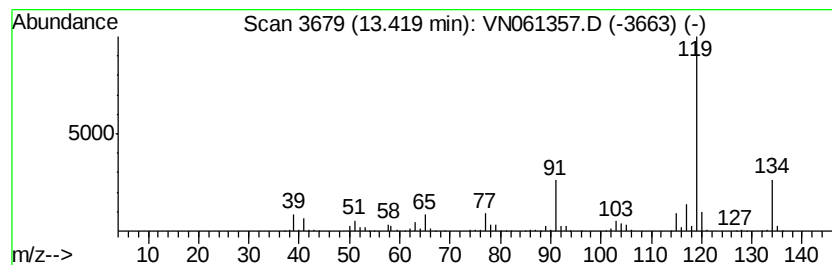
Instrument :
MSVOA_N
ClientSampled :
ST008MW037-200504

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Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-methyl-3-(1-meth... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.42	45.01 ug/l	524313	1,4-Dichlorobenzene-d4	13.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1-methyl-3-(1-methyleth...	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, 1-ethyl-2,3-dimethyl-	134 C10H14	000933-98-2	91
5	Benzene, 4-ethyl-1,2-dimethyl-	134 C10H14	000934-80-5	91



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Instrument :
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ClientSampleId :
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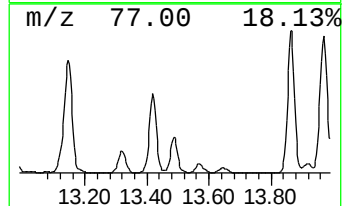
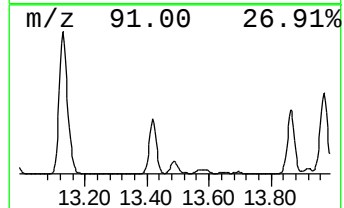
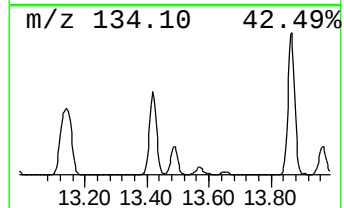
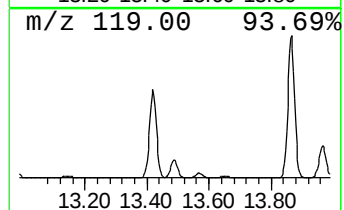
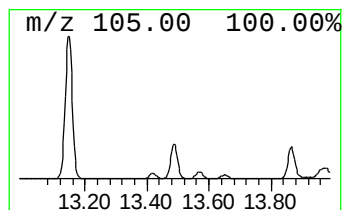
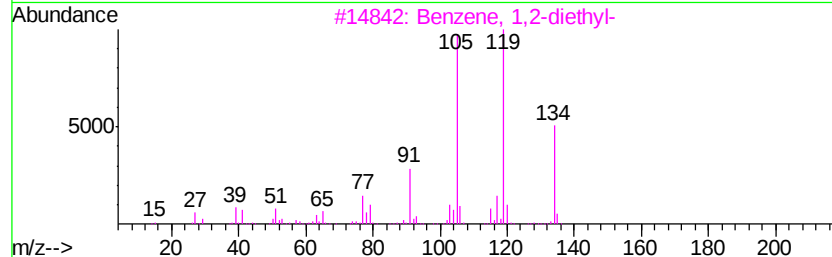
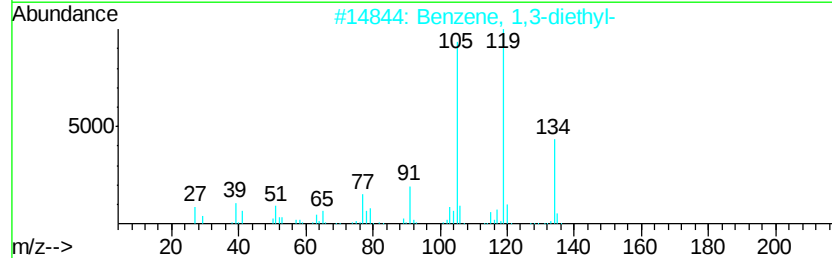
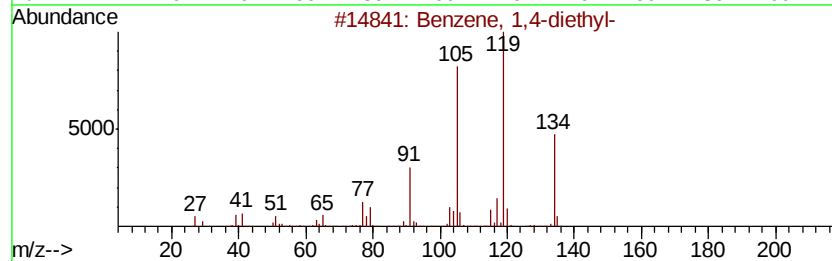
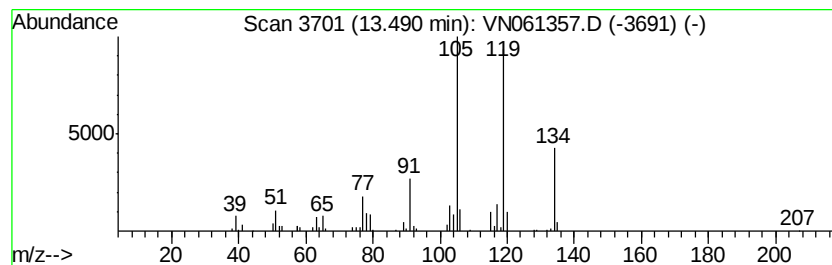
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TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1,4-diethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.49	16.10 ug/l	187611	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	97
2		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
3		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



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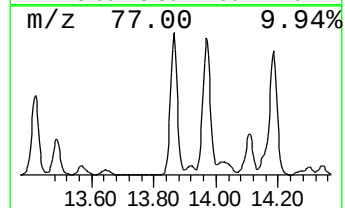
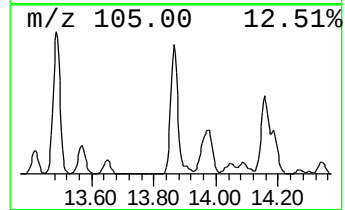
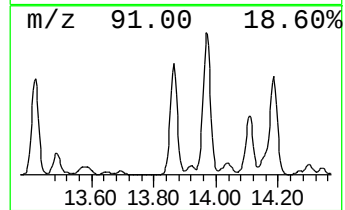
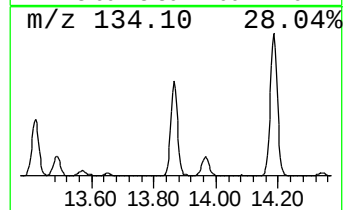
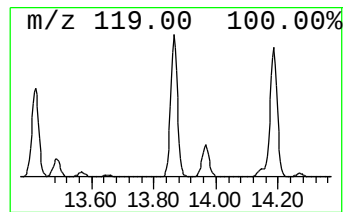
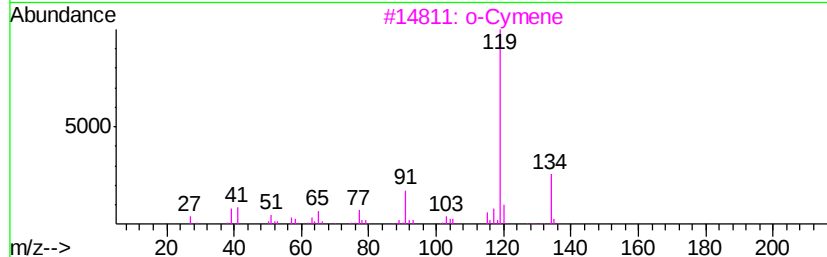
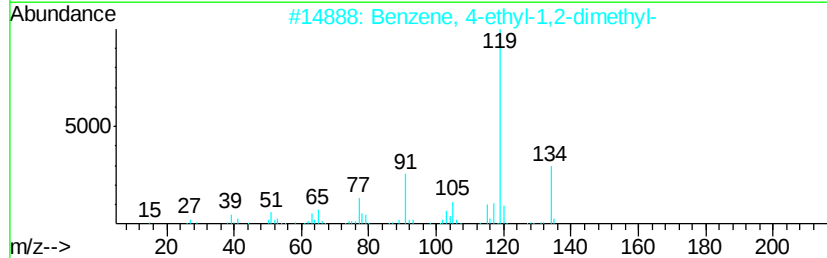
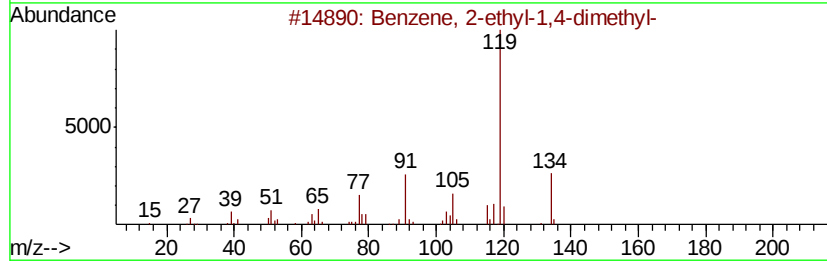
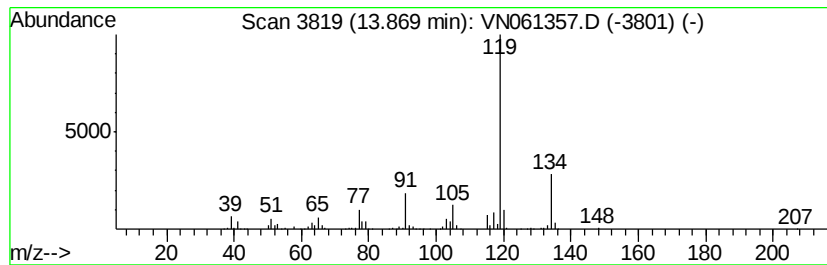
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	67.32 ug/l	784260	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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		p-Cymene	134	C10H14	000099-87-6	95
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95



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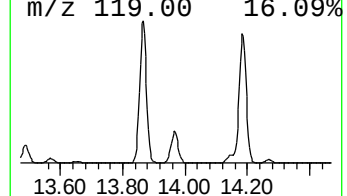
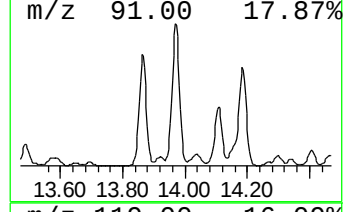
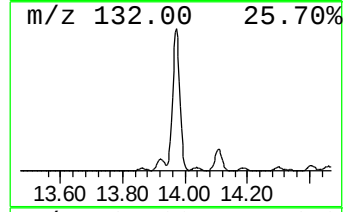
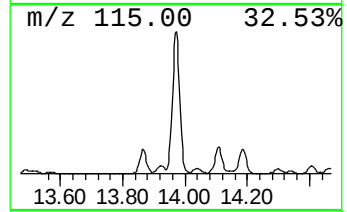
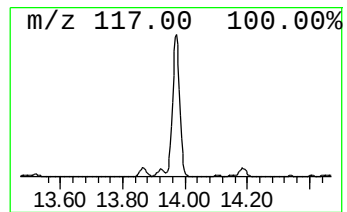
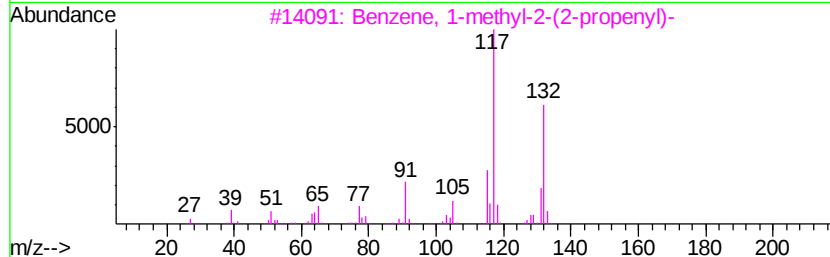
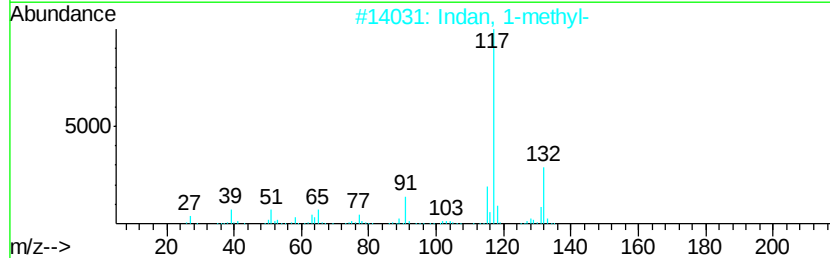
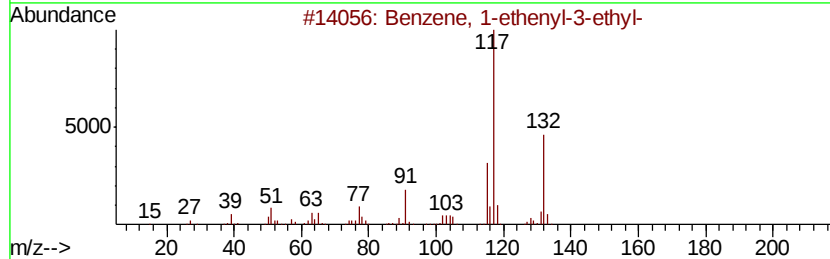
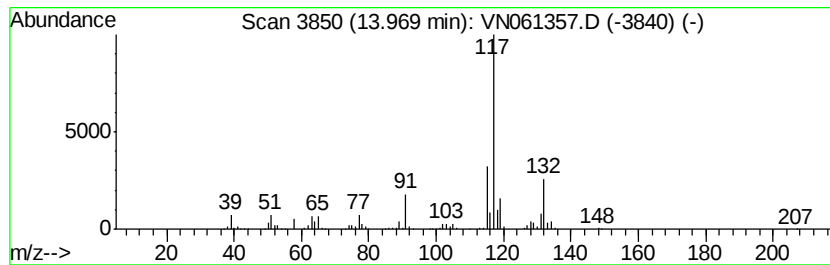
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	104.83 ug/l	1221200	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
2		Indan, 1-methyl-	132	C10H12	000767-58-8	81
3		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	81
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	80
5		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	74



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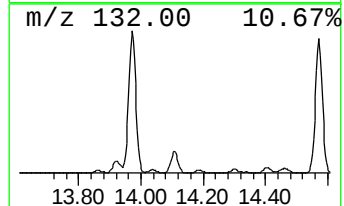
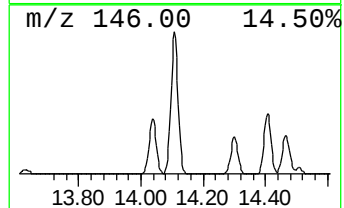
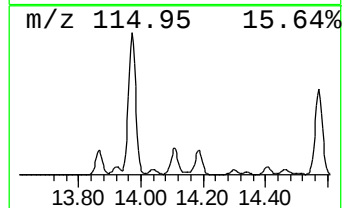
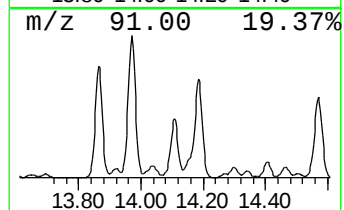
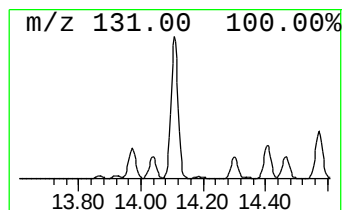
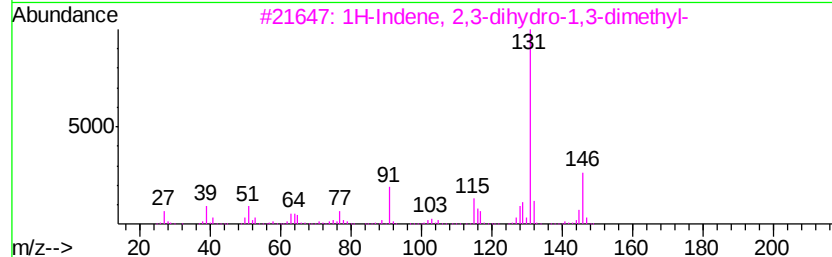
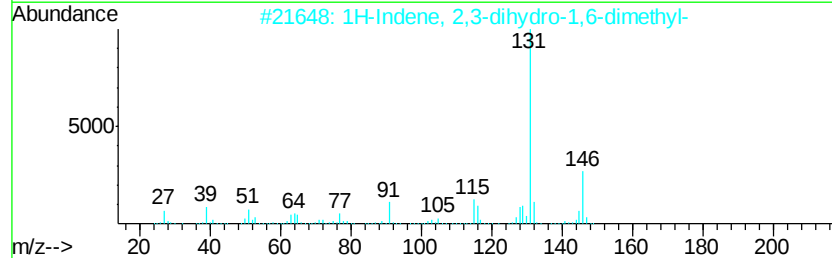
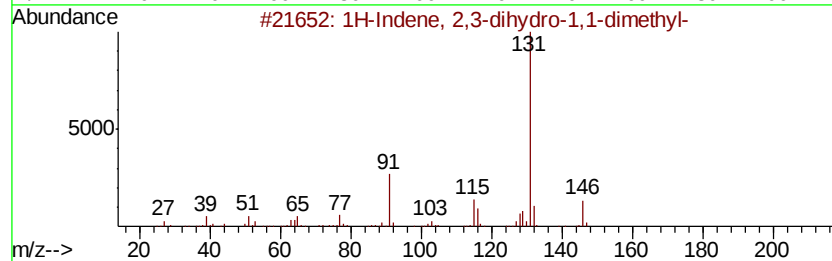
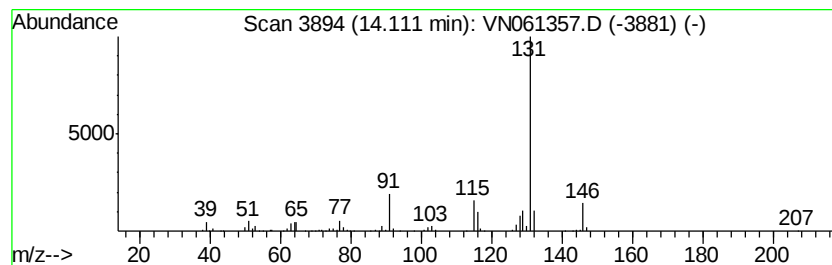
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Quant Title : SW846 8260

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

Peak Number 7 1H-Indene, 2,3-dihydro-1,1-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.11	32.60 ug/l	379799	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	97
2		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	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		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN050820\
Data File : VN061357.D
Acq On : 8 May 2020 13:26
Operator : JC/MD
Sample : L2544-05
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
ST008MW037-200504

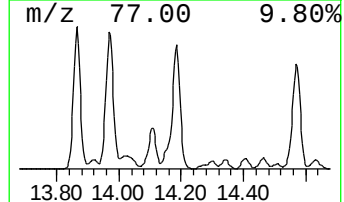
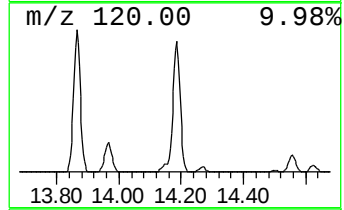
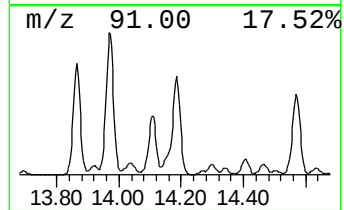
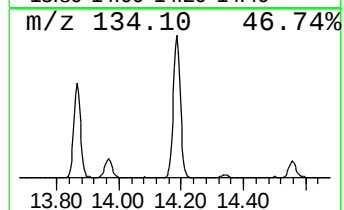
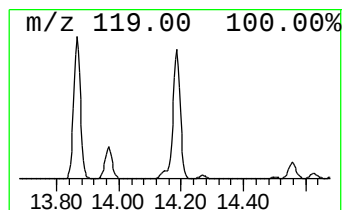
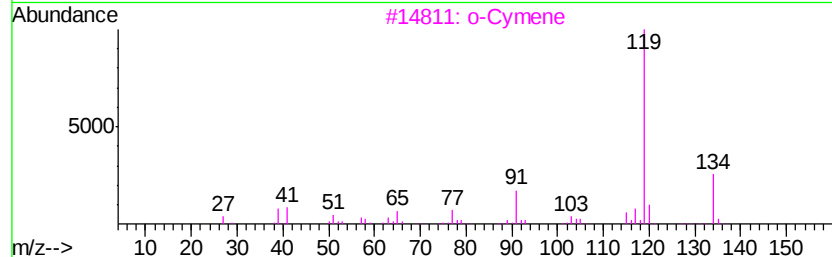
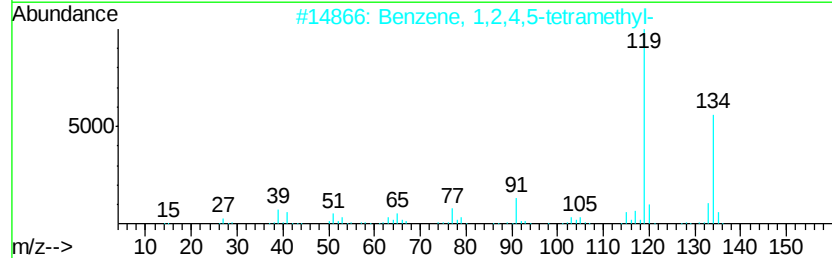
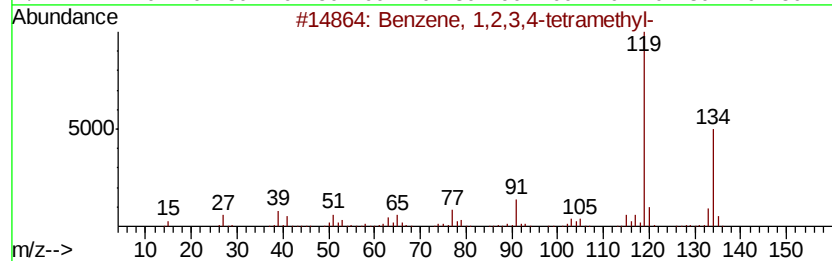
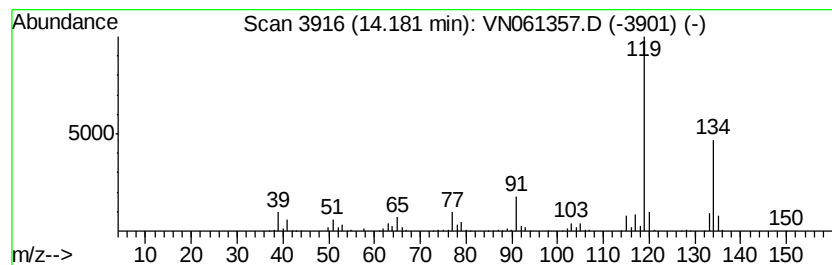
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.18	78.67 ug/l	916435	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
3		o-Cymene	134	C10H14	000527-84-4	95
4		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
5		p-Cymene	134	C10H14	000099-87-6	94



Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN050820\
Data File : VN061357.D
Acq On : 8 May 2020 13:26
Operator : JC/MD
Sample : L2544-05
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampled :
ST008MW037-200504

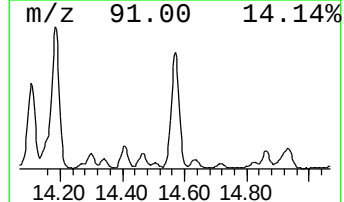
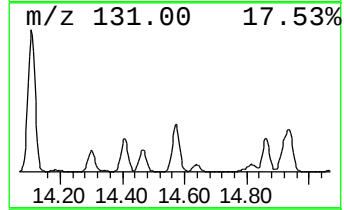
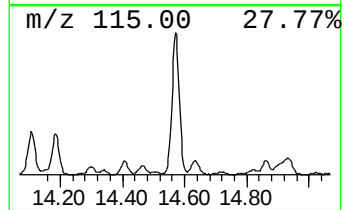
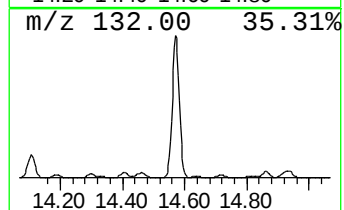
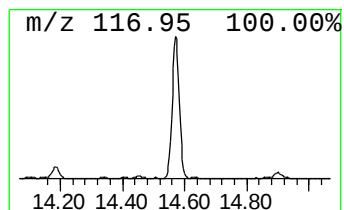
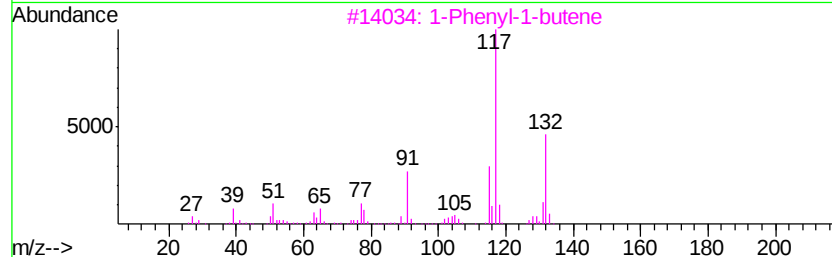
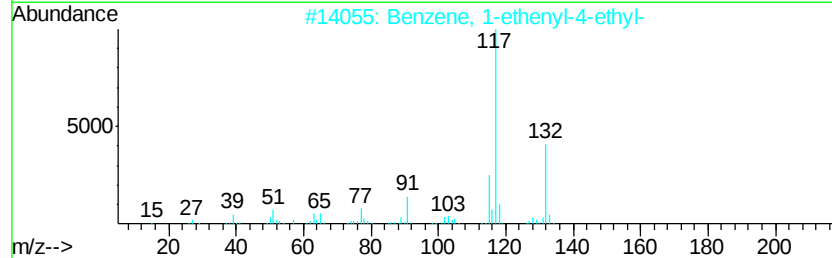
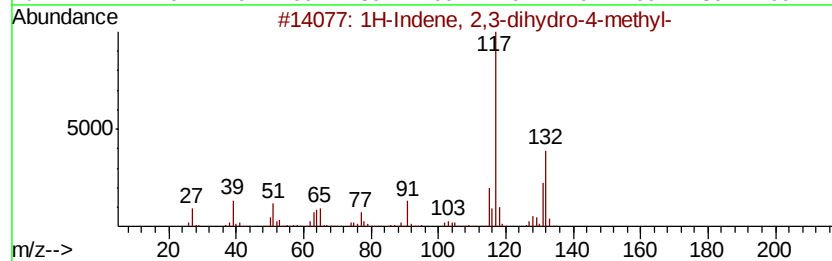
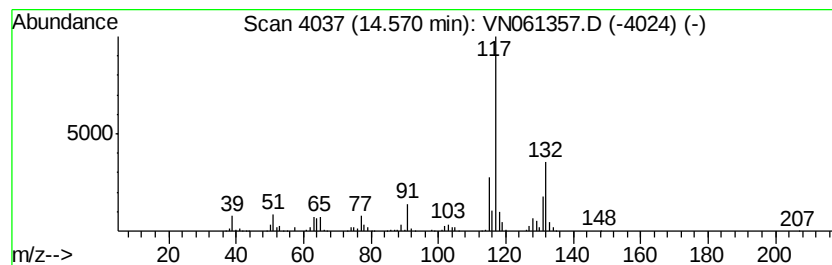
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Quant Title : SW846 8260

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

Peak Number 9 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	84.00 ug/l	978573	1,4-Dichlorobenzene-d4	13.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3		1-Phenyl-1-butene	132	C10H12	000824-90-8	94
4		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN050820\
Data File : VN061357.D
Acq On : 8 May 2020 13:26
Operator : JC/MD
Sample : L2544-05
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 9 Sample Multiplier: 1

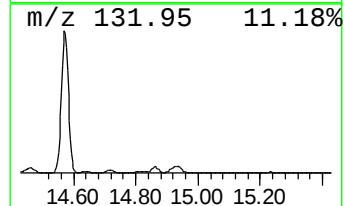
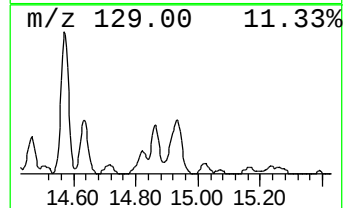
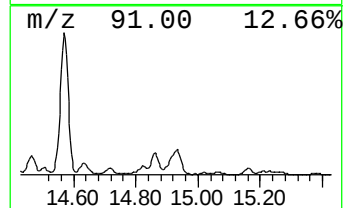
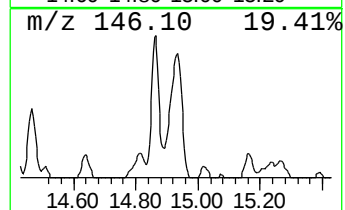
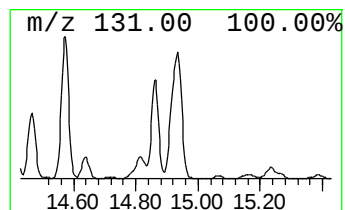
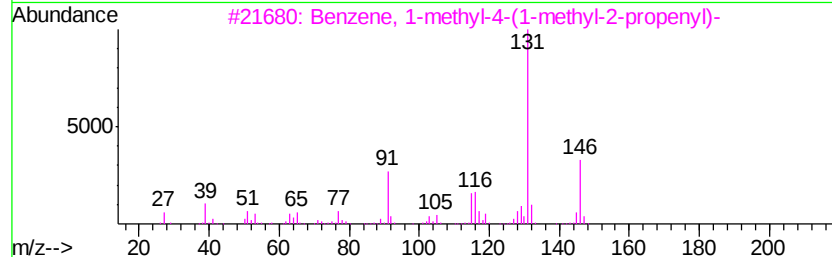
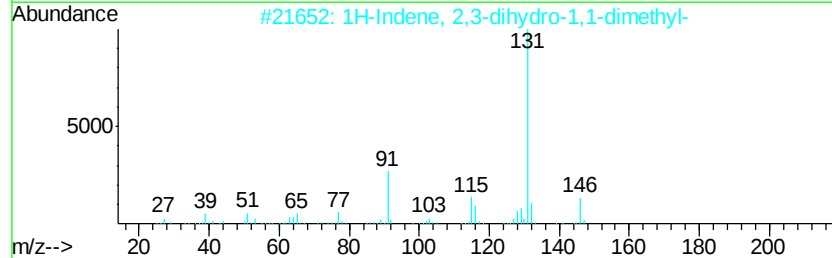
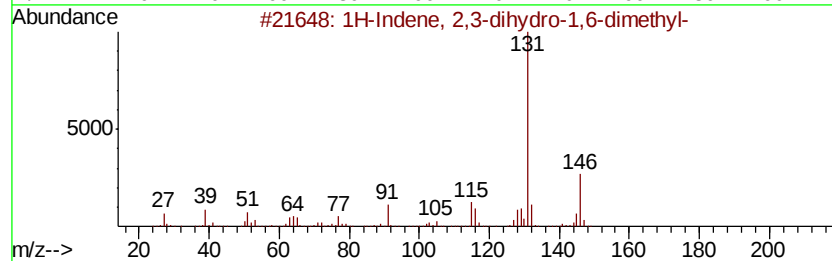
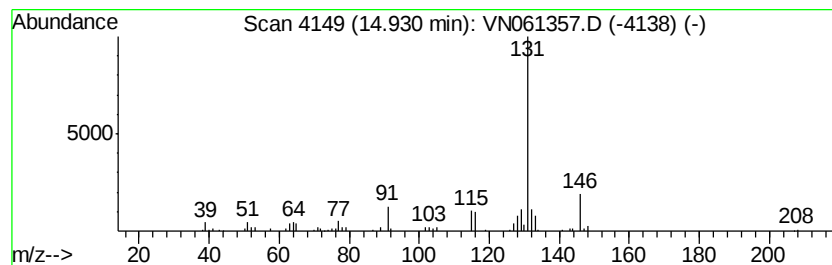
Instrument :
MSVOA_N
ClientSampleId :
ST008MW037-200504

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 10 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.93	17.89 ug/l	208457	1,4-Dichlorobenzene-d4	13.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146 C11H14	017059-48-2	91
2	1H-Indene, 2,3-dihydro-1,1-dimet...	146 C11H14	004912-92-9	90
3	Benzene, 1-methyl-4-(1-methyl-2-...	146 C11H14	097664-18-1	90
4	2,2-Dimethylindene, 2,3-dihydro-	146 C11H14	020836-11-7	90
5	1H-Indene, 2,3-dihydro-1,3-dimet...	146 C11H14	004175-53-5	90



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN050820\
Data File : VN061357.D
Acq On : 8 May 2020 13:26
Operator : JC/MD
Sample : L2544-05
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
ST008MW037-200504

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
Pentane, 2,2,4-tr...	8.17	18.4	ug/l	168504	2	8.55	458461	50.0
Benzene, 1-methyl...	13.42	45.0	ug/l	524313	4	13.32	582492	50.0
Benzene, 1,4-diet...	13.49	16.1	ug/l	187611	4	13.32	582492	50.0
Benzene, 2-ethyl-...	13.87	67.3	ug/l	784260	4	13.32	582492	50.0
Benzene, 1-etheny...	13.97	104.8	ug/l	1221200	4	13.32	582492	50.0
1H-Indene, 2,3-di...	14.11	32.6	ug/l	379799	4	13.32	582492	50.0
Benzene, 1,2,3,4-...	14.18	78.7	ug/l	916435	4	13.32	582492	50.0
1H-Indene, 2,3-di...	14.57	84.0	ug/l	978573	4	13.32	582492	50.0
1H-Indene, 2,3-di...	14.93	17.9	ug/l	208457	4	13.32	582492	50.0