

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN080218\
 Data File : VN050292.D
 Acq On : 3 Aug 2018 3:48
 Operator : MD\SY
 Sample : J4304-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 EW-A07

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\82N072418W.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.661	3	22	55	rBV	2892846	4478306	100.00%	11.451%
2	7.593	1848	1867	1879	rBV2	504940	1282060	28.63%	3.278%
3	7.667	1879	1890	1926	rVB	711969	1699545	37.95%	4.346%
4	8.030	1984	2003	2029	rBV	507536	1201469	26.83%	3.072%
5	8.242	2039	2069	2072	rBV3	171485	650187	14.52%	1.663%
6	8.590	2161	2177	2211	rVB	1083772	2303312	51.43%	5.890%
7	10.094	2625	2645	2678	rBV	1930777	3568043	79.67%	9.123%
8	10.432	2738	2750	2762	rVB4	35116	67637	1.51%	0.173%
9	11.409	3039	3054	3073	rBV	1425098	2470927	55.18%	6.318%
10	11.496	3073	3081	3090	rVB2	37177	61833	1.38%	0.158%
11	11.657	3120	3131	3152	rVB5	51199	137114	3.06%	0.351%
12	11.924	3200	3214	3222	rBV7	41909	95395	2.13%	0.244%
13	11.972	3223	3229	3252	rVB3	33534	84718	1.89%	0.217%
14	12.120	3267	3275	3285	rVB2	61478	102488	2.29%	0.262%
15	12.252	3307	3316	3327	rVB	85044	151299	3.38%	0.387%
16	12.403	3350	3363	3376	rBV	1215838	2050089	45.78%	5.242%
17	12.474	3377	3385	3397	rVV3	47602	88083	1.97%	0.225%
18	12.567	3399	3414	3429	rVV4	71649	180820	4.04%	0.462%
19	12.728	3456	3464	3475	rVB9	29438	59332	1.32%	0.152%
20	12.962	3525	3537	3542	rBV3	54303	98203	2.19%	0.251%
21	13.171	3590	3602	3610	rBV3	74898	148697	3.32%	0.380%
22	13.345	3643	3656	3670	rBV	1262537	2073808	46.31%	5.303%
23	13.445	3670	3687	3698	rVV	183136	344652	7.70%	0.881%
24	13.512	3698	3708	3730	rVB	513778	855737	19.11%	2.188%
25	13.676	3745	3759	3778	rVB	506584	908903	20.30%	2.324%
26	13.888	3816	3825	3835	rVB2	255379	388792	8.68%	0.994%
27	13.995	3846	3858	3868	rBV4	593410	1044946	23.33%	2.672%
28	14.133	3891	3901	3908	rBV	228360	369688	8.26%	0.945%
29	14.210	3909	3925	3941	rVB2	843469	1662777	37.13%	4.252%
30	14.294	3942	3951	3956	rBV2	129866	202185	4.51%	0.517%
31	14.326	3957	3961	3967	rVV	115713	170185	3.80%	0.435%
32	14.364	3967	3973	3981	rVV	159535	248473	5.55%	0.635%
33	14.419	3981	3990	4002	rVB3	285519	602429	13.45%	1.540%
34	14.490	4002	4012	4018	rBV3	117597	197191	4.40%	0.504%

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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 >
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35	14.532	4018	4025	4032	rVV3	119524	187774	4.19%	0.480%
36	14.583	4032	4041	4055	rVV	255031	445986	9.96%	1.140%
37	14.657	4055	4064	4073	rVV3	169501	332279	7.42%	0.850%
38	14.705	4073	4079	4088	rVB	68251	101959	2.28%	0.261%
39	14.789	4096	4105	4113	rBV5	85294	167192	3.73%	0.428%
40	14.853	4115	4125	4132	rBV3	363599	610734	13.64%	1.562%
41	14.888	4132	4136	4142	rVV	64573	89661	2.00%	0.229%
42	14.959	4148	4158	4177	rVB2	684268	1496852	33.42%	3.827%
43	15.049	4177	4186	4193	rBV2	89362	164810	3.68%	0.421%
44	15.094	4194	4200	4210	rVB4	105536	182654	4.08%	0.467%
45	15.191	4219	4230	4240	rBV	615212	1087117	24.28%	2.780%
46	15.265	4241	4253	4288	rVB3	676596	1671546	37.33%	4.274%
47	15.419	4289	4301	4311	rBV2	85866	175744	3.92%	0.449%
48	15.486	4314	4322	4332	rBV5	93915	180172	4.02%	0.461%
49	15.702	4380	4389	4399	rBV2	160441	281673	6.29%	0.720%
50	15.914	4442	4455	4467	rBV5	324135	670039	14.96%	1.713%
51	16.104	4499	4514	4525	rBV3	233632	540010	12.06%	1.381%
52	16.168	4526	4534	4543	rVV3	118706	223124	4.98%	0.571%
53	16.358	4580	4593	4604	rBV6	164229	449908	10.05%	1.150%

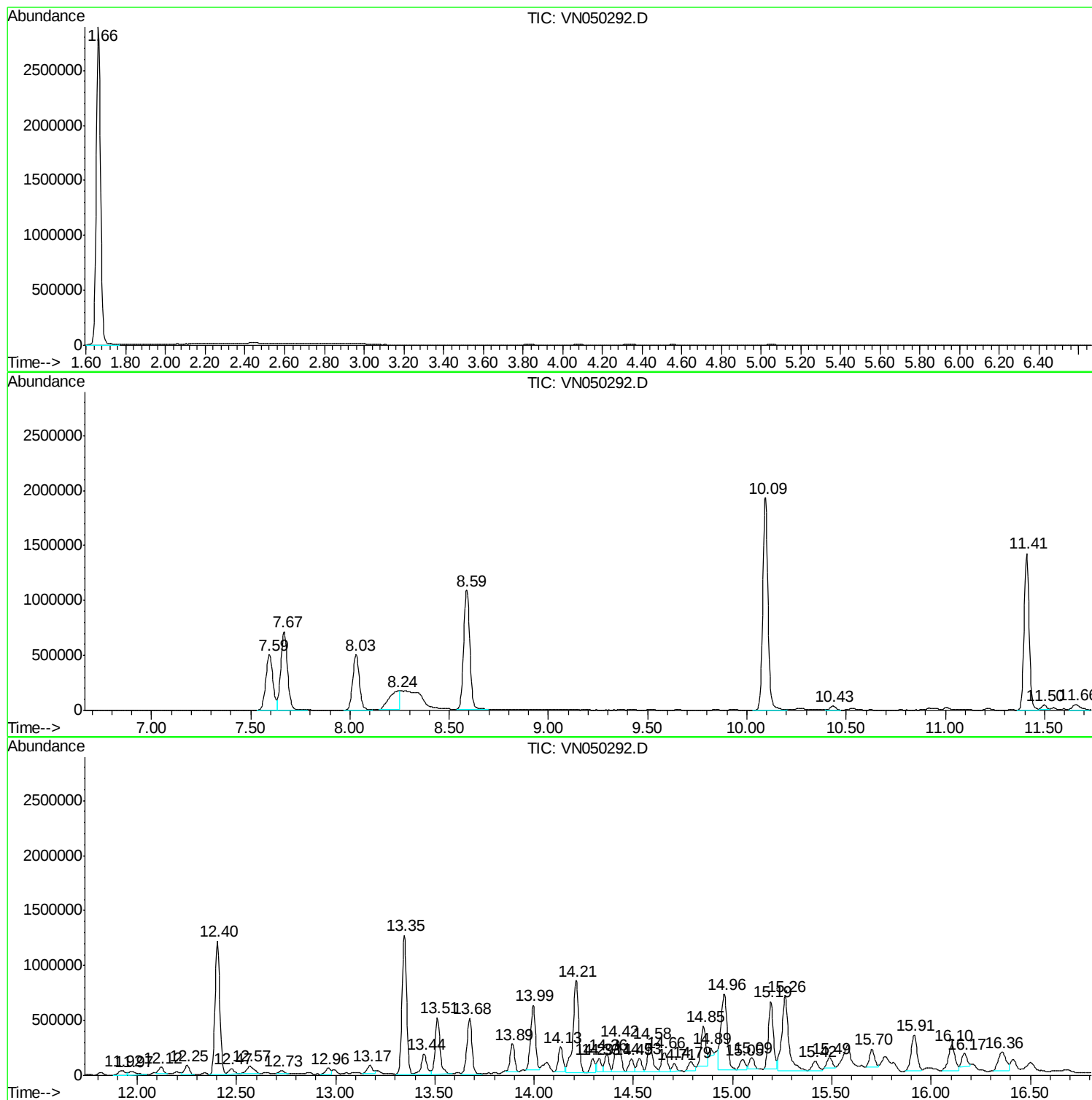
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Quant Title : SW846 8260

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



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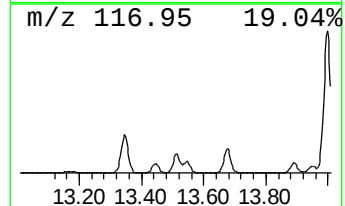
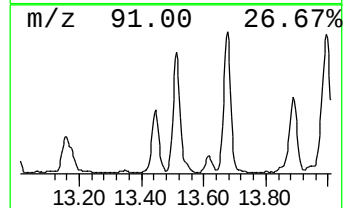
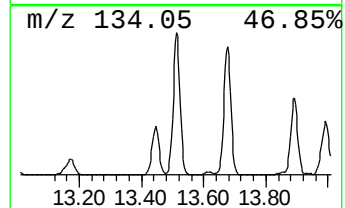
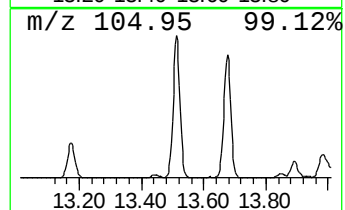
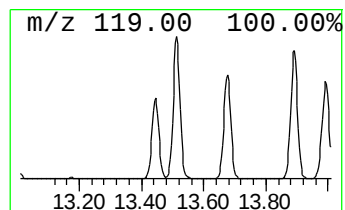
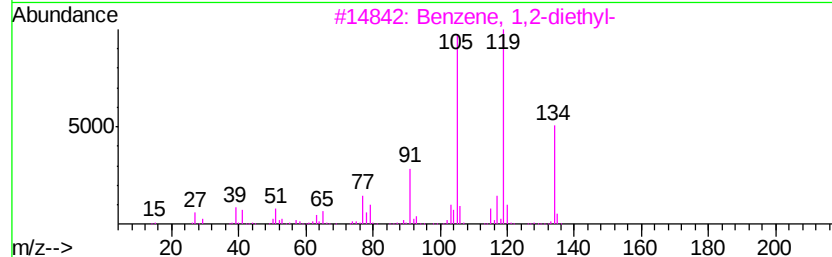
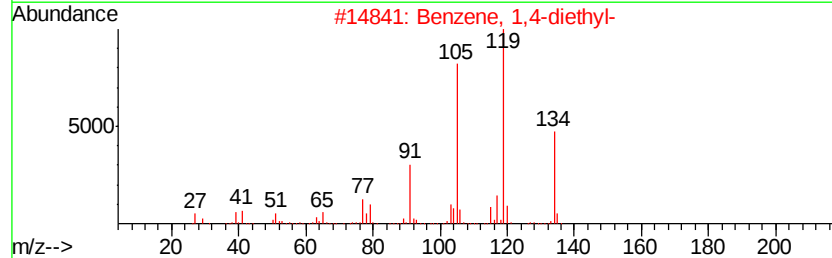
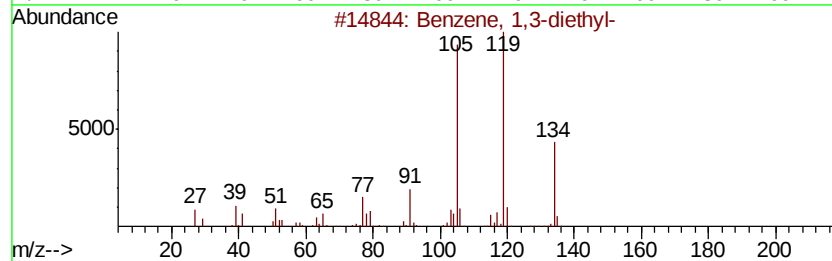
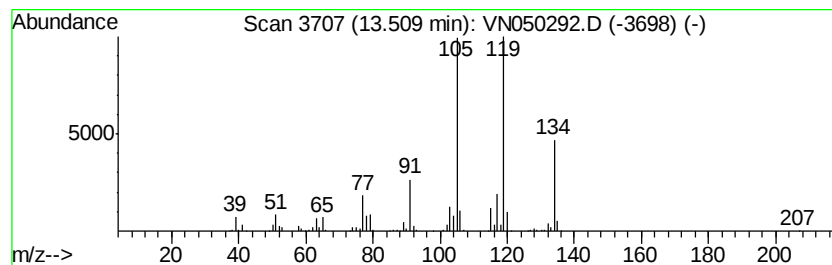
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1,3-diethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.51	20.63 ug/l	855737	1,4-Dichlorobenzene-d4	13.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	97
2		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	97
3		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	90
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	87



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Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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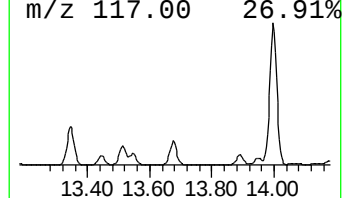
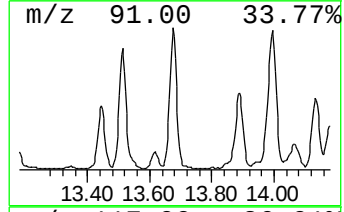
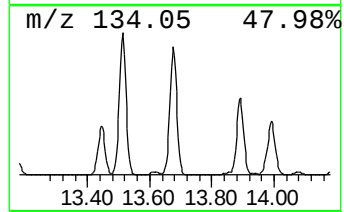
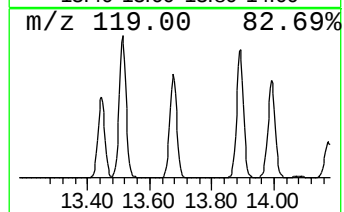
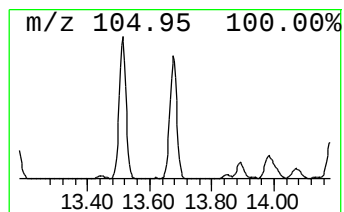
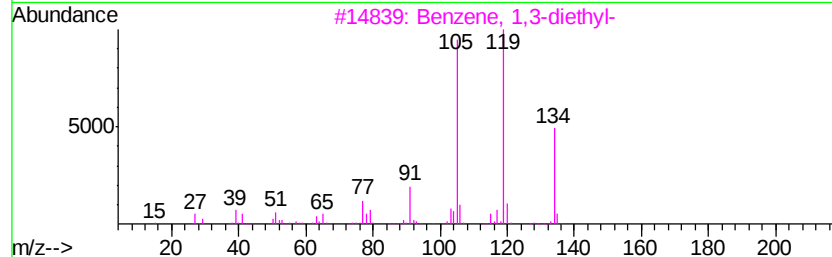
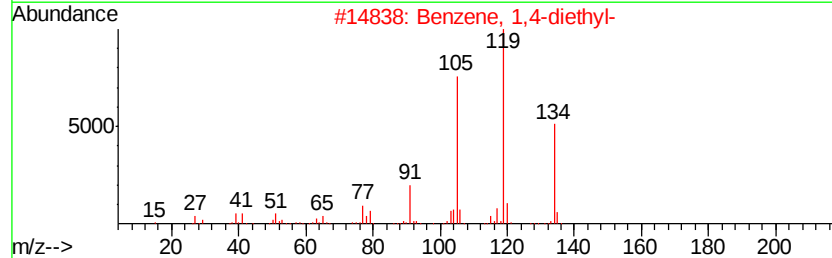
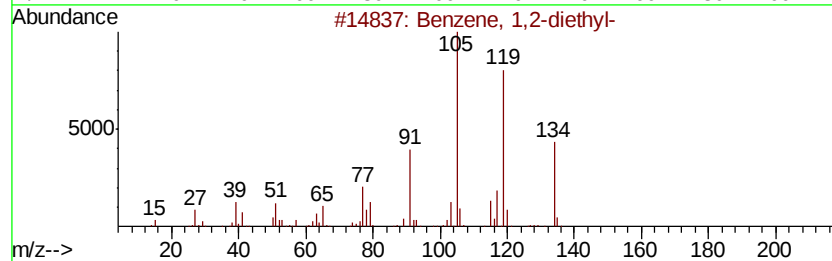
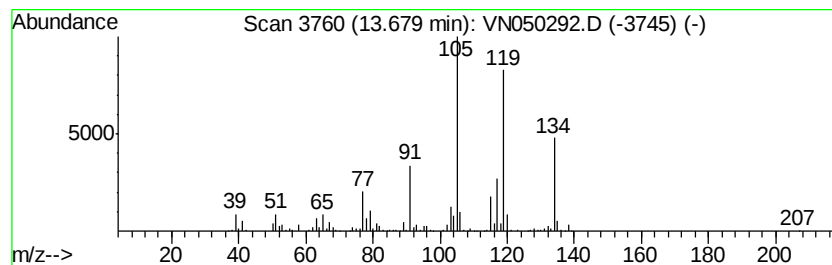
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1,2-diethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.68	21.91 ug/l	908903	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	90
3		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	87
4		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	62
5		Ethanone, 1-(4-methylphenyl)-	134	C9H10O	000122-00-9	60



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Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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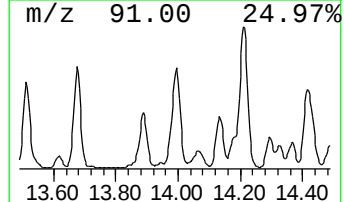
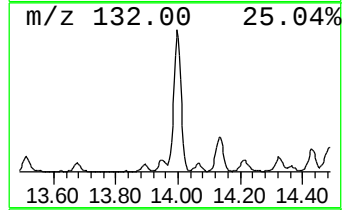
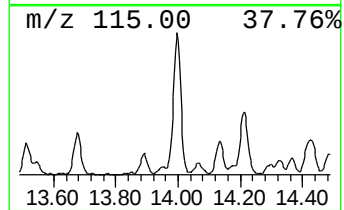
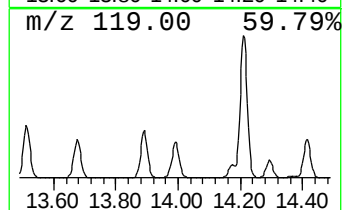
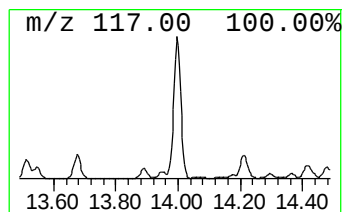
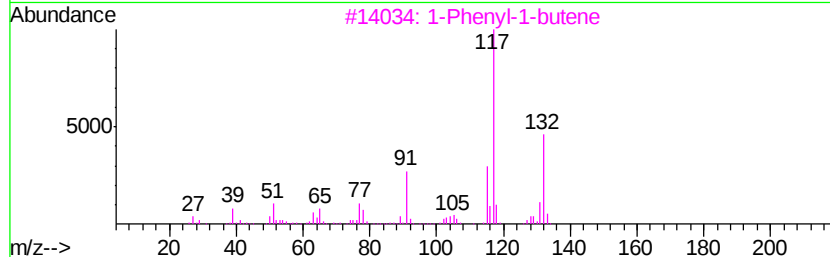
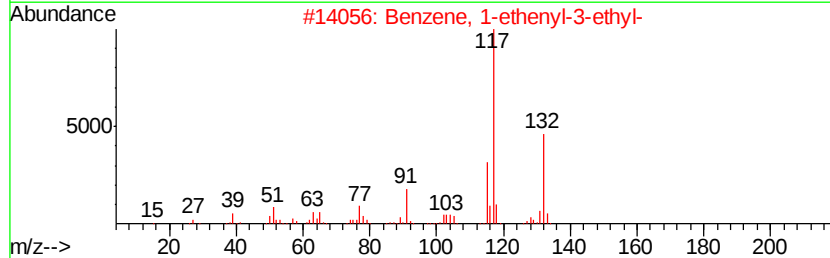
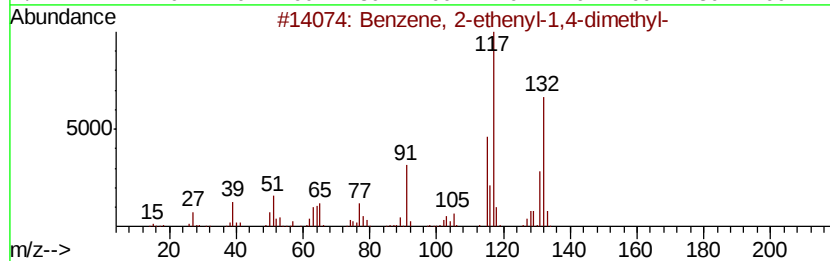
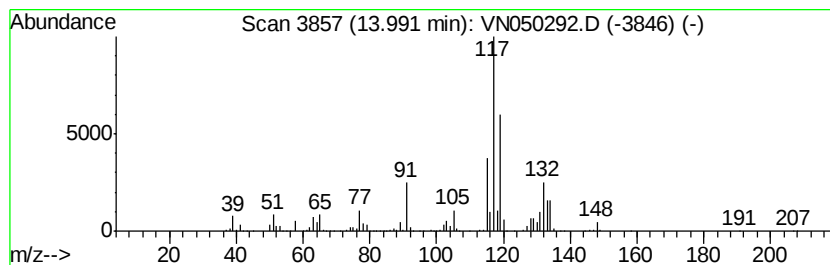
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	25.19 ug/l	1044950	1,4-Dichlorobenzene-d4	13.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	64
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	58
3		1-Phenyl-1-butene	132	C10H12	000824-90-8	58
4		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	58
5		Indan, 1-methyl-	132	C10H12	000767-58-8	58



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ALS Vial : 40 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampled :
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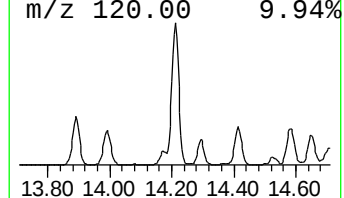
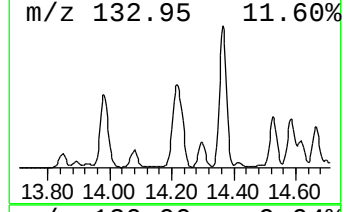
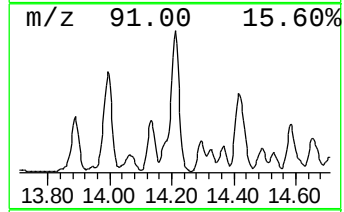
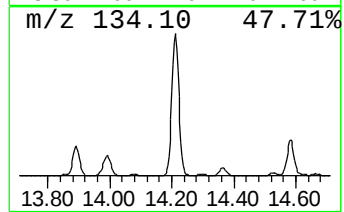
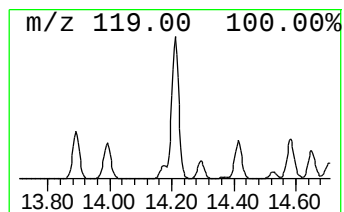
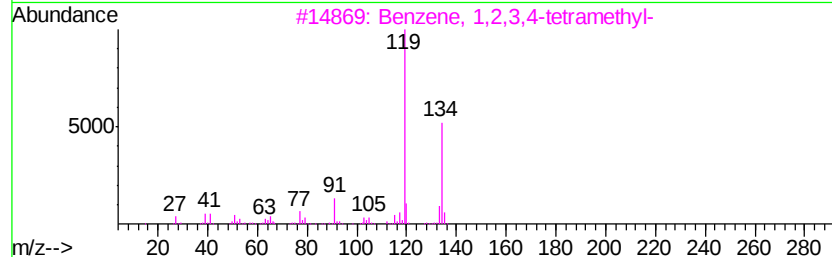
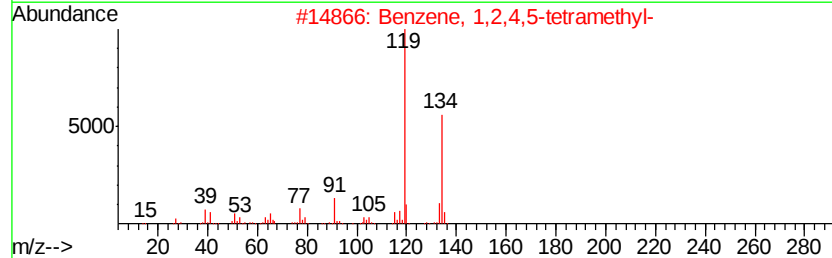
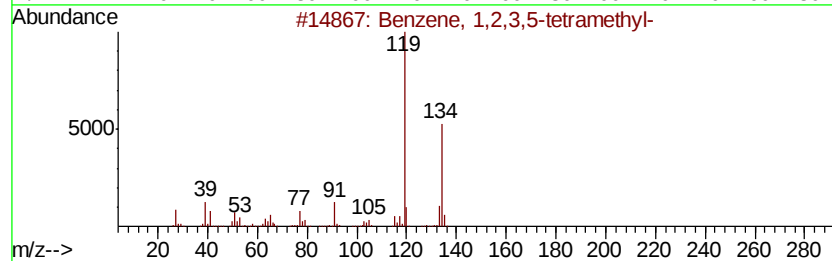
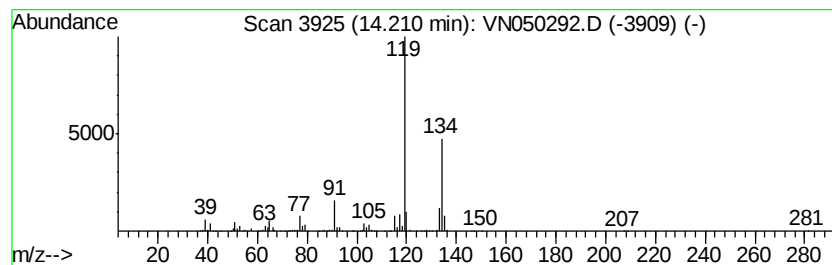
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	40.09 ug/l	1662780	1,4-Dichlorobenzene-d4	13.35

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	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-2,3-dimethyl-	134	C10H14	000933-98-2	94
5	o-Cymene	134	C10H14	000527-84-4	91



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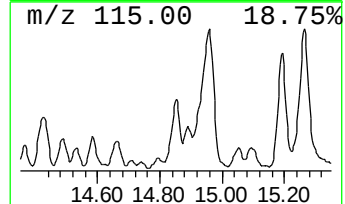
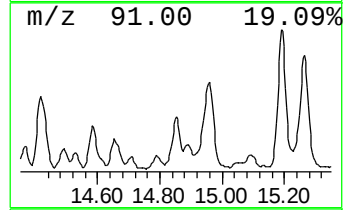
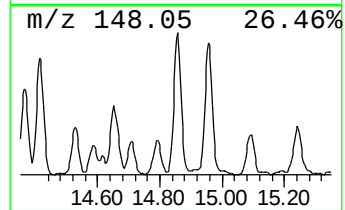
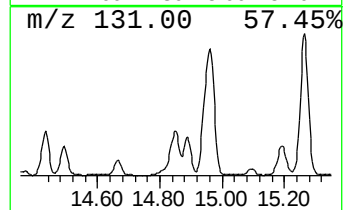
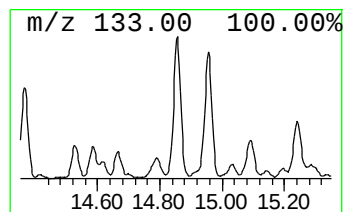
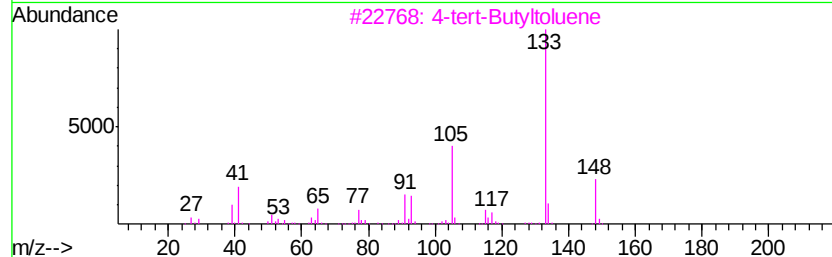
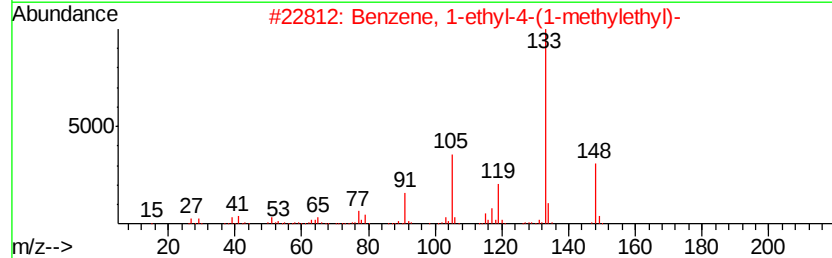
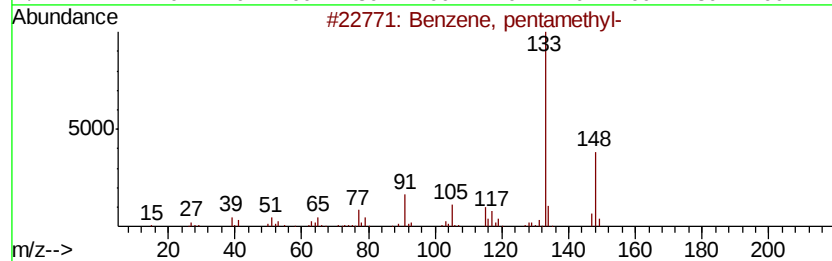
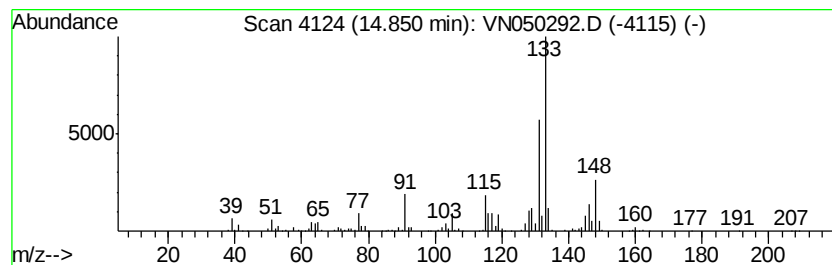
Instrument :
MSVOA_N
ClientSampleId :
EW-A07

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Quant Title : SW846 8260

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

Peak Number 6 Benzene, pentamethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	14.72 ug/l	610734	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, pentamethyl-	148 C11H16	000700-12-9 64
2		Benzene, 1-ethyl-4-(1-methylethyl)-	148 C11H16	004218-48-8 53
3		4-tert-Butyltoluene	148 C11H16	000098-51-1 53
4		Benzene, 1-ethyl-2,4,5-trimethyl-	148 C11H16	017851-27-3 52
5		1,5,6,7-Tetramethylbicyclo[3.2.0...]	148 C11H16	134329-46-7 52



Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN080218\
Data File : VN050292.D
Acq On : 3 Aug 2018 3:48
Operator : MD\SY
Sample : J4304-01
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 40 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
EW-A07

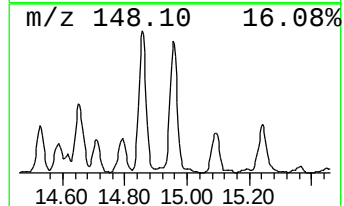
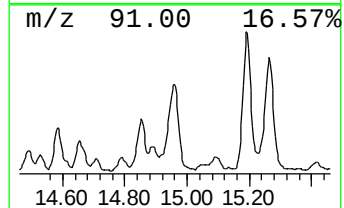
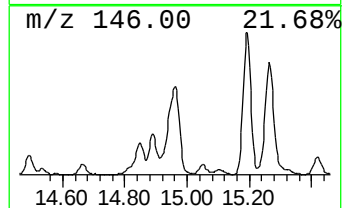
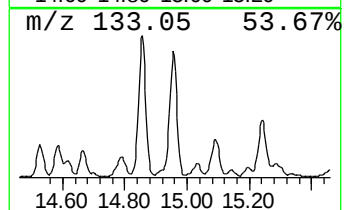
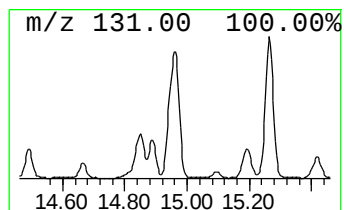
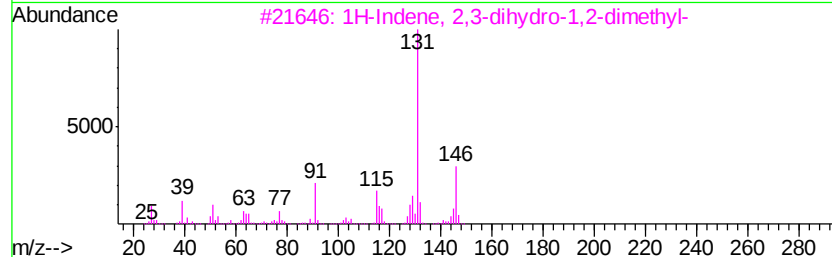
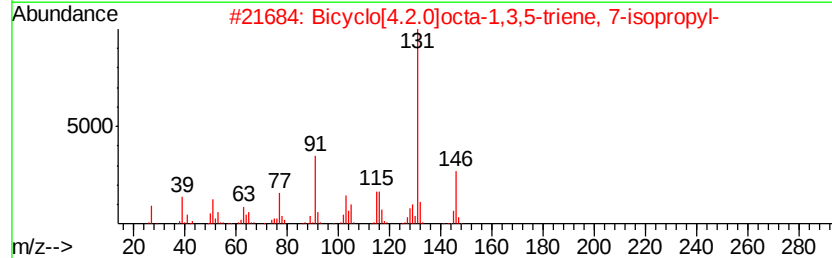
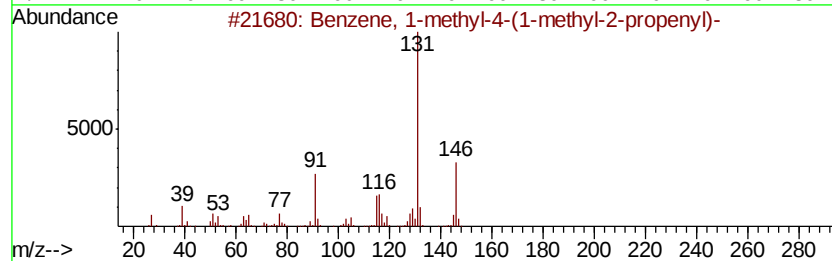
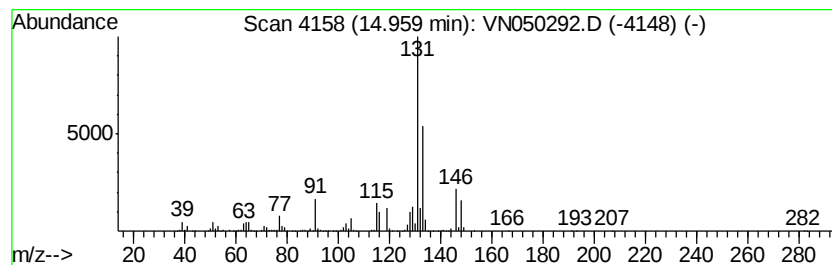
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	36.09 ug/l	1496850	1,4-Dichlorobenzene-d4	13.35

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	70
2		Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	64
3		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	64
4		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	64
5		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN080218\
Data File : VN050292.D
Acq On : 3 Aug 2018 3:48
Operator : MD\SY
Sample : J4304-01
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 40 Sample Multiplier: 1

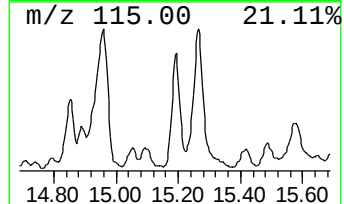
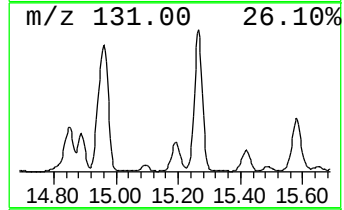
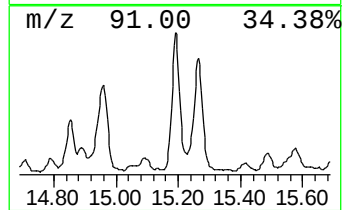
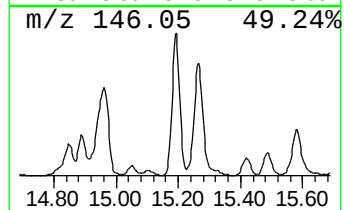
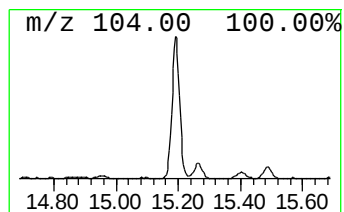
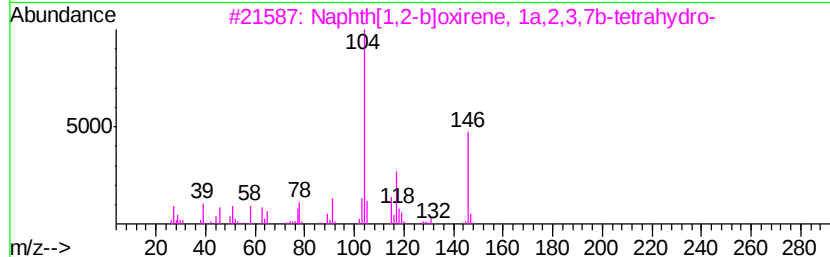
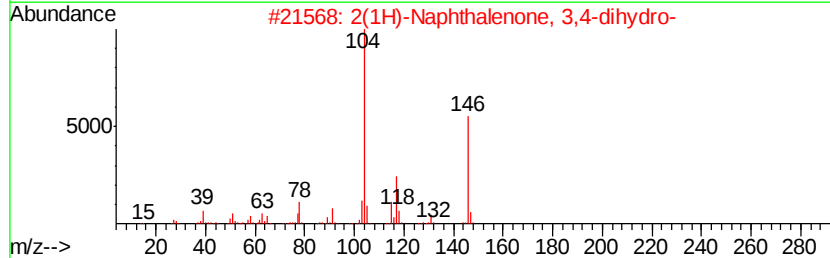
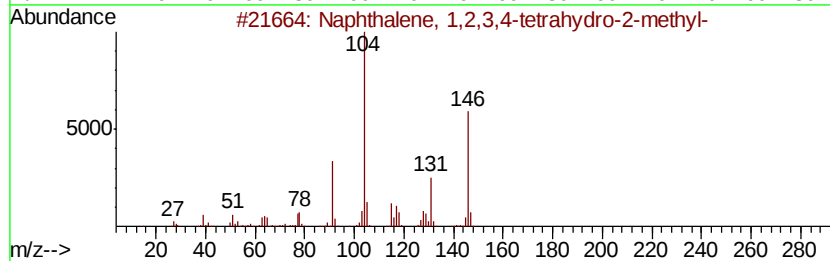
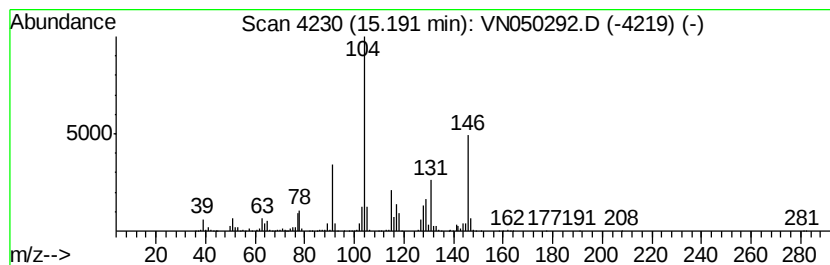
Instrument :
MSVOA_N
ClientSampleId :
EW-A07

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetra... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.19	26.21 ug/l	1087120	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	003877-19-8 94
2		2(1H)-Naphthalenone, 3,4-dihydro-	146 C10H10O	000530-93-8 64
3		Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146 C10H10O	002461-34-9 62
4		2,6-Piperidinedione, 3-phenyl-	189 C11H11N02	014149-34-9 50
5		Naphthalene, 2-ethyl-1,2,3,4-tet...	160 C12H16	032367-54-7 47



Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN080218\
Data File : VN050292.D
Acq On : 3 Aug 2018 3:48
Operator : MD\SY
Sample : J4304-01
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 40 Sample Multiplier: 1

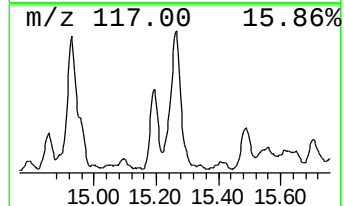
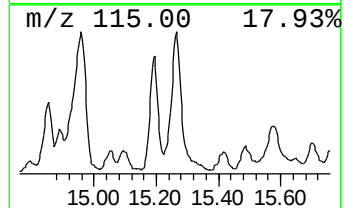
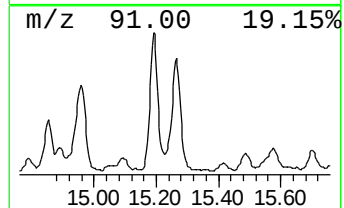
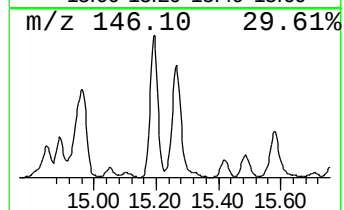
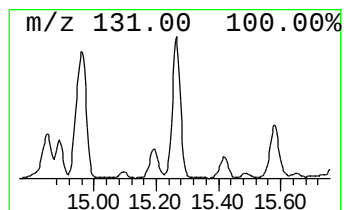
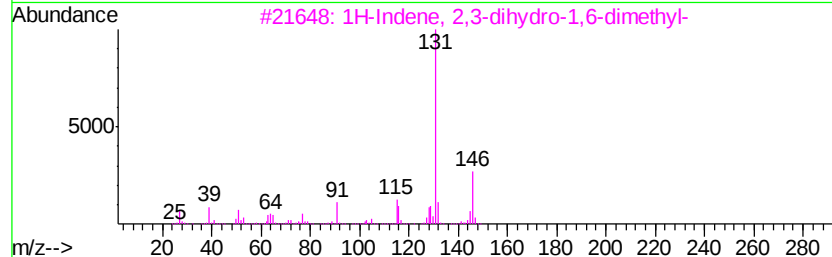
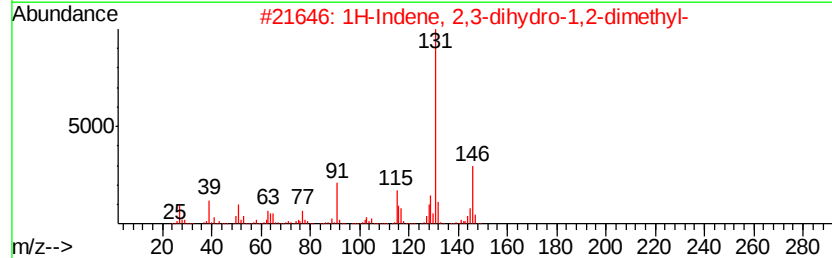
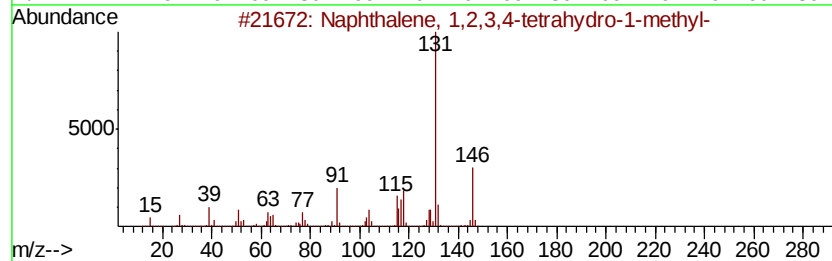
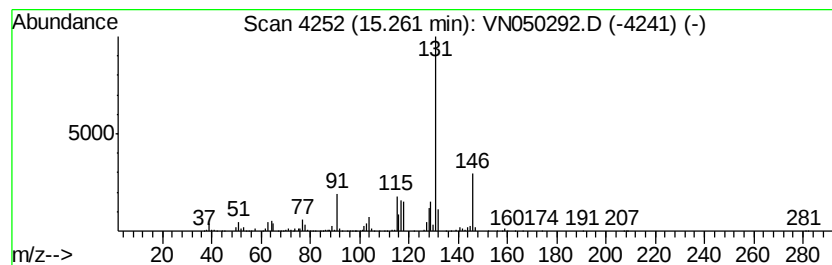
Instrument :
MSVOA_N
ClientSampleId :
EW-A07

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetra... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.26	40.30 ug/l	1671550	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001559-81-5 94
2		1H-Indene, 2,3-dihydro-1,2-dimet...	146 C11H14	017057-82-8 90
3		1H-Indene, 2,3-dihydro-1,6-dimet...	146 C11H14	017059-48-2 90
4		Benzene, 1-methyl-2-(1-methyl-2-...	146 C11H14	097664-19-2 87
5		1H-Indene, 2,3-dihydro-1,3-dimet...	146 C11H14	004175-53-5 87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN080218\
Data File : VN050292.D
Acq On : 3 Aug 2018 3:48
Operator : MD\SY
Sample : J4304-01
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 40 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
EW-A07

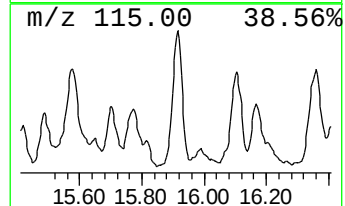
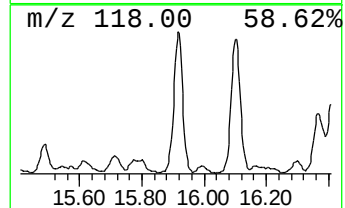
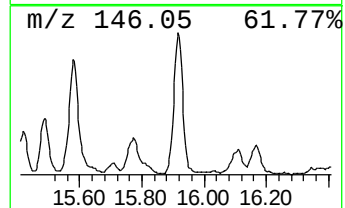
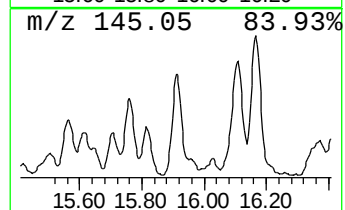
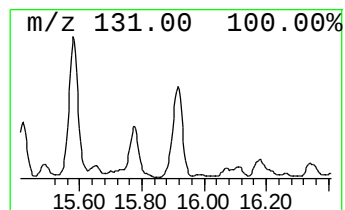
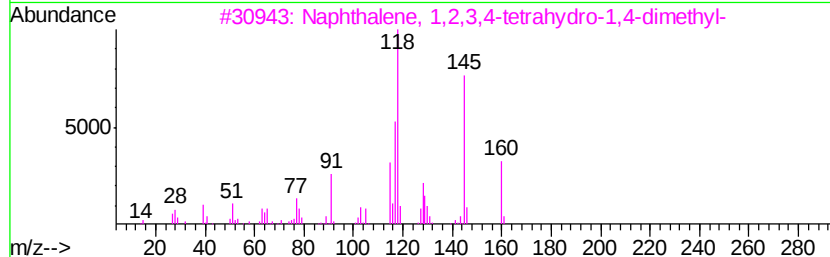
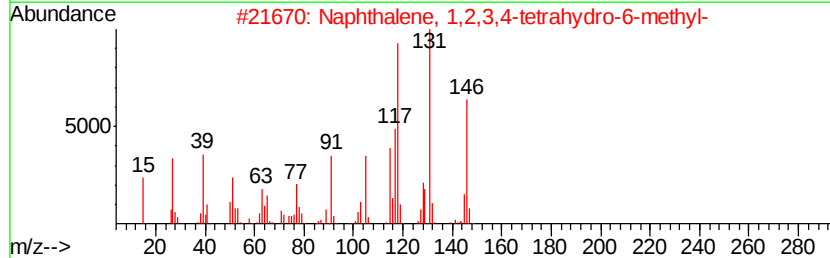
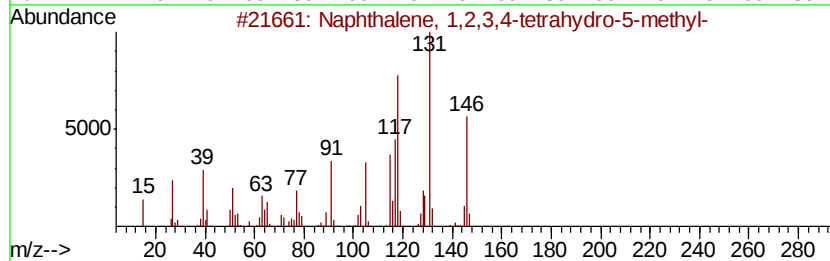
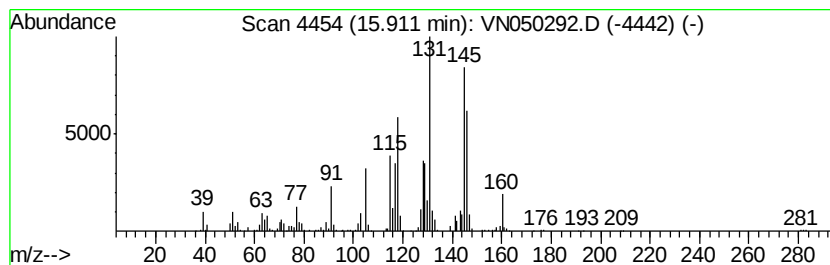
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.91	16.15 ug/l	670039	1,4-Dichlorobenzene-d4	13.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	83
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	60
3		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	55
4		2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	49
5		Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	49



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN080218\
Data File : VN050292.D
Acq On : 3 Aug 2018 3:48
Operator : MD\SY
Sample : J4304-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 40 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
EW-A07

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.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, 1,3-diet...	13.51	20.6	ug/l	855737	4	13.35	2073810	50.0
Benzene, 1,2-diet...	13.68	21.9	ug/l	908903	4	13.35	2073810	50.0
Benzene, 2-etheny...	13.99	25.2	ug/l	1044950	4	13.35	2073810	50.0
Benzene, 1,2,3,5-...	14.21	40.1	ug/l	1662780	4	13.35	2073810	50.0
Benzene, pentamet...	14.85	14.7	ug/l	610734	4	13.35	2073810	50.0
Benzene, 1-methyl...	14.96	36.1	ug/l	1496850	4	13.35	2073810	50.0
Naphthalene, 1,2,...	15.19	26.2	ug/l	1087120	4	13.35	2073810	50.0
Naphthalene, 1,2,...	15.26	40.3	ug/l	1671550	4	13.35	2073810	50.0
Naphthalene, 1,2,...	15.91	16.1	ug/l	670039	4	13.35	2073810	50.0