

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091118\
 Data File : VN051124.D
 Acq On : 11 Sep 2018 13:58
 Operator : MD\SY
 Sample : J4849-04 5X
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 EP1-MW-D

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\82N091018W.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	7.593	1790	1810	1821	rBV	485205	1212567	4.45%	1.602%
2	7.667	1821	1833	1871	rVB	738765	1797291	6.60%	2.374%
3	8.030	1927	1946	1971	rBV	451626	1074243	3.95%	1.419%
4	8.586	2103	2119	2153	rBV	1054489	2229540	8.19%	2.945%
5	10.091	2571	2587	2621	rBV	1928983	3614337	13.28%	4.774%
6	11.406	2983	2996	3015	rBV	1432391	2545686	9.35%	3.362%
7	11.509	3015	3028	3046	rVB2	148786	302590	1.11%	0.400%
8	11.628	3046	3065	3083	rBV2	196848	494253	1.82%	0.653%
9	12.249	3247	3258	3273	rVB	1210858	1960010	7.20%	2.589%
10	12.400	3293	3305	3319	rBV	1327100	2283718	8.39%	3.016%
11	12.590	3341	3364	3376	rBV	2302265	3770589	13.85%	4.980%
12	12.689	3386	3395	3401	rBV	848890	1294577	4.76%	1.710%
13	12.731	3401	3408	3419	rVB	980210	1563022	5.74%	2.064%
14	13.040	3492	3504	3529	rVB	16827917	27218618	100.00%	35.951%
15	13.168	3530	3544	3555	rBV3	645914	1348314	4.95%	1.781%
16	13.287	3572	3581	3589	rBV	470980	708473	2.60%	0.936%
17	13.342	3589	3598	3606	rBV	1323972	2384684	8.76%	3.150%
18	13.442	3619	3629	3640	rBV	266572	418321	1.54%	0.553%
19	13.512	3640	3651	3654	rVV	722170	1049627	3.86%	1.386%
20	13.541	3655	3660	3669	rVV	1249246	1970309	7.24%	2.602%
21	13.593	3669	3676	3691	rVV3	641473	1426433	5.24%	1.884%
22	13.673	3691	3701	3721	rVB	378863	681460	2.50%	0.900%
23	13.802	3727	3741	3746	rBV	978511	1614231	5.93%	2.132%
24	13.831	3746	3750	3759	rVV	1025066	1454452	5.34%	1.921%
25	13.889	3759	3768	3777	rVV	1564190	2361272	8.68%	3.119%
26	13.946	3777	3786	3792	rVV	504679	825513	3.03%	1.090%
27	13.991	3792	3800	3812	rVB2	1445985	2430100	8.93%	3.210%
28	14.130	3833	3843	3851	rBV2	232287	357149	1.31%	0.472%
29	14.210	3851	3868	3873	rBV	701788	1160127	4.26%	1.532%
30	14.249	3873	3880	3889	rVB	999969	1551809	5.70%	2.050%
31	14.477	3943	3951	3960	rBV2	174672	280097	1.03%	0.370%
32	14.583	3973	3984	3999	rBV4	1162079	2326345	8.55%	3.073%

LSC Area Percent Report

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ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EP1-MW-D

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\82N091018W.M
Title : SW846 8260

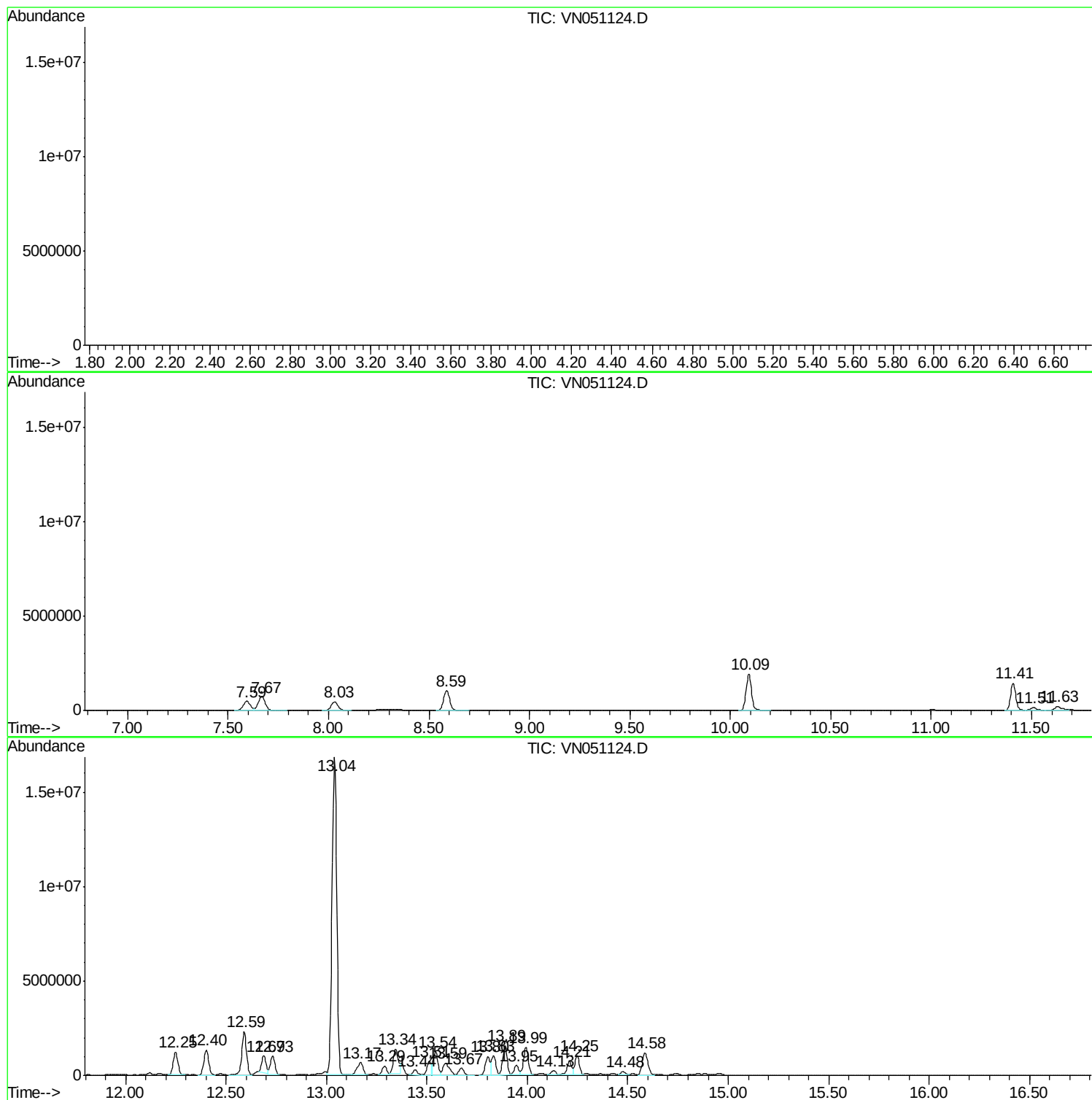
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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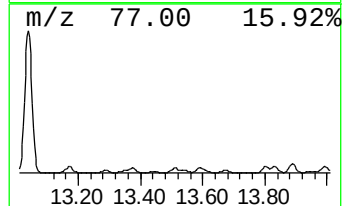
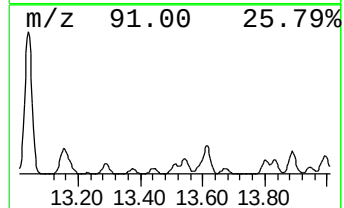
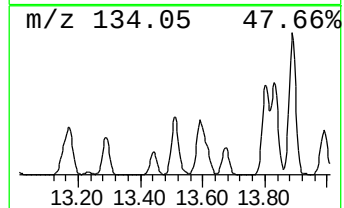
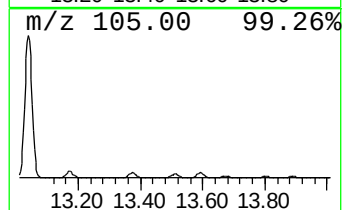
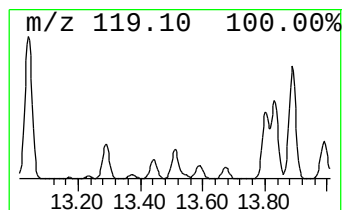
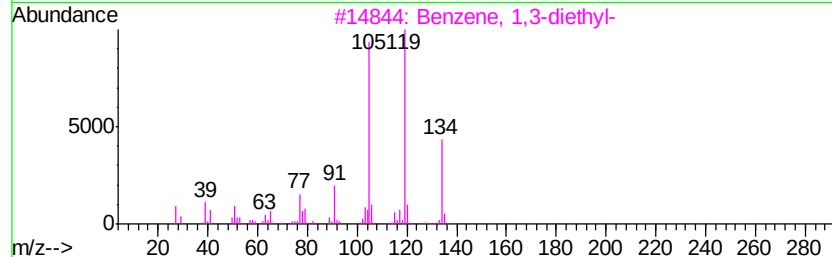
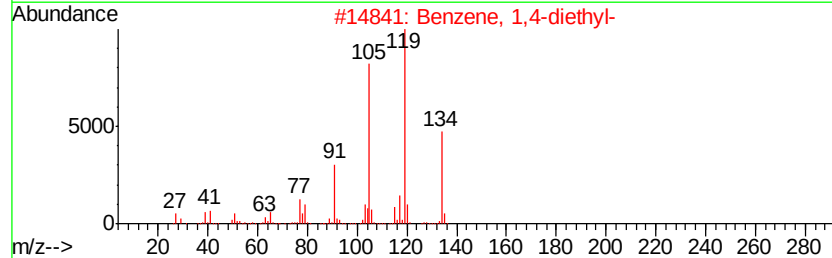
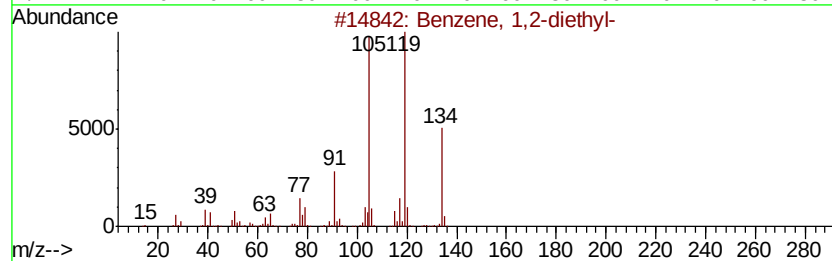
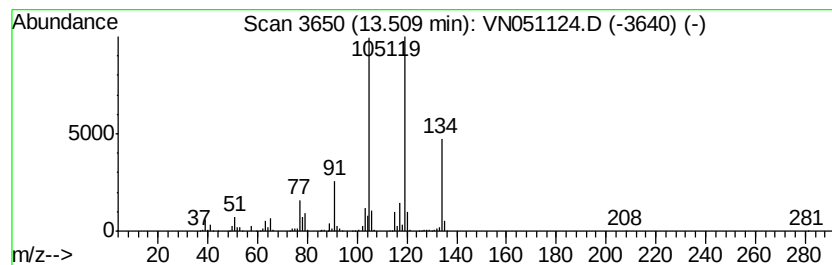
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.51	22.01 ug/l	1049630	1,4-Dichlorobenzene-d4	13.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
2		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	96
3		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	93
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	90
5		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	87



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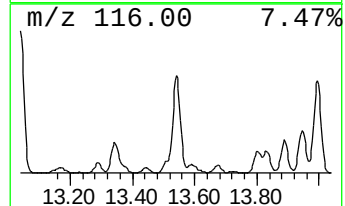
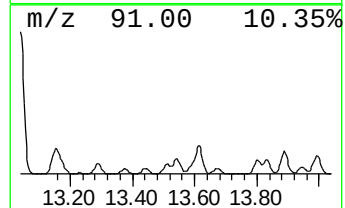
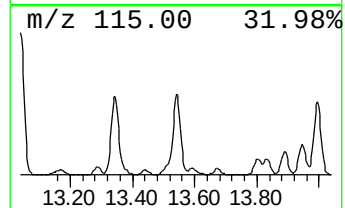
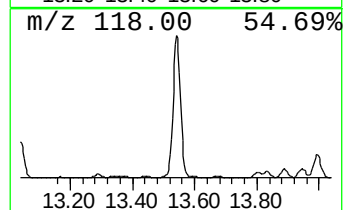
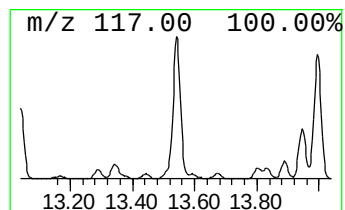
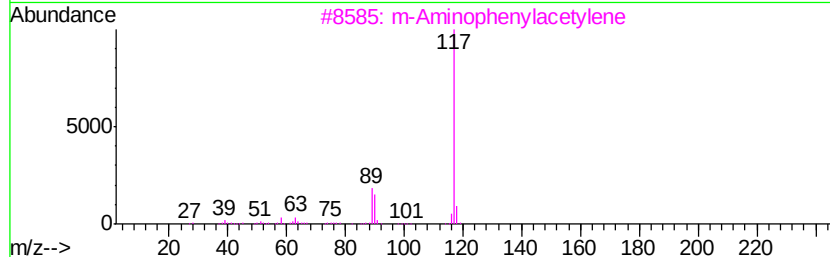
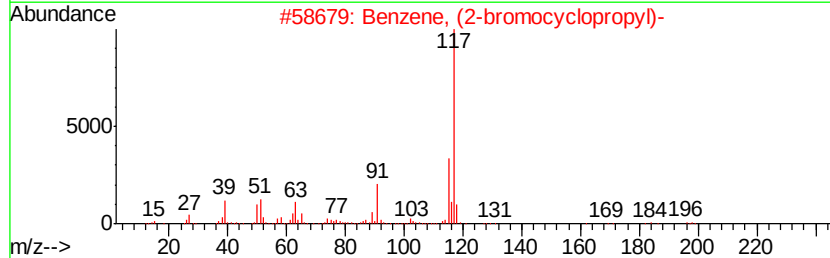
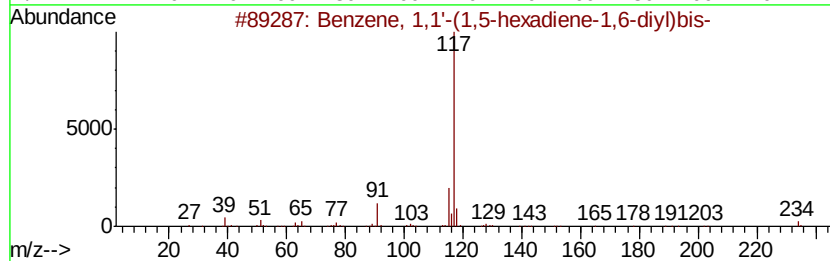
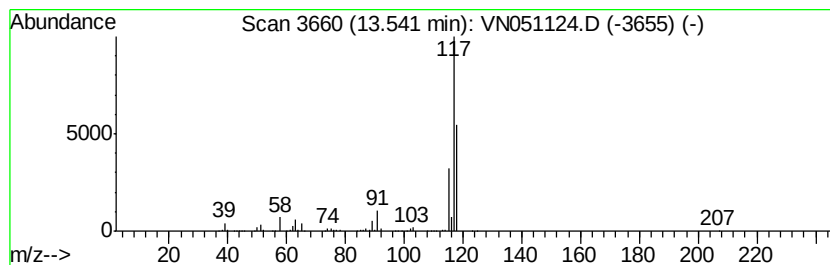
Instrument :
MSVOA_N
ClientSampleId :
EP1-MW-D

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091018W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1,1'-(1,5-hexadien... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.54	41.31 ug/l	1970310	1,4-Dichlorobenzene-d4	13.35
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		Benzene, (2-bromocyclopropyl)-	196 C9H9Br	036617-02-4 45
3		m-Aminophenylacetylene	117 C8H7N	054060-30-9 37
4		trans-Cinnamyl bromide	196 C9H9Br	026146-77-0 25
5		3-Bromo-1-phenyl-1-propene	196 C9H9Br	004392-24-9 17



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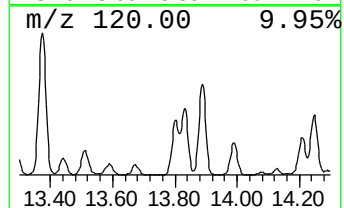
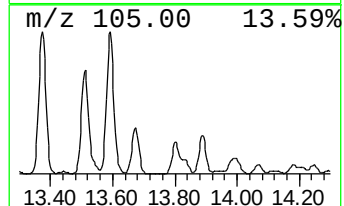
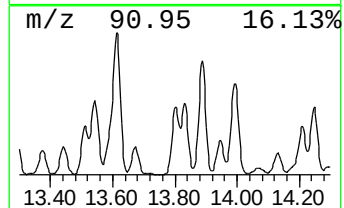
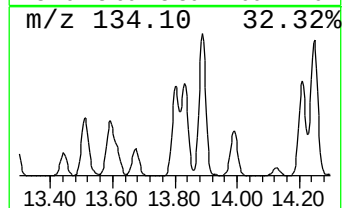
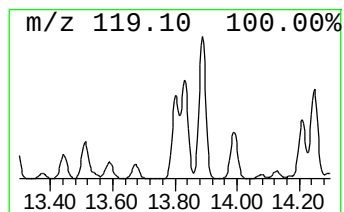
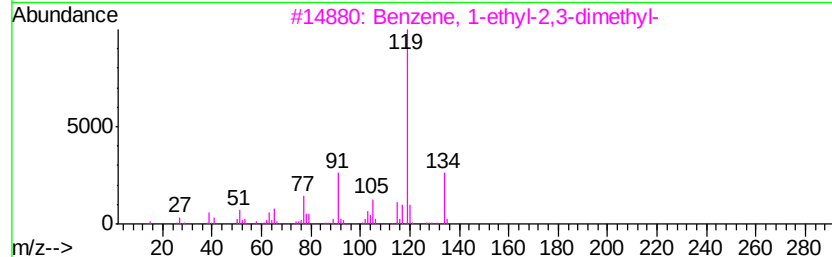
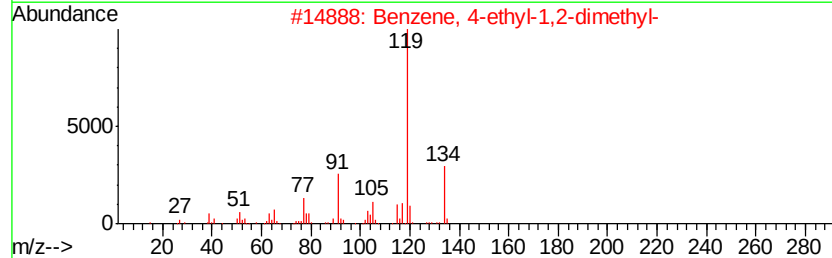
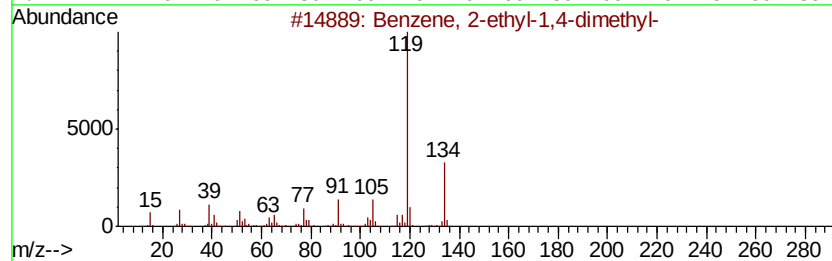
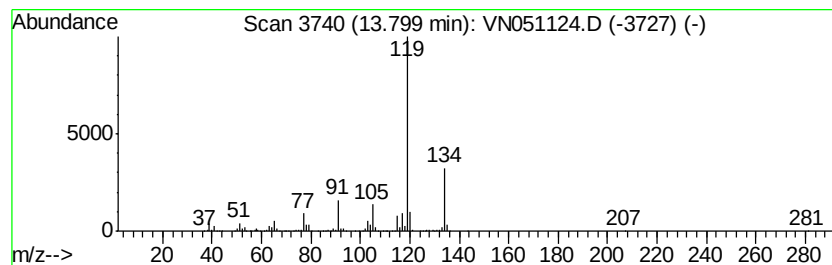
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.80	33.85 ug/l	1614230	1,4-Dichlorobenzene-d4	13.35

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
5	o-Cymene	134	C10H14	000527-84-4	95



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Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 16 Sample Multiplier: 1

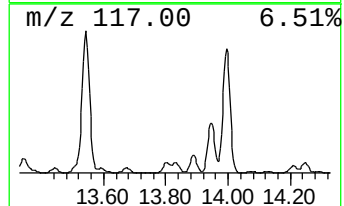
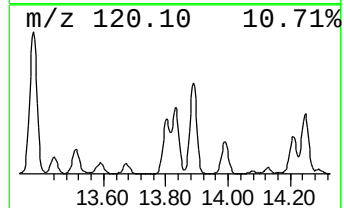
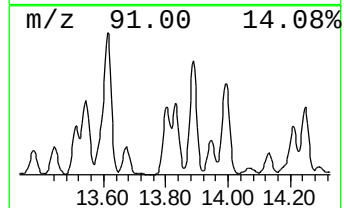
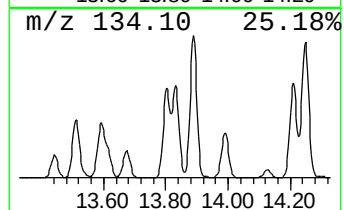
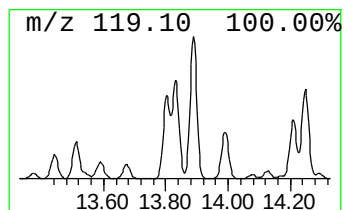
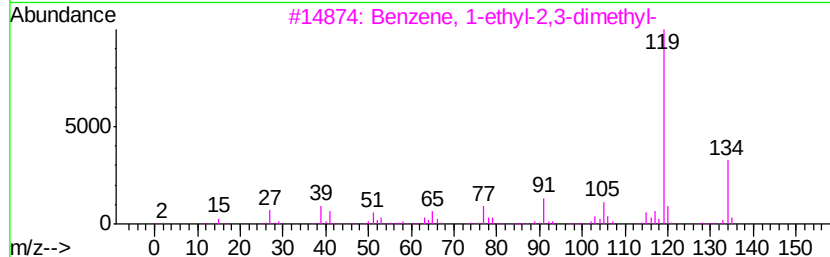
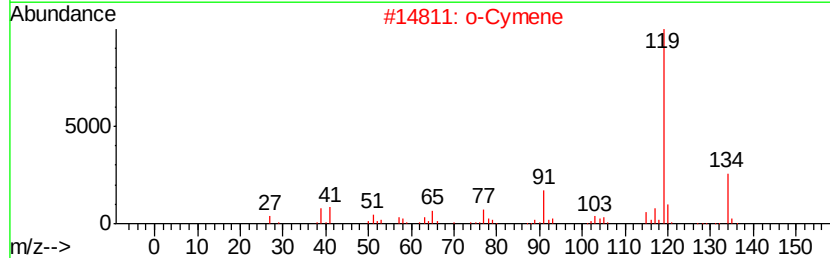
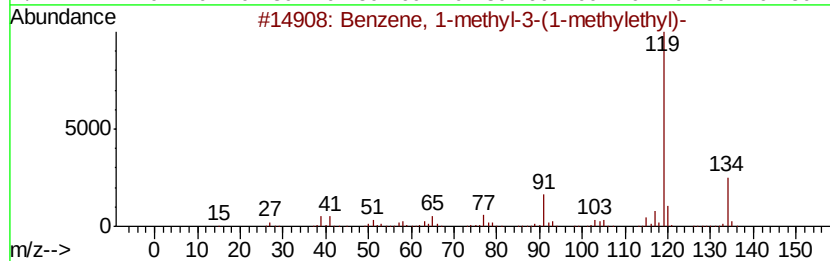
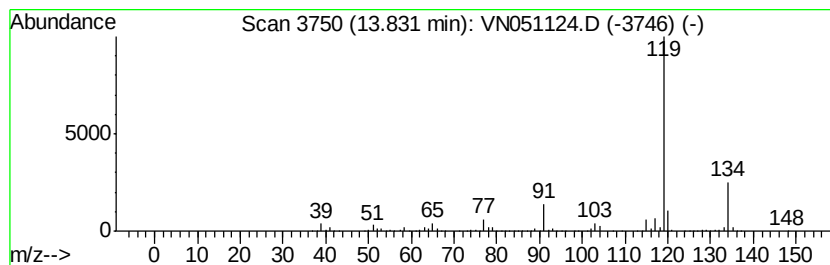
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ClientSampled :
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-methyl-3-(1-meth... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	30.50 ug/l	1454450	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1-methyl-3-(1-methyleth...	134 C10H14	000535-77-3	94
2	o-Cymene	134 C10H14	000527-84-4	94
3	Benzene, 1-ethyl-2,3-dimethyl-	134 C10H14	000933-98-2	91
4	p-Cymene	134 C10H14	000099-87-6	91
5	Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9	91



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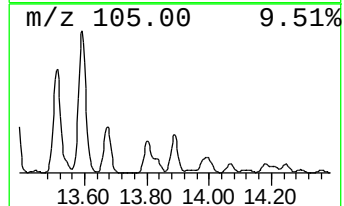
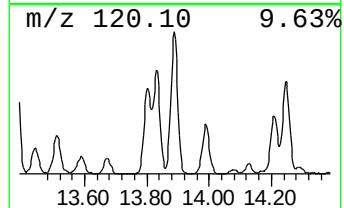
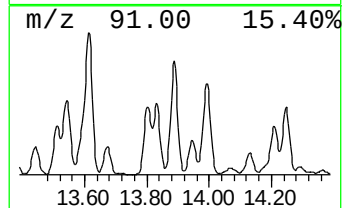
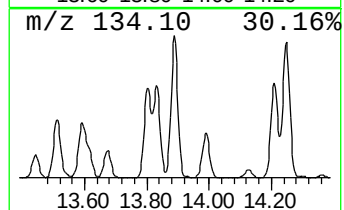
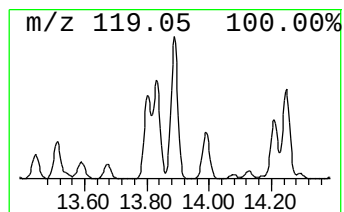
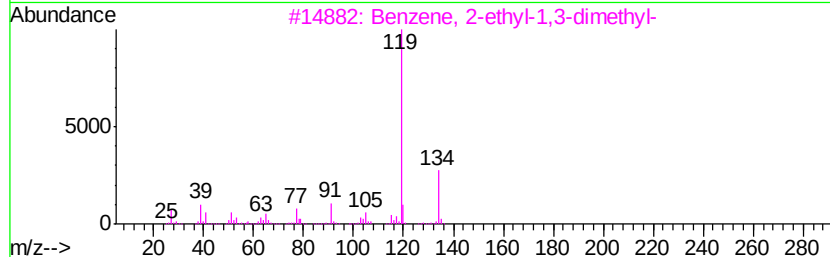
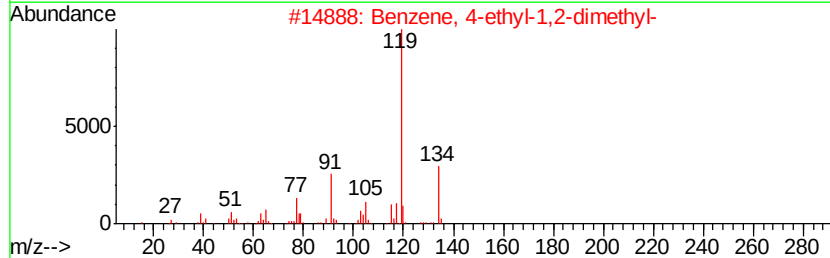
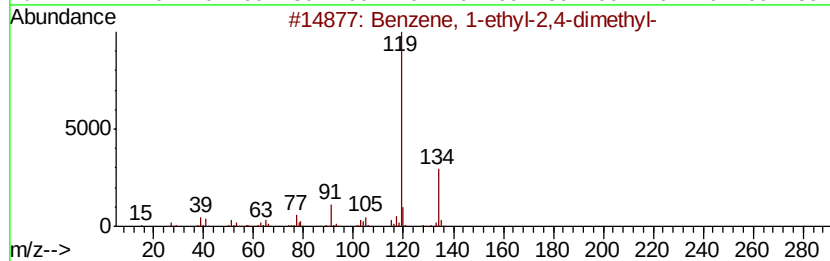
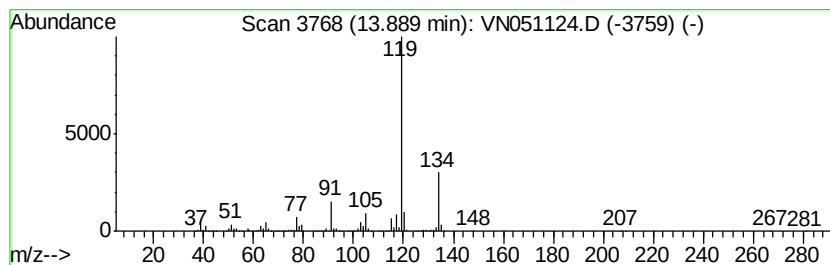
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	49.51 ug/l	2361270	1,4-Dichlorobenzene-d4	13.35

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



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Sample : J4849-04 5X
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
EP1-MW-D

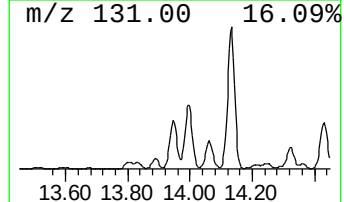
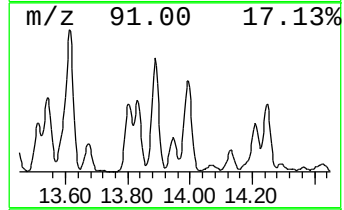
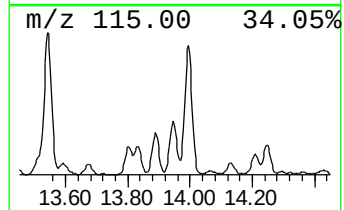
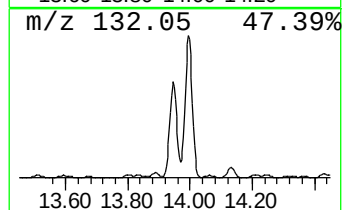
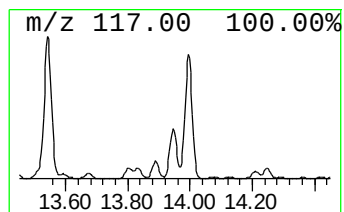
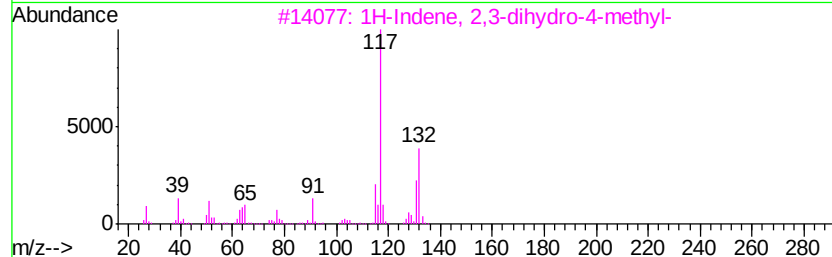
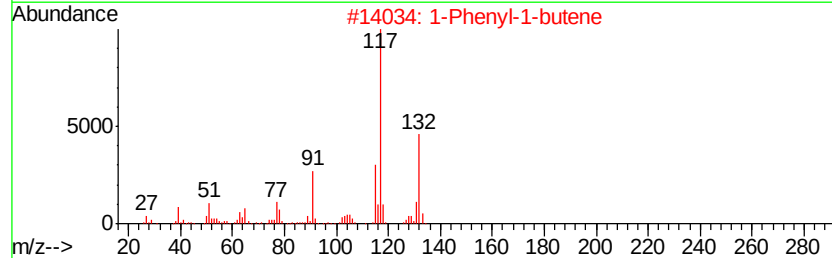
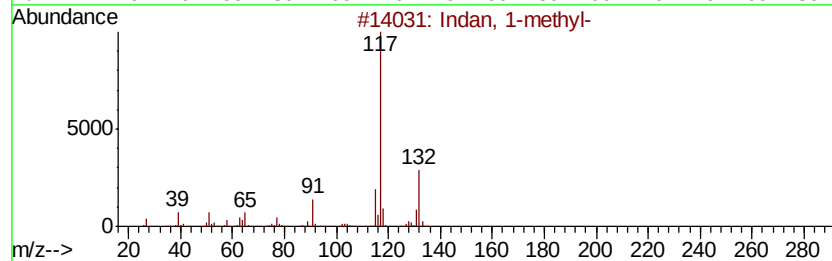
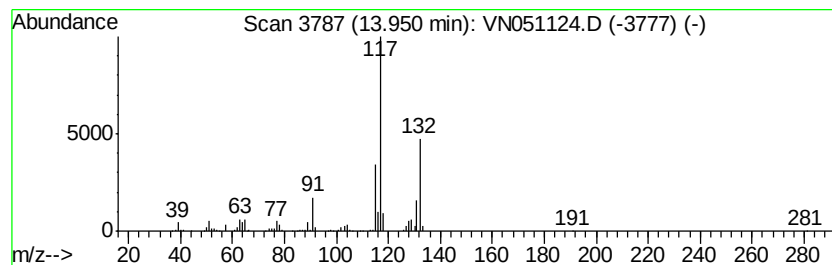
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Quant Title : SW846 8260

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

Peak Number 6 Indan, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.95	17.31 ug/l	825513	1,4-Dichlorobenzene-d4	13.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indan, 1-methyl-	132	C10H12	000767-58-8	90
2		1-Phenyl-1-butene	132	C10H12	000824-90-8	87
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	86
4		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	86
5		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	86



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN091118\
Data File : VN051124.D
Acq On : 11 Sep 2018 13:58
Operator : MD\SY
Sample : J4849-04 5X
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 16 Sample Multiplier: 1

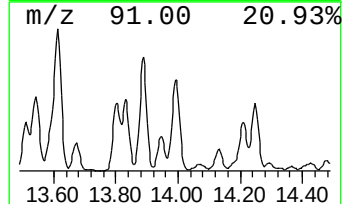
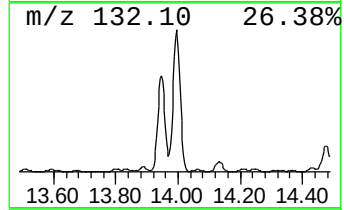
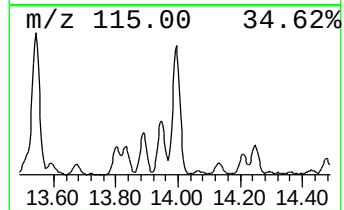
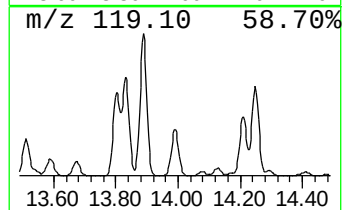
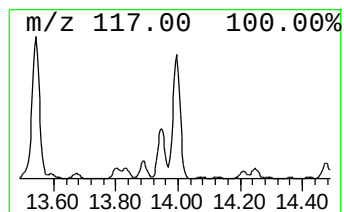
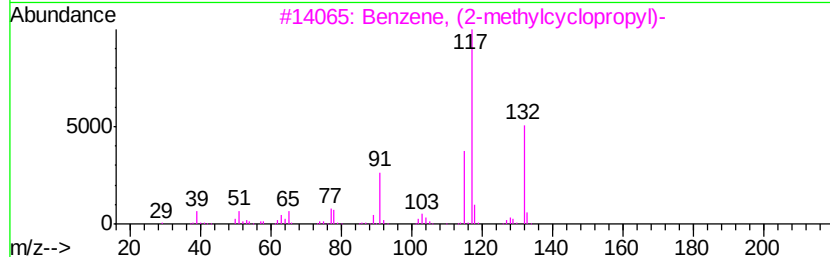
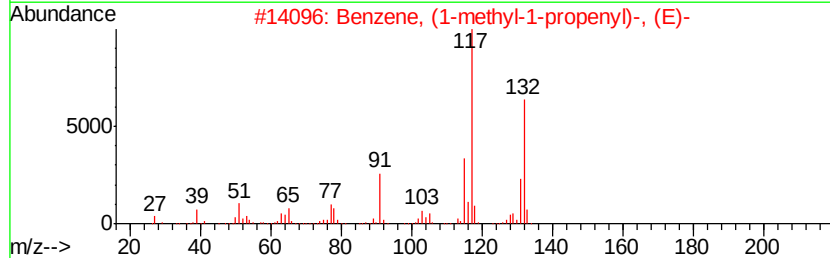
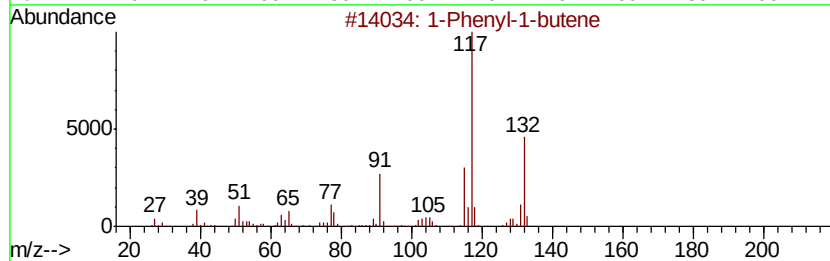
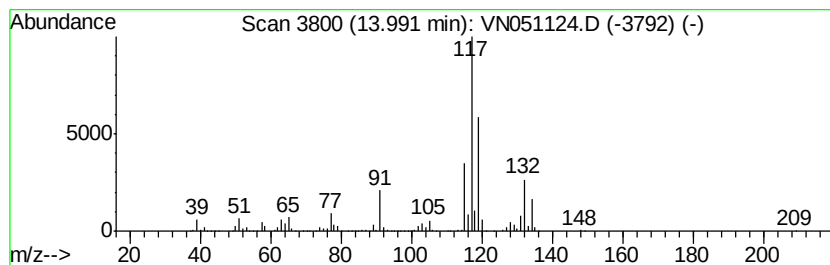
Instrument :
MSVOA_N
ClientSampleId :
EP1-MW-D

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091018W.M
Quant Title : SW846 8260

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

Peak Number 7 1-Phenyl-1-butene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	50.95 ug/l	2430100	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Phenyl-1-butene	132 C10H12	000824-90-8	70
2	Benzene, (1-methyl-1-propenyl)-,...	132 C10H12	000768-00-3	70
3	Benzene, (2-methylcyclopropyl)-	132 C10H12	1000327-39-0	64
4	Indan, 1-methyl-	132 C10H12	000767-58-8	64
5	Benzene, (2-methyl-1-propenyl)-	132 C10H12	000768-49-0	62



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN091118\
Data File : VN051124.D
Acq On : 11 Sep 2018 13:58
Operator : MD\SY
Sample : J4849-04 5X
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampled :
EP1-MW-D

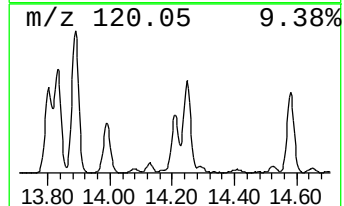
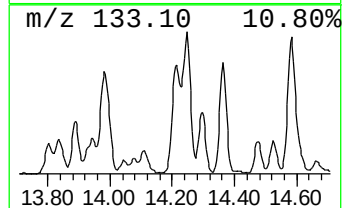
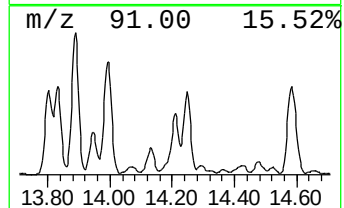
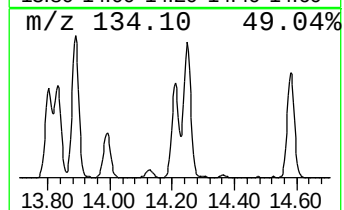
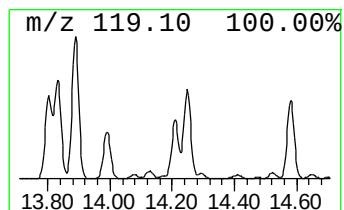
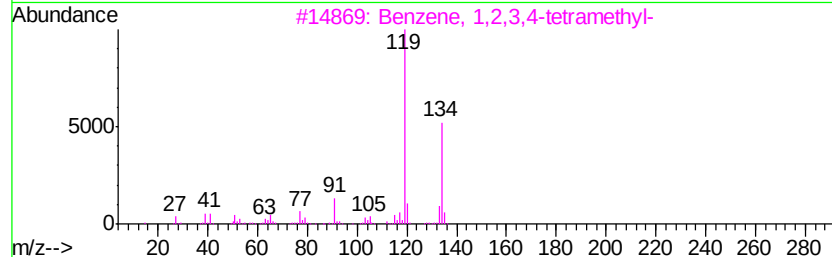
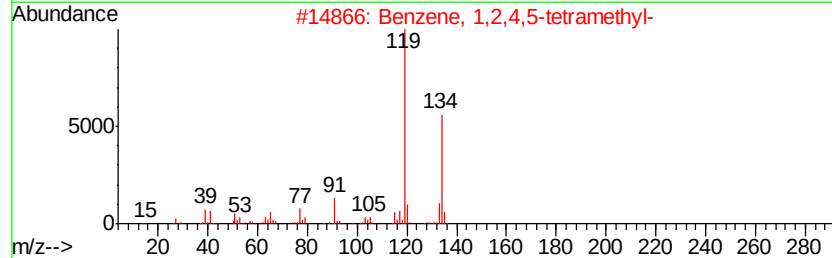
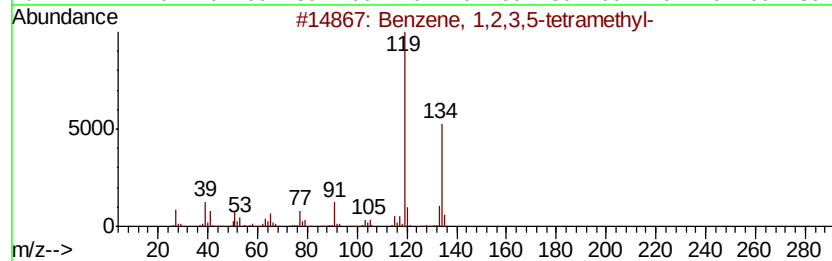
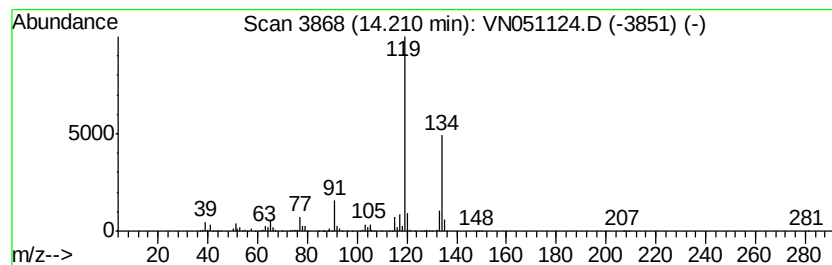
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	24.32 ug/l	1160130	1,4-Dichlorobenzene-d4	13.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN091118\
Data File : VN051124.D
Acq On : 11 Sep 2018 13:58
Operator : MD\SY
Sample : J4849-04 5X
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 16 Sample Multiplier: 1

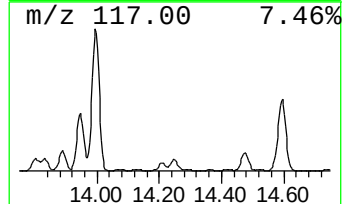
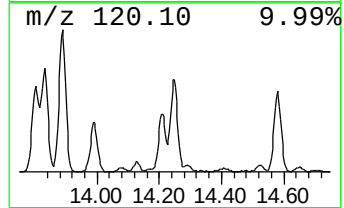
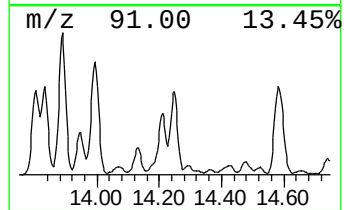
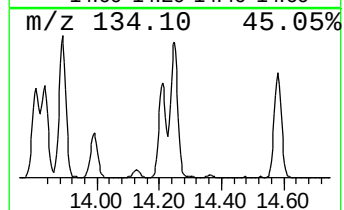
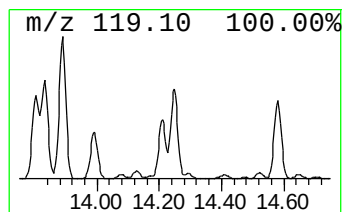
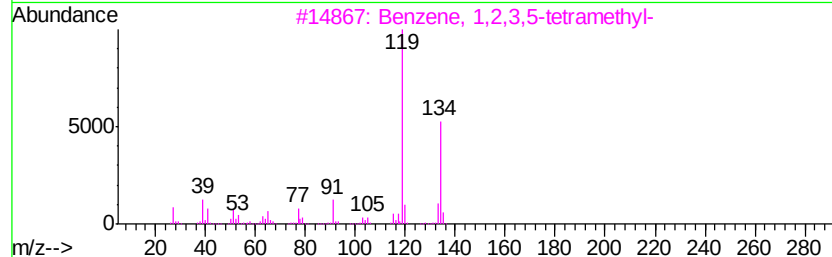
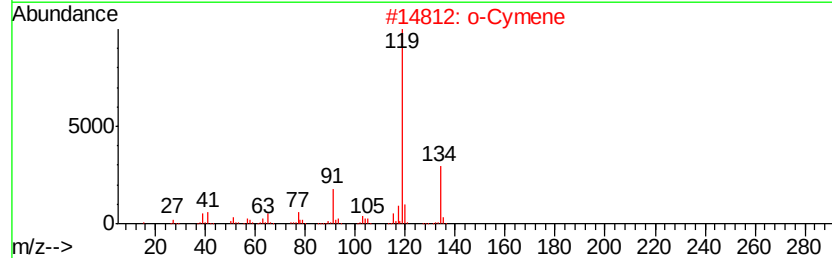
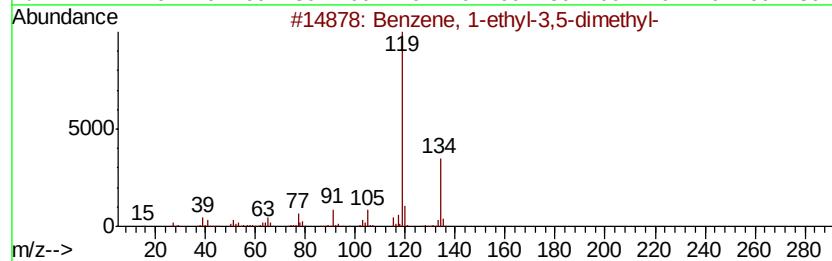
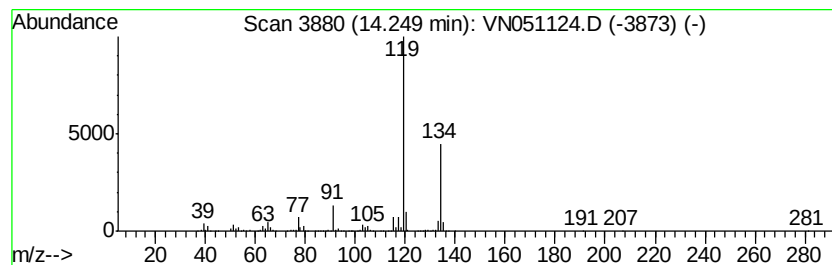
Instrument :
MSVOA_N
ClientSampleId :
EP1-MW-D

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091018W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.25	32.54 ug/l	1551810	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1-ethyl-3,5-dimethyl-	134 C10H14	000934-74-7	94
2	o-Cymene	134 C10H14	000527-84-4	94
3	Benzene, 1,2,3,5-tetramethyl-	134 C10H14	000527-53-7	94
4	Benzene, 1,2,4,5-tetramethyl-	134 C10H14	000095-93-2	94
5	Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN091118\
Data File : VN051124.D
Acq On : 11 Sep 2018 13:58
Operator : MD\SY
Sample : J4849-04 5X
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampled :
EP1-MW-D

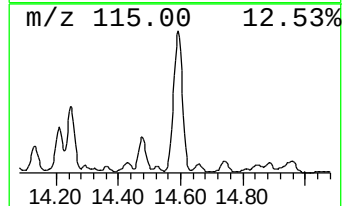
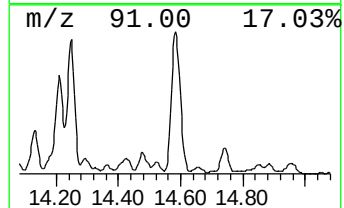
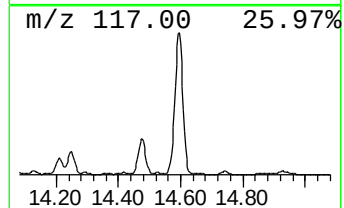
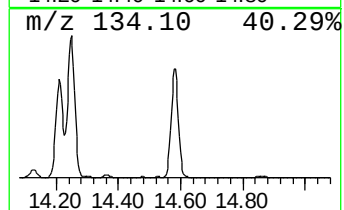
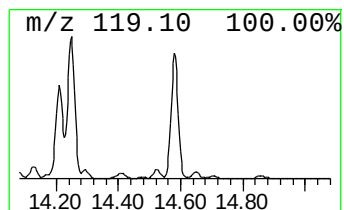
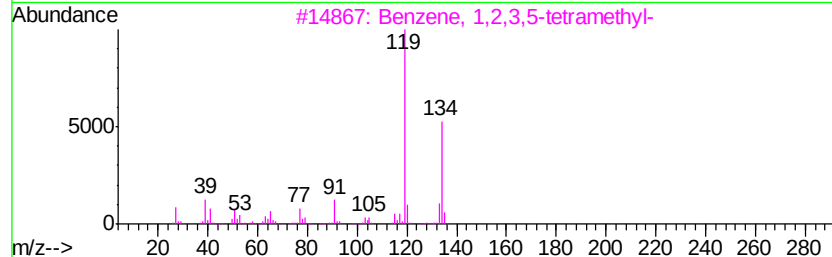
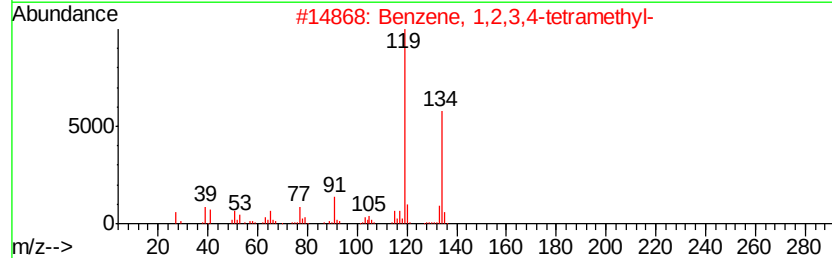
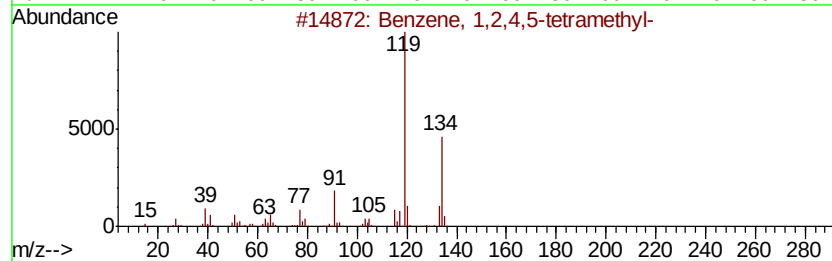
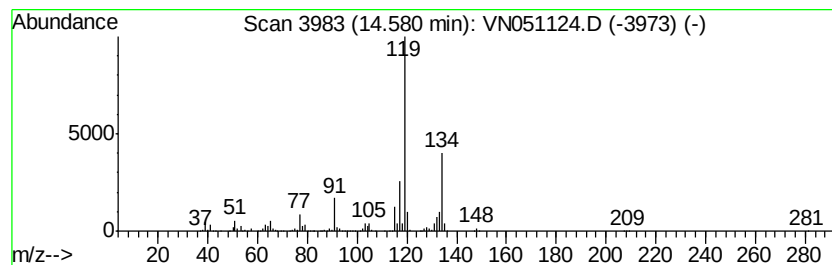
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091018W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.58	48.78 ug/l	2326350	1,4-Dichlorobenzene-d4	13.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	83
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	81
4		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	81
5		o-Cymene	134	C10H14	000527-84-4	81



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN091118\
 Data File : VN051124.D
 Acq On : 11 Sep 2018 13:58
 Operator : MD\SY
 Sample : J4849-04 5X
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 EP1-MW-D

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091018W.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, 1,2-diet...	13.51	22.0	ug/l	1049630	4	13.35	2384680	50.0
Benzene, 1,1'-(1,...	13.54	41.3	ug/l	1970310	4	13.35	2384680	50.0
Benzene, 2-ethyl-...	13.80	33.9	ug/l	1614230	4	13.35	2384680	50.0
Benzene, 1-methyl...	13.83	30.5	ug/l	1454450	4	13.35	2384680	50.0
Benzene, 1-ethyl-...	13.89	49.5	ug/l	2361270	4	13.35	2384680	50.0
Indan, 1-methyl-	13.95	17.3	ug/l	825513	4	13.35	2384680	50.0
1-Phenyl-1-butene	13.99	51.0	ug/l	2430100	4	13.35	2384680	50.0
Benzene, 1,2,3,5-...	14.21	24.3	ug/l	1160130	4	13.35	2384680	50.0
Benzene, 1-ethyl-...	14.25	32.5	ug/l	1551810	4	13.35	2384680	50.0
Benzene, 1,2,4,5-...	14.58	48.8	ug/l	2326350	4	13.35	2384680	50.0