

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN071020\
 Data File : VN062504.D
 Acq On : 10 Jul 2020 19:58
 Operator : JC/MD
 Sample : L3217-09 100X
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 E3-41-NHL

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\82N070820W.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	7.551	1778	1797	1807	rBV	132508	333791	1.58%	0.374%
2	7.625	1808	1820	1842	rVB	232410	548050	2.59%	0.614%
3	7.989	1914	1933	1960	rBV2	155311	371665	1.76%	0.416%
4	8.551	2092	2108	2126	rBV	366510	771313	3.64%	0.864%
5	10.059	2563	2577	2590	rBV	603221	1087130	5.14%	1.218%
6	11.381	2975	2988	3009	rBV	554256	959773	4.53%	1.075%
7	11.590	3039	3053	3077	rVB	195607	360688	1.70%	0.404%
8	11.921	3144	3156	3177	rVB2	133588	257392	1.22%	0.288%
9	12.226	3239	3251	3268	rBV	301488	516829	2.44%	0.579%
10	12.377	3280	3298	3312	rBV	440561	751097	3.55%	0.842%
11	12.631	3365	3377	3384	rBV	236527	406182	1.92%	0.455%
12	12.712	3393	3402	3404	rBV	247568	334495	1.58%	0.375%
13	13.014	3481	3496	3510	rBV	474975	851626	4.02%	0.954%
14	13.236	3549	3565	3580	rBV	12220202	19311133	91.23%	21.636%
15	13.320	3580	3591	3608	rVB2	619816	1215654	5.74%	1.362%
16	13.519	3634	3653	3663	rBV	1395509	2355357	11.13%	2.639%
17	13.699	3697	3709	3722	rVB	878110	1435553	6.78%	1.608%
18	13.969	3784	3793	3804	rBV	177729	271436	1.28%	0.304%
19	14.040	3804	3815	3827	rVB	390620	588689	2.78%	0.660%
20	14.156	3840	3851	3866	rBV3	171562	372360	1.76%	0.417%
21	14.249	3866	3880	3895	rVV2	303616	728112	3.44%	0.816%
22	14.451	3933	3943	3953	rBV	241687	394763	1.86%	0.442%
23	14.567	3966	3979	3990	rBV2	410604	765167	3.61%	0.857%
24	14.631	3990	3999	4012	rVB	235614	394912	1.87%	0.442%
25	14.715	4013	4025	4040	rVB	321869	538165	2.54%	0.603%
26	15.033	4109	4124	4130	rBV5	109140	242854	1.15%	0.272%
27	15.104	4133	4146	4163	rVB	9412082	17372094	82.07%	19.463%
28	15.197	4163	4175	4190	rBV	827023	1543281	7.29%	1.729%
29	16.062	4429	4444	4449	rBV2	206209	428807	2.03%	0.480%
30	16.123	4449	4463	4482	rVB	10008898	21168300	100.00%	23.717%
31	16.239	4489	4499	4508	rBV	181879	385647	1.82%	0.432%
32	16.313	4508	4522	4544	rVB	5587812	12193028	57.60%	13.661%

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ClientSampleId :
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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_N\METHODS\82N070820W.M
Title : SW846 8260

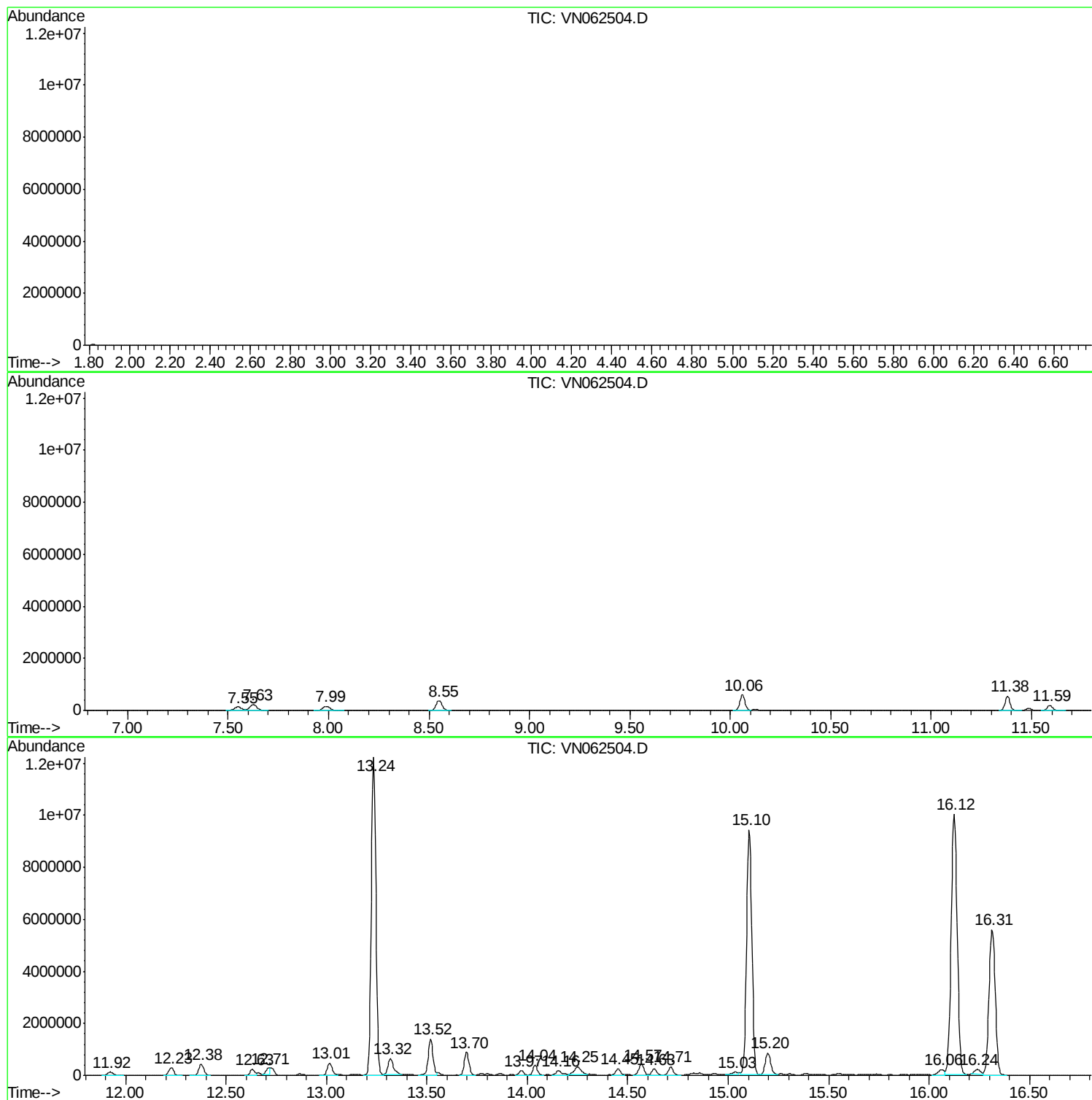
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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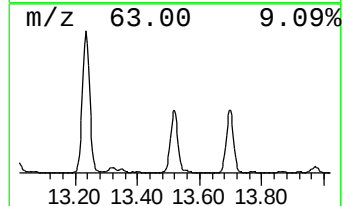
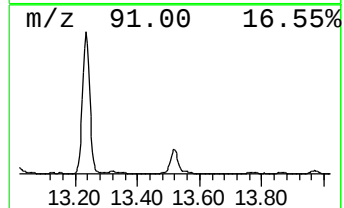
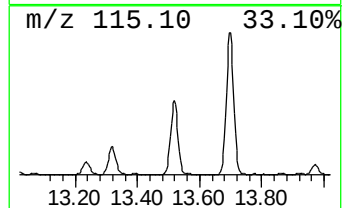
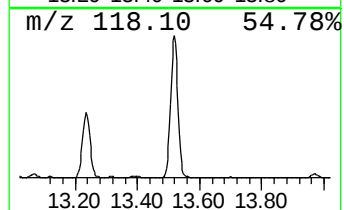
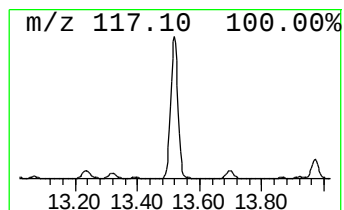
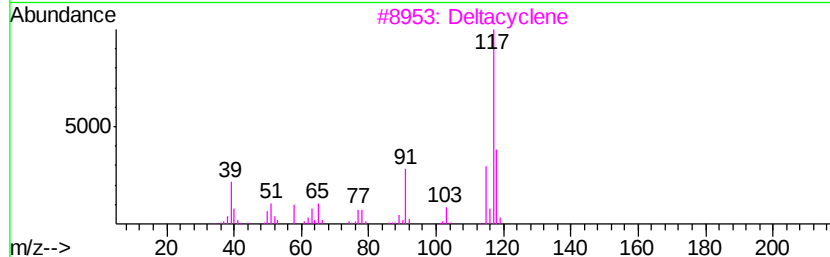
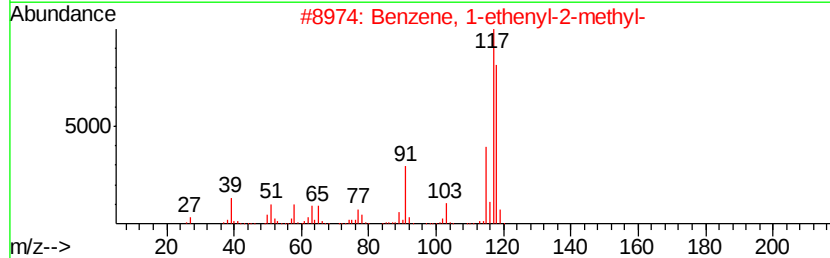
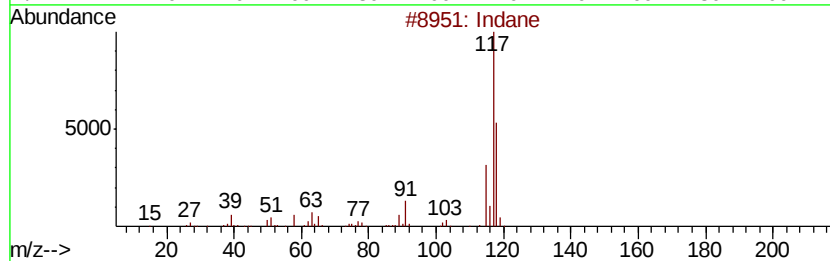
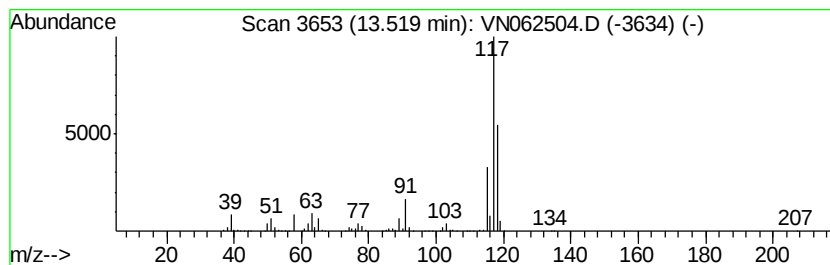
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.M
Quant Title : SW846 8260

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

Peak Number 2 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.52	96.88 ug/l	2355360	1,4-Dichlorobenzene-d4	13.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	94
2	Benzene, 1-ethenyl-2-methyl-		118	C9H10	000611-15-4	78
3	Deltacyclene		118	C9H10	007785-10-6	68
4	Benzene, cyclopropyl-		118	C9H10	000873-49-4	64
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...		118	C9H10	1000191-13-7	64



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Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampled :
E3-41-NHL

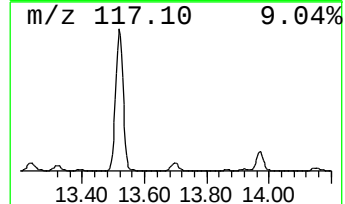
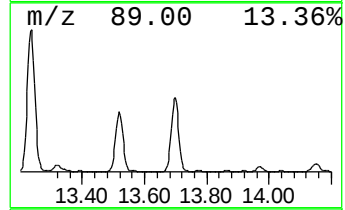
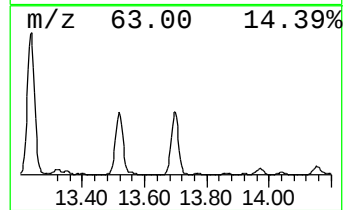
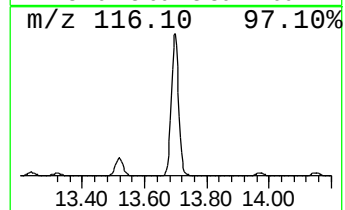
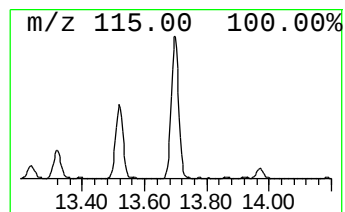
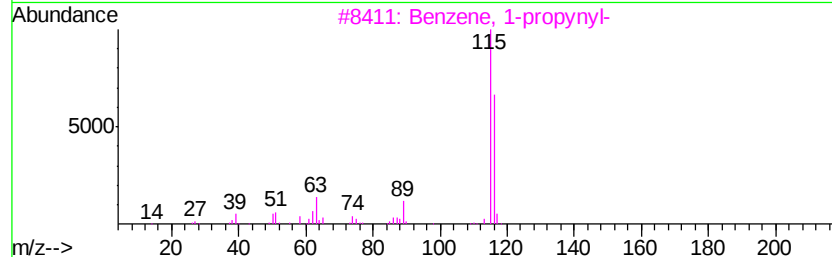
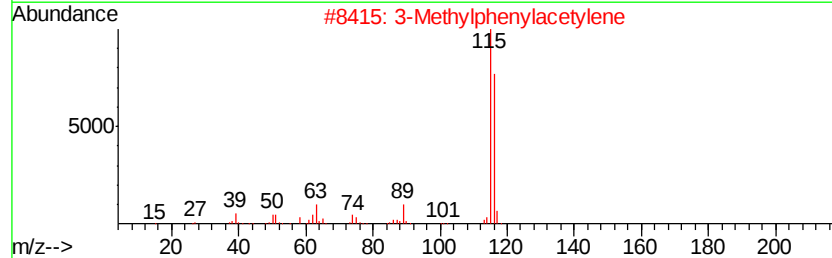
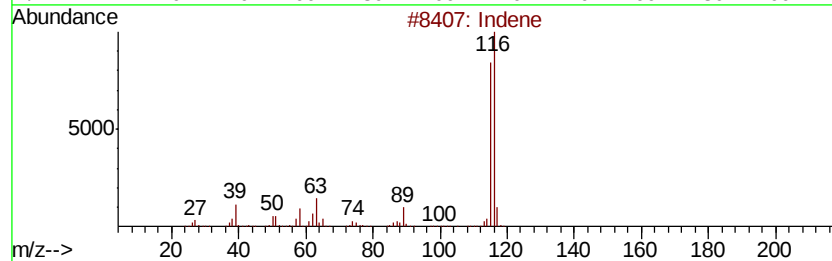
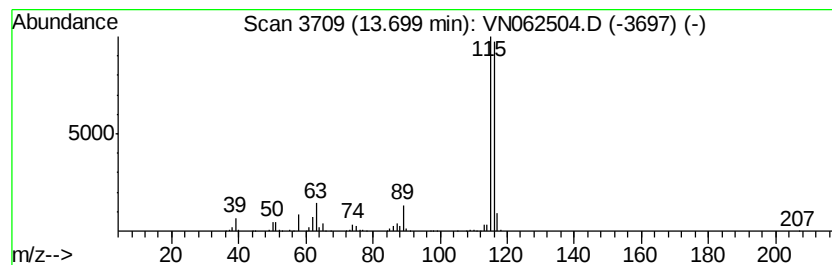
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.M
Quant Title : SW846 8260

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

Peak Number 3 Indene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.70	59.04 ug/l	1435550	1,4-Dichlorobenzene-d4	13.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene		116	C9H8	000095-13-6	96
2	3-Methylphenylacetylene		116	C9H8	000766-82-5	94
3	Benzene, 1-propynyl-		116	C9H8	000673-32-5	91
4	Benzene, 1,2-propadienyl-		116	C9H8	002327-99-3	91
5	Benzene, 1-ethynyl-4-methyl-		116	C9H8	000766-97-2	91



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Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 24 Sample Multiplier: 1

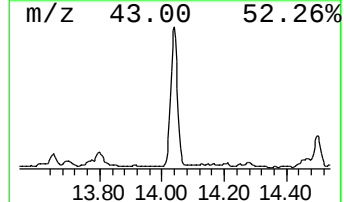
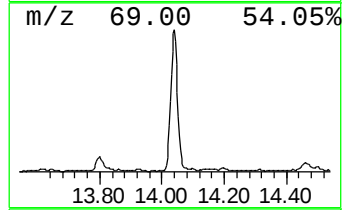
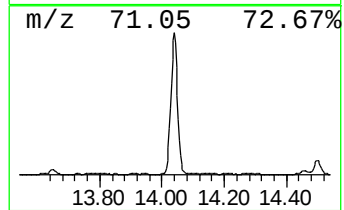
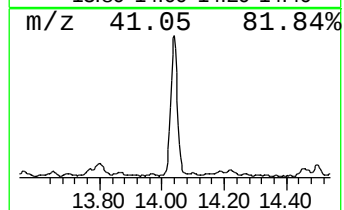
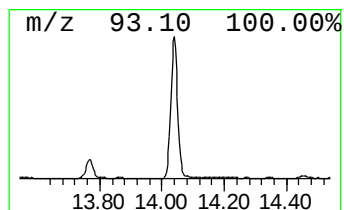
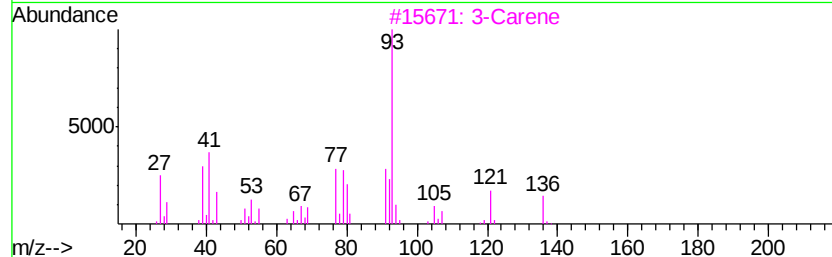
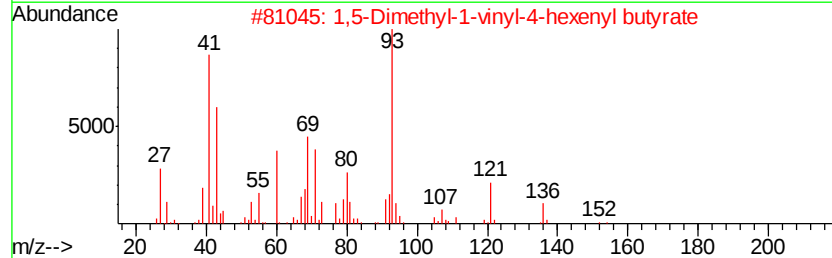
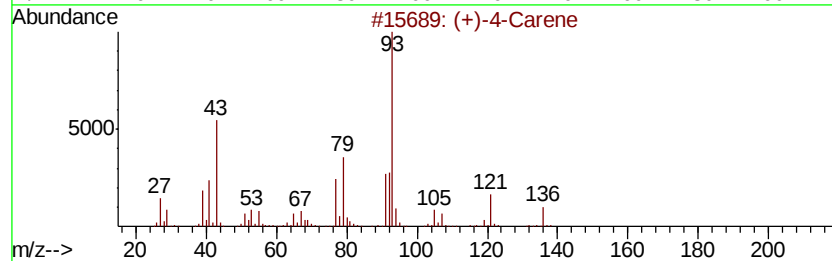
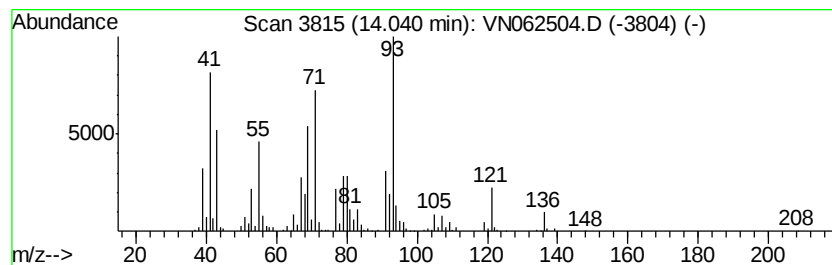
Instrument :
MSVOA_N
ClientSampleId :
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Quant Title : SW846 8260

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

Peak Number 4 (+)-4-Carene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.04	24.21 ug/l	588689	1,4-Dichlorobenzene-d4	13.32	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(+)-4-Carene	136	C10H16	029050-33-7	90
2	1,5-Dimethyl-1-vinyl-4-hexenyl b...	224	C14H24O2	000078-36-4	72
3	3-Carene	136	C10H16	013466-78-9	64
4	.beta.-Myrcene	136	C10H16	000123-35-3	64
5	Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000508-32-7	64



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Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
E3-41-NHL

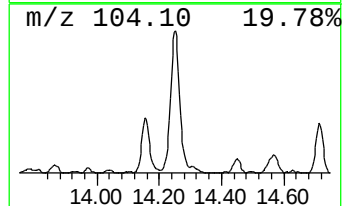
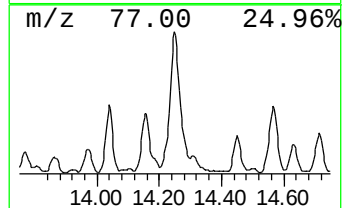
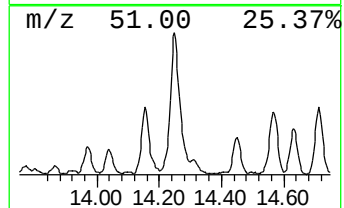
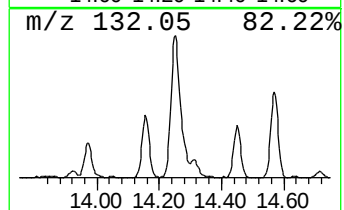
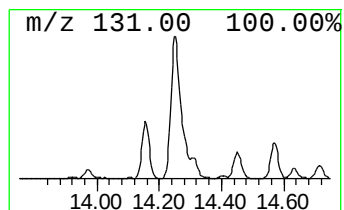
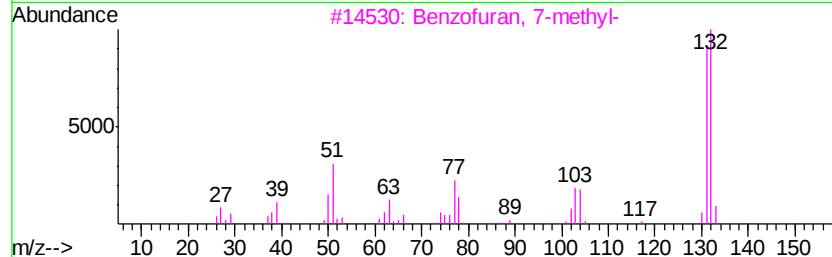
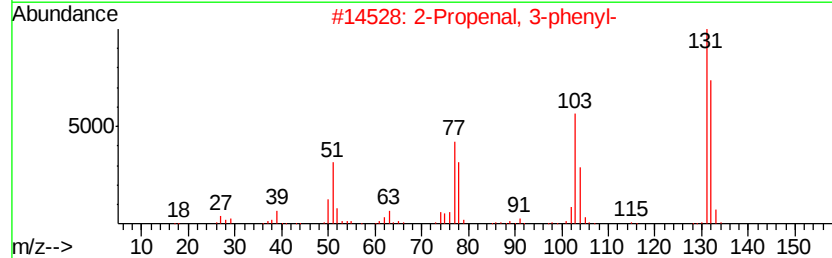
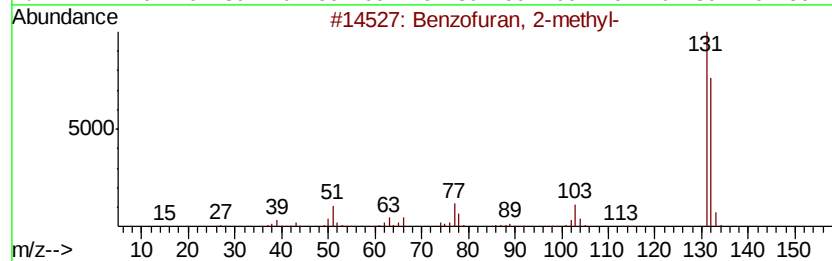
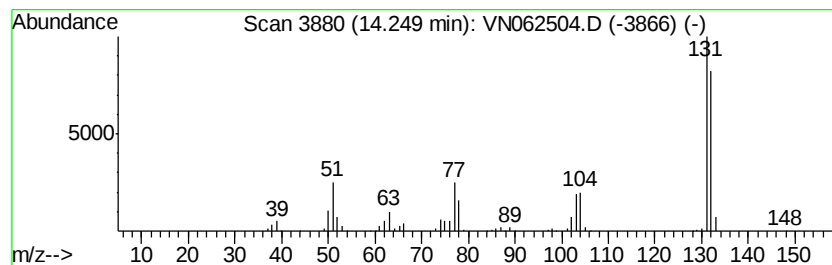
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzofuran, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.25	29.95 ug/l	728112	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	95
2		2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	94
3		Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	94
4		3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	90
5		Benzene, (2-propynyloxy)-	132	C9H8O	013610-02-1	90



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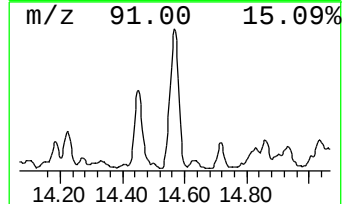
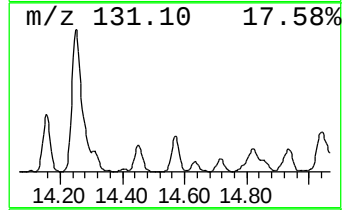
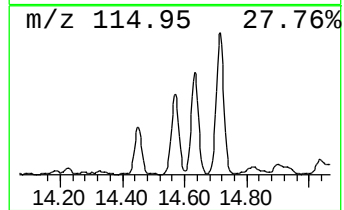
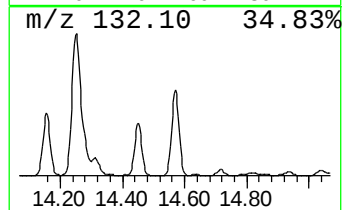
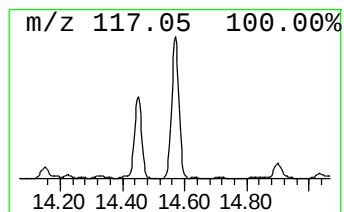
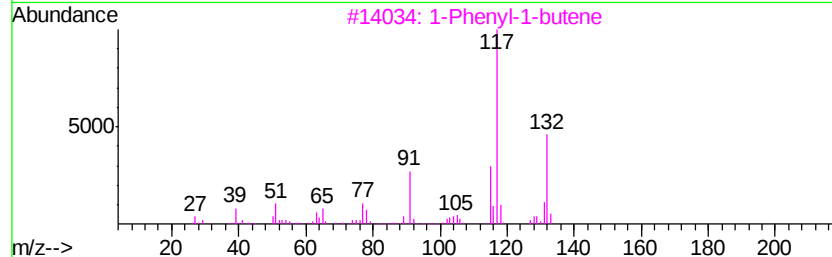
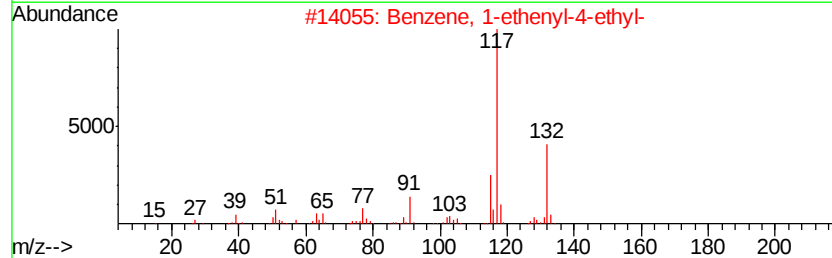
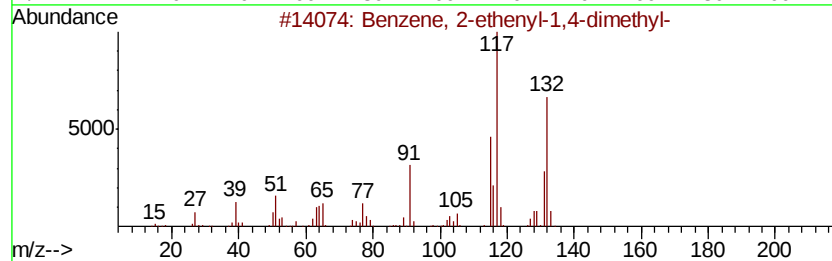
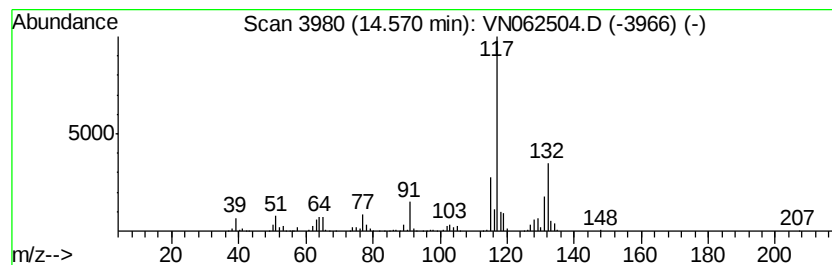
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	31.47 ug/l	765167	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	93
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN071020\
Data File : VN062504.D
Acq On : 10 Jul 2020 19:58
Operator : JC/MD
Sample : L3217-09 100X
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
E3-41-NHL

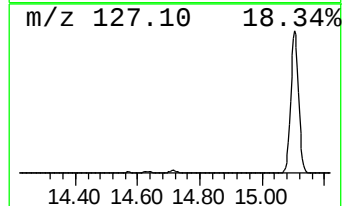
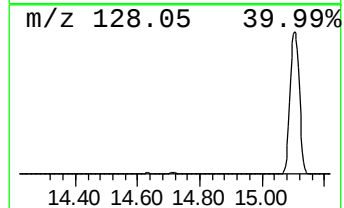
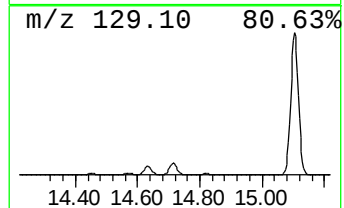
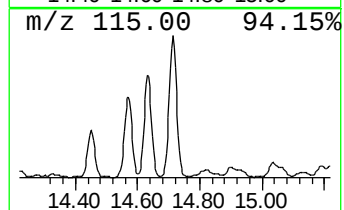
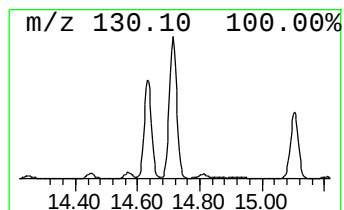
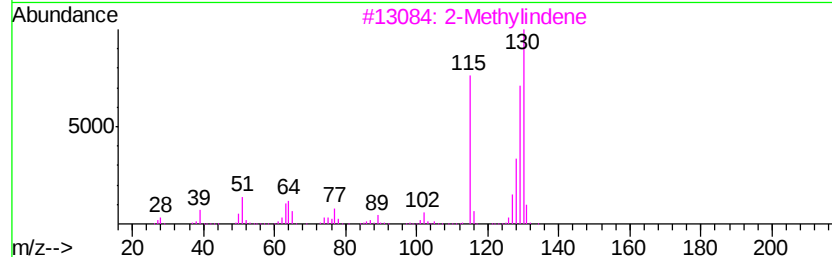
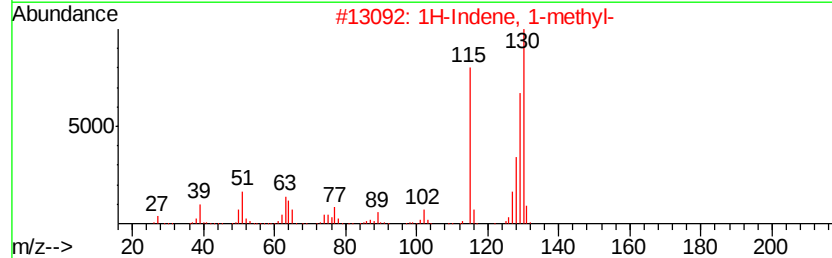
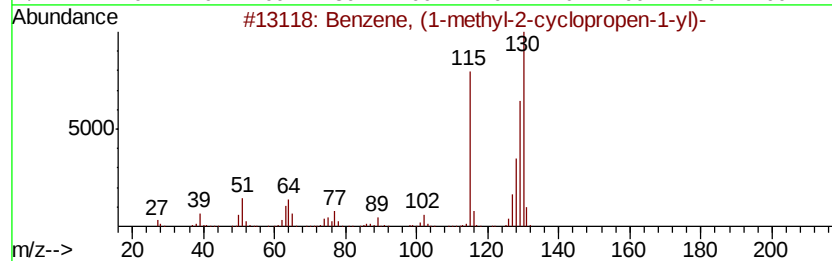
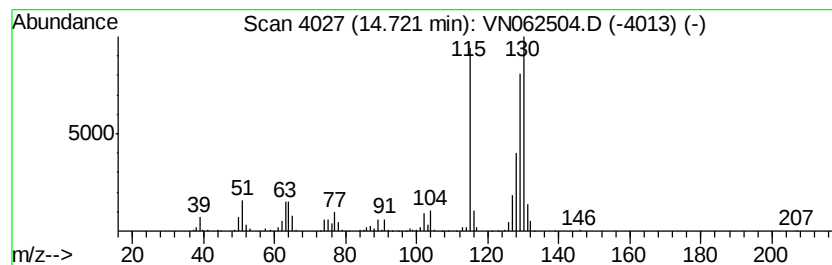
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, (1-methyl-2-cyclop... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.72	22.13 ug/l	538165	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	96
2		1H-Indene, 1-methyl-	130	C10H10	000767-59-9	96
3		2-Methylindene	130	C10H10	002177-47-1	95
4		Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	95
5		1,4-Dihydronaphthalene	130	C10H10	000612-17-9	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN071020\
Data File : VN062504.D
Acq On : 10 Jul 2020 19:58
Operator : JC/MD
Sample : L3217-09 100X
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
E3-41-NHL

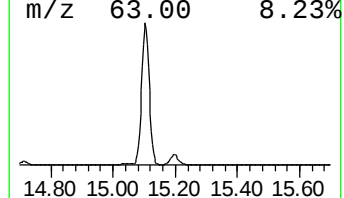
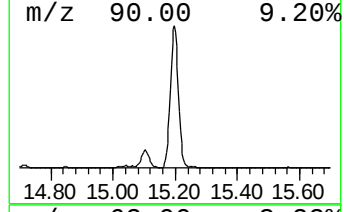
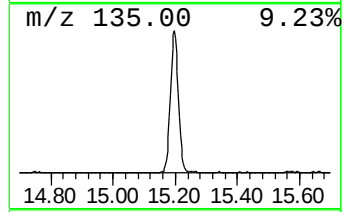
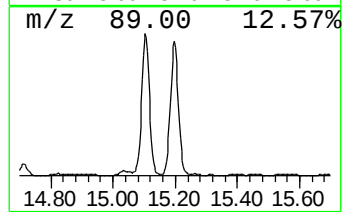
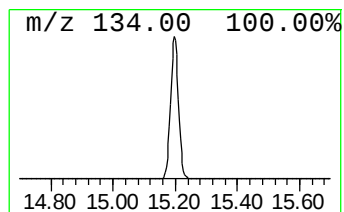
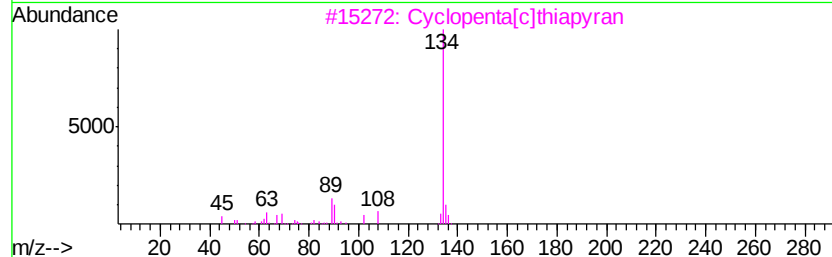
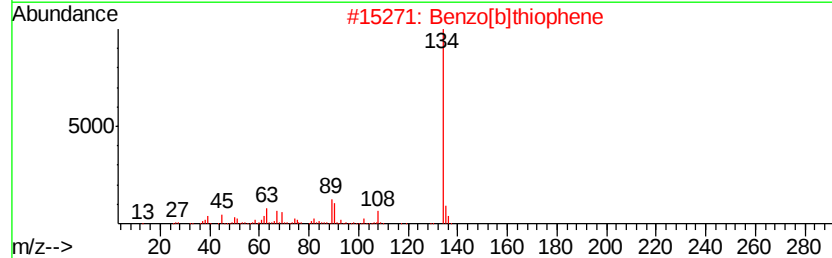
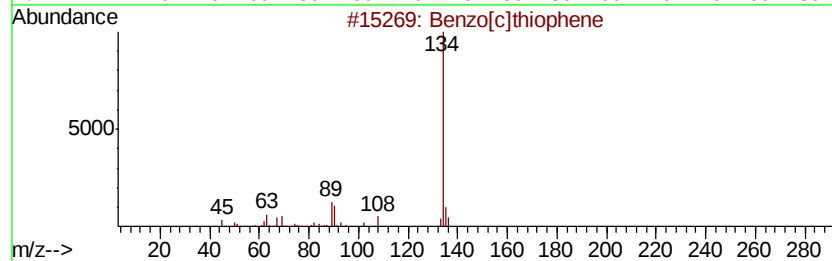
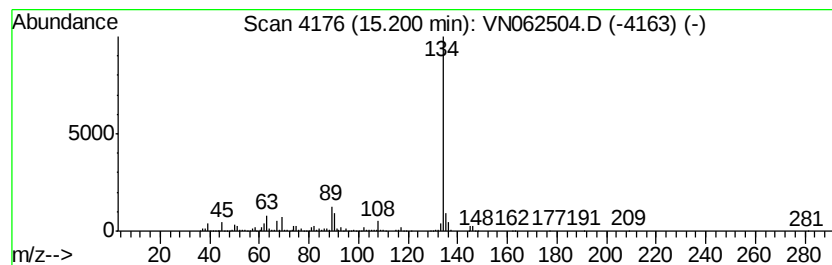
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Quant Title : SW846 8260

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

Peak Number 8 Benzo[c]thiophene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.20	63.48 ug/l	1543280	1,4-Dichlorobenzene-d4	13.32

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[c]thiophene	134	C8H6S	000270-82-6	97
2	Benzo[b]thiophene	134	C8H6S	000095-15-8	95
3	Cyclopenta[c]thiopyran	134	C8H6S	000270-63-3	94
4	5H-Pyrrolo(3,2-d)pyrimidin-4-amine	134	C6H6N4	002227-98-7	42
5	[1]Benzothieno[2',3':3,4]cyclobu...	246	C13H10O3S	034002-19-2	39



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN071020\
Data File : VN062504.D
Acq On : 10 Jul 2020 19:58
Operator : JC/MD
Sample : L3217-09 100X
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 24 Sample Multiplier: 1

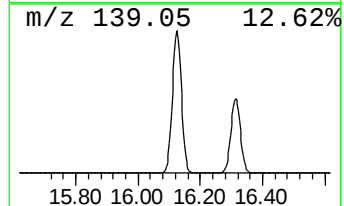
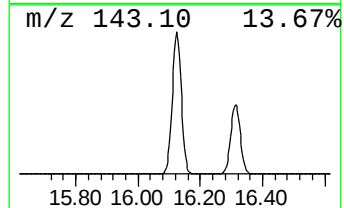
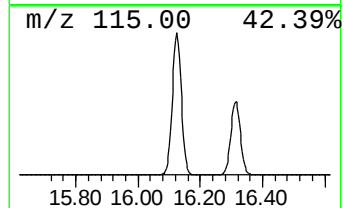
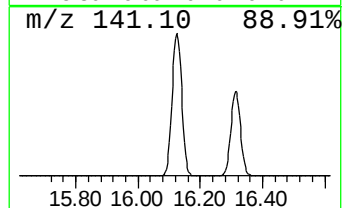
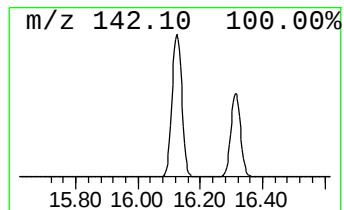
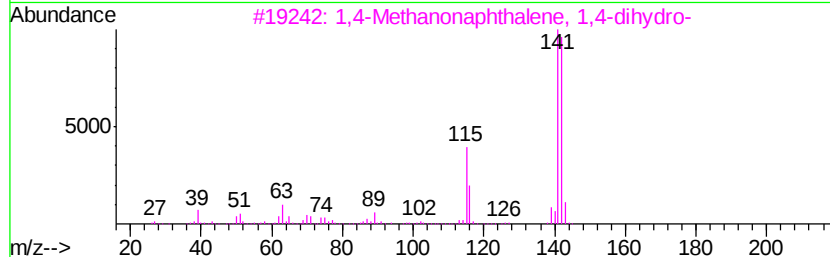
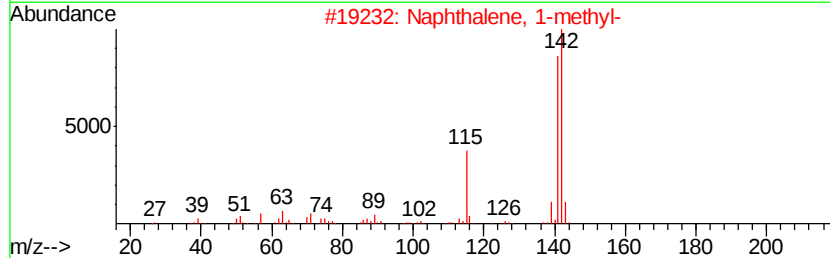
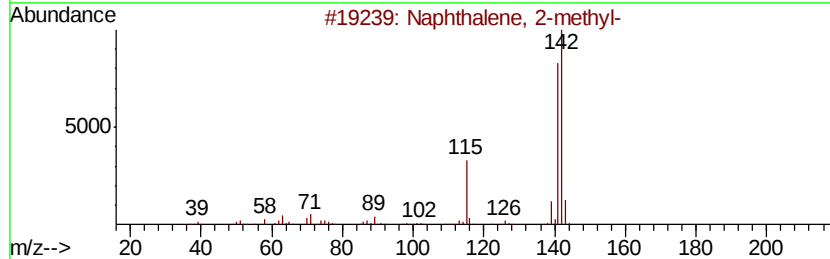
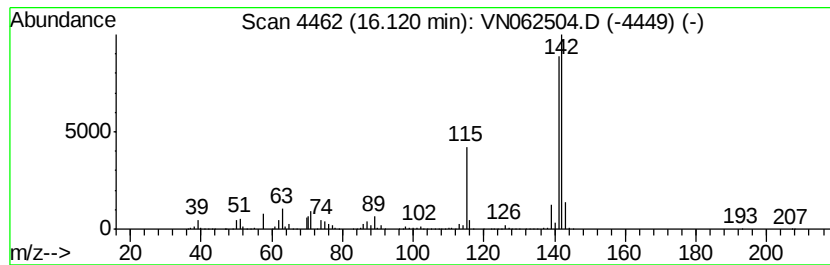
Instrument :
MSVOA_N
ClientSampleId :
E3-41-NHL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.12	870.66 ug/l	21168300	1,4-Dichlorobenzene-d4	13.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2-methyl-	142 C11H10	000091-57-6 96
2		Naphthalene, 1-methyl-	142 C11H10	000090-12-0 96
3		1,4-Methanonaphthalene, 1,4-dihv...	142 C11H10	004453-90-1 93
4		Benzocycloheptatriene	142 C11H10	000264-09-5 93
5		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142 C11H10	002443-46-1 91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN071020\
Data File : VN062504.D
Acq On : 10 Jul 2020 19:58
Operator : JC/MD
Sample : L3217-09 100X
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
E3-41-NHL

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.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
D-Limonene	13.24	794.3	ug/l	19311100	4	13.32	1215650	50.0
Indane	13.52	96.9	ug/l	2355360	4	13.32	1215650	50.0
Indene	13.70	59.0	ug/l	1435550	4	13.32	1215650	50.0
(+)-4-Carene	14.04	24.2	ug/l	588689	4	13.32	1215650	50.0
Benzofuran, 2-met...	14.25	29.9	ug/l	728112	4	13.32	1215650	50.0
Benzene, 2-etheny...	14.57	31.5	ug/l	765167	4	13.32	1215650	50.0
Benzene, (1-methy...	14.72	22.1	ug/l	538165	4	13.32	1215650	50.0
Benzo[c]thiophene	15.20	63.5	ug/l	1543280	4	13.32	1215650	50.0
Naphthalene, 1-me...	16.12	870.7	ug/l	21168300	4	13.32	1215650	50.0