

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091119\
 Data File : VN057887.D
 Acq On : 11 Sep 2019 00:23
 Operator : JC/SP
 Sample : K4818-07
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 T11-MW-13-8.0-091019

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\82N091019W.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	2.426	195	203	212	rVB3	24801	39678	1.68%	0.223%
2	3.805	617	632	644	rBV2	12470	31548	1.34%	0.177%
3	4.050	691	708	731	rBV	51855	153834	6.52%	0.864%
4	5.021	984	1010	1036	rBV3	193827	710156	30.12%	3.988%
5	7.577	1786	1805	1816	rBV	288794	717982	30.45%	4.032%
6	7.651	1816	1828	1848	rVB	365021	870907	36.94%	4.891%
7	8.014	1921	1941	1964	rBV2	386168	919267	38.99%	5.163%
8	8.577	2100	2116	2142	rVB	552276	1163274	49.34%	6.533%
9	9.069	2253	2269	2285	rVB6	14283	35343	1.50%	0.198%
10	10.088	2571	2586	2606	rBV	1278968	2357805	100.00%	13.242%
11	10.603	2736	2746	2757	rVB2	17839	32663	1.39%	0.183%
12	10.792	2798	2805	2815	rVB	15491	24941	1.06%	0.140%
13	11.406	2981	2996	3013	rBV	816566	1442440	61.18%	8.101%
14	12.120	3200	3218	3227	rBV8	14895	34050	1.44%	0.191%
15	12.252	3248	3259	3270	rVB2	112599	188569	8.00%	1.059%
16	12.403	3293	3306	3321	rBV	880746	1489581	63.18%	8.366%
17	12.596	3350	3366	3378	rVB	120717	216748	9.19%	1.217%
18	12.998	3485	3491	3498	rVV2	18035	29261	1.24%	0.164%
19	13.175	3533	3546	3555	rBV3	60489	122805	5.21%	0.690%
20	13.349	3588	3600	3621	rBV	962382	1626873	69.00%	9.137%
21	13.448	3621	3631	3642	rVV2	37621	60953	2.59%	0.342%
22	13.548	3643	3662	3672	rVV2	373402	760010	32.23%	4.268%
23	13.599	3673	3678	3691	rVV5	32667	67137	2.85%	0.377%
24	13.676	3692	3702	3713	rVB3	74341	133012	5.64%	0.747%
25	13.895	3759	3770	3779	rBV	230603	363095	15.40%	2.039%
26	13.953	3779	3788	3794	rBV	83199	133846	5.68%	0.752%
27	14.001	3794	3803	3814	rVB2	323242	528346	22.41%	2.967%
28	14.136	3837	3845	3853	rVB	48284	71139	3.02%	0.400%
29	14.217	3854	3870	3888	rVB	194030	355867	15.09%	1.999%
30	14.435	3927	3938	3945	rVV5	31076	62248	2.64%	0.350%
31	14.484	3945	3953	3963	rVV2	66581	115817	4.91%	0.650%
32	14.596	3975	3988	4001	rVV3	553388	1147231	48.66%	6.443%
33	14.664	4001	4009	4020	rVB5	24367	41929	1.78%	0.235%
34	14.747	4026	4035	4046	rVB	126916	193117	8.19%	1.085%

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35	14.857	4062	4069	4075	rBV4	42298	64352	2.73%	0.361%
36	14.895	4075	4081	4089	rVV3	43922	66037	2.80%	0.371%
37	14.963	4091	4102	4111	rVV2	95447	196618	8.34%	1.104%
38	15.091	4136	4142	4147	rBV4	22691	32664	1.39%	0.183%
39	15.130	4147	4154	4163	rVV	77077	125219	5.31%	0.703%
40	15.184	4164	4171	4178	rVV2	43440	67654	2.87%	0.380%
41	15.252	4179	4192	4215	rVB5	62593	212003	8.99%	1.191%
42	15.390	4226	4235	4245	rBV5	22098	38293	1.62%	0.215%
43	15.451	4246	4254	4264	rBV5	22901	39345	1.67%	0.221%
44	15.535	4270	4280	4284	rBV3	40402	69772	2.96%	0.392%
45	15.641	4305	4313	4321	rBV3	22648	38214	1.62%	0.215%
46	15.705	4325	4333	4344	rVB4	21104	41187	1.75%	0.231%
47	15.831	4362	4372	4384	rBV3	103377	188032	7.97%	1.056%
48	16.049	4431	4440	4461	rVB2	59192	136637	5.80%	0.767%
49	16.213	4479	4491	4503	rBV2	126044	248043	10.52%	1.393%

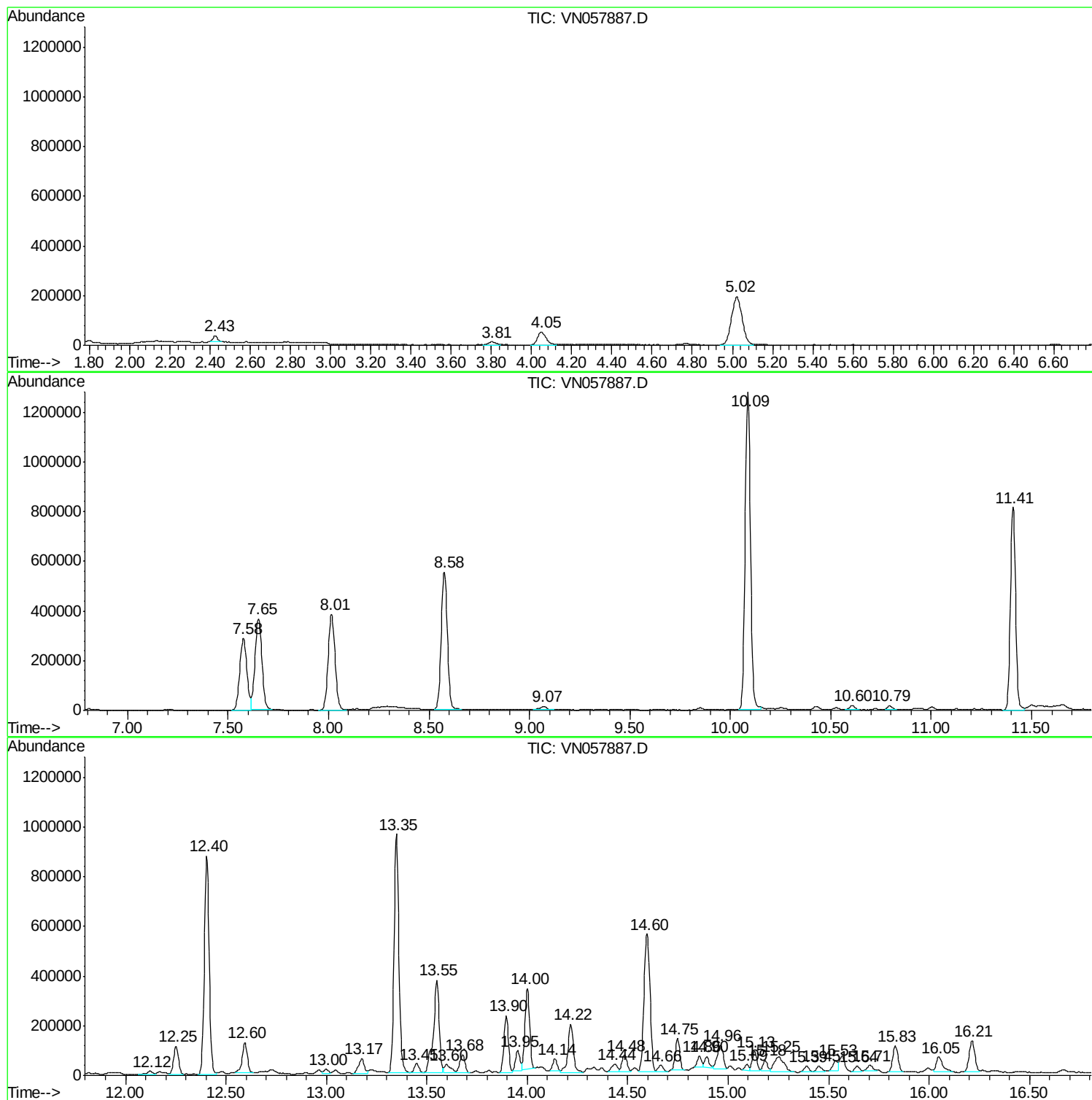
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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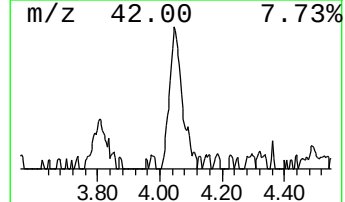
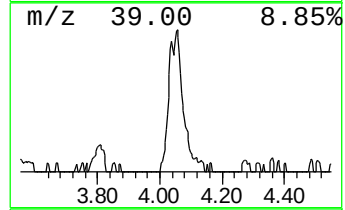
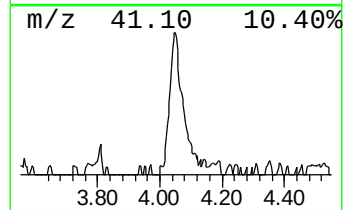
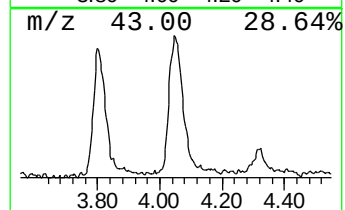
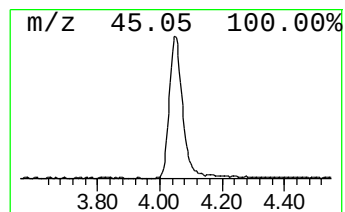
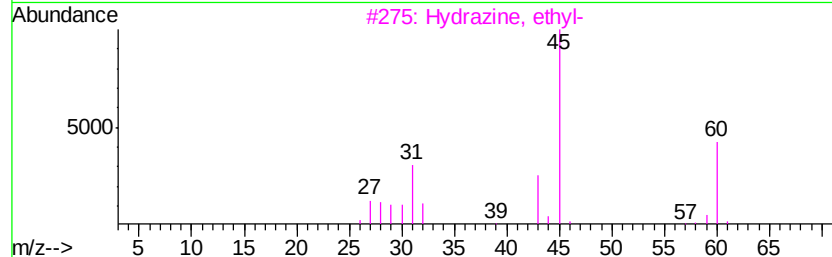
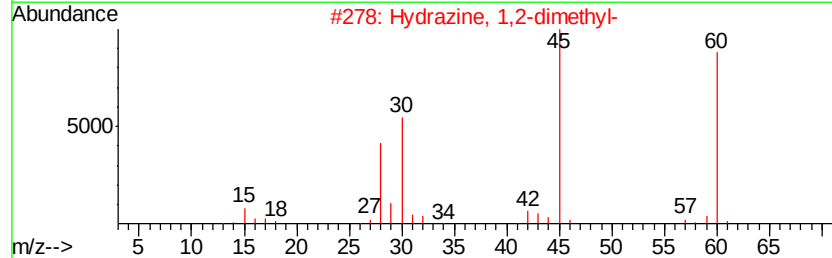
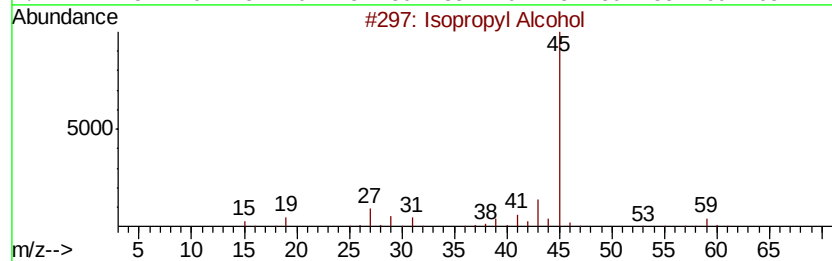
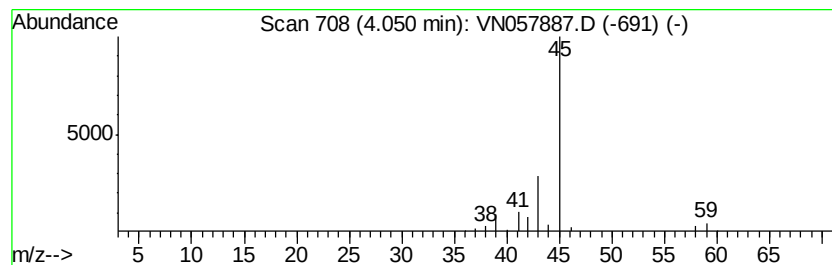
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Quant Title : SW846 8260

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

Peak Number 1 Isopropyl Alcohol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.05	8.83 ug/l	153834	Pentafluorobenzene	7.65

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Isopropyl Alcohol	60	C3H8O	000067-63-0	86
2		Hydrazine, 1,2-dimethyl-	60	C2H8N2	000540-73-8	9
3		Hydrazine, ethyl-	60	C2H8N2	000624-80-6	9
4		Ethylamine	45	C2H7N	000075-04-7	5
5		Oxirane, (methoxymethyl)-	88	C4H8O2	000930-37-0	4



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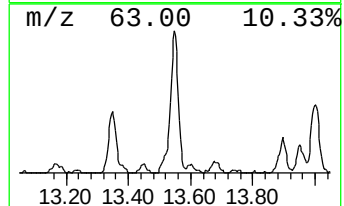
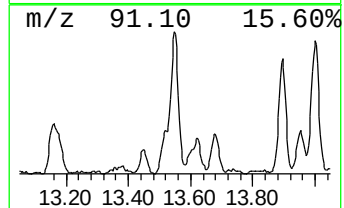
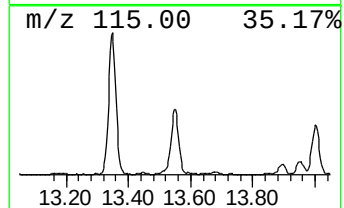
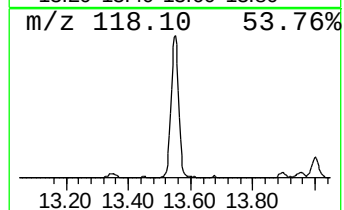
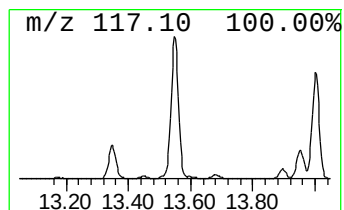
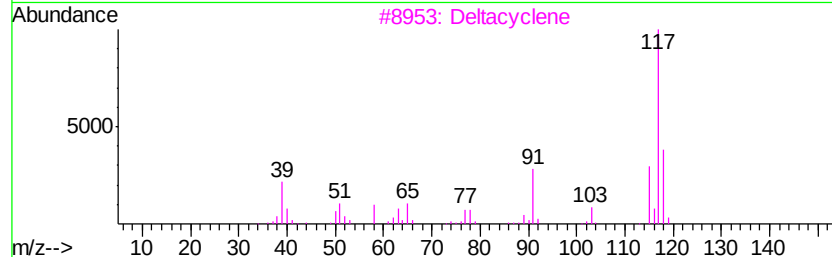
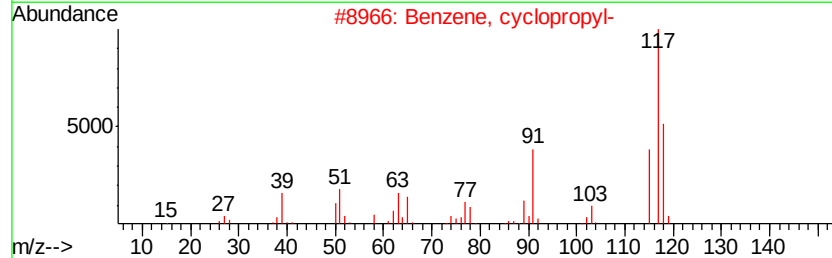
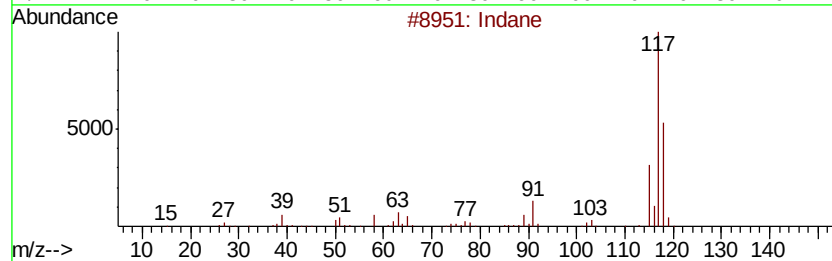
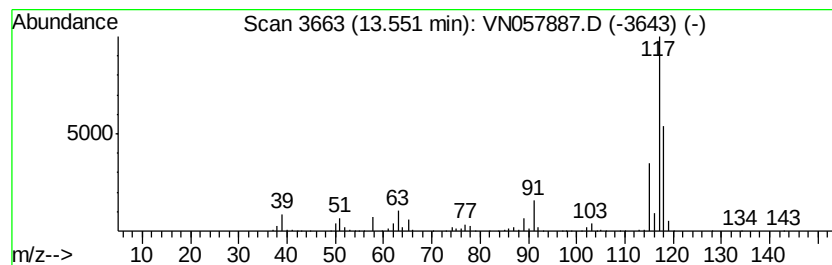
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.55	23.36 ug/l	760010	1,4-Dichlorobenzene-d4		13.35	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Indane	118	C9H10	000496-11-7	95
2		Benzene, cyclopropyl-	118	C9H10	000873-49-4	81
3		Deltacyclene	118	C9H10	007785-10-6	76
4		Benzene, 2-propenyl-	118	C9H10	000300-57-2	74
5		Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	64



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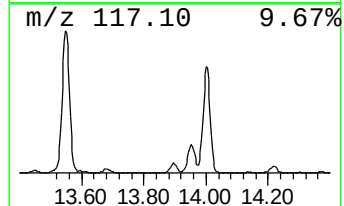
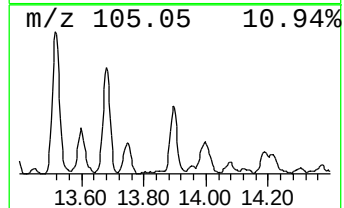
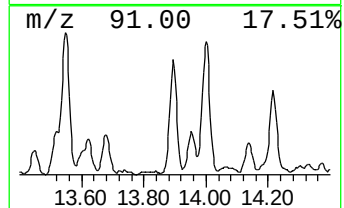
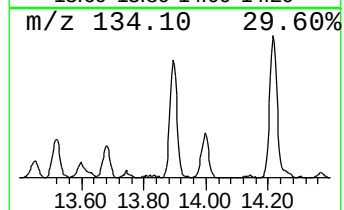
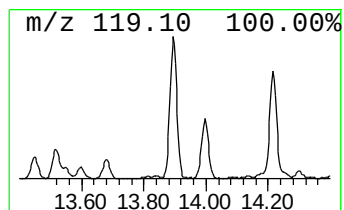
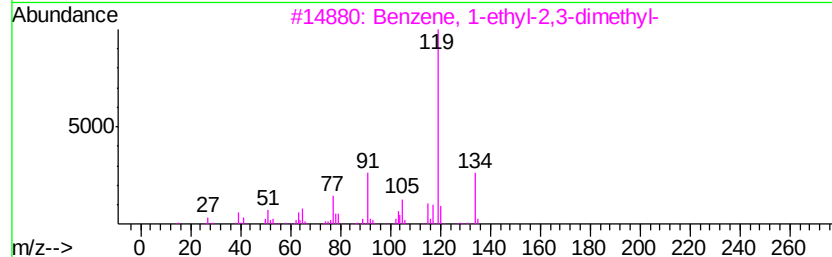
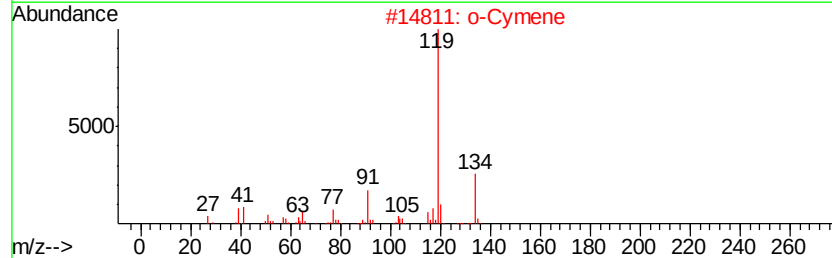
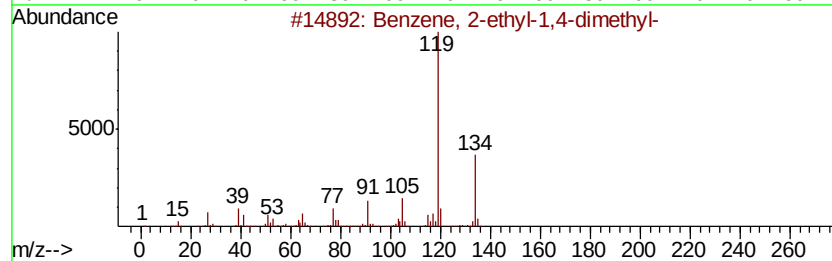
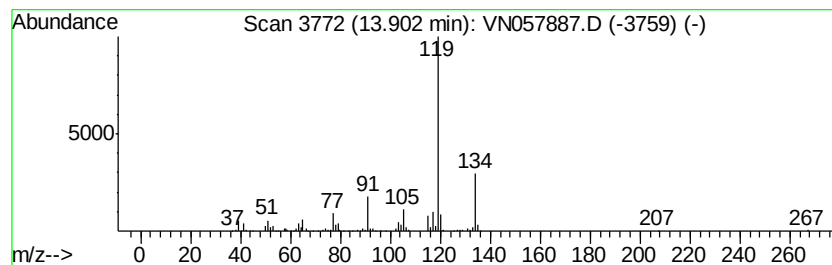
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Peak Number 3 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.90	11.16 ug/l	363095	1,4-Dichlorobenzene-d4	13.35	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2	o-Cymene	134	C10H14	000527-84-4	95
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
5	p-Cymene	134	C10H14	000099-87-6	94



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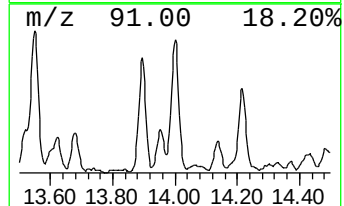
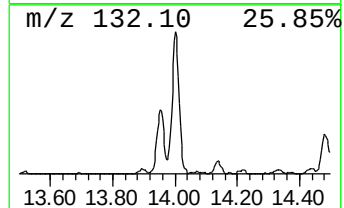
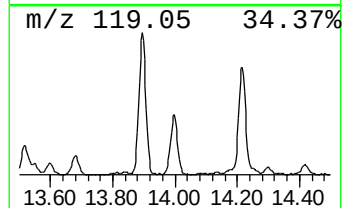
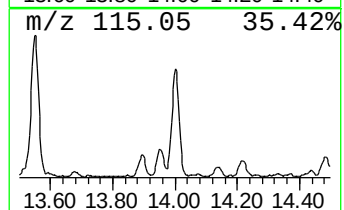
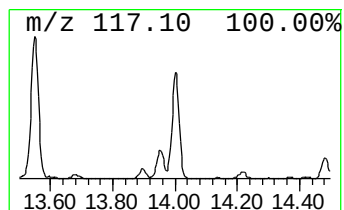
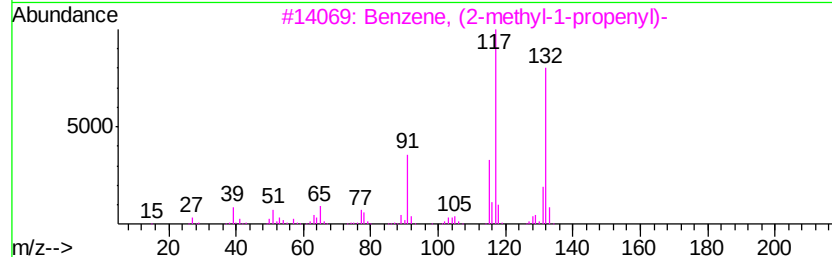
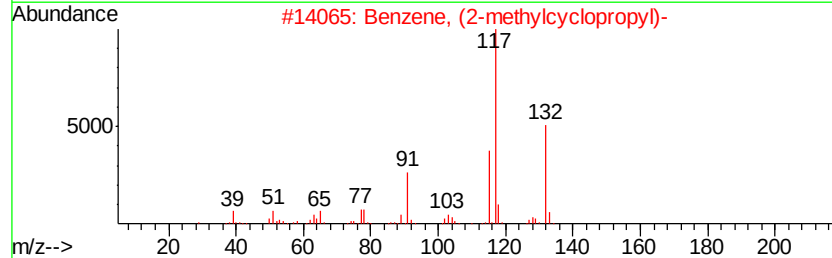
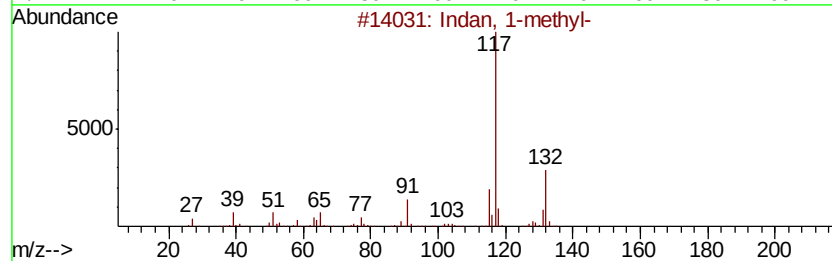
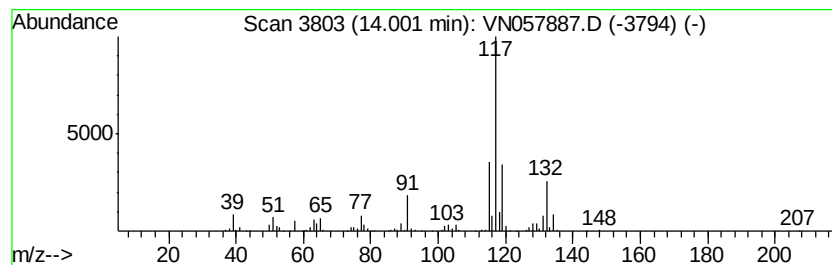
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Indan, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.00	16.24 ug/l	528346	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indan, 1-methyl-	132 C10H12	000767-58-8	81
2	Benzene, (2-methylcyclopropyl)-	132 C10H12	1000327-39-0	76
3	Benzene, (2-methyl-1-propenyl)-	132 C10H12	000768-49-0	74
4	3-Phenylbut-1-ene	132 C10H12	000934-10-1	74
5	Benzene, 2-ethenyl-1,3-dimethyl-	132 C10H12	002039-90-9	72



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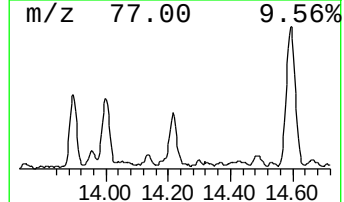
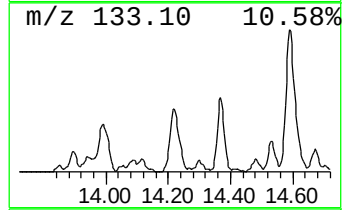
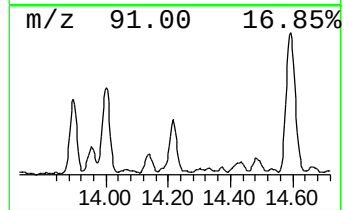
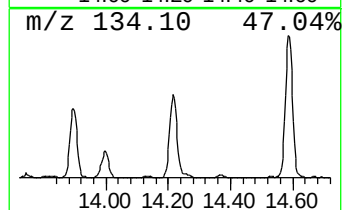
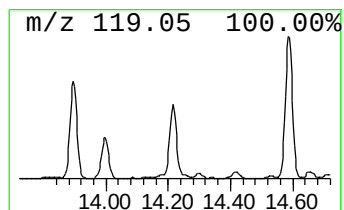
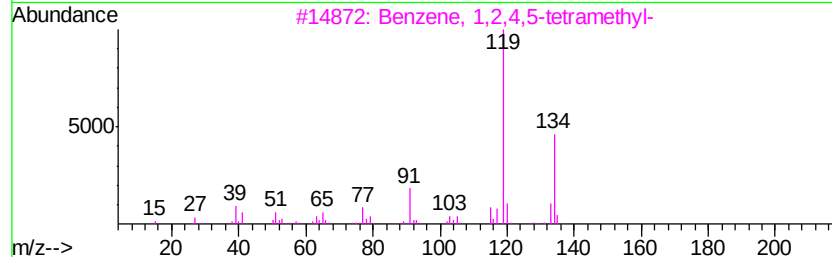
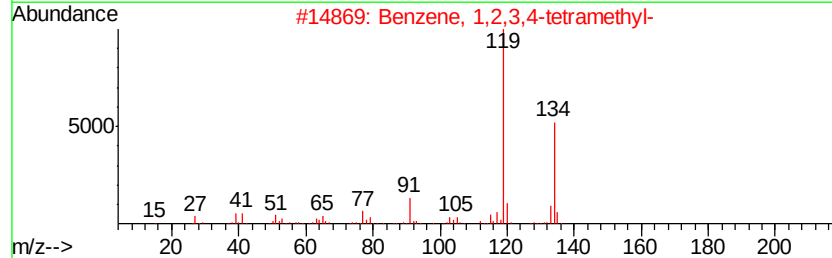
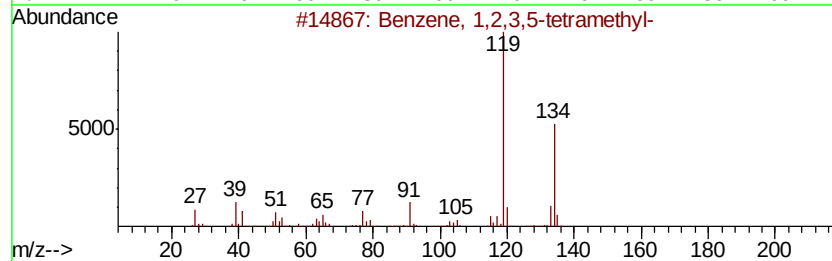
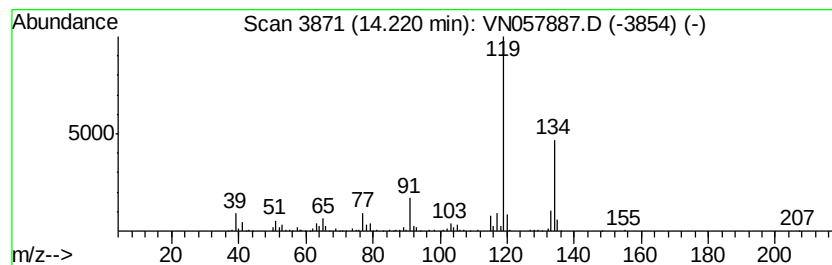
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.22	10.94 ug/l	355867	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	96
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4	p-Cymene	134	C10H14	000099-87-6	93
5	o-Cymene	134	C10H14	000527-84-4	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN091119\
Data File : VN057887.D
Acq On : 11 Sep 2019 00:23
Operator : JC/SP
Sample : K4818-07
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 28 Sample Multiplier: 1

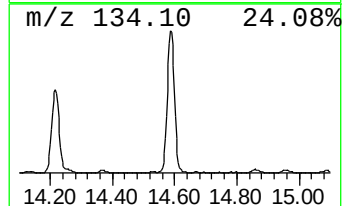
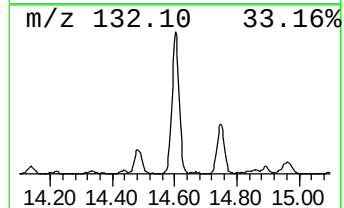
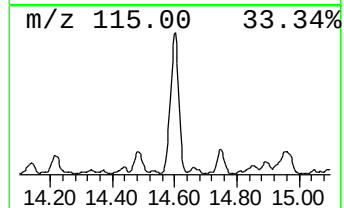
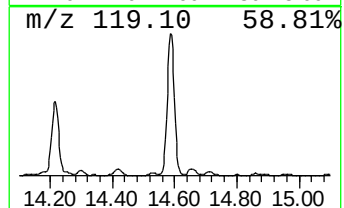
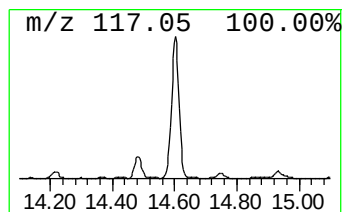
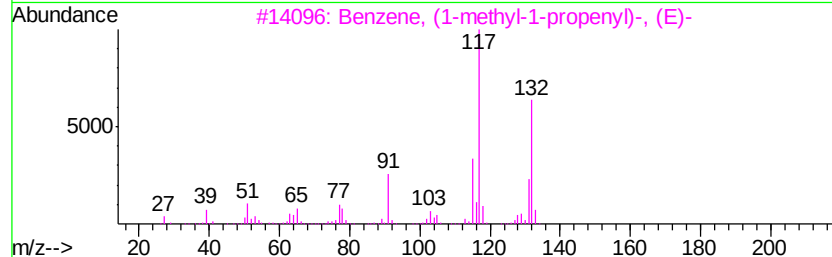
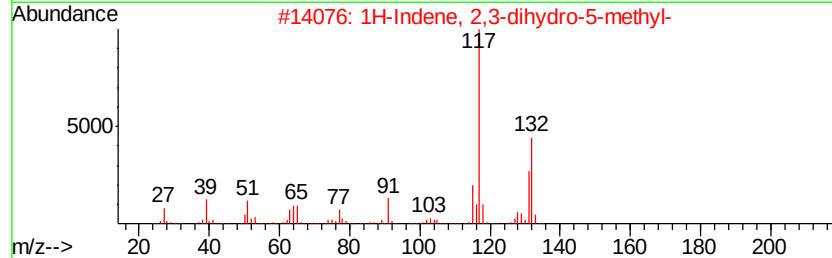
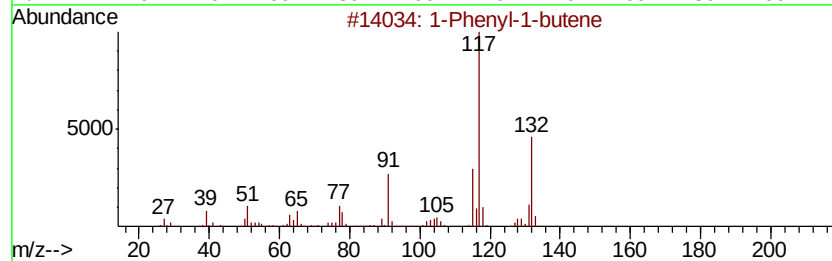
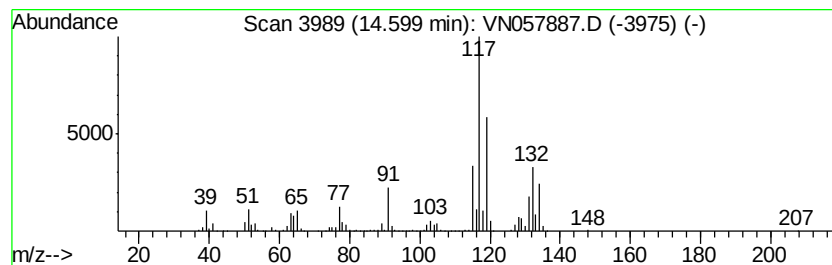
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ClientSampleId :
T11-MW-13-8.0-091019

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Quant Title : SW846 8260

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

Peak Number 6 1-Phenyl-1-butene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.60	35.26 ug/l	1147230	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	55
2	1H-Indene, 2,3-dihydro-5-methyl-	132 C10H12	000874-35-1	55
3	Benzene, (1-methyl-1-propenyl)-,...	132 C10H12	000768-00-3	55
4	3-Phenylbut-1-ene	132 C10H12	000934-10-1	55
5	Benzene, 1,2,3,4-tetramethyl-	134 C10H14	000488-23-3	50



Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091119\
Data File : VN057887.D
Acq On : 11 Sep 2019 00:23
Operator : JC/SP
Sample : K4818-07
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
T11-MW-13-8.0-091019

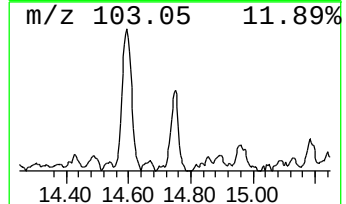
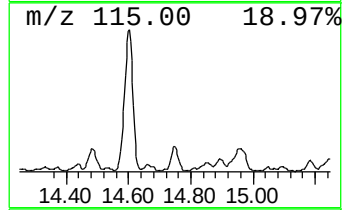
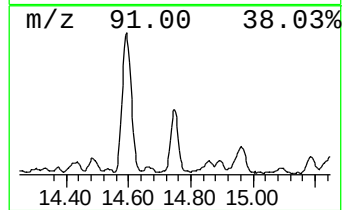
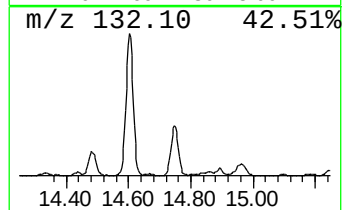
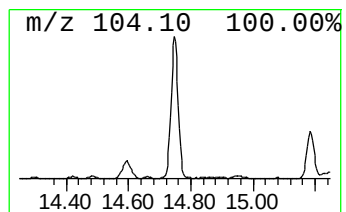
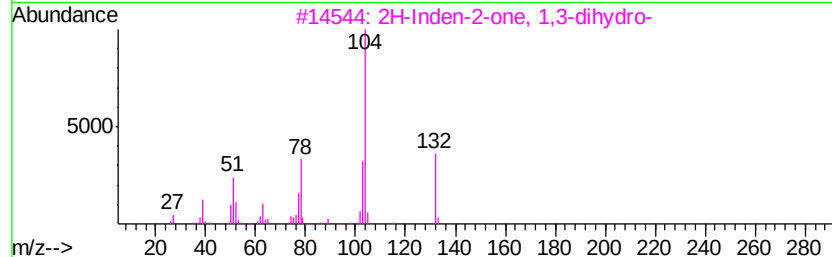
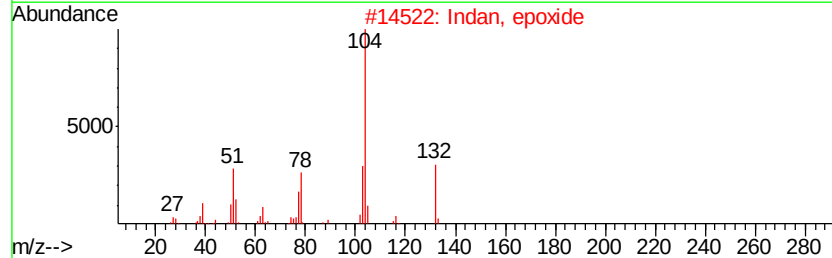
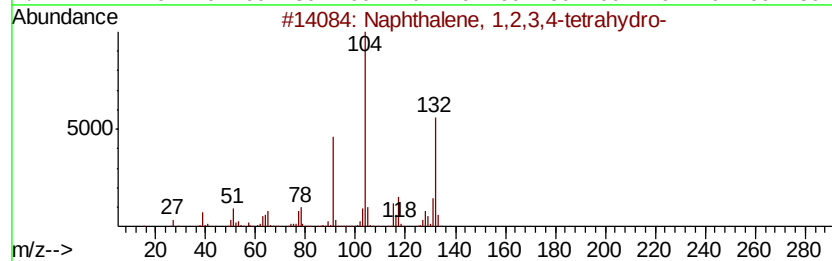
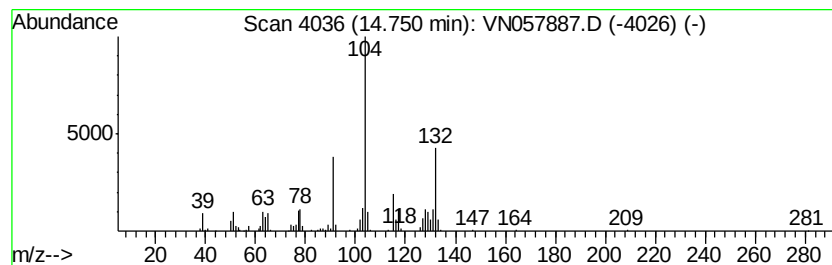
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Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.75	5.94 ug/l	193117	1,4-Dichlorobenzene-d4	13.35

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	96
2			Indan, epoxide	132	C9H8O	000768-22-9	64
3			2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	58
4			2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	43
5			2-Furanone, 4-phenyltetrahydro	162	C10H10O2	1000197-38-4	38



Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091119\
Data File : VN057887.D
Acq On : 11 Sep 2019 00:23
Operator : JC/SP
Sample : K4818-07
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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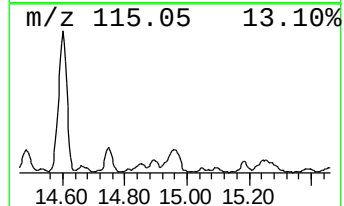
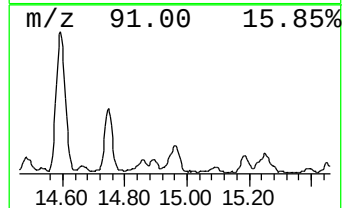
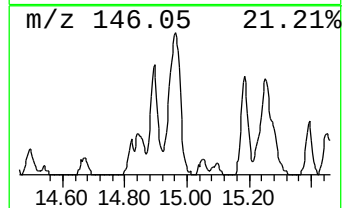
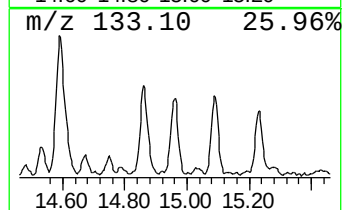
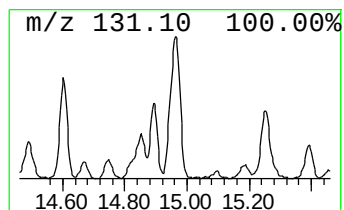
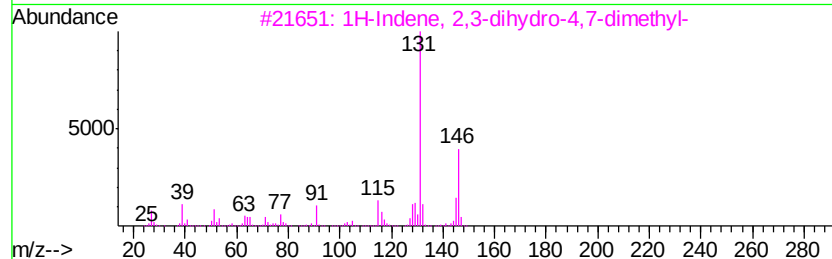
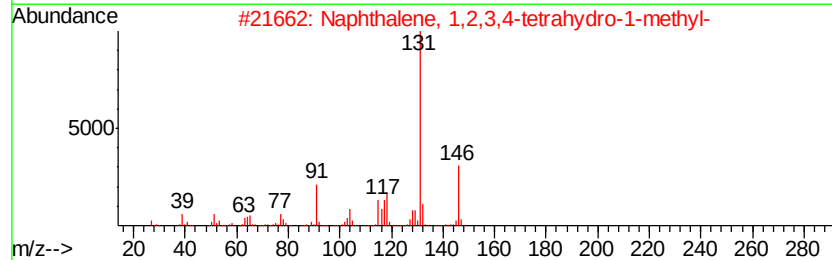
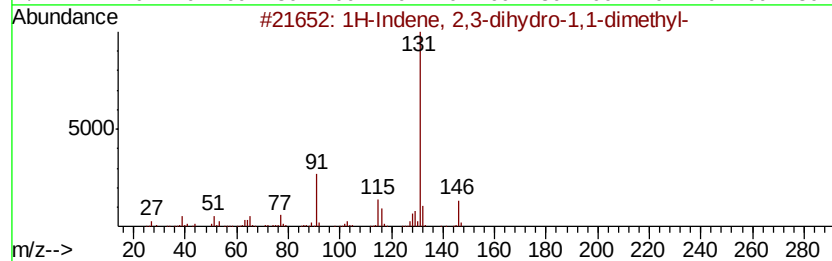
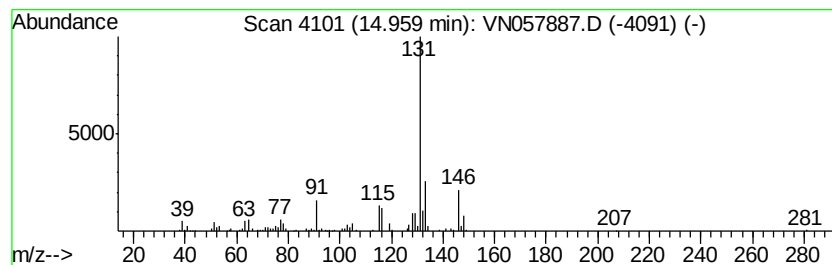
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Quant Title : SW846 8260

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

Peak Number 8 1H-Indene, 2,3-dihydro-1,1-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	6.04 ug/l	196618	1,4-Dichlorobenzene-d4	13.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	87
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	87
3		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	87
4		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	87
5		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	87



Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091119\
Data File : VN057887.D
Acq On : 11 Sep 2019 00:23
Operator : JC/SP
Sample : K4818-07
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 28 Sample Multiplier: 1

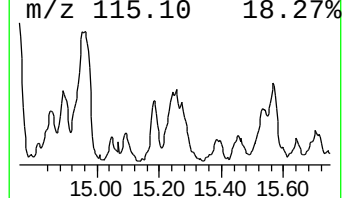
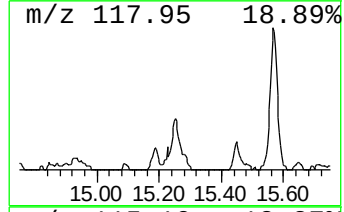
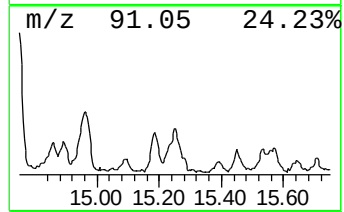
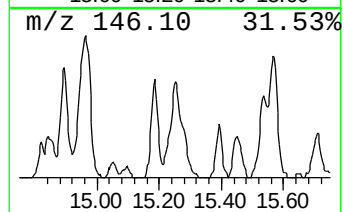
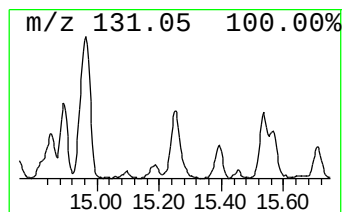
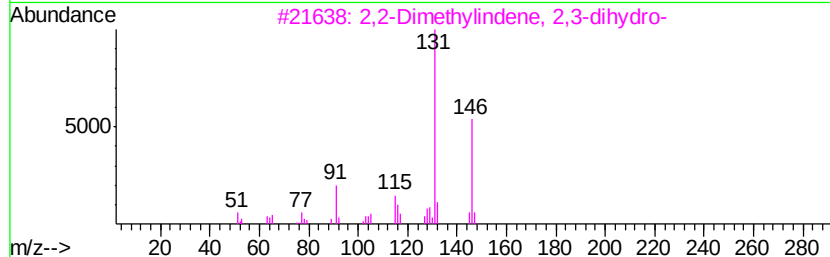
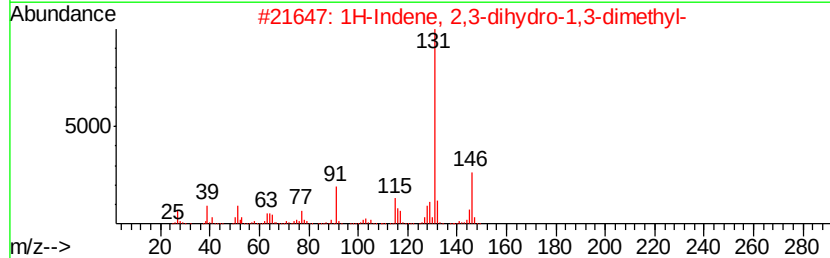
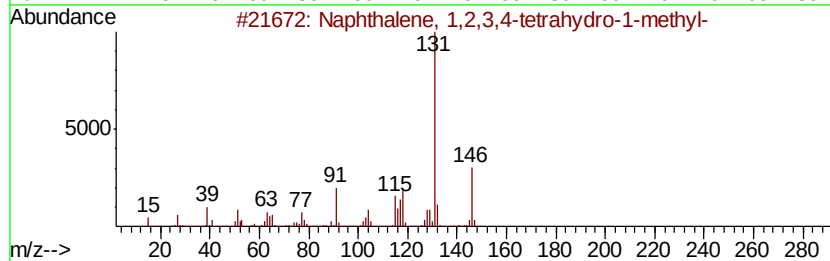
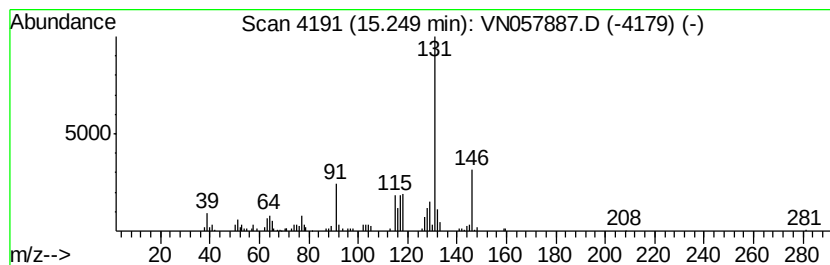
Instrument :
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ClientSampleId :
T11-MW-13-8.0-091019

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091019W.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 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.25	6.52 ug/l	212003	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-1,3-dimet...	146 C11H14	004175-53-5 81
3		2,2-Dimethylindene, 2,3-dihydro-	146 C11H14	020836-11-7 81
4		Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3 81
5		Benzene, (1,2-dimethyl-1-propenyl)-	146 C11H14	000769-57-3 76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN091119\
Data File : VN057887.D
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Sample : K4818-07
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 28 Sample Multiplier: 1

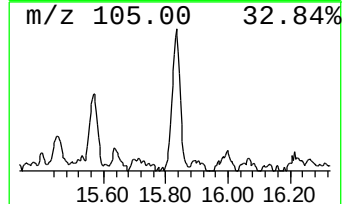
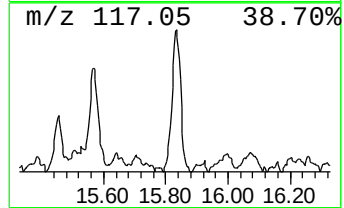
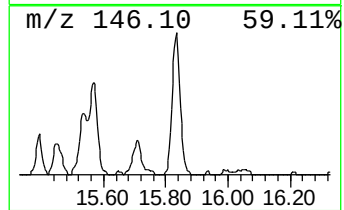
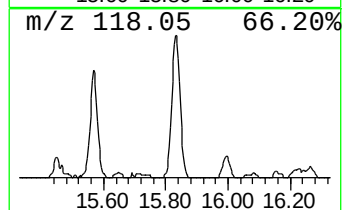
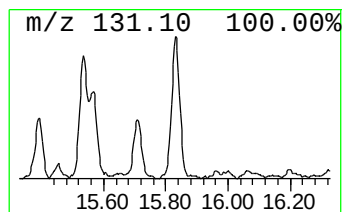
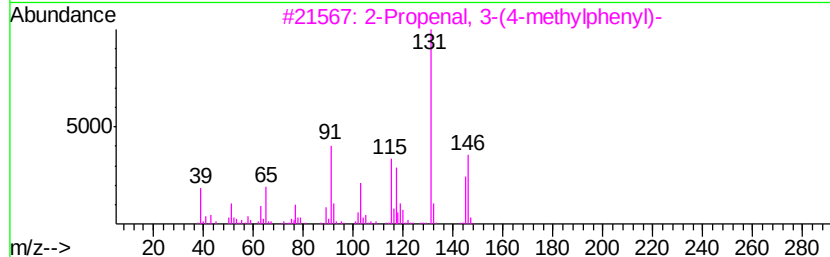
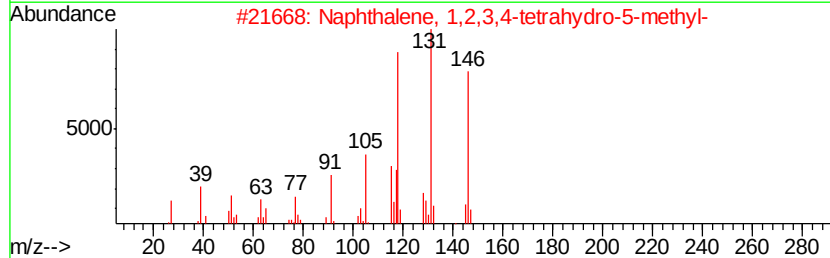
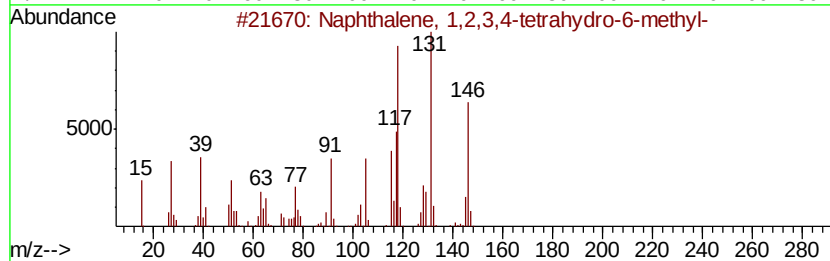
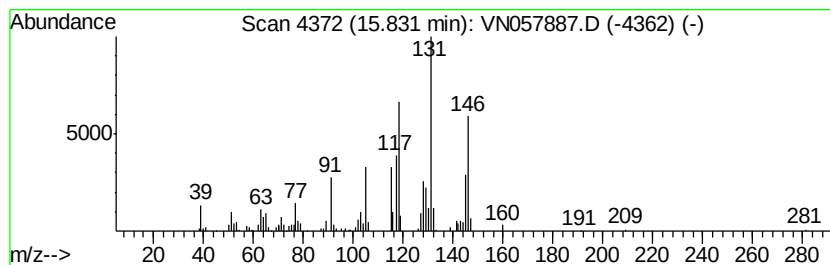
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetra... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.83	5.78 ug/l	188032	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 93
3		2-Propenal, 3-(4-methylphenyl)-	146 C10H10O	001504-75-2 60
4		Benzene, 2-ethenyl-1,3,5-trimethyl-	146 C11H14	000769-25-5 60
5		Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3 60



Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091119\
Data File : VN057887.D
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Sample : K4818-07
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 28 Sample Multiplier: 1

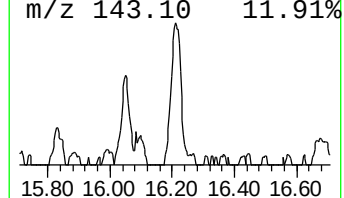
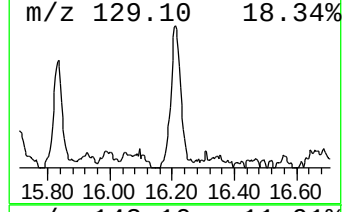
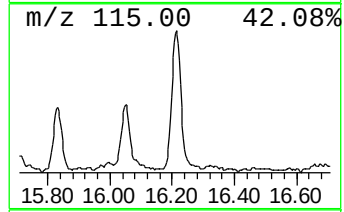
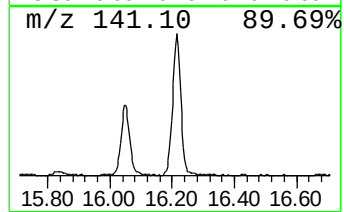
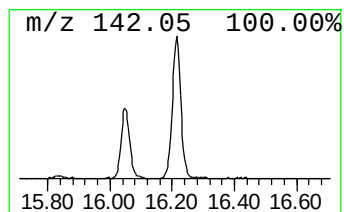
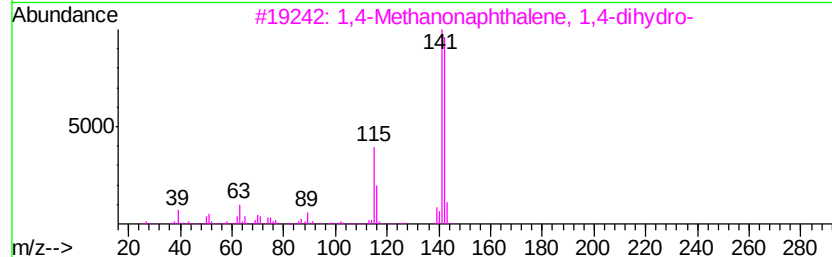
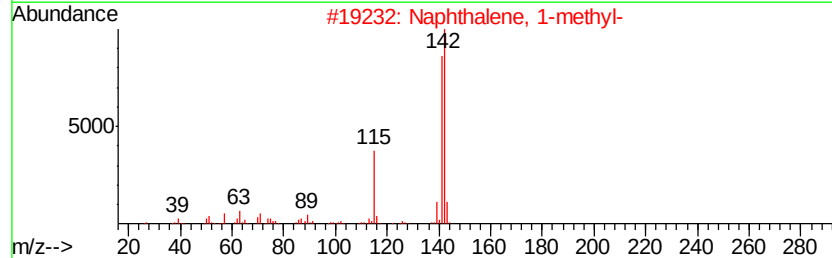
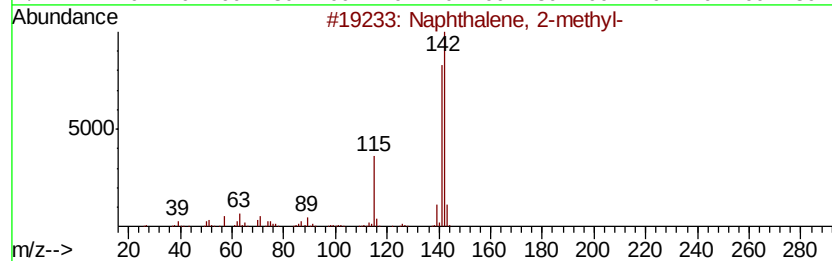
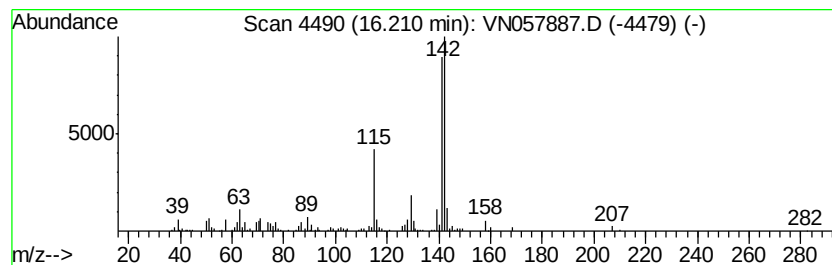
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ClientSampleId :
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Quant Title : SW846 8260

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

Peak Number 11 Naphthalene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.21	7.62 ug/l	248043	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2-methyl-	142 C11H10	000091-57-6 96
2		Naphthalene, 1-methyl-	142 C11H10	000090-12-0 96
3		1,4-Methanonaphthalene, 1,4-dihv...	142 C11H10	004453-90-1 91
4		Benzocycloheptatriene	142 C11H10	000264-09-5 91
5		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142 C11H10	002443-46-1 87



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN091119\
Data File : VN057887.D
Acq On : 11 Sep 2019 00:23
Operator : JC/SP
Sample : K4818-07
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
T11-MW-13-8.0-091019

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091019W.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
Isopropyl Alcohol	4.05	8.8	ug/l	153834	1	7.65	870907	50.0
Indane	13.55	23.4	ug/l	760010	4	13.35	1626870	50.0
Benzene, 2-ethyl-...	13.90	11.2	ug/l	363095	4	13.35	1626870	50.0
Indan, 1-methyl-	14.00	16.2	ug/l	528346	4	13.35	1626870	50.0
Benzene, 1,2,3,5-...	14.22	10.9	ug/l	355867	4	13.35	1626870	50.0
1-Phenyl-1-butene	14.60	35.3	ug/l	1147230	4	13.35	1626870	50.0
Naphthalene, 1,2,...	14.75	5.9	ug/l	193117	4	13.35	1626870	50.0
1H-Indene, 2,3-di...	14.96	6.0	ug/l	196618	4	13.35	1626870	50.0
Naphthalene, 1,2,...	15.25	6.5	ug/l	212003	4	13.35	1626870	50.0
Naphthalene, 1,2,...	15.83	5.8	ug/l	188032	4	13.35	1626870	50.0
Naphthalene, 2-me...	16.21	7.6	ug/l	248043	4	13.35	1626870	50.0