

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN031819\
 Data File : VN054386.D
 Acq On : 18 Mar 2019 11:42
 Operator : JC/SP
 Sample : K1960-11
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 T11-MW-4-9.0-031519

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	3.818	617	636	661	rBV	978213	2527756	14.51%	2.085%
2	5.050	999	1019	1047	rVB2	79794	292957	1.68%	0.242%
3	7.590	1787	1809	1820	rBV2	1059464	2674547	15.36%	2.207%
4	7.664	1820	1832	1883	rVB	2000423	5175933	29.72%	4.270%
5	8.027	1928	1945	1974	rBV	1051474	2439802	14.01%	2.013%
6	8.587	2103	2119	2173	rVB	2621824	5918777	33.98%	4.883%
7	9.079	2249	2272	2288	rBV3	198139	470114	2.70%	0.388%
8	10.091	2568	2587	2629	rBV	3789716	7803432	44.80%	6.438%
9	11.407	2981	2996	3016	rBV	3357803	6175155	35.45%	5.095%
10	11.497	3018	3024	3044	rVB3	110872	239099	1.37%	0.197%
11	11.622	3048	3063	3085	rBV3	127521	383286	2.20%	0.316%
12	11.950	3146	3165	3196	rVB5	81299	248073	1.42%	0.205%
13	12.249	3246	3258	3272	rBV	898411	1494269	8.58%	1.233%
14	12.403	3292	3306	3322	rBV	2642495	4674993	26.84%	3.857%
15	12.593	3351	3365	3378	rBV	800044	1298641	7.46%	1.071%
16	12.992	3482	3489	3497	rVV	197517	327019	1.88%	0.270%
17	13.040	3497	3504	3516	rVB	146224	237829	1.37%	0.196%
18	13.172	3529	3545	3555	rBV2	674018	1387463	7.97%	1.145%
19	13.342	3586	3598	3617	rVB	3000176	5064492	29.08%	4.178%
20	13.442	3619	3629	3639	rBV	454882	715452	4.11%	0.590%
21	13.513	3639	3651	3653	rVV	1448985	1912332	10.98%	1.578%
22	13.545	3654	3661	3671	rVV	4147505	6939252	39.84%	5.725%
23	13.590	3671	3675	3690	rVV2	337931	645672	3.71%	0.533%
24	13.673	3691	3701	3714	rVB	887715	1411667	8.11%	1.165%
25	13.889	3756	3768	3777	rBV	3681835	5545369	31.84%	4.575%
26	13.947	3777	3786	3792	rVV	1209780	1985068	11.40%	1.638%
27	13.995	3792	3801	3813	rVV2	4731078	7951115	45.65%	6.560%
28	14.062	3817	3822	3831	rVB3	123424	189922	1.09%	0.157%
29	14.130	3831	3843	3852	rBV	526934	879055	5.05%	0.725%
30	14.207	3852	3867	3885	rVB	2883682	4798522	27.55%	3.959%
31	14.294	3886	3894	3899	rBV	164305	253005	1.45%	0.209%
32	14.361	3909	3915	3923	rVB	247141	334829	1.92%	0.276%
33	14.426	3923	3935	3942	rBV3	355542	699561	4.02%	0.577%
34	14.477	3942	3951	3960	rVV2	1163853	1951743	11.21%	1.610%

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Integration Parameters: RTEINT.P

Integrator: RTE
Smoothing : ON
Sampling : 1
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Max Peaks: 100
Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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35	14.525	3961	3966	3972	rVV2	188421	263477	1.51%	0.217%
36	14.583	3972	3984	3999	rVV4	8661947	17417224	100.00%	14.370%
37	14.654	3999	4006	4016	rVB4	290964	506386	2.91%	0.418%
38	14.850	4042	4067	4073	rBV3	807435	2046125	11.75%	1.688%
39	14.885	4073	4078	4086	rVV	697447	1131505	6.50%	0.934%
40	14.956	4088	4100	4111	rVB2	1308134	2901493	16.66%	2.394%
41	15.088	4134	4141	4151	rVB	333483	495177	2.84%	0.409%
42	15.188	4162	4172	4180	rBV	539735	927950	5.33%	0.766%
43	15.258	4180	4194	4229	rVB5	703010	2631758	15.11%	2.171%
44	15.416	4232	4243	4254	rBV	277749	518635	2.98%	0.428%
45	15.484	4254	4264	4275	rBV3	239820	410828	2.36%	0.339%
46	15.577	4281	4293	4297	rBV3	513706	918690	5.27%	0.758%
47	15.612	4298	4304	4320	rVB	1072457	1904942	10.94%	1.572%
48	15.696	4321	4330	4340	rBV	224987	401403	2.30%	0.331%
49	15.770	4345	4353	4363	rVB3	206690	367721	2.11%	0.303%
50	15.914	4384	4398	4414	rBV2	1077457	2117321	12.16%	1.747%
51	16.101	4448	4456	4466	rVB5	112879	214591	1.23%	0.177%
52	16.165	4468	4476	4506	rVB8	120725	447134	2.57%	0.369%
53	16.348	4521	4533	4546	rBV5	219428	538036	3.09%	0.444%

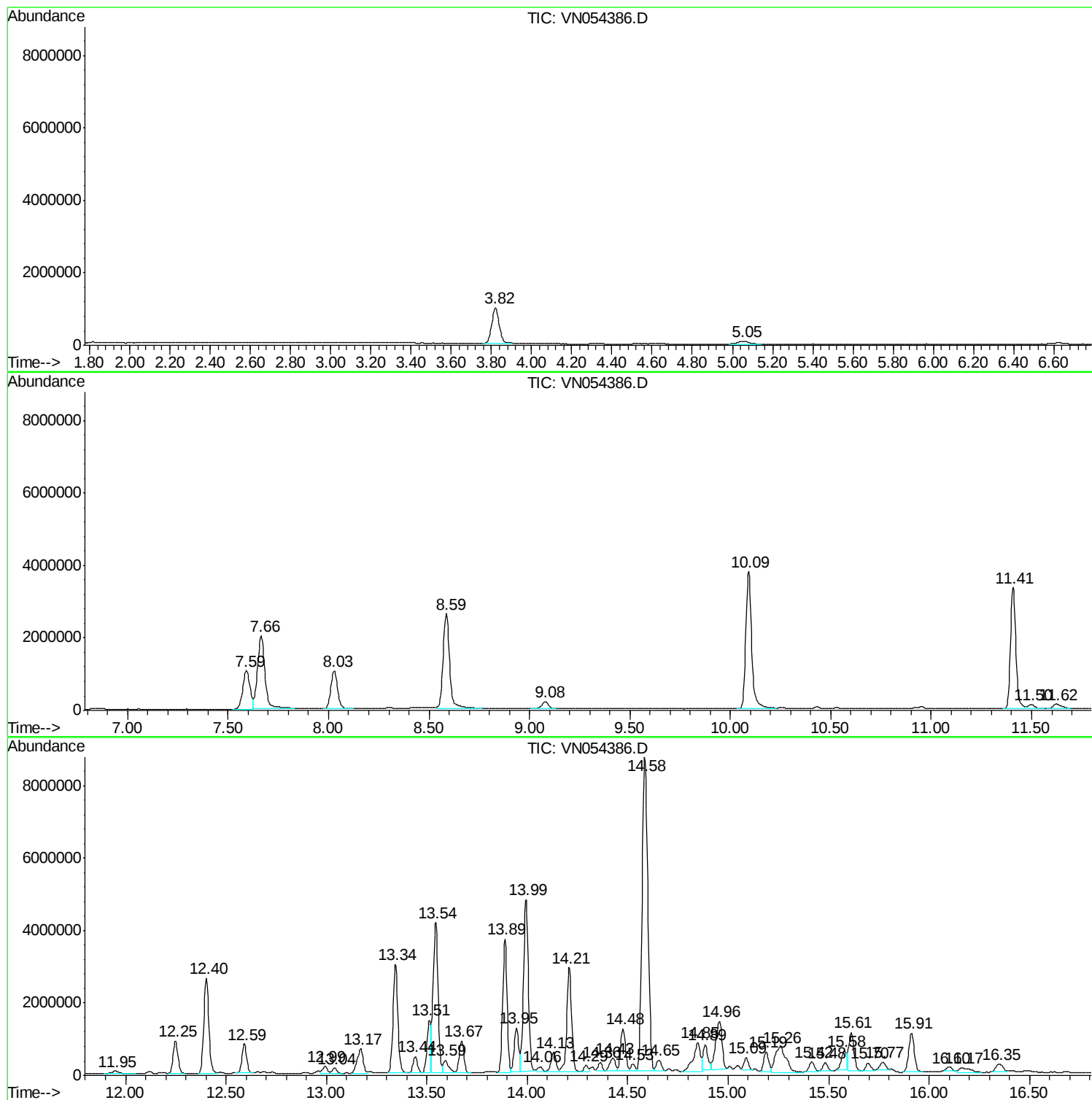
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

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



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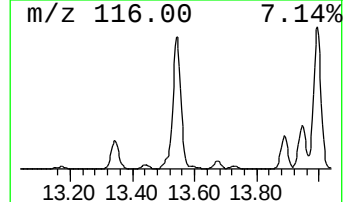
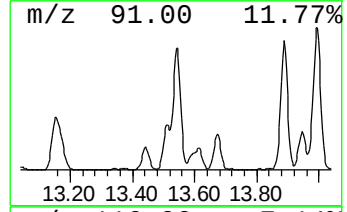
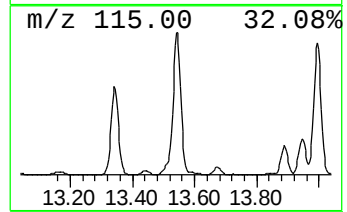
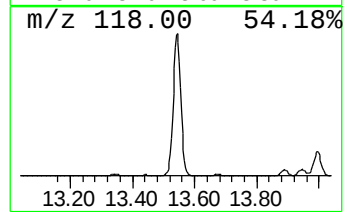
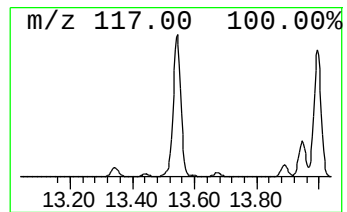
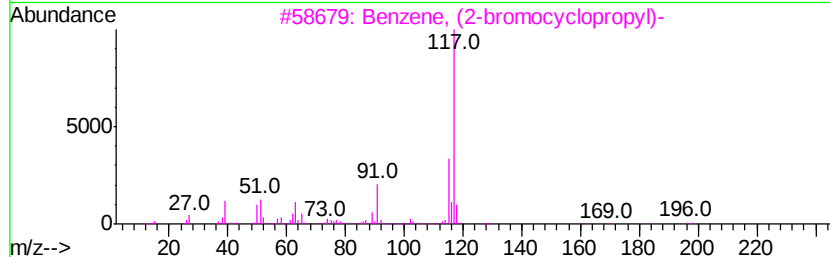
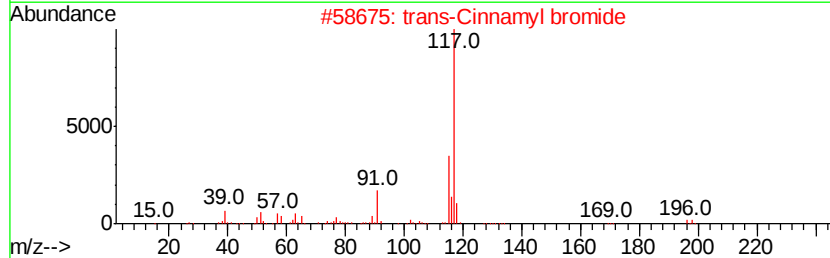
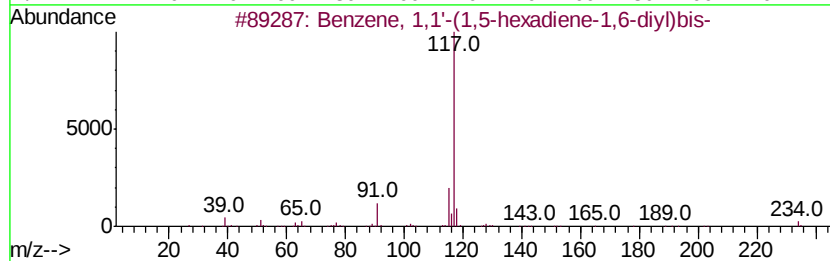
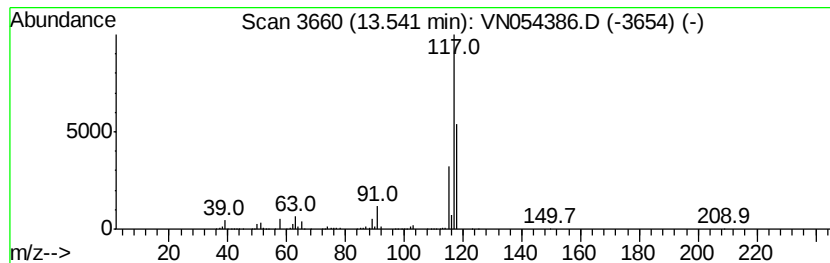
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1,1'-(1,5-hexadien... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.54	68.51 ug/l	6939250	1,4-Dichlorobenzene-d4	13.34

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	53
2		trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	42
3		Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	38
4		m-Aminophenylacetylene	117	C8H7N	054060-30-9	37
5		Benzocyclobutene, 1-bromo-1-methyl-	196	C9H9Br	1000161-76-4	10



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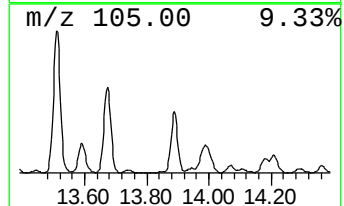
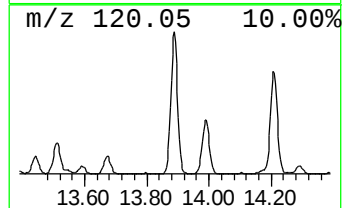
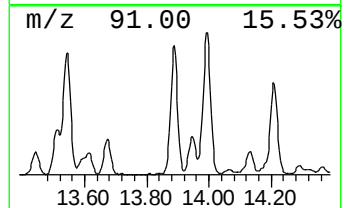
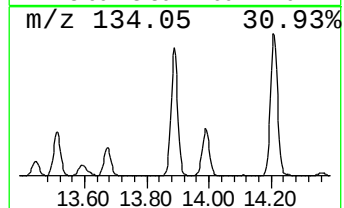
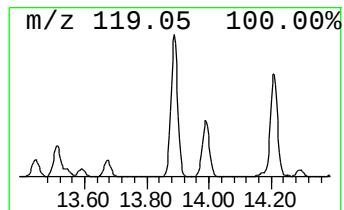
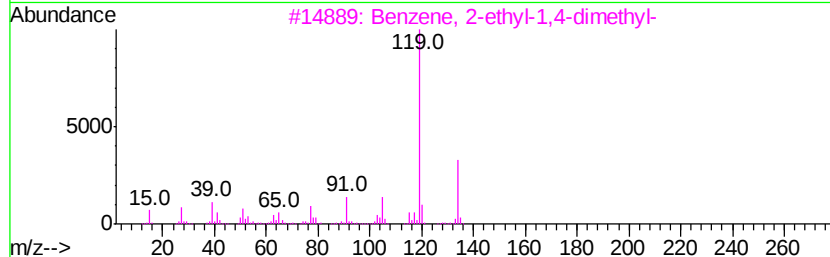
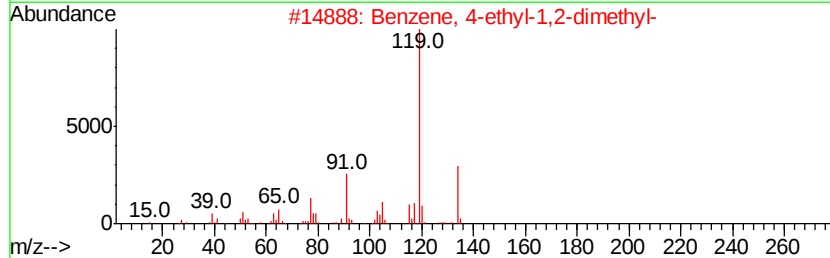
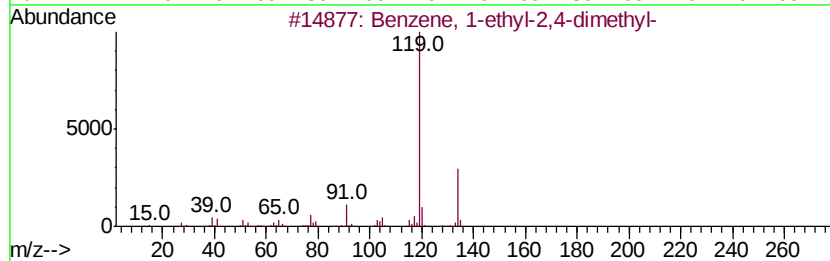
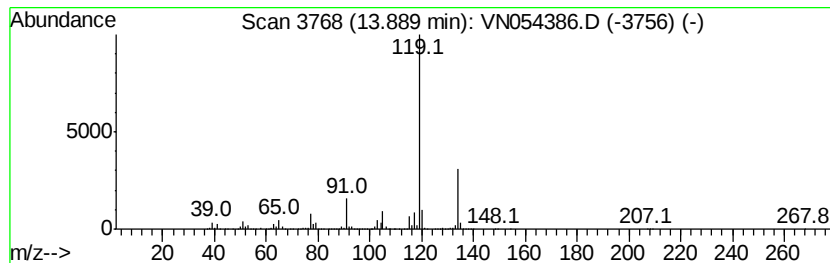
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	54.75 ug/l	5545370	1,4-Dichlorobenzene-d4	13.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5		o-Cymene	134	C10H14	000527-84-4	95



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ALS Vial : 8 Sample Multiplier: 1

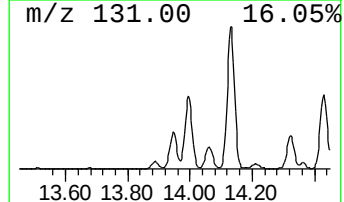
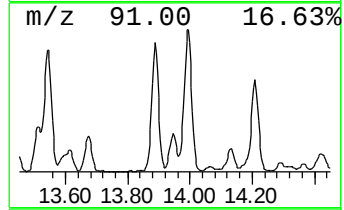
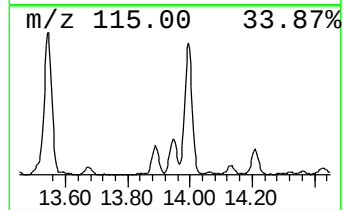
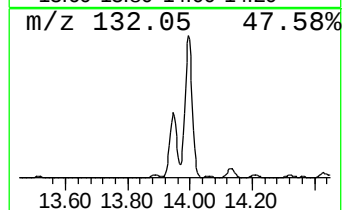
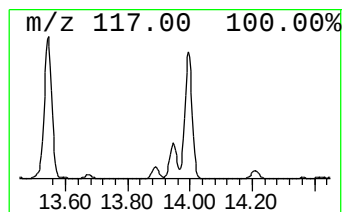
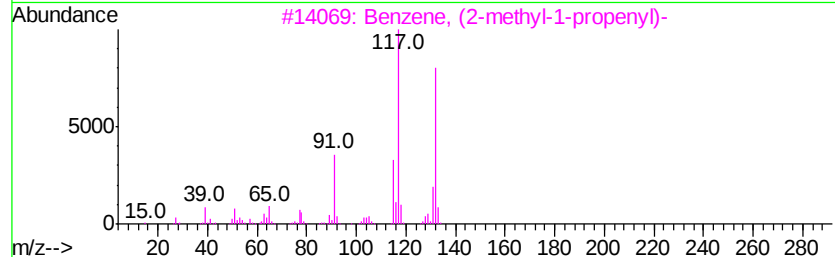
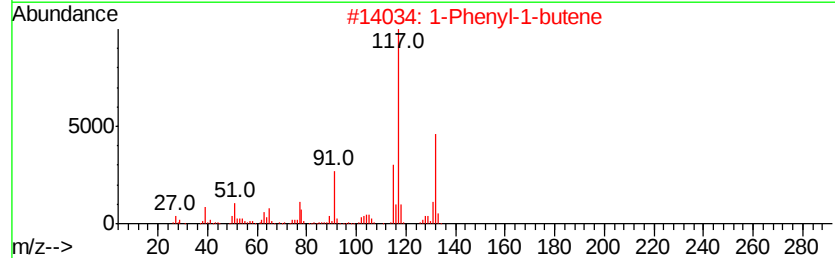
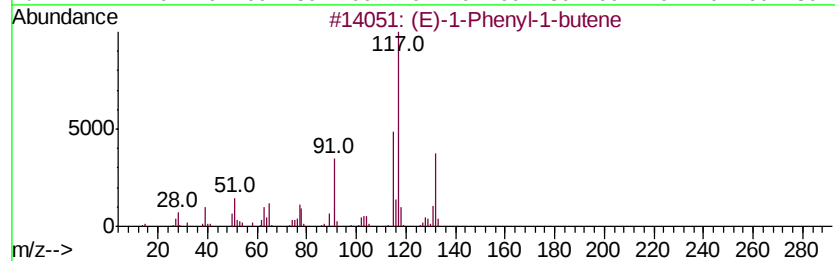
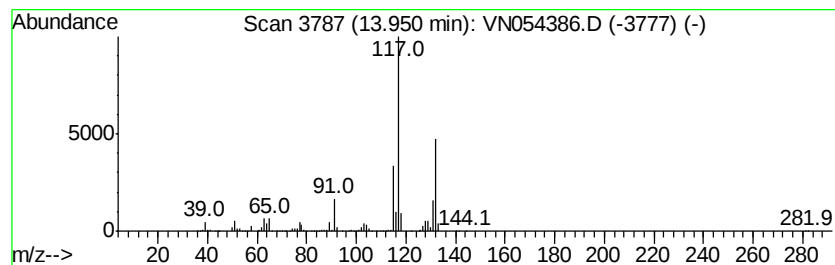
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (E)-1-Phenyl-1-butene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.95	19.60 ug/l	1985070	1,4-Dichlorobenzene-d4	13.34	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	94
2	1-Phenyl-1-butene	132	C10H12	000824-90-8	94
3	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	91
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	91
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	91



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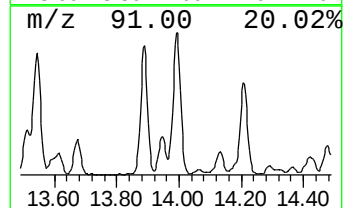
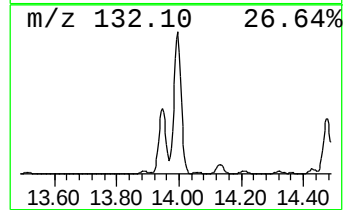
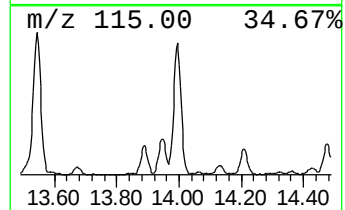
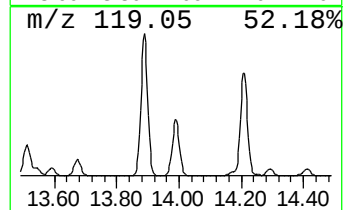
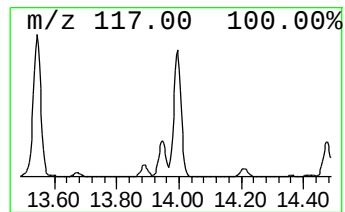
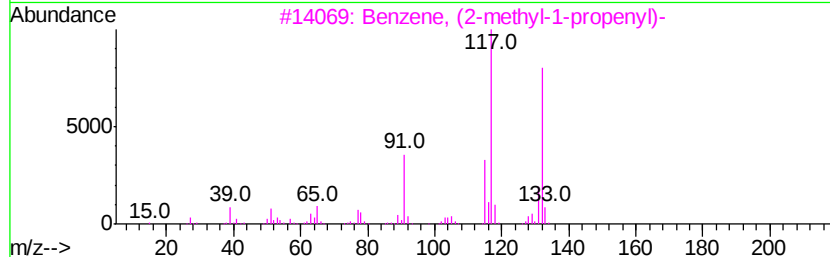
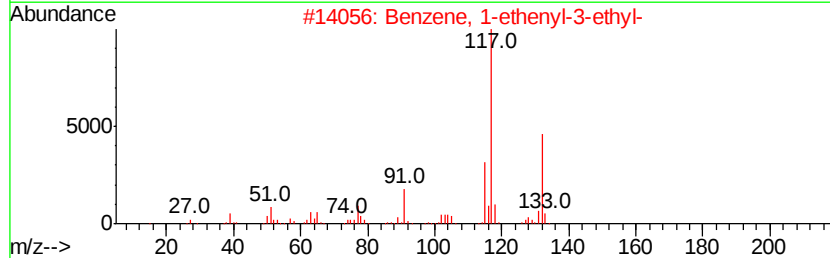
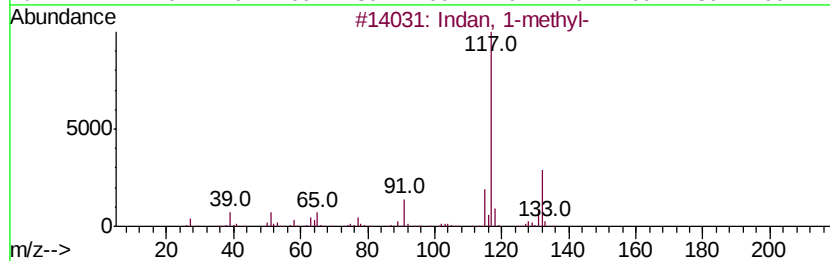
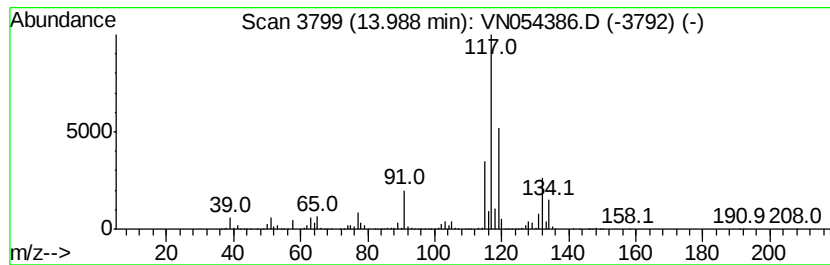
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Indan, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	78.50 ug/l	7951120	1,4-Dichlorobenzene-d4	13.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indan, 1-methyl-	132	C10H12	000767-58-8	70
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	68
3		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	68
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	64
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	64



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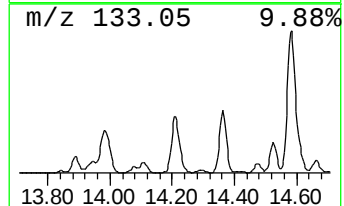
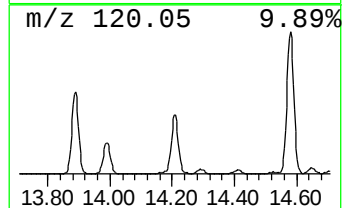
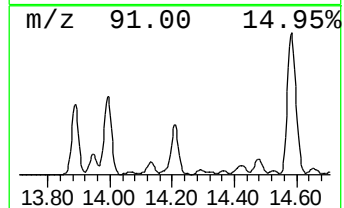
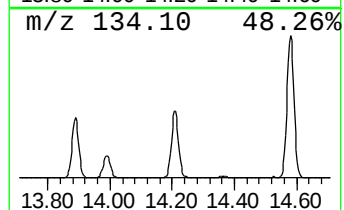
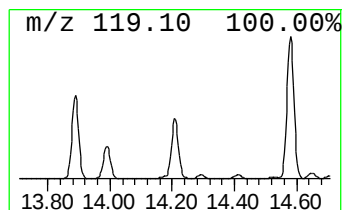
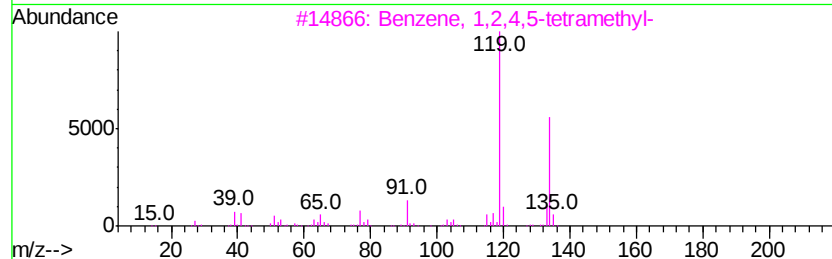
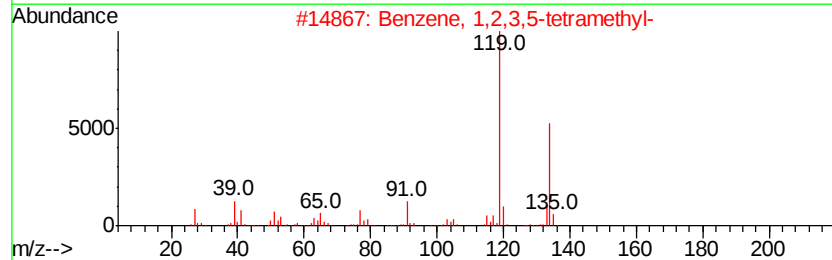
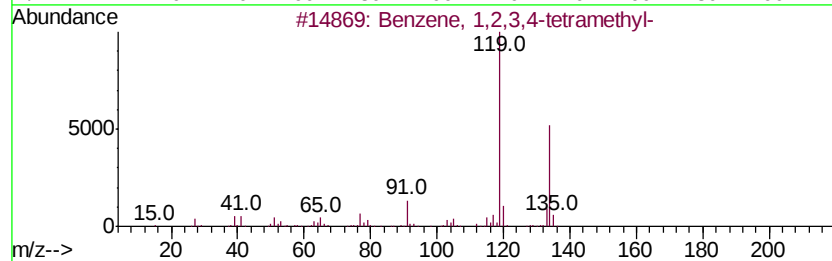
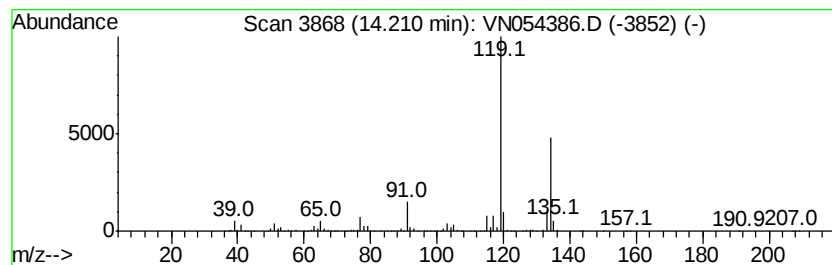
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	47.37 ug/l	4798520	1,4-Dichlorobenzene-d4	13.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN031819\
Data File : VN054386.D
Acq On : 18 Mar 2019 11:42
Operator : JC/SP
Sample : K1960-11
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
T11-MW-4-9.0-031519

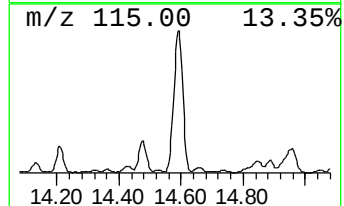
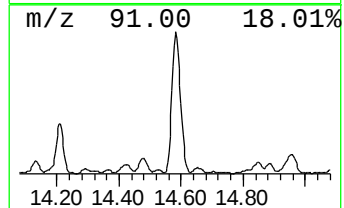
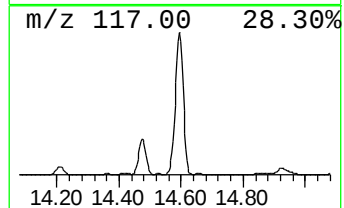
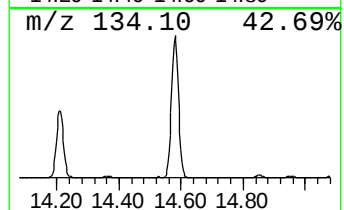
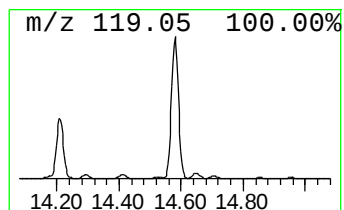
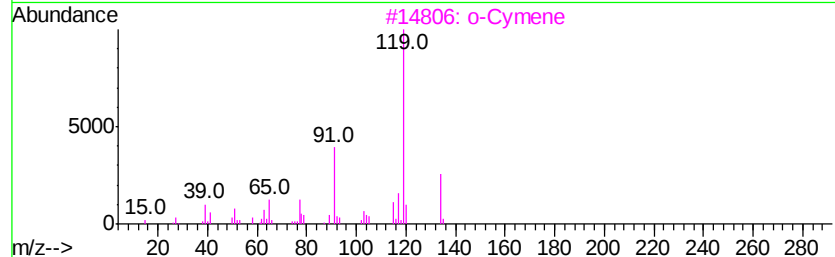
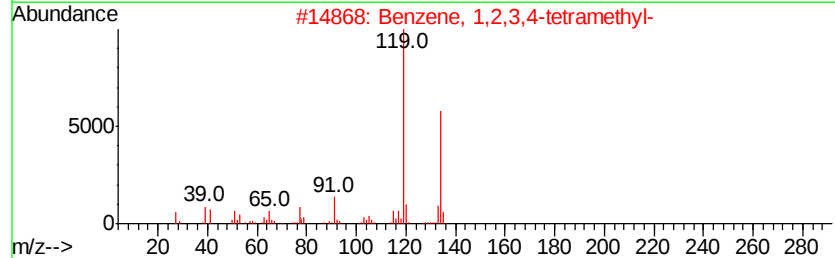
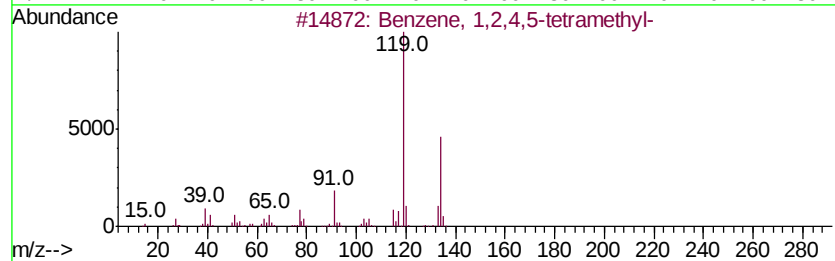
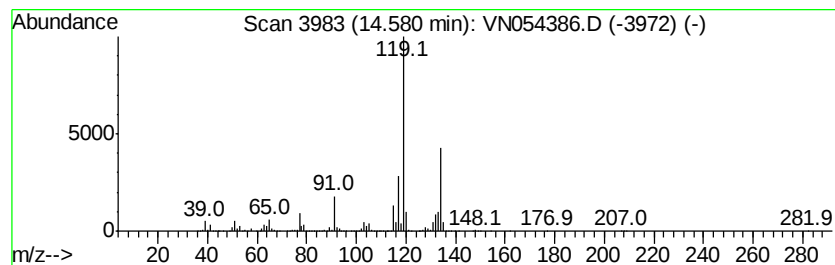
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.58	171.95 ug/l	17417200	1,4-Dichlorobenzene-d4	13.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	93
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	76
3		o-Cymene	134	C10H14	000527-84-4	70
4		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	70
5		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	70



Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN031819\
Data File : VN054386.D
Acq On : 18 Mar 2019 11:42
Operator : JC/SP
Sample : K1960-11
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 8 Sample Multiplier: 1

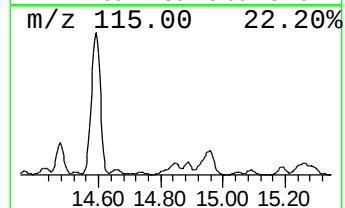
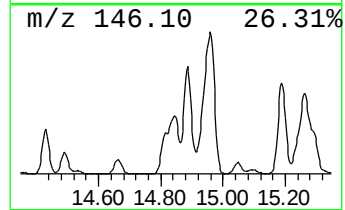
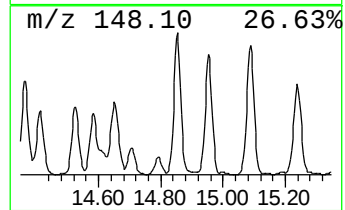
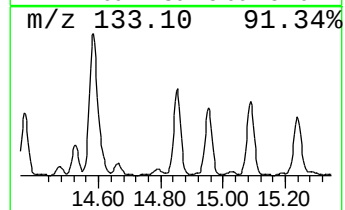
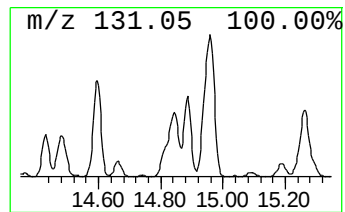
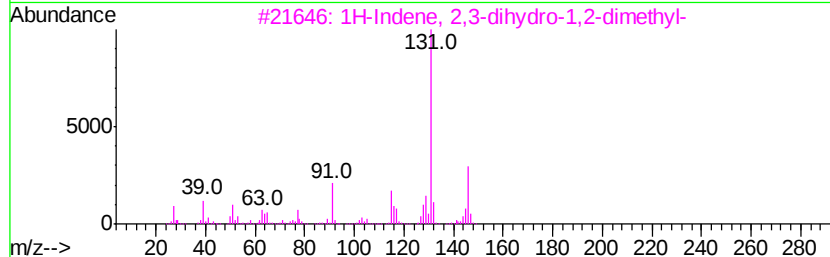
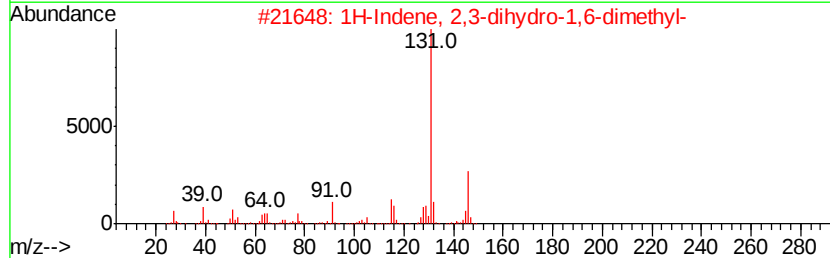
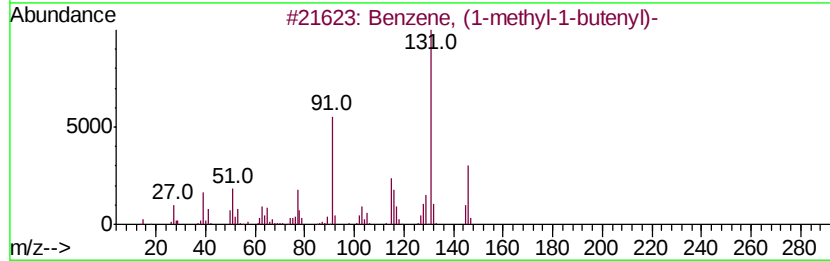
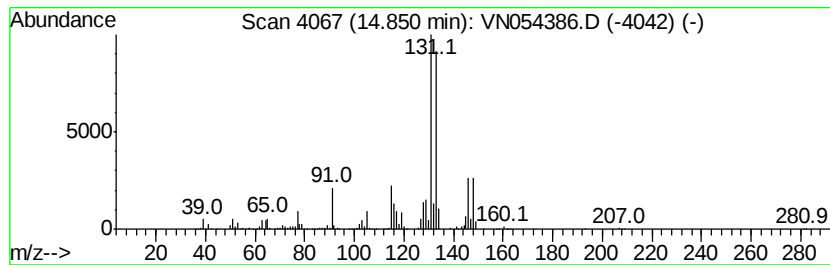
Instrument :
MSVOA_N
ClientSampleId :
T11-MW-4-9.0-031519

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 Benzene, (1-methyl-1-butenyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	20.20 ug/l	2046130	1,4-Dichlorobenzene-d4	13.34
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (1-methyl-1-butenyl)-	146 C11H14	053172-84-2 90
2		1H-Indene, 2,3-dihydro-1,6-dimet...	146 C11H14	017059-48-2 87
3		1H-Indene, 2,3-dihydro-1,2-dimet...	146 C11H14	017057-82-8 70
4		Benzene, 1-methyl-4-(1-methyl-2-...	146 C11H14	097664-18-1 70
5		Benzene, (2-methyl-1-butenyl)-	146 C11H14	056253-64-6 64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN031819\
Data File : VN054386.D
Acq On : 18 Mar 2019 11:42
Operator : JC/SP
Sample : K1960-11
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 8 Sample Multiplier: 1

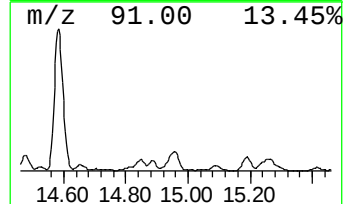
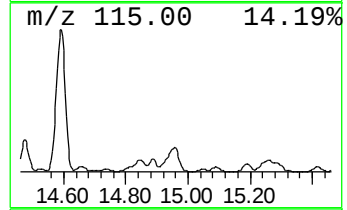
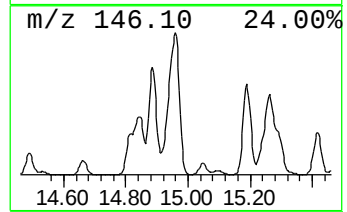
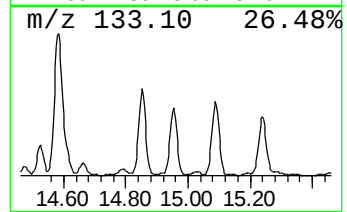
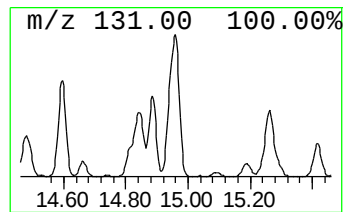
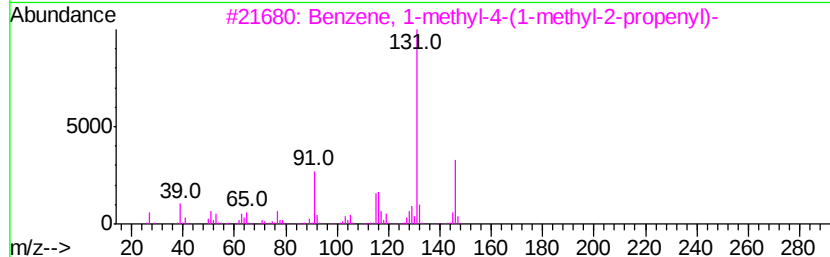
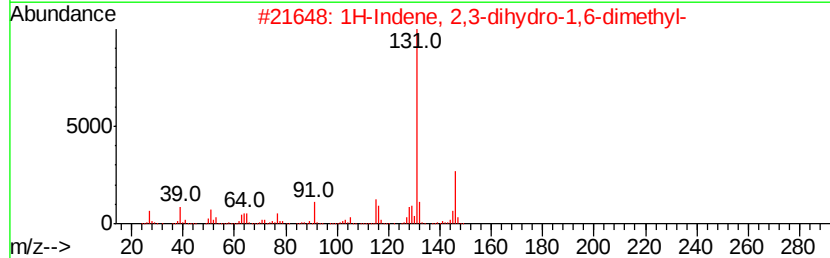
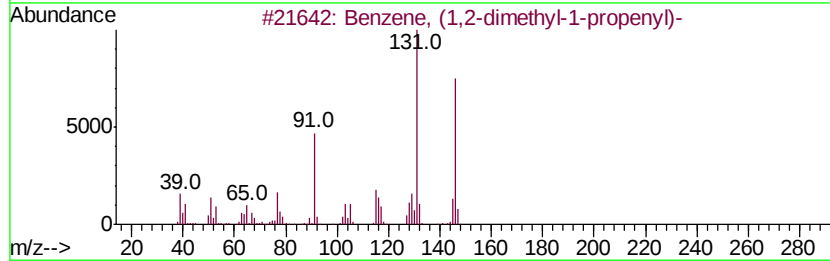
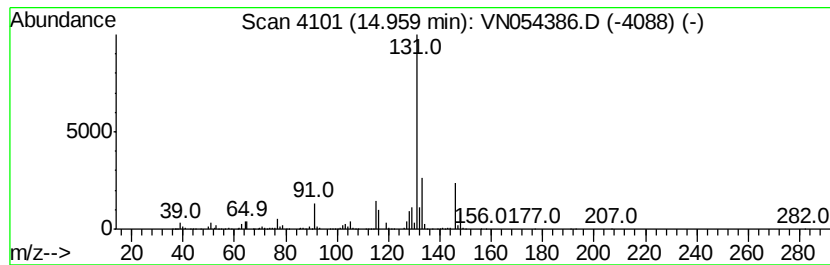
Instrument :
MSVOA_N
ClientSampleId :
T11-MW-4-9.0-031519

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 8 Benzene, (1,2-dimethyl-1-propenyl)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	28.65 ug/l	2901490	1,4-Dichlorobenzene-d4	13.34
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (1,2-dimethyl-1-propenyl)-	146 C11H14	000769-57-3 81
2		1H-Indene, 2,3-dihydro-1,6-dimet...	146 C11H14	017059-48-2 81
3		Benzene, 1-methyl-4-(1-methyl-2-...	146 C11H14	097664-18-1 81
4		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 81
5		1H-Indene, 2,3-dihydro-1,3-dimet...	146 C11H14	004175-53-5 81



Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN031819\
Data File : VN054386.D
Acq On : 18 Mar 2019 11:42
Operator : JC/SP
Sample : K1960-11
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 8 Sample Multiplier: 1

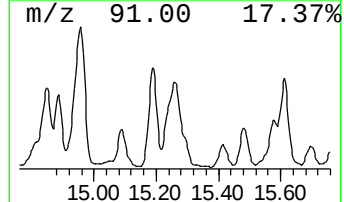
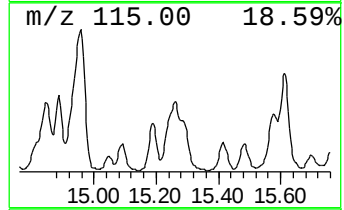
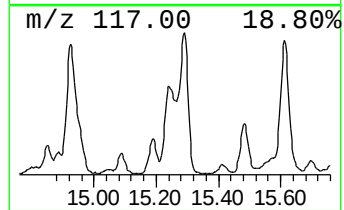
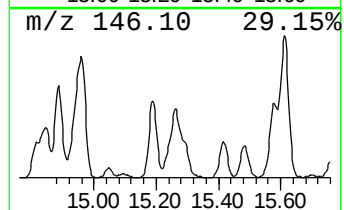
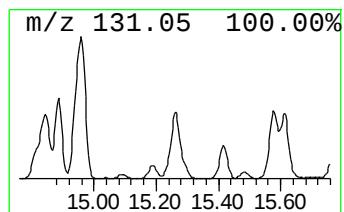
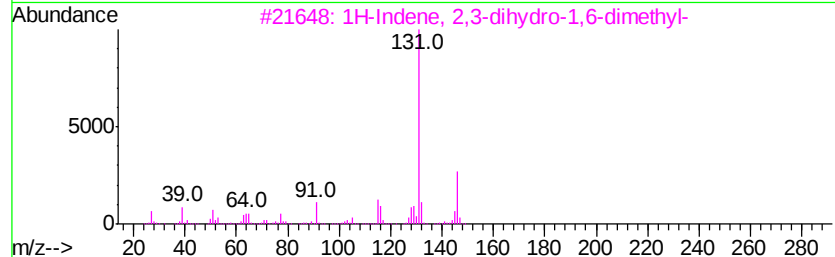
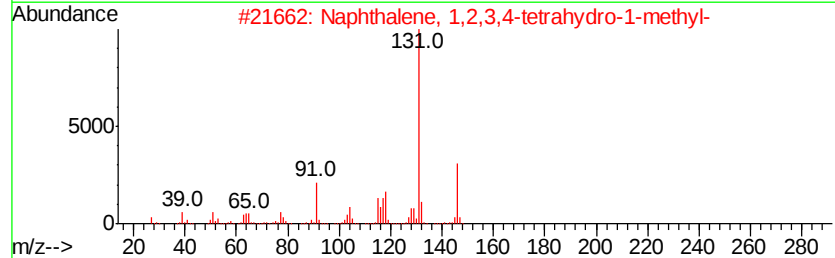
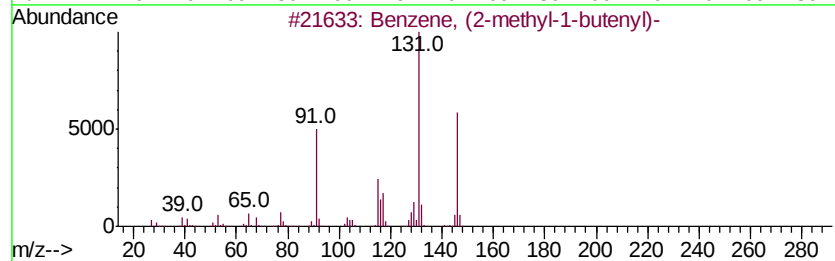
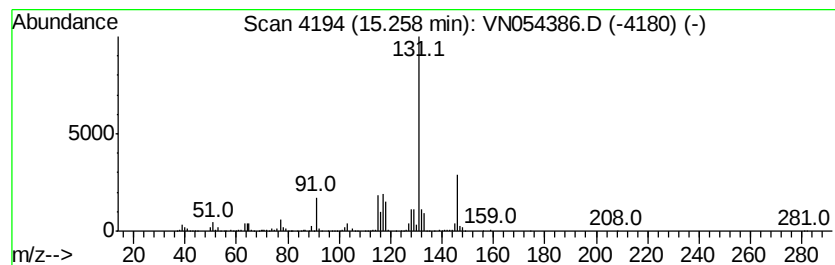
Instrument :
MSVOA_N
ClientSampleId :
T11-MW-4-9.0-031519

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 Benzene, (2-methyl-1-butenyl)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.26	25.98 ug/l	2631760	1,4-Dichlorobenzene-d4	13.34
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (2-methyl-1-butenyl)-	146 C11H14	056253-64-6 95
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001559-81-5 94
3		1H-Indene, 2,3-dihydro-1,6-dimet...	146 C11H14	017059-48-2 81
4		1H-Indene, 2,3-dihydro-1,3-dimet...	146 C11H14	004175-53-5 81
5		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 74



Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN031819\
Data File : VN054386.D
Acq On : 18 Mar 2019 11:42
Operator : JC/SP
Sample : K1960-11
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 8 Sample Multiplier: 1

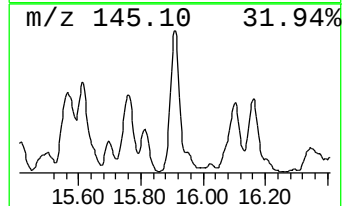
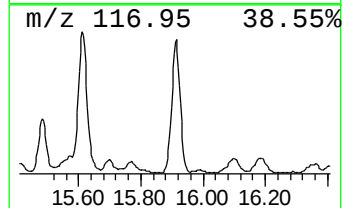
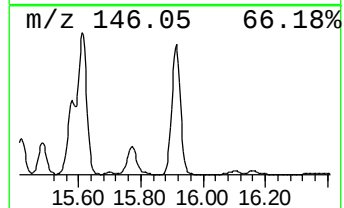
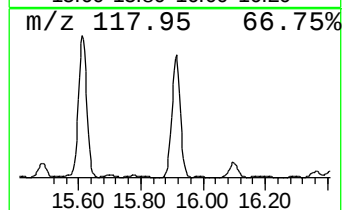
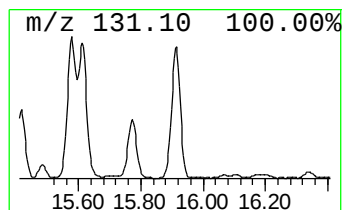
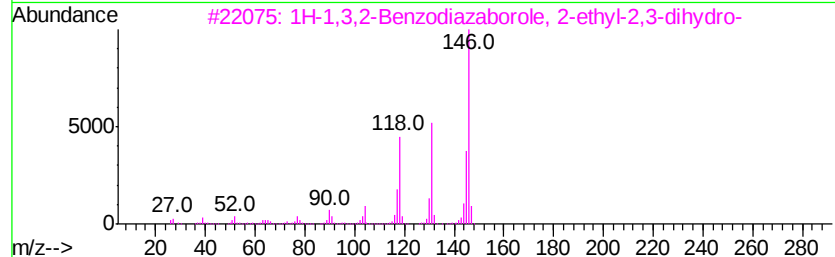
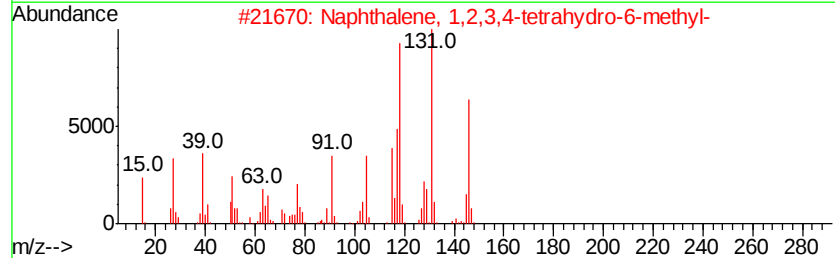
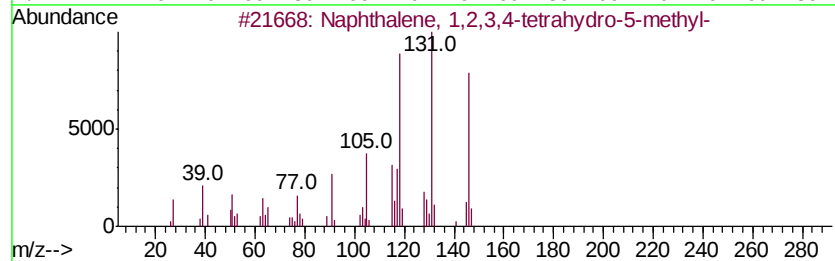
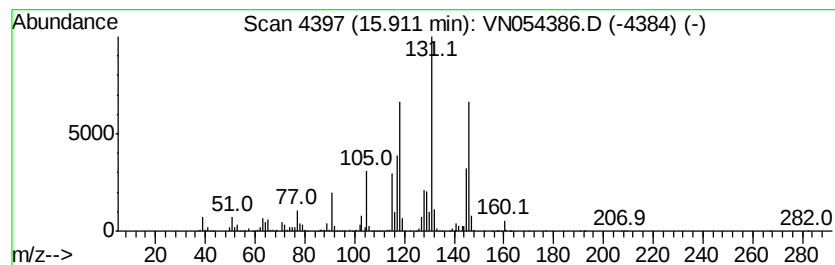
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetra... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.91	20.90 ug/l	2117320	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 95
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 94
3		1H-1,3,2-Benzodiazaborole, 2-eth...	146 C8H11BN2	074663-81-3 90
4		Benzene, 2-ethenyl-1,3,5-trimethyl-	146 C11H14	000769-25-5 78
5		Benzene, (1-methyl-1-butenyl)-	146 C11H14	053172-84-2 64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN031819\
 Data File : VN054386.D
 Acq On : 18 Mar 2019 11:42
 Operator : JC/SP
 Sample : K1960-11
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 T11-MW-4-9.0-031519

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, 1,1'-(1,...	13.54	68.5	ug/l	6939250	4	13.34	5064490	50.0
Benzene, 1-ethyl-...	13.89	54.8	ug/l	5545370	4	13.34	5064490	50.0
(E)-1-Phenyl-1-bu...	13.95	19.6	ug/l	1985070	4	13.34	5064490	50.0
Indan, 1-methyl-	13.99	78.5	ug/l	7951120	4	13.34	5064490	50.0
Benzene, 1,2,3,4-...	14.21	47.4	ug/l	4798520	4	13.34	5064490	50.0
Benzene, 1,2,4,5-...	14.58	171.9	ug/l	17417200	4	13.34	5064490	50.0
Benzene, (1-methy...	14.85	20.2	ug/l	2046130	4	13.34	5064490	50.0
Benzene, (1,2-dim...	14.96	28.6	ug/l	2901490	4	13.34	5064490	50.0
Benzene, (2-methy...	15.26	26.0	ug/l	2631760	4	13.34	5064490	50.0
Naphthalene, 1,2,...	15.91	20.9	ug/l	2117320	4	13.34	5064490	50.0