

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN080218\
 Data File : VN050298.D
 Acq On : 3 Aug 2018 6:17
 Operator : MD\SY
 Sample : J4303-03
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-A19R

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	1.664	4	23	48	rBV	2941089	4405075	40.11%	4.635%
2	7.593	1846	1867	1878	rBV2	546538	1362977	12.41%	1.434%
3	7.667	1879	1890	1909	rVB	803073	1897966	17.28%	1.997%
4	8.030	1982	2003	2030	rBV	548502	1295168	11.79%	1.363%
5	8.249	2036	2071	2074	rBV2	179874	765460	6.97%	0.805%
6	8.587	2161	2176	2204	rBV	1177544	2477906	22.56%	2.607%
7	9.082	2308	2330	2344	rBV3	78708	189452	1.73%	0.199%
8	10.091	2629	2644	2674	rBV	2131468	4002747	36.45%	4.212%
9	10.432	2738	2750	2766	rVB2	72764	141345	1.29%	0.149%
10	10.953	2905	2912	2920	rVB	87709	137509	1.25%	0.145%
11	11.008	2920	2929	2939	rVB2	72574	122372	1.11%	0.129%
12	11.410	3039	3054	3072	rBV	1507894	2700611	24.59%	2.842%
13	11.500	3072	3082	3091	rBV3	83713	143464	1.31%	0.151%
14	11.654	3119	3130	3143	rBV6	90438	203750	1.86%	0.214%
15	11.921	3199	3213	3221	rBV4	65515	147374	1.34%	0.155%
16	11.972	3224	3229	3248	rVB4	49061	110045	1.00%	0.116%
17	12.120	3267	3275	3283	rVB2	99110	150391	1.37%	0.158%
18	12.252	3305	3316	3330	rVB	950354	1567757	14.28%	1.650%
19	12.403	3351	3363	3376	rBV	1393019	2331646	21.23%	2.453%
20	12.474	3376	3385	3396	rVB3	71245	124655	1.14%	0.131%
21	12.593	3399	3422	3435	rBV	1213274	2055301	18.72%	2.163%
22	12.734	3454	3466	3477	rVB	577026	950096	8.65%	1.000%
23	12.963	3527	3537	3541	rBV2	80489	132327	1.20%	0.139%
24	12.995	3542	3547	3553	rVV	94055	137189	1.25%	0.144%
25	13.043	3553	3562	3575	rVB	444398	695927	6.34%	0.732%
26	13.172	3588	3602	3621	rBV2	659561	1310529	11.93%	1.379%
27	13.345	3644	3656	3662	rBV	1481210	2536964	23.10%	2.670%
28	13.445	3678	3687	3697	rBV	314181	479365	4.37%	0.504%
29	13.512	3698	3708	3712	rBV	1198958	1828765	16.65%	1.924%
30	13.545	3712	3718	3726	rVV	2077392	3386046	30.83%	3.563%
31	13.590	3726	3732	3748	rVB3	630073	1321123	12.03%	1.390%
32	13.673	3748	3758	3769	rVB	914192	1505944	13.71%	1.585%
33	13.741	3769	3779	3789	rBV	463614	718233	6.54%	0.756%
34	13.805	3789	3799	3805	rBV	400450	680140	6.19%	0.716%

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35	13.892	3815	3826	3835	rBV	2584211	3846118	35.02%	4.047%
36	13.947	3835	3843	3849	rVV	714562	1078169	9.82%	1.135%
37	13.995	3849	3858	3869	rVB3	2758911	4582525	41.73%	4.822%
38	14.066	3869	3880	3889	rVB8	149396	347118	3.16%	0.365%
39	14.133	3889	3901	3909	rBV2	729118	1178335	10.73%	1.240%
40	14.210	3909	3925	3933	rBV	1609370	2741246	24.96%	2.884%
41	14.249	3933	3937	3945	rVB	646827	865189	7.88%	0.910%
42	14.297	3945	3952	3958	rBV2	284725	399389	3.64%	0.420%
43	14.365	3968	3973	3980	rVB	181003	233334	2.12%	0.246%
44	14.422	3981	3991	4000	rVV4	328358	667312	6.08%	0.702%
45	14.480	4000	4009	4017	rVV2	628686	1018823	9.28%	1.072%
46	14.528	4017	4024	4031	rVV2	348870	489983	4.46%	0.516%
47	14.586	4031	4042	4056	rVV4	5487305	10981824	100.00%	11.556%
48	14.657	4057	4064	4073	rVB4	269257	455559	4.15%	0.479%
49	14.744	4082	4091	4100	rVB	1163964	1732921	15.78%	1.823%
50	14.853	4117	4125	4131	rBV3	572156	956682	8.71%	1.007%
51	14.885	4132	4135	4145	rVV2	486205	661309	6.02%	0.696%
52	14.959	4147	4158	4177	rVB3	1167745	2664529	24.26%	2.804%
53	15.049	4178	4186	4192	rBV	158623	263957	2.40%	0.278%
54	15.091	4192	4199	4212	rVB3	303376	457868	4.17%	0.482%
55	15.191	4220	4230	4239	rBV	1088429	1833674	16.70%	1.929%
56	15.265	4239	4253	4287	rVB3	1123141	3349603	30.50%	3.525%
57	15.419	4289	4301	4312	rBV3	214655	442666	4.03%	0.466%
58	15.487	4313	4322	4333	rBV5	175083	316106	2.88%	0.333%
59	15.580	4335	4351	4353	rVV4	443146	783364	7.13%	0.824%
60	15.615	4354	4362	4378	rVV	1675362	3168260	28.85%	3.334%
61	15.702	4380	4389	4399	rVV2	234704	460584	4.19%	0.485%
62	15.773	4401	4411	4438	rVB7	244250	807233	7.35%	0.849%
63	15.918	4441	4456	4472	rBV2	1640730	3267269	29.75%	3.438%
64	16.104	4499	4514	4525	rBV2	402532	909359	8.28%	0.957%
65	16.168	4525	4534	4559	rVB7	251793	766264	6.98%	0.806%
66	16.358	4578	4593	4604	rBV6	390545	981990	8.94%	1.033%
67	16.499	4628	4637	4652	rVB6	128187	310534	2.83%	0.327%

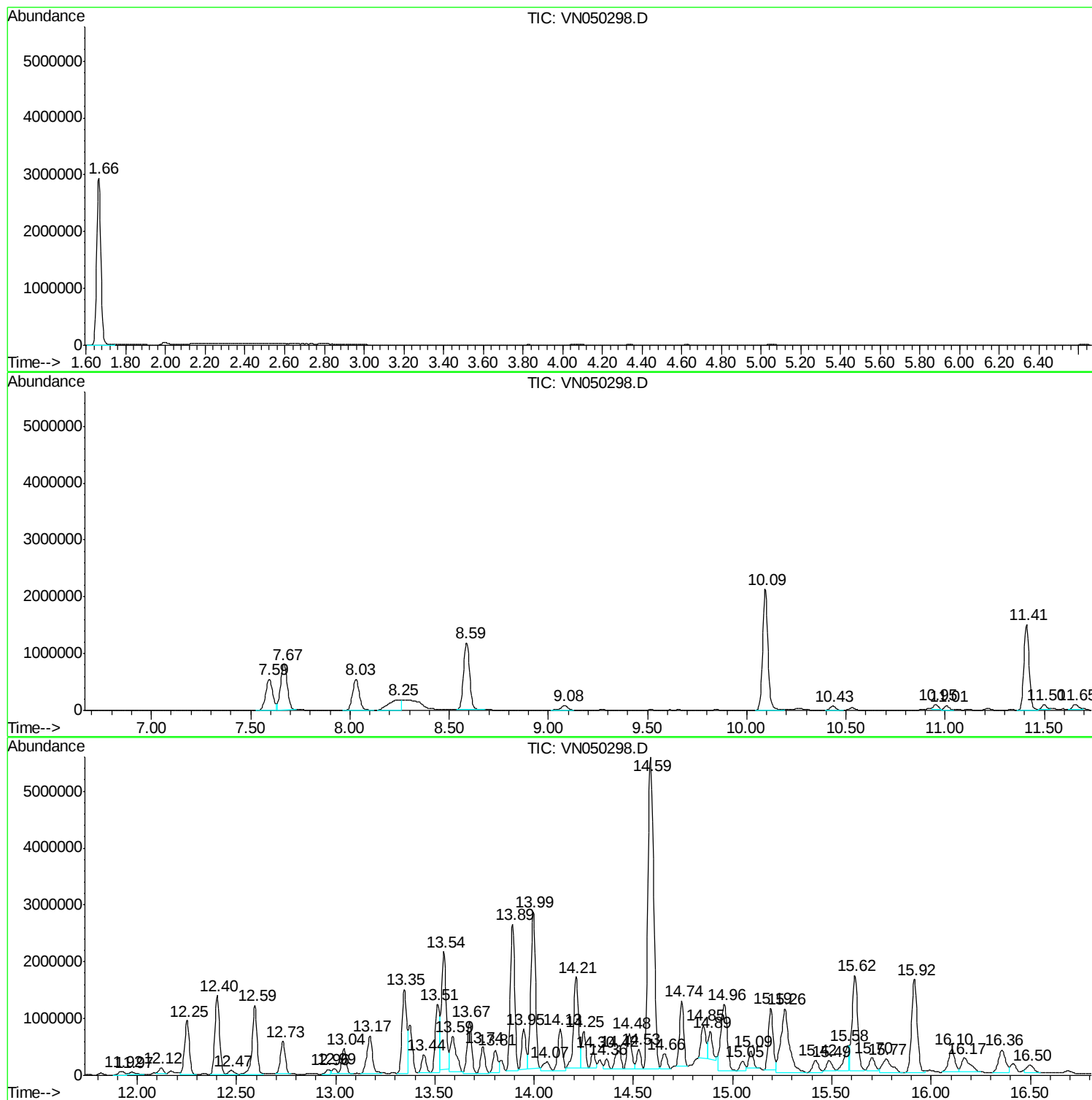
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Quant Title : SW846 8260

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



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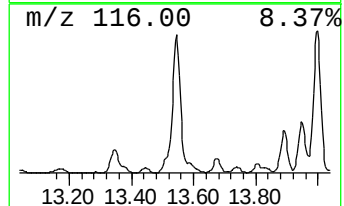
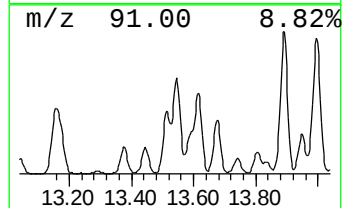
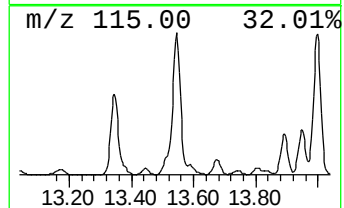
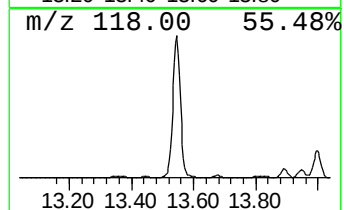
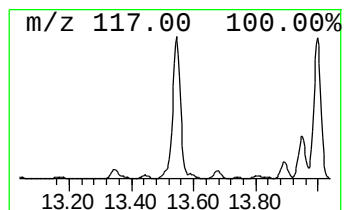
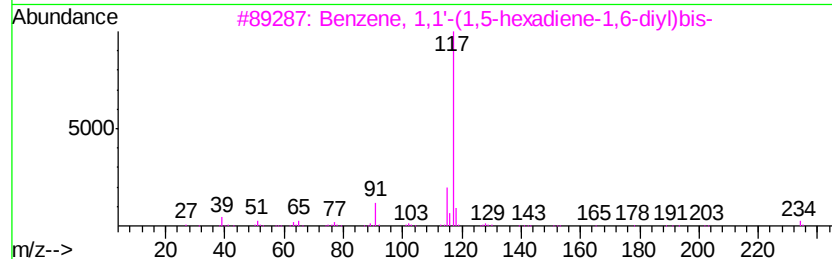
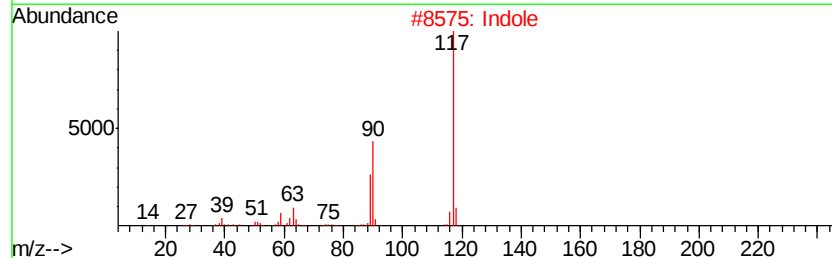
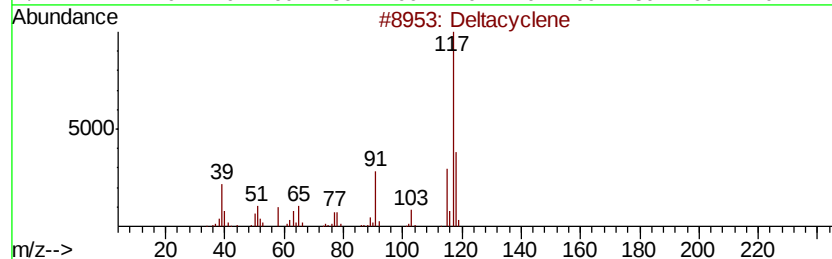
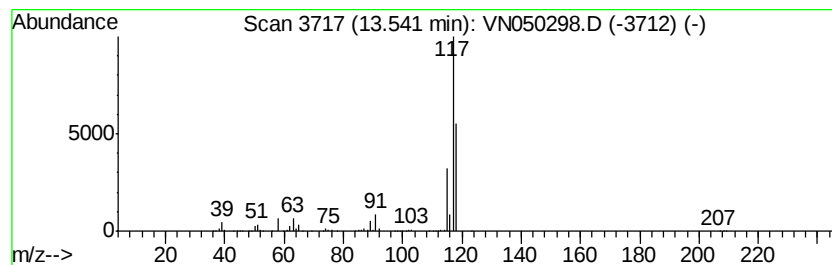
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Quant Title : SW846 8260

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

Peak Number 2 Deltacyclene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.54	66.73 ug/l	3386050	1,4-Dichlorobenzene-d4	13.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Deltacyclene	118	C9H10	007785-10-6	72
2	Indole	117	C8H7N	000120-72-9	43
3	Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	42
4	Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	38
5	Benzyl nitrile	117	C8H7N	000140-29-4	38



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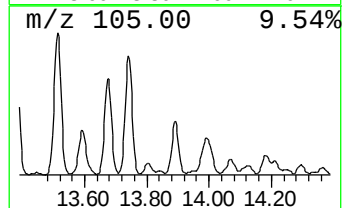
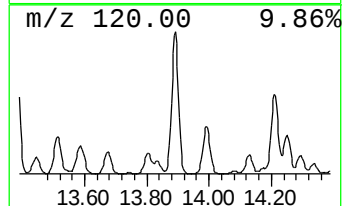
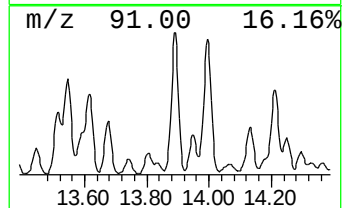
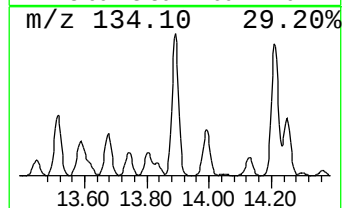
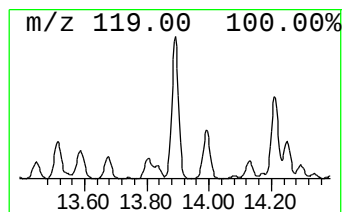
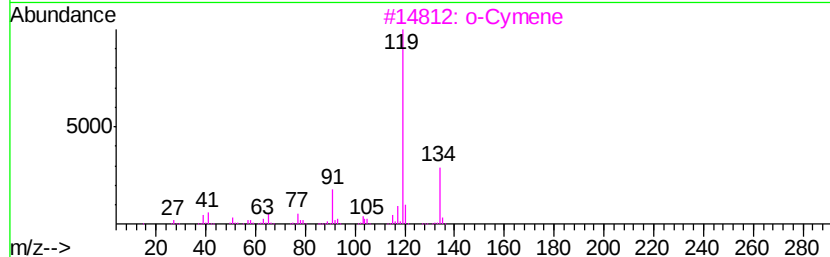
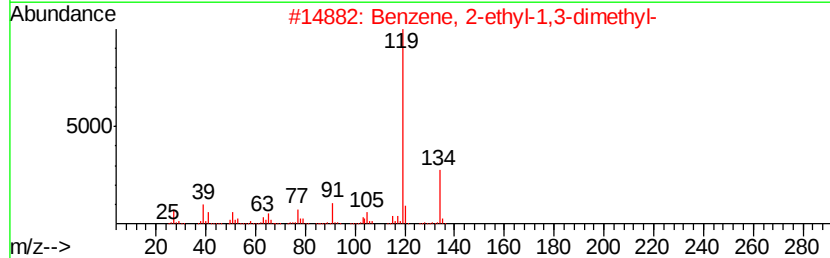
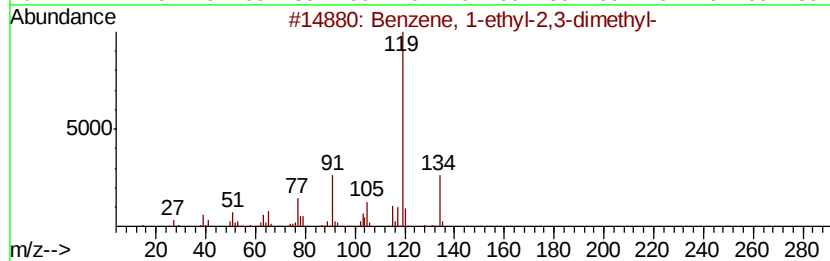
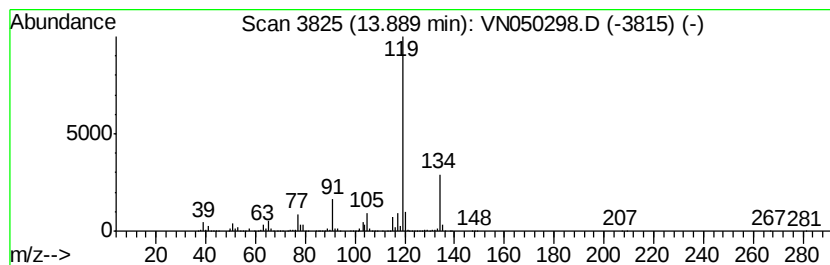
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	75.80 ug/l	3846120	1,4-Dichlorobenzene-d4	13.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3		o-Cymene	134	C10H14	000527-84-4	95
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



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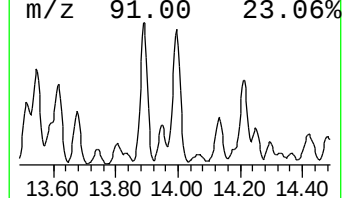
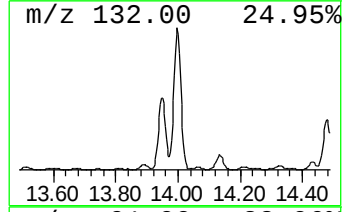
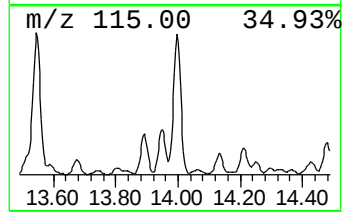
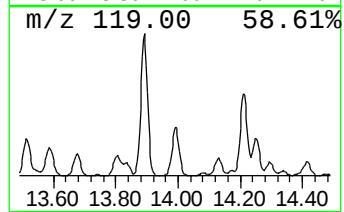
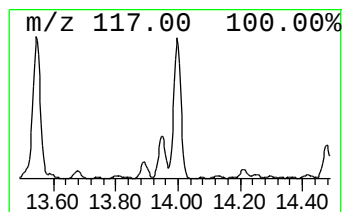
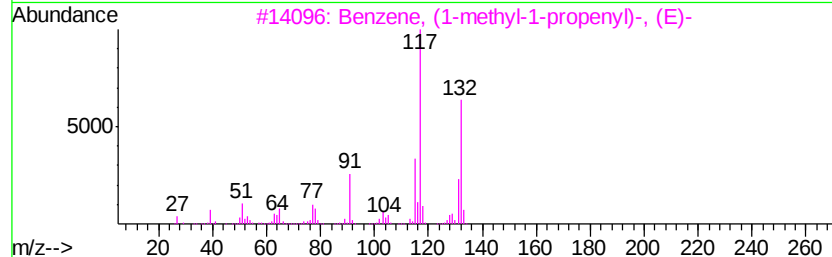
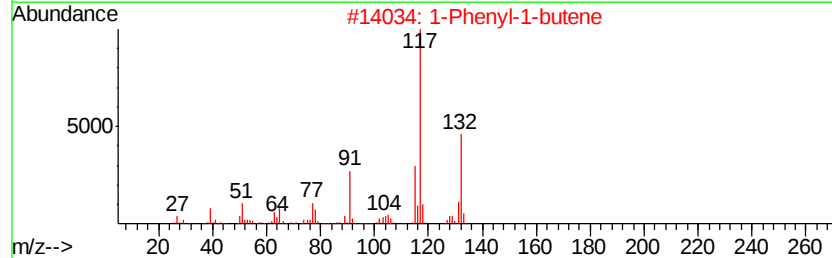
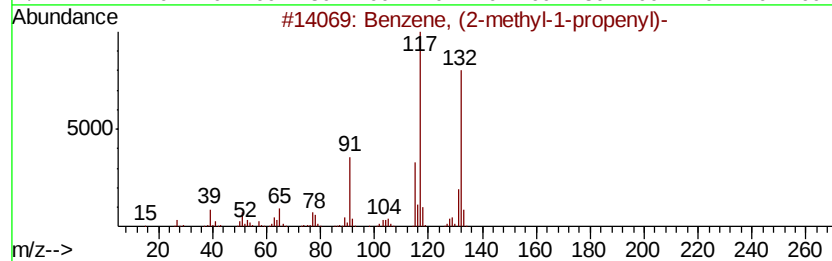
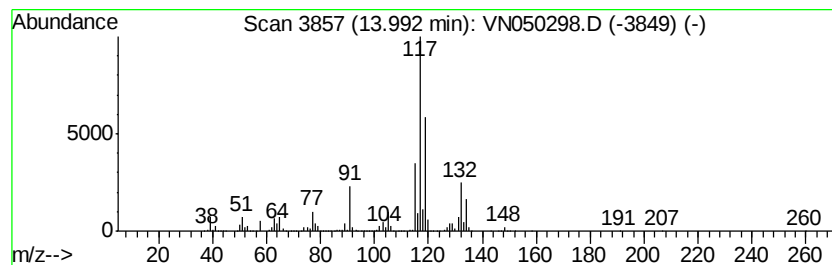
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Peak Number 4 Benzene, (2-methyl-1-propen... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	90.32 ug/l	4582530	1,4-Dichlorobenzene-d4	13.35

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	70
2		1-Phenyl-1-butene	132	C10H12	000824-90-8	70
3		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	70
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	64
5		Indan, 1-methyl-	132	C10H12	000767-58-8	64



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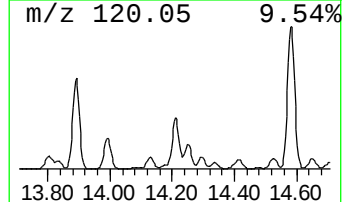
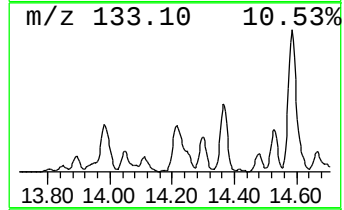
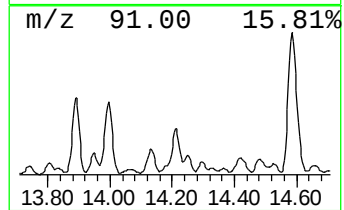
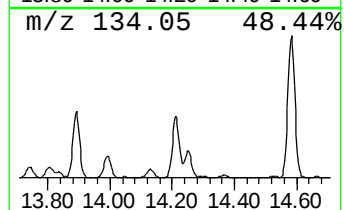
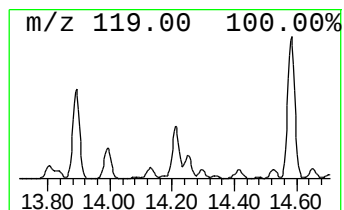
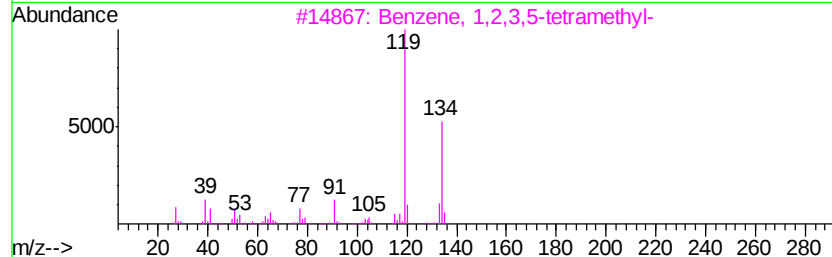
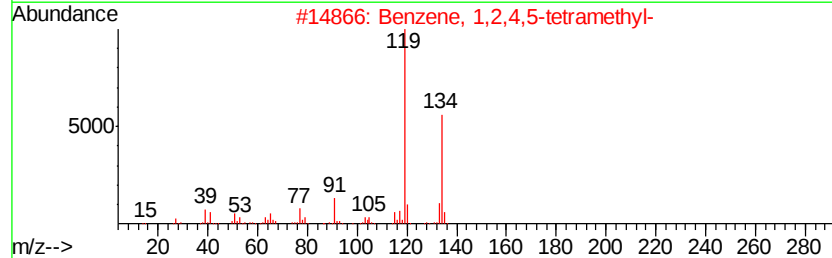
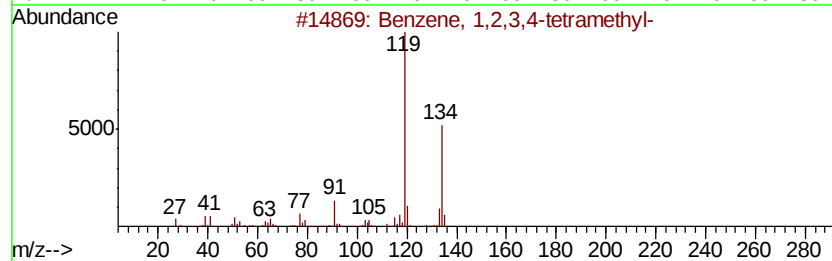
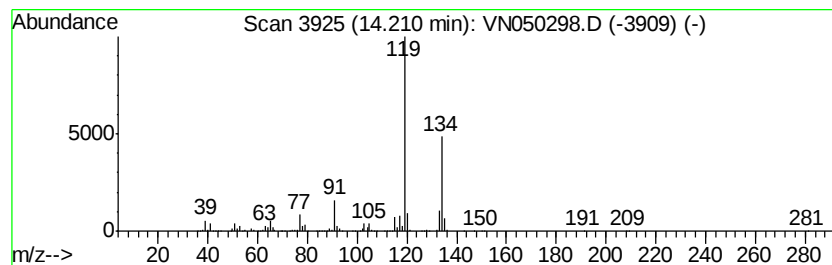
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Peak Number 5 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	54.03 ug/l	2741250	1,4-Dichlorobenzene-d4	13.35

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
4	o-Cymene	134	C10H14	000527-84-4	94
5	p-Cymene	134	C10H14	000099-87-6	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN080218\
Data File : VN050298.D
Acq On : 3 Aug 2018 6:17
Operator : MD\SY
Sample : J4303-03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-A19R

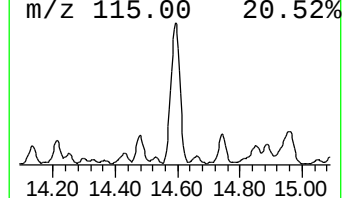
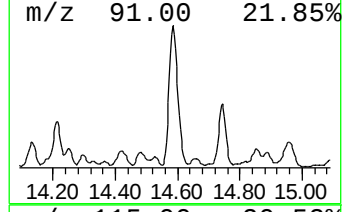
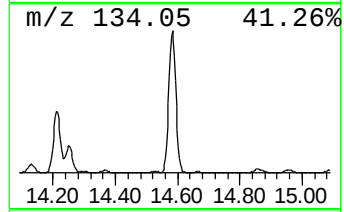
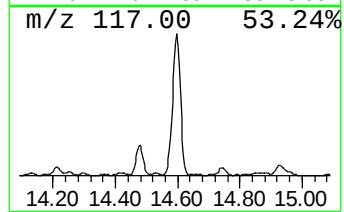
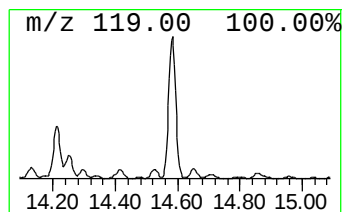
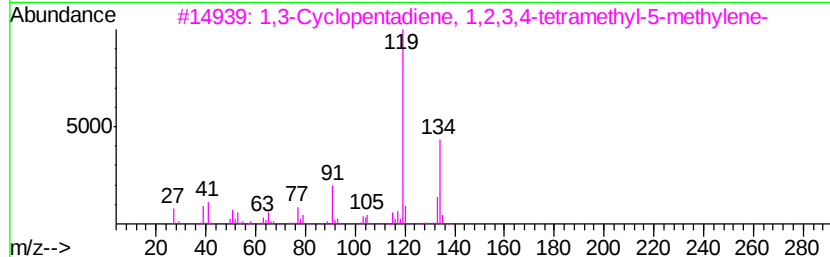
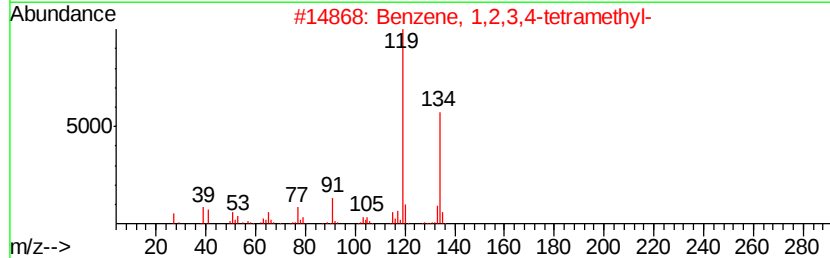
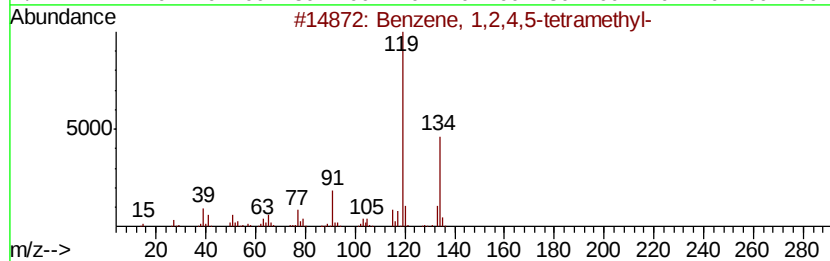
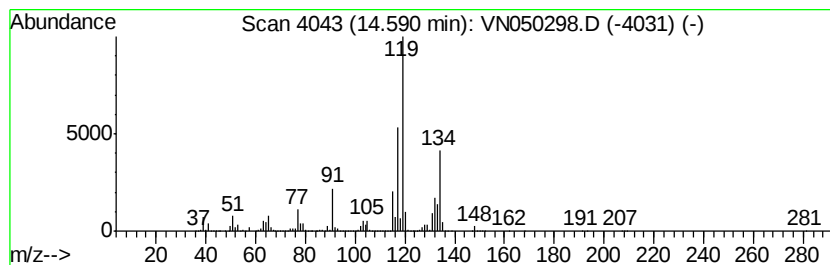
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Quant Title : SW846 8260

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.
14.59	216.44 ug/l	10981800	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	93
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	87
3		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	87
4		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	81
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN080218\
Data File : VN050298.D
Acq On : 3 Aug 2018 6:17
Operator : MD\SY
Sample : J4303-03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-A19R

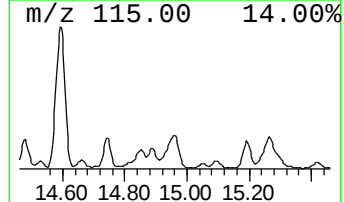
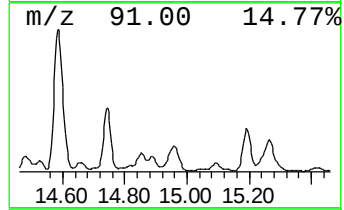
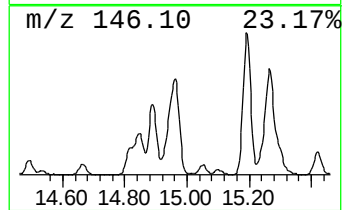
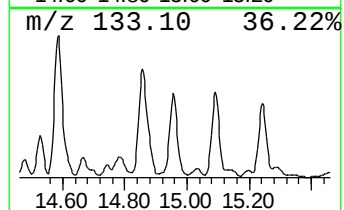
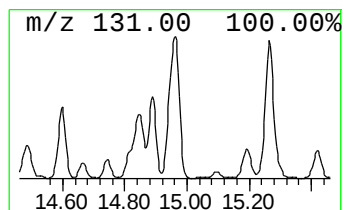
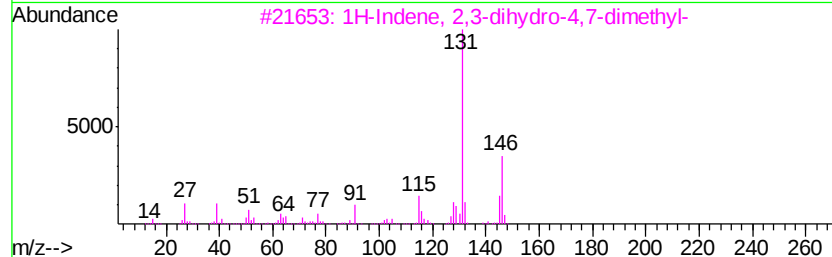
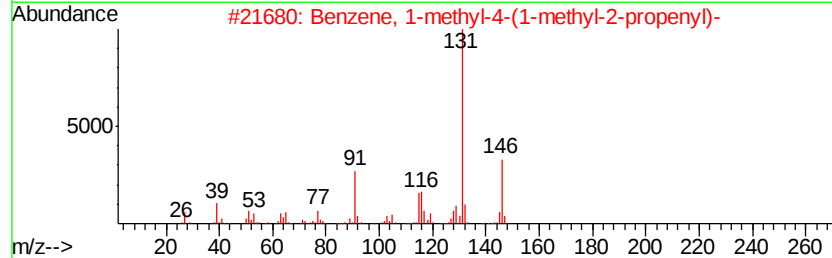
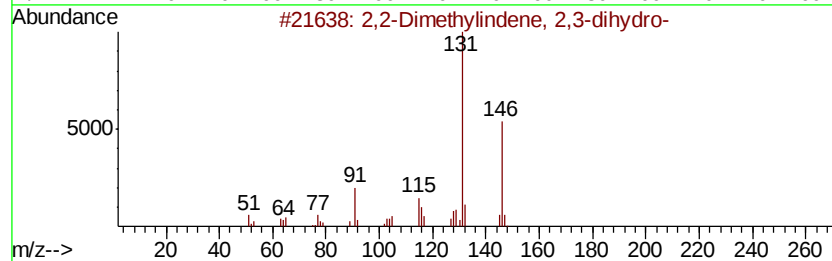
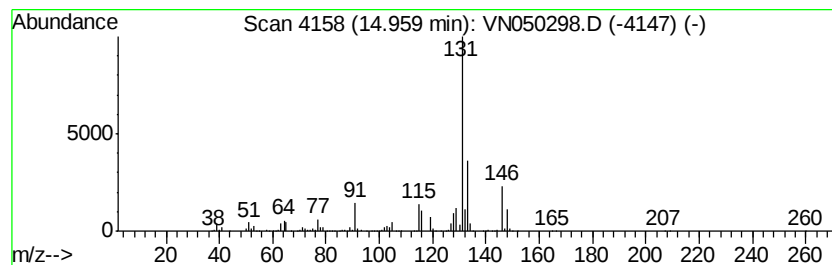
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Quant Title : SW846 8260

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

Peak Number 7 2,2-Dimethylindene, 2,3-dih... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	52.51 ug/l	2664530	1,4-Dichlorobenzene-d4	13.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	81
2		Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	81
3		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	81
4		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	81
5		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN080218\
Data File : VN050298.D
Acq On : 3 Aug 2018 6:17
Operator : MD\SY
Sample : J4303-03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-A19R

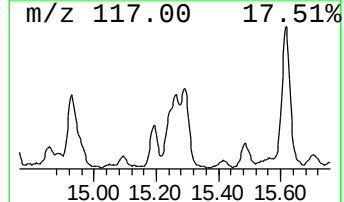
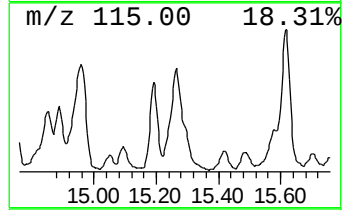
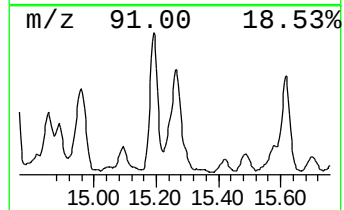
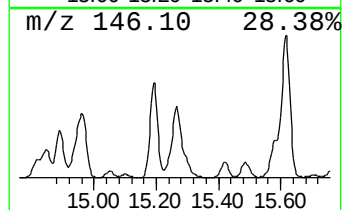
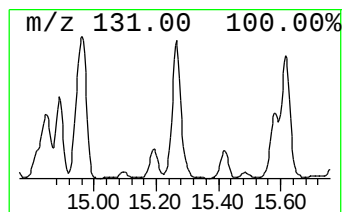
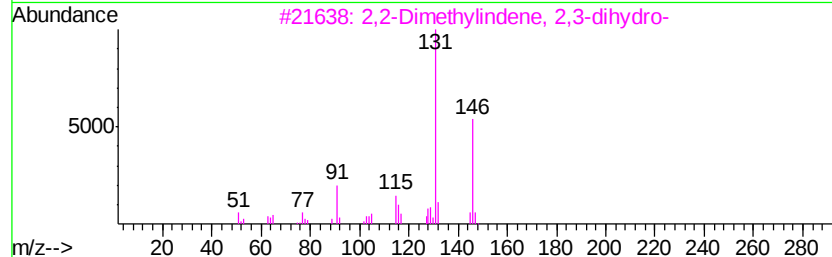
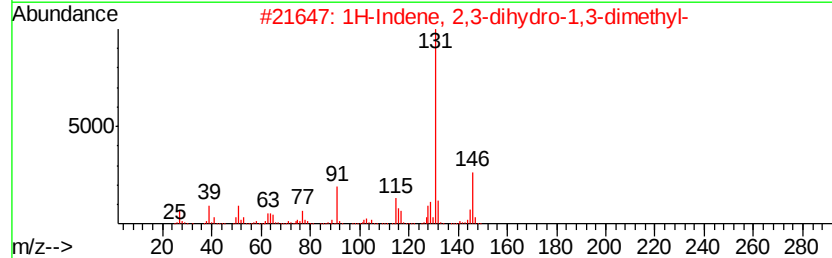
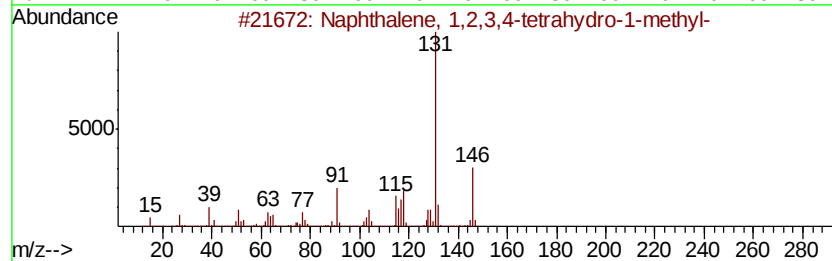
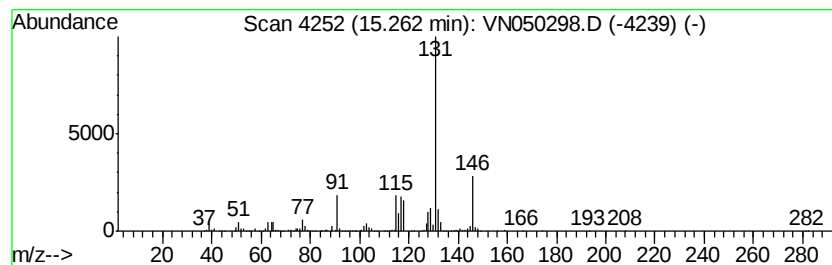
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Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.26	66.02 ug/l	3349600	1,4-Dichlorobenzene-d4	13.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	94
2		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90
3		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	87
4		Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	87
5		1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN080218\
Data File : VN050298.D
Acq On : 3 Aug 2018 6:17
Operator : MD\SY
Sample : J4303-03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 46 Sample Multiplier: 1

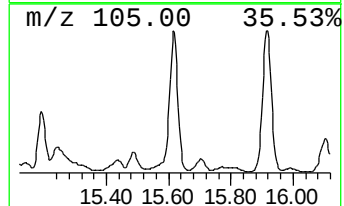
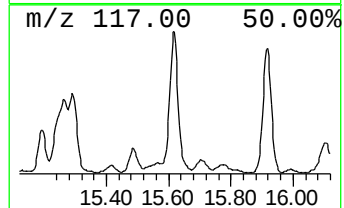
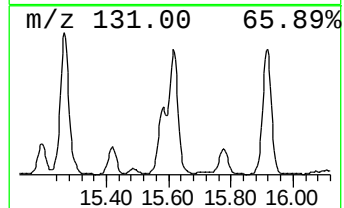
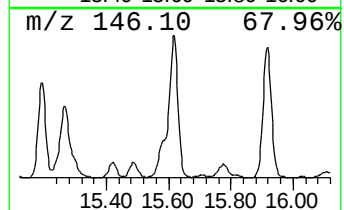
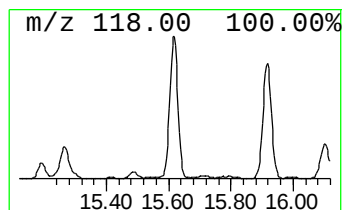
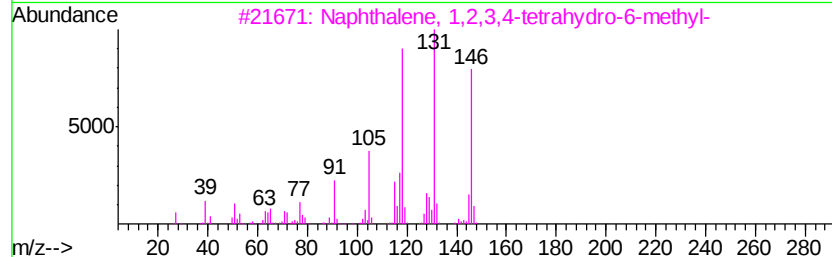
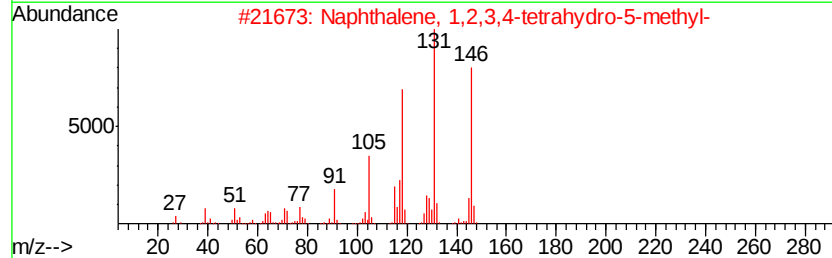
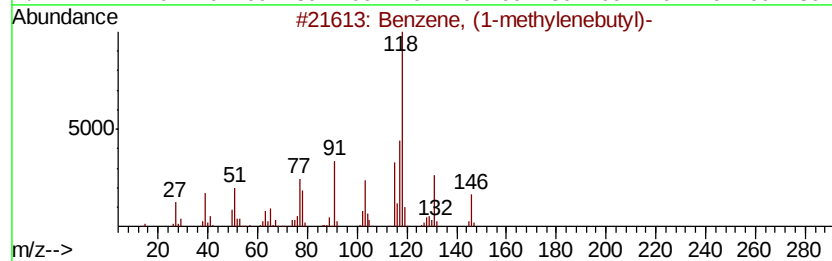
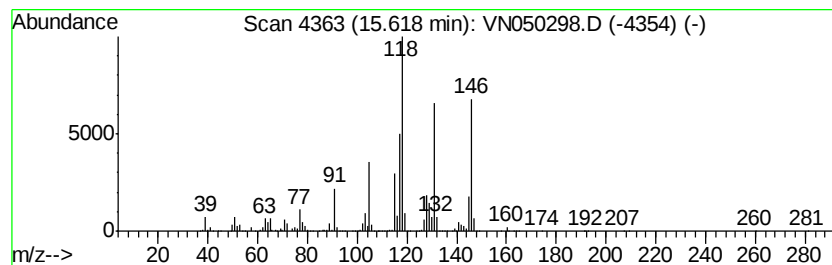
Instrument :
MSVOA_N
ClientSampleId :
MW-A19R

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 9 Benzene, (1-methylenebutyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.62	62.44 ug/l	3168260	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (1-methylenebutyl)-	146 C11H14	005676-32-4 83
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 70
3		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 70
4		2,3,4,5,6,7-Hexahydro-1H-cyclope...	146 C11H14	1000189-31-0 49
5		1(2H)-Naphthalenone, 3,4-dihydro-	146 C10H10O	000529-34-0 46



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN080218\
Data File : VN050298.D
Acq On : 3 Aug 2018 6:17
Operator : MD\SY
Sample : J4303-03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 46 Sample Multiplier: 1

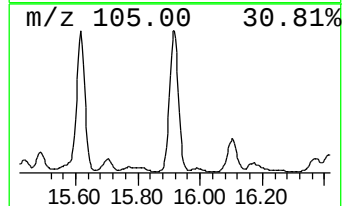
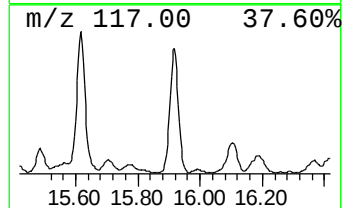
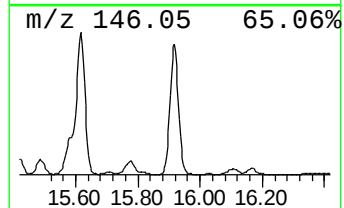
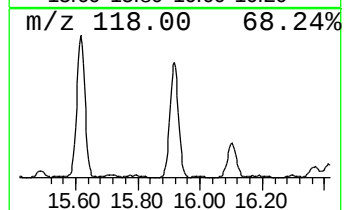
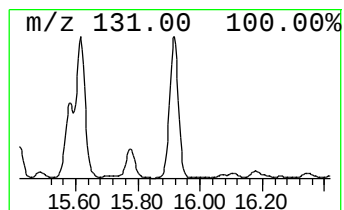
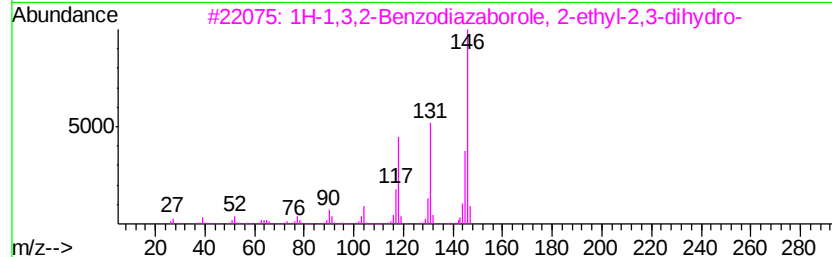
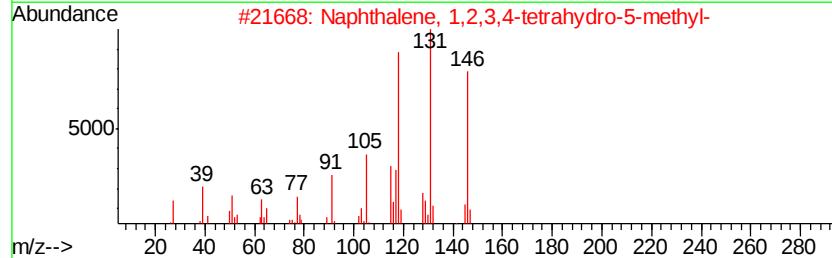
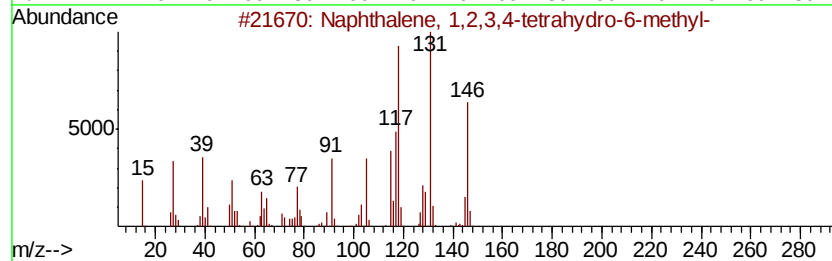
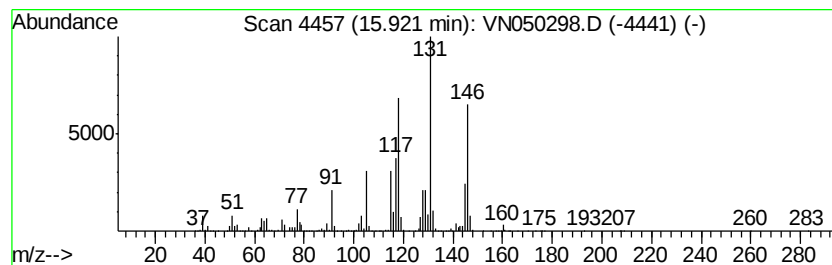
Instrument :
MSVOA_N
ClientSampleId :
MW-A19R

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-tetra... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.92	64.39 ug/l	3267270	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 95
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 95
3		1H-1,3,2-Benzodiazaborole, 2-eth...	146 C8H11BN2	074663-81-3 90
4		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 64
5		1H-Inden-1-one, 2,3-dihydro-2-me...	146 C10H10O	017496-14-9 62



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN080218\
Data File : VN050298.D
Acq On : 3 Aug 2018 6:17
Operator : MD\SY
Sample : J4303-03
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-A19R

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
Deltacyclene	13.54	66.7	ug/l	3386050	4	13.35	2536960	50.0
Benzene, 1-ethyl-...	13.89	75.8	ug/l	3846120	4	13.35	2536960	50.0
Benzene, (2-methy...	13.99	90.3	ug/l	4582530	4	13.35	2536960	50.0
Benzene, 1,2,3,4-...	14.21	54.0	ug/l	2741250	4	13.35	2536960	50.0
Benzene, 1,2,4,5-...	14.59	216.4	ug/l	10981800	4	13.35	2536960	50.0
2,2-Dimethylinden...	14.96	52.5	ug/l	2664530	4	13.35	2536960	50.0
Naphthalene, 1,2,...	15.26	66.0	ug/l	3349600	4	13.35	2536960	50.0
Benzene, (1-methy...	15.62	62.4	ug/l	3168260	4	13.35	2536960	50.0
Naphthalene, 1,2,...	15.92	64.4	ug/l	3267270	4	13.35	2536960	50.0