

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN031819\
 Data File : VN054423.D
 Acq On : 19 Mar 2019 5:29
 Operator : JC/SP
 Sample : K1970-04 10X
 Misc : 6.33g/10mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 P-4-E.WALL-S.CENTER

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\82N031219W.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	7.590	1788	1809	1820	rBV	1031019	2587409	33.29%	3.339%
2	7.664	1820	1832	1874	rVB	1860639	4664253	60.01%	6.020%
3	8.027	1928	1945	1975	rBV	1002987	2351406	30.25%	3.035%
4	8.587	2103	2119	2173	rVB	2573969	5828846	74.99%	7.523%
5	10.091	2572	2587	2633	rBV	3766127	7772847	100.00%	10.032%
6	11.005	2862	2871	2882	rVB4	78468	139923	1.80%	0.181%
7	11.201	2923	2932	2953	rVB8	75673	183086	2.36%	0.236%
8	11.297	2953	2962	2976	rBV	70799	129844	1.67%	0.168%
9	11.406	2982	2996	3032	rBV	3347661	6313875	81.23%	8.149%
10	11.654	3065	3073	3082	rBV4	55079	104628	1.35%	0.135%
11	11.924	3142	3157	3169	rBV7	138534	390923	5.03%	0.505%
12	12.043	3187	3194	3203	rVB	154288	212832	2.74%	0.275%
13	12.120	3210	3218	3222	rBV3	118090	180098	2.32%	0.232%
14	12.156	3224	3229	3240	rVB4	180218	314351	4.04%	0.406%
15	12.336	3280	3285	3294	rVB2	173459	239683	3.08%	0.309%
16	12.400	3294	3305	3327	rBV	2788982	4897176	63.00%	6.321%
17	12.561	3348	3355	3371	rVB4	272193	574058	7.39%	0.741%
18	12.644	3372	3381	3391	rBV8	90286	201476	2.59%	0.260%
19	12.728	3394	3407	3417	rBV7	294667	780565	10.04%	1.007%
20	12.866	3445	3450	3457	rVB6	111422	142766	1.84%	0.184%
21	12.963	3473	3480	3493	rVB	519554	849064	10.92%	1.096%
22	13.236	3553	3565	3573	rBV5	359520	763466	9.82%	0.985%
23	13.342	3587	3598	3616	rVB	3323155	5672793	72.98%	7.322%
24	13.599	3671	3678	3687	rBV3	149394	354052	4.55%	0.457%
25	13.664	3689	3698	3707	rVV5	625581	1188407	15.29%	1.534%
26	13.889	3760	3768	3777	rVB2	968303	1510939	19.44%	1.950%
27	13.992	3792	3800	3811	rVB3	780160	1313597	16.90%	1.695%
28	14.172	3847	3856	3863	rBV6	659996	1381132	17.77%	1.783%
29	14.291	3886	3893	3899	rBV3	373908	576836	7.42%	0.744%
30	14.332	3900	3906	3913	rVB3	419072	602698	7.75%	0.778%
31	14.477	3943	3951	3960	rBV3	465985	853745	10.98%	1.102%
32	14.590	3975	3986	3995	rBV4	1905129	3756395	48.33%	4.848%
33	14.644	3996	4003	4015	rVB3	871614	1597841	20.56%	2.062%
34	14.850	4061	4067	4073	rVB4	832258	1143013	14.71%	1.475%

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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

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35	14.956	4092	4100	4112	rVB2	1027885	1993791	25.65%	2.573%
36	15.188	4165	4172	4180	rBV	726311	1072843	13.80%	1.385%
37	15.242	4183	4189	4199	rBV5	459114	1065641	13.71%	1.375%
38	15.416	4234	4243	4256	rBV	766084	1534833	19.75%	1.981%
39	15.612	4298	4304	4316	rVB	1489035	2476050	31.86%	3.196%
40	15.908	4387	4396	4407	rBV3	1060163	2009554	25.85%	2.594%
41	16.101	4448	4456	4465	rVB4	1036254	1848071	23.78%	2.385%
42	16.159	4466	4474	4485	rVB2	1389505	2464909	31.71%	3.181%
43	16.348	4522	4533	4547	rBV2	1382076	3440177	44.26%	4.440%

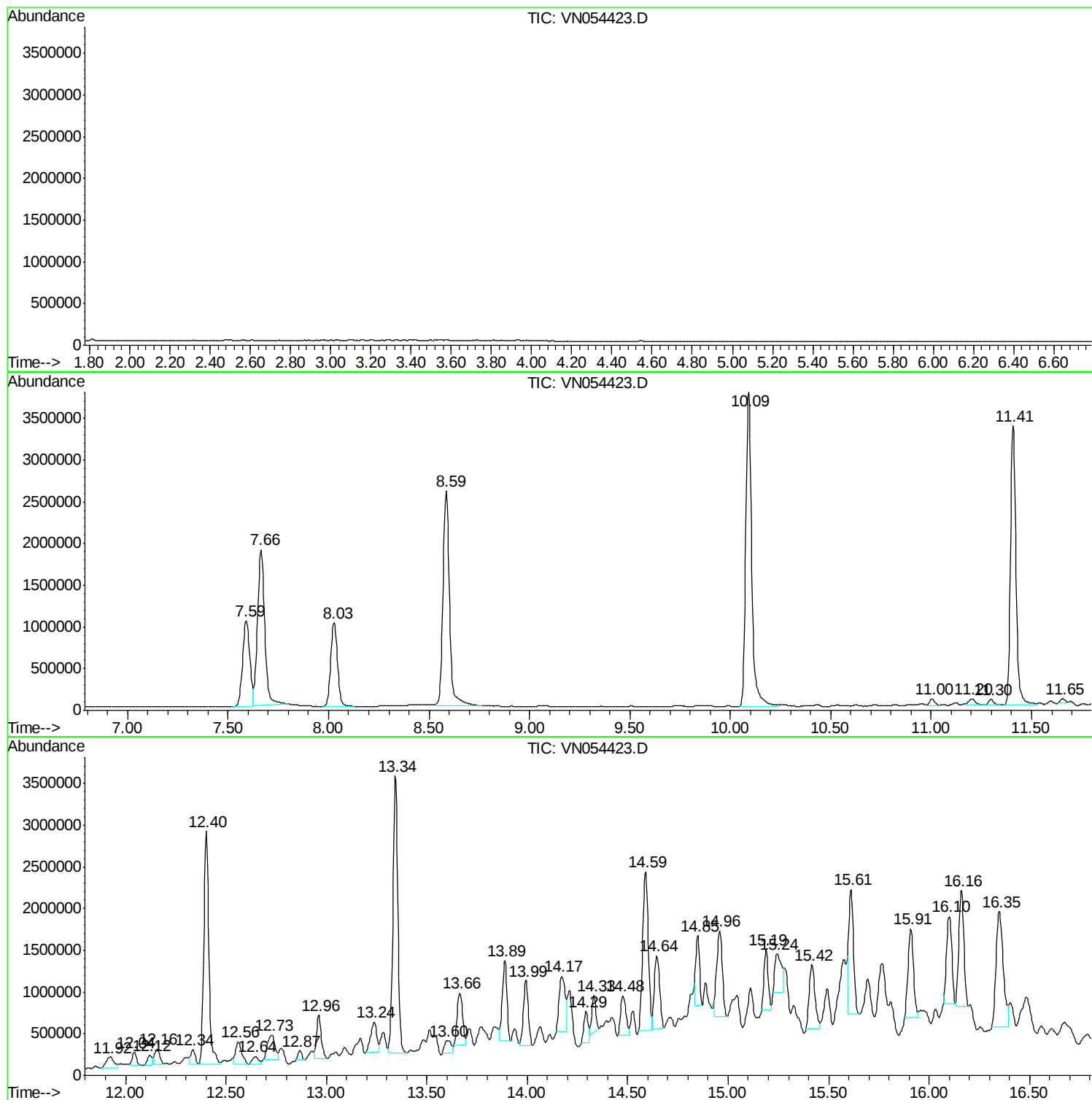
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Quant Title : SW846 8260

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



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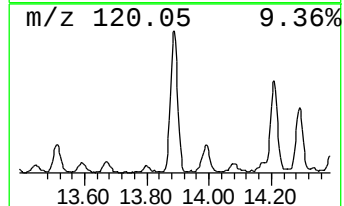
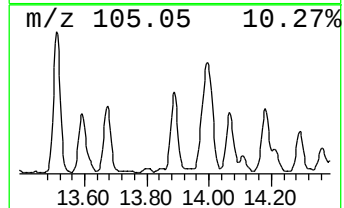
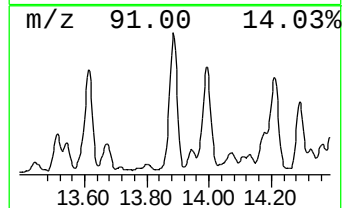
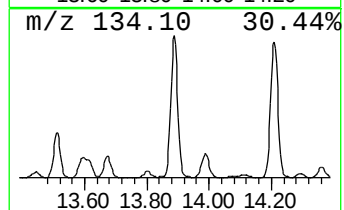
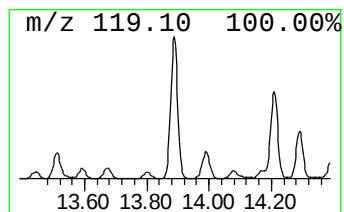
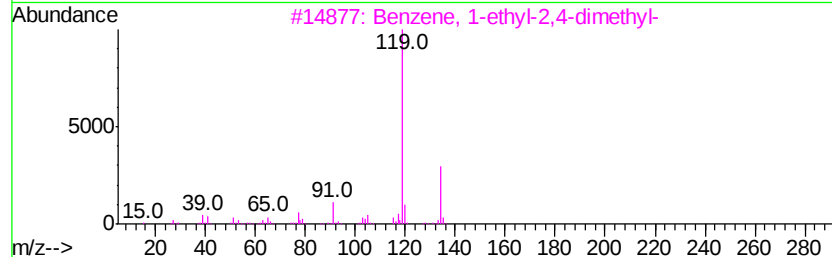
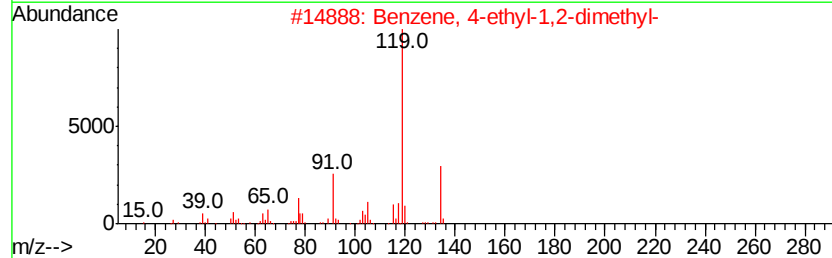
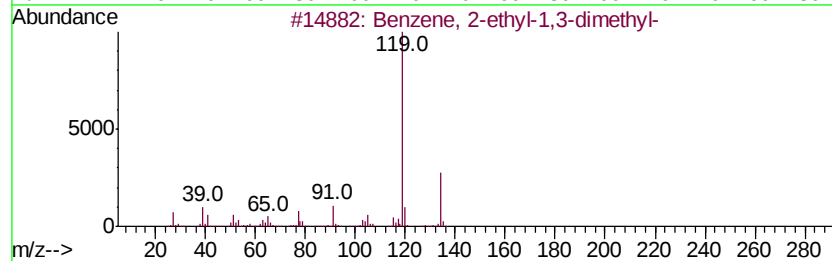
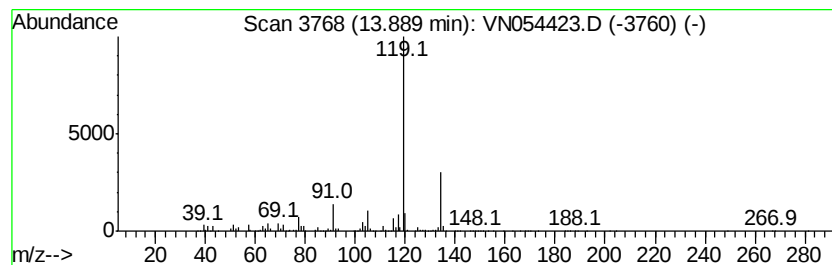
Instrument :
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ClientSampleId :
P-4-E.WALL-S.CENTER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.89	13.32 ug/l	1510940	1,4-Dichlorobenzene-d4	13.34	
Hit# of	5	Tentative ID	MW MolForm	CAS#	Qual
1		Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4	95
2		Benzene, 4-ethyl-1,2-dimethyl-	134 C10H14	000934-80-5	95
3		Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9	94
4		Benzene, 1-ethyl-3,5-dimethyl-	134 C10H14	000934-74-7	94
5		Benzene, 1-methyl-3-(1-methyleth...	134 C10H14	000535-77-3	93



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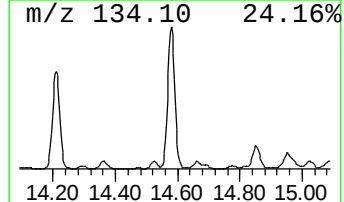
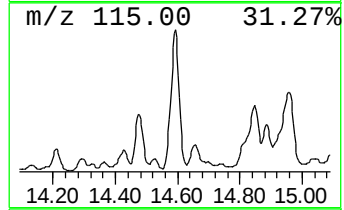
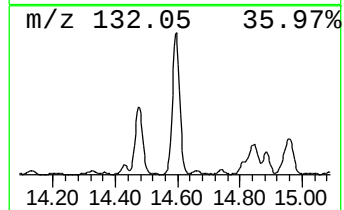
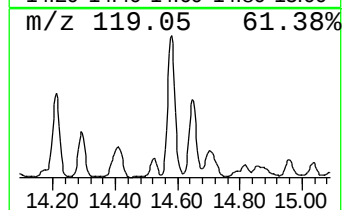
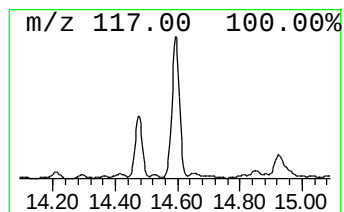
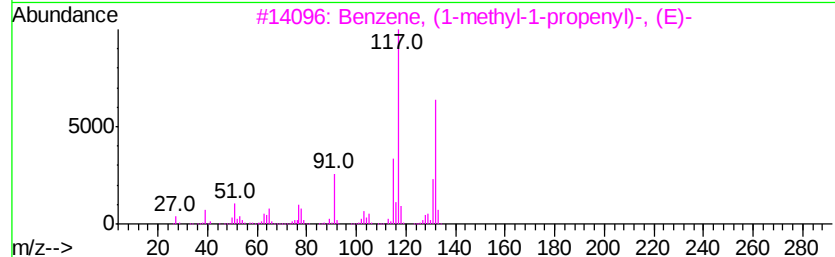
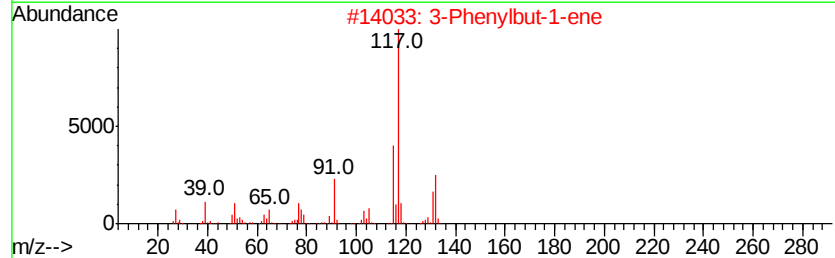
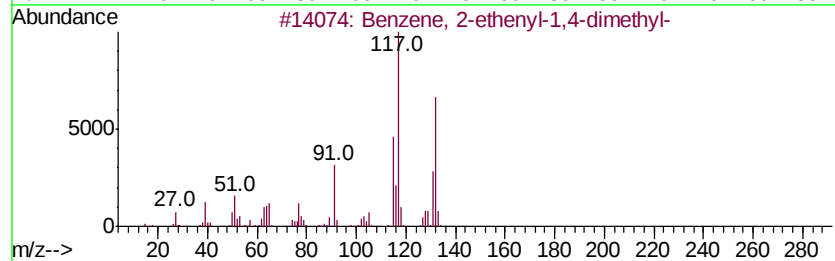
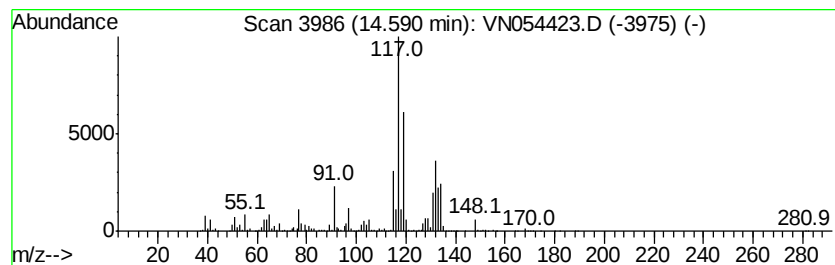
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.59	33.11 ug/l	3756400	1,4-Dichlorobenzene-d4	13.34
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 2-ethenyl-1,4-dimethyl-	132 C10H12	002039-89-6 86
2		3-Phenylbut-1-ene	132 C10H12	000934-10-1 64
3		Benzene, (1-methyl-1-propenyl)-,...	132 C10H12	000768-00-3 64
4		Benzene, 1-methyl-2-(2-propenyl)-	132 C10H12	001587-04-8 60
5		1-Phenyl-1-butene	132 C10H12	000824-90-8 60



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Misc : 6.33g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 47 Sample Multiplier: 1

Instrument :
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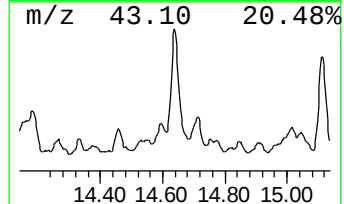
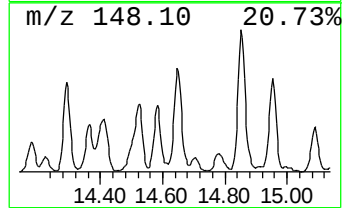
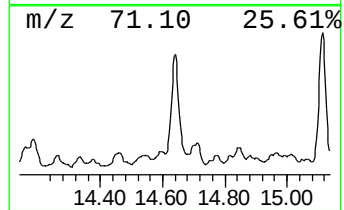
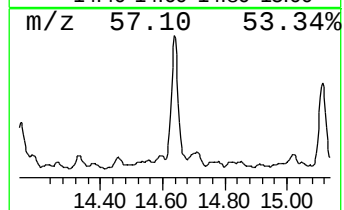
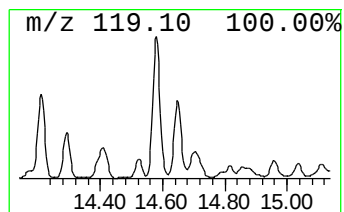
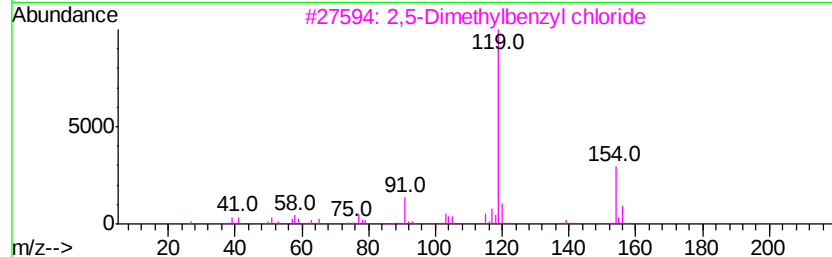
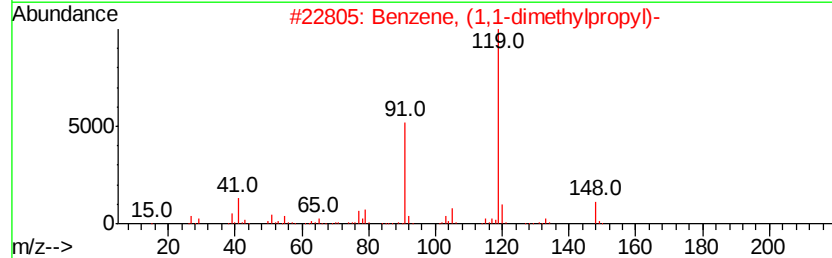
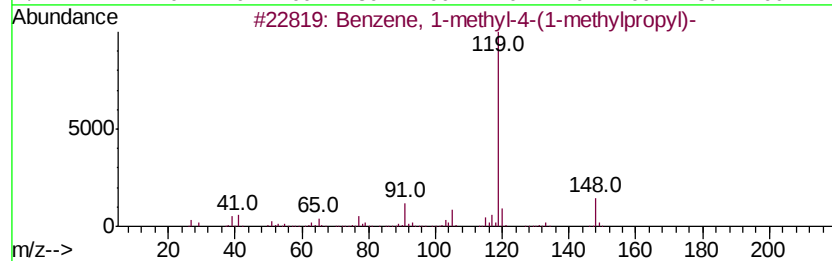
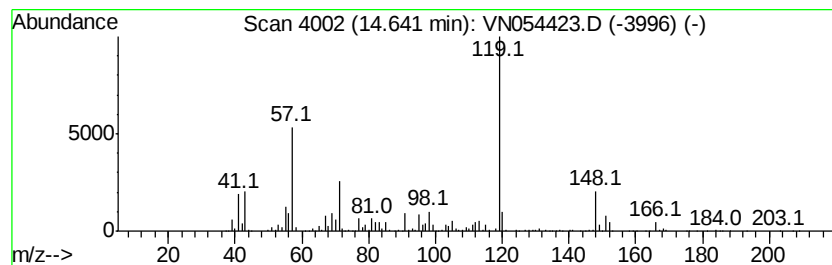
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-methyl-4-(1-meth... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.64	14.08 ug/l	1597840	1,4-Dichlorobenzene-d4	13.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	68
2		Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	59
3		2,5-Dimethylbenzyl chloride	154	C9H11Cl	000824-45-3	47
4		Benzene, 1,1'-(1,1,2,2-tetrameth...	238	C18H22	001889-67-4	47
5		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	47



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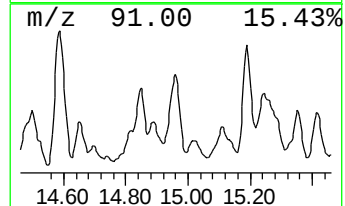
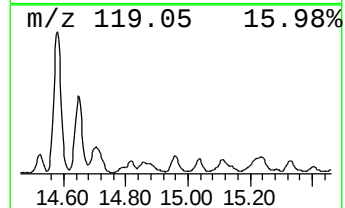
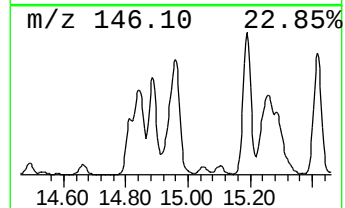
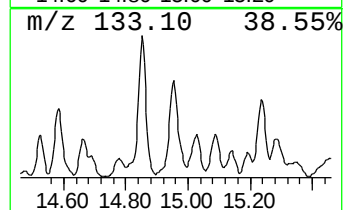
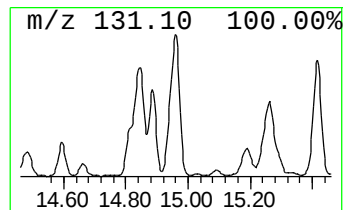
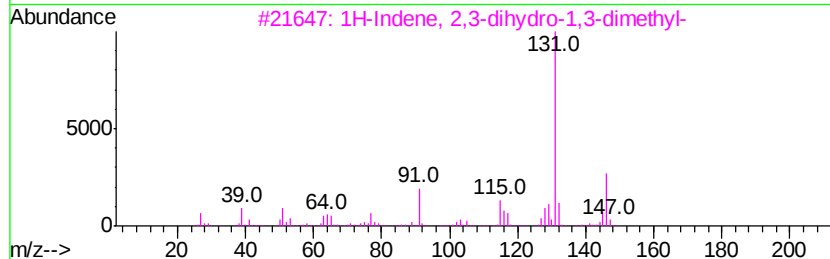
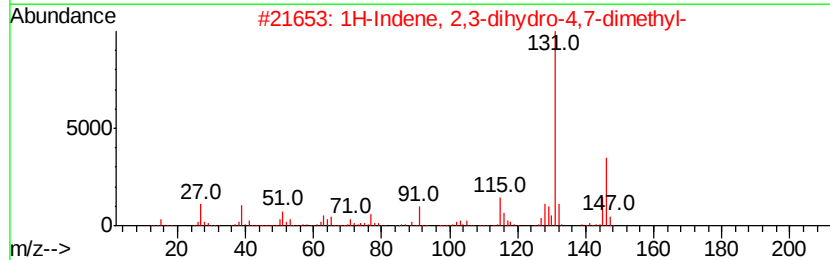
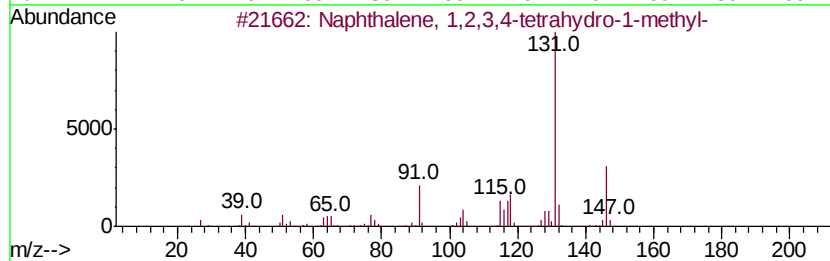
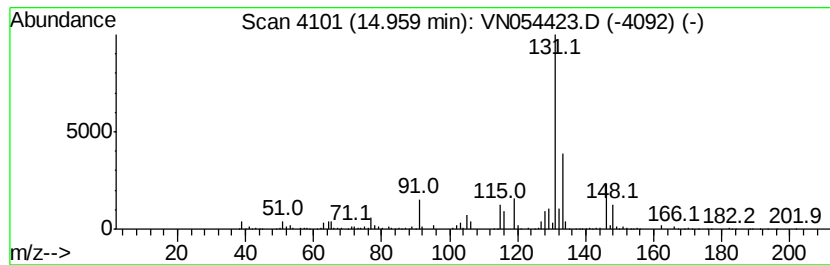
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TIC Integration Parameters: LSCINT.P

Peak Number 4 Naphthalene, 1,2,3,4-tetra... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	17.57 ug/l	1993790	1,4-Dichlorobenzene-d4	13.34
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001559-81-5 64
2		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 64
3		1H-Indene, 2,3-dihydro-1,3-dimet...	146 C11H14	004175-53-5 64
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	146 C11H14	027087-54-3 64
5		1H-Indene, 2,3-dihydro-1,1-dimet...	146 C11H14	004912-92-9 64



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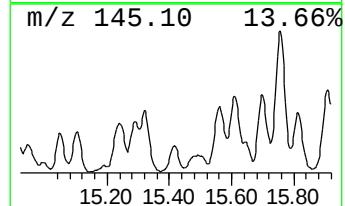
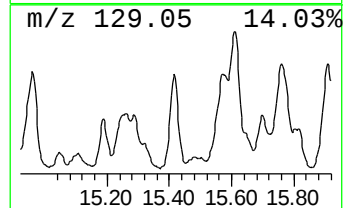
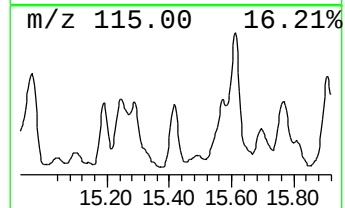
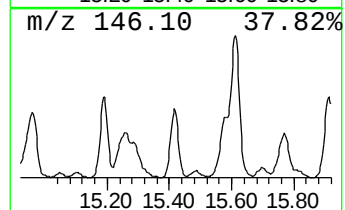
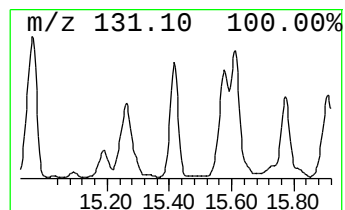
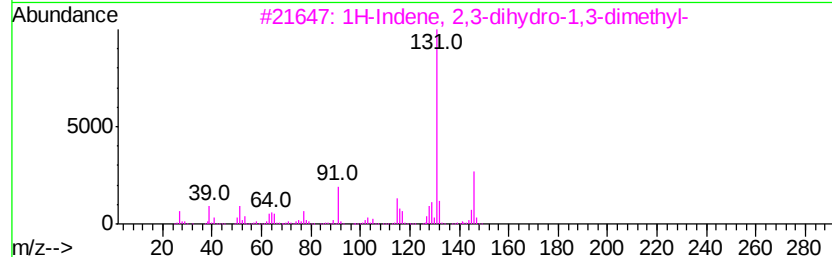
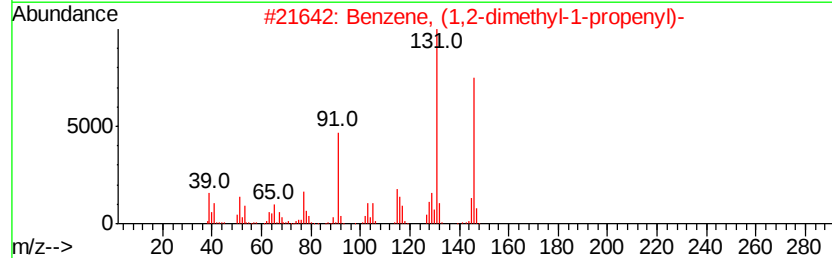
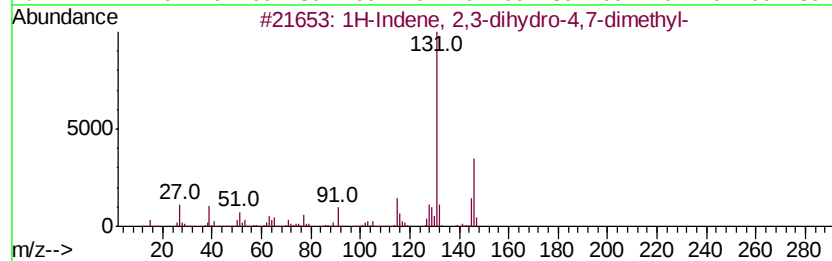
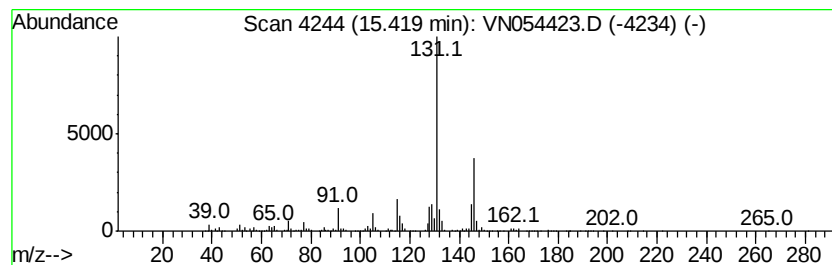
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Quant Title : SW846 8260

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

Peak Number 5 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.42	13.53 ug/l	1534830	1,4-Dichlorobenzene-d4	13.34

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14		006682-71-9	97
2	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14		000769-57-3	94
3	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14		004175-53-5	91
4	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14		017057-82-8	91
5	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14		001075-22-5	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_N\DATA\VN031819\
Data File : VN054423.D
Acq On : 19 Mar 2019 5:29
Operator : JC/SP
Sample : K1970-04 10X
Misc : 6.33g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 47 Sample Multiplier: 1

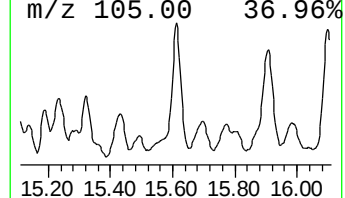
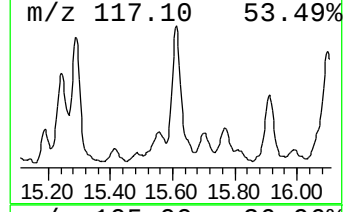
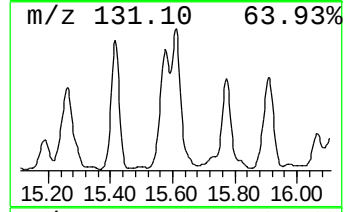
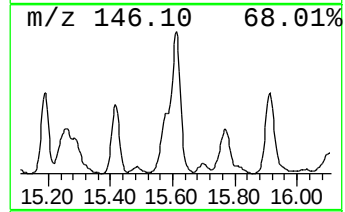
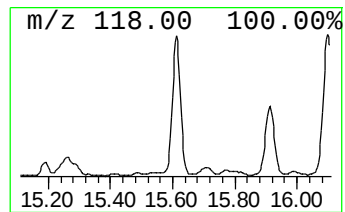
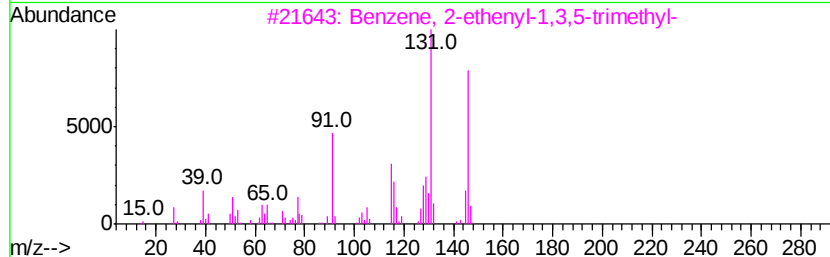
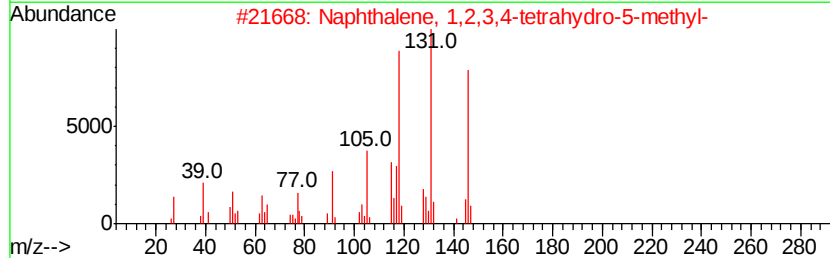
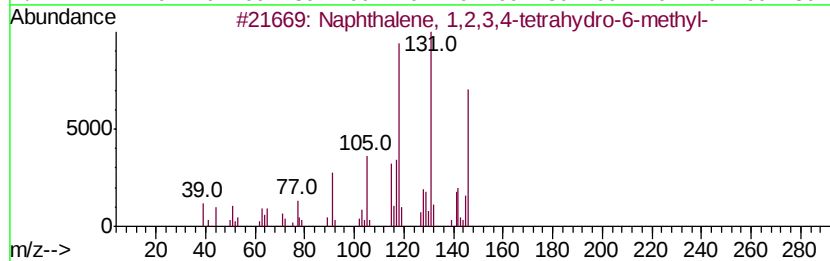
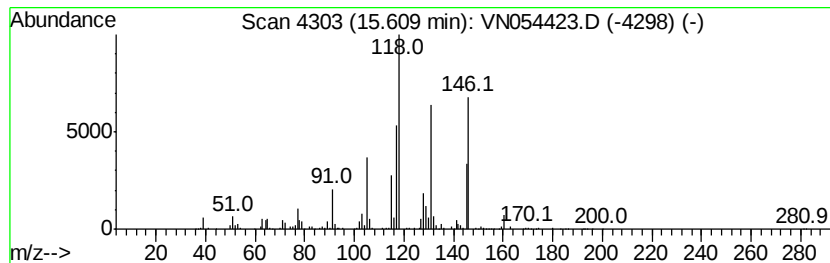
Instrument :
MSVOA_N
ClientSampleId :
P-4-E.WALL-S.CENTER

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

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

Peak Number 6 Naphthalene, 1,2,3,4-tetra... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.61	21.82 ug/l	2476050	1,4-Dichlorobenzene-d4	13.34
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 68
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 64
3		Benzene, 2-ethenyl-1,3,5-trimethyl-	146 C11H14	000769-25-5 41
4		Benzene, (1-ethyl-1-propenyl)-	146 C11H14	004701-36-4 38
5		Cyclobutanone, 3-phenylmethyl-	160 C11H12O	055262-02-7 38



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN031819\
Data File : VN054423.D
Acq On : 19 Mar 2019 5:29
Operator : JC/SP
Sample : K1970-04 10X
Misc : 6.33q/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 47 Sample Multiplier: 1

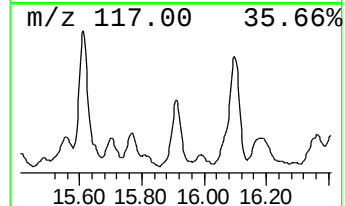
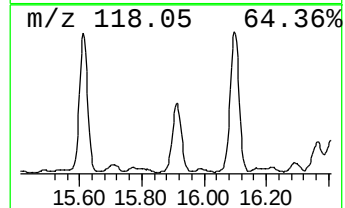
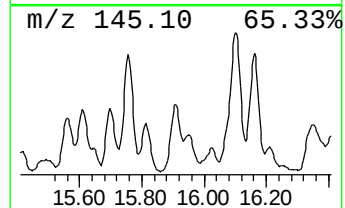
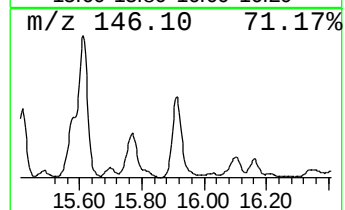
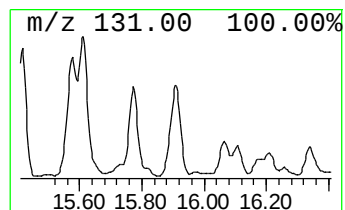
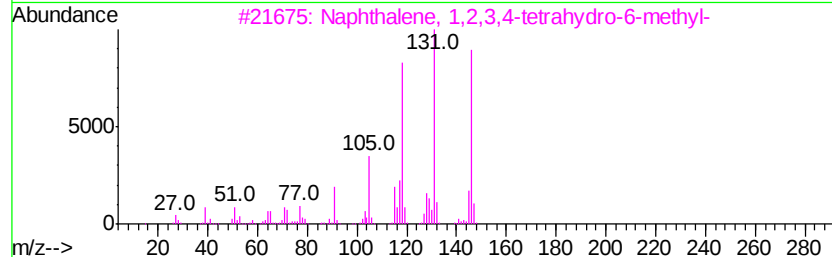
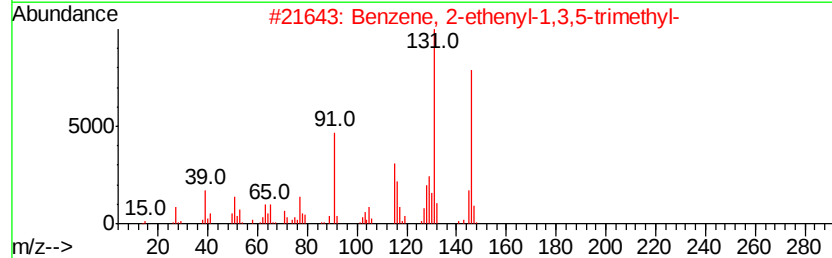
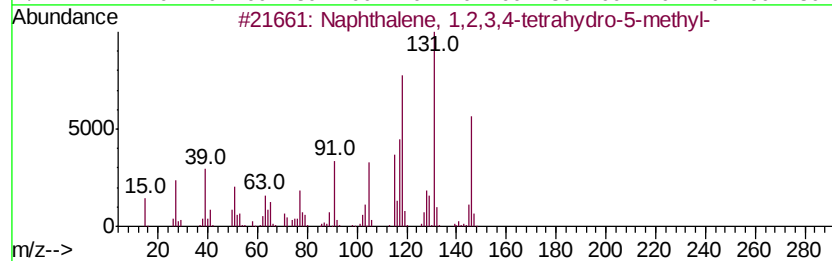
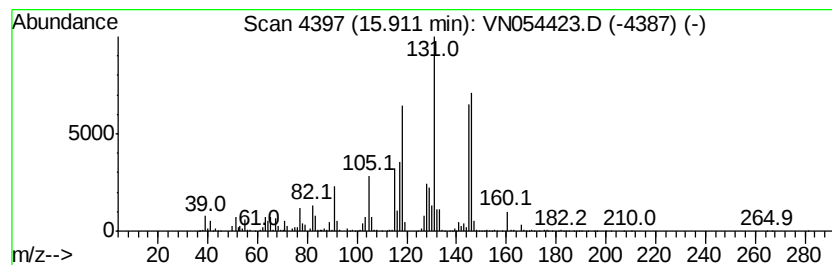
Instrument :
MSVOA_N
ClientSampleId :
P-4-E.WALL-S.CENTER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1,2,3,4-tetra... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.91	17.71 ug/l	2009550	1,4-Dichlorobenzene-d4	13.34
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 90
2		Benzene, 2-ethenyl-1,3,5-trimethyl-	146 C11H14	000769-25-5 78
3		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 58
4		1H-Benzimidazole, 5,6-dimethyl-	146 C9H10N2	000582-60-5 53
5		2,3,4,5,6,7-Hexahydro-1H-cyclope...	146 C11H14	1000189-31-0 49



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_N\DATA\VN031819\
Data File : VN054423.D
Acq On : 19 Mar 2019 5:29
Operator : JC/SP
Sample : K1970-04 10X
Misc : 6.33q/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 47 Sample Multiplier: 1

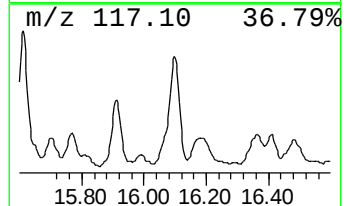
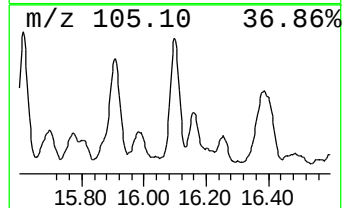
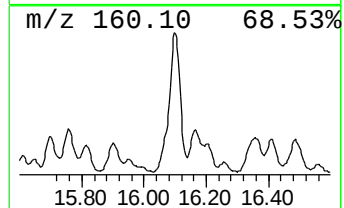
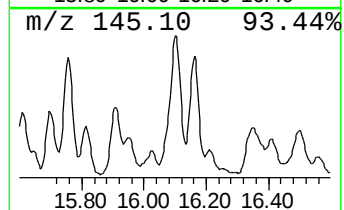
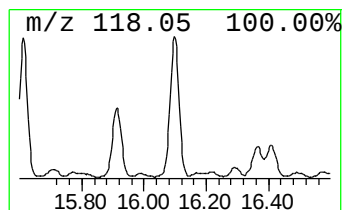
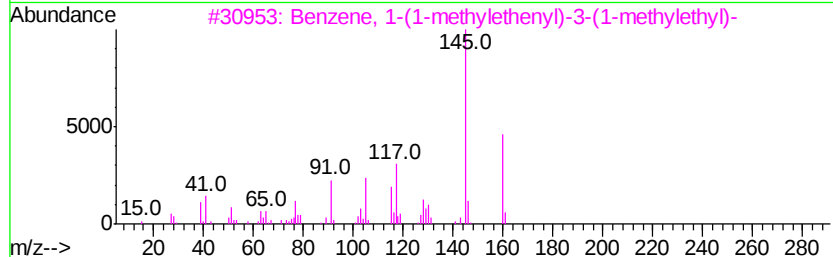
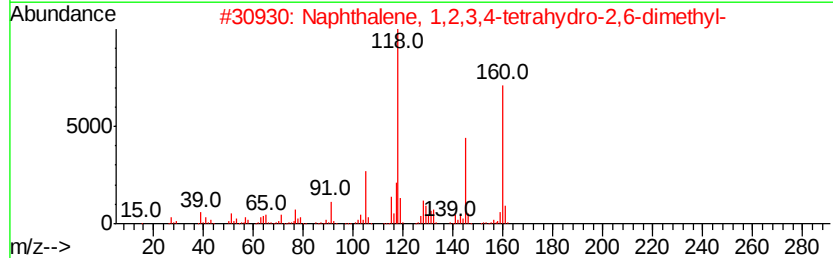
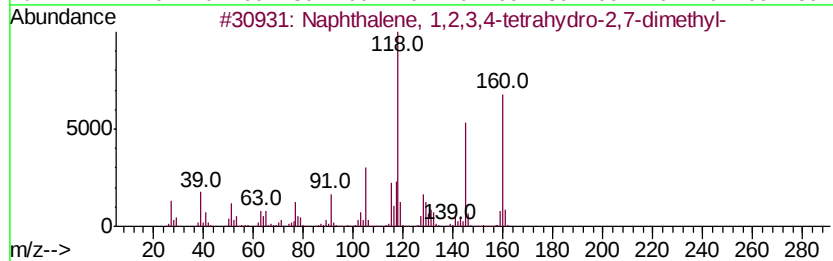
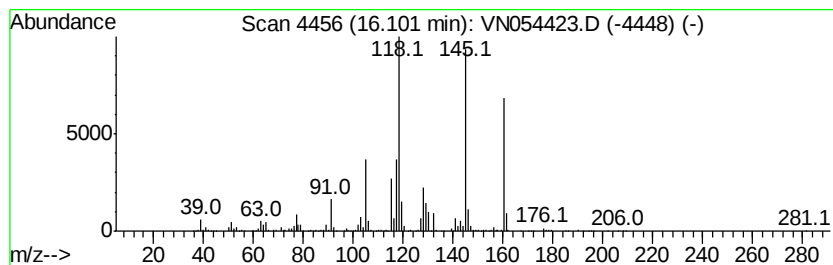
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ClientSampleId :
P-4-E.WALL-S.CENTER

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.10	16.29 ug/l	1848070	1,4-Dichlorobenzene-d4	13.34
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	013065-07-1 90
2		Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	007524-63-2 76
3		Benzene, 1-(1-methylethenyl)-3-(...	160 C12H16	001129-29-9 64
4		2,5-Dimethylbenzyl cyanide	145 C10H11N	016213-85-7 64
5		Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	020027-77-4 60



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN031819\
Data File : VN054423.D
Acq On : 19 Mar 2019 5:29
Operator : JC/SP
Sample : K1970-04 10X
Misc : 6.33q/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
P-4-E.WALL-S.CENTER

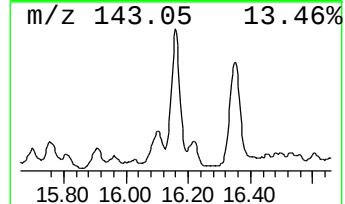
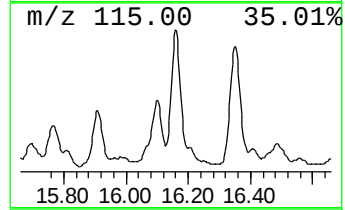
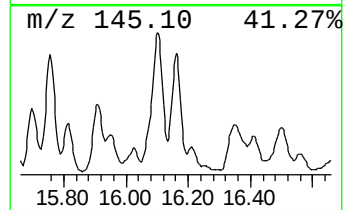
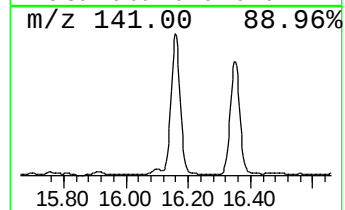
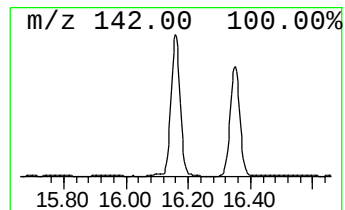
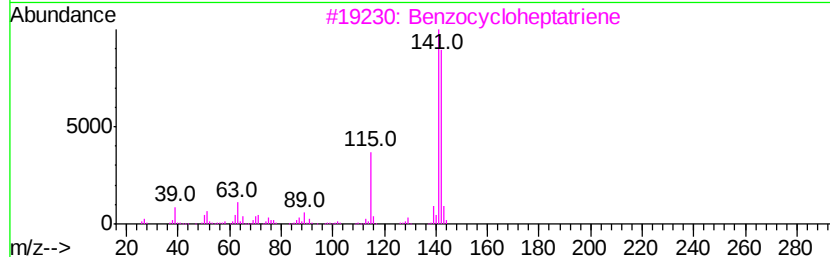
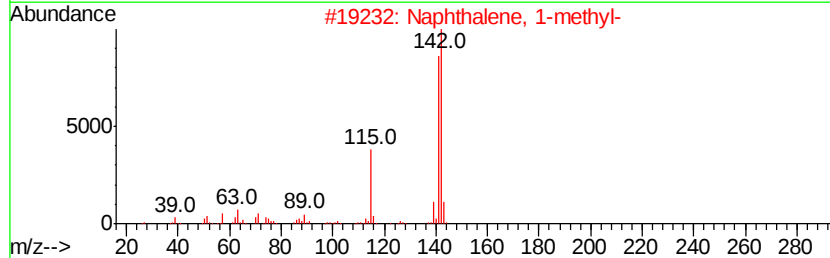
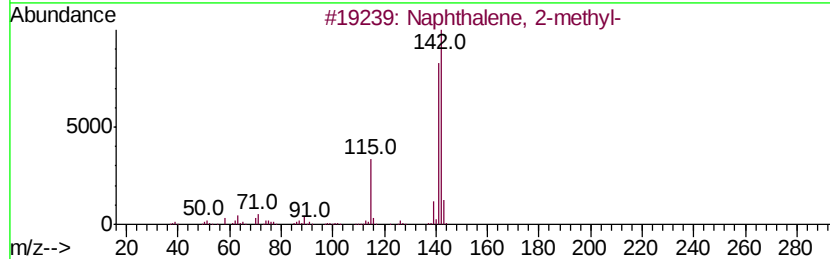
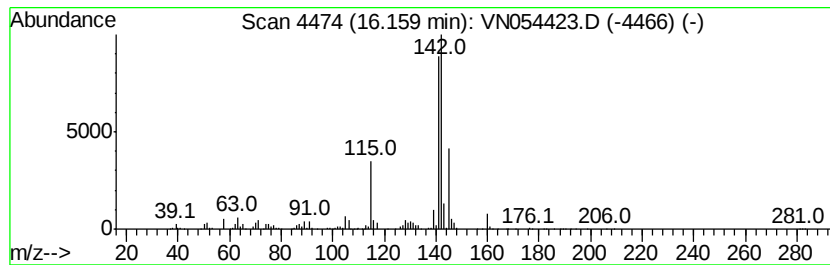
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.16	21.73 ug/l	2464910	1,4-Dichlorobenzene-d4	13.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3		Benzocycloheptatriene	142	C11H10	000264-09-5	76
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	76
5		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	70



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN031819\
Data File : VN054423.D
Acq On : 19 Mar 2019 5:29
Operator : JC/SP
Sample : K1970-04 10X
Misc : 6.33g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
P-4-E.WALL-S.CENTER

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.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, 2-ethyl-...	13.89	13.3	ug/l	1510940	4	13.34	5672790	50.0
Benzene, 2-etheny...	14.59	33.1	ug/l	3756400	4	13.34	5672790	50.0
Benzene, 1-methyl...	14.64	14.1	ug/l	1597840	4	13.34	5672790	50.0
Naphthalene, 1,2,...	14.96	17.6	ug/l	1993790	4	13.34	5672790	50.0
1H-Indene, 2,3-di...	15.42	13.5	ug/l	1534830	4	13.34	5672790	50.0
Naphthalene, 1,2,...	15.61	21.8	ug/l	2476050	4	13.34	5672790	50.0
Naphthalene, 1,2,...	15.91	17.7	ug/l	2009550	4	13.34	5672790	50.0
Naphthalene, 1,2,...	16.10	16.3	ug/l	1848070	4	13.34	5672790	50.0
Naphthalene, 1-me...	16.16	21.7	ug/l	2464910	4	13.34	5672790	50.0