

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX030819\
 Data File : VX007946.D
 Acq On : 08 Mar 2019 22:15
 Operator : JC/SP
 Sample : K1800-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 ERM-33

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_X\METHOD\82X030619W.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	4.403	522	533	541	rBV3	5806	16299	1.28%	0.121%
2	5.507	701	714	723	rBV	157364	448636	35.10%	3.337%
3	5.665	730	740	757	rVB	250022	699362	54.71%	5.202%
4	6.067	795	806	814	rBV	159918	417533	32.66%	3.106%
5	6.147	814	819	830	rVB2	24410	63652	4.98%	0.473%
6	6.860	926	936	953	rBV	390269	926406	72.47%	6.891%
7	7.464	1025	1035	1044	rVB3	18276	44161	3.45%	0.328%
8	8.713	1234	1240	1256	rVB	818688	1278334	100.00%	9.509%
9	10.109	1463	1469	1478	rBV	797133	1110151	86.84%	8.258%
10	10.250	1487	1492	1497	rVB	10754	15218	1.19%	0.113%
11	10.359	1502	1510	1516	rVB4	6716	15849	1.24%	0.118%
12	11.018	1612	1618	1623	rBV	46909	59632	4.66%	0.444%
13	11.134	1632	1637	1644	rBV	682917	890711	69.68%	6.625%
14	11.359	1669	1674	1684	rVB	45033	57250	4.48%	0.426%
15	11.670	1719	1725	1733	rVB2	12623	19363	1.51%	0.144%
16	11.804	1744	1747	1752	rVB	20490	25194	1.97%	0.187%
17	11.944	1767	1770	1779	rVB	36305	44488	3.48%	0.331%
18	12.079	1786	1792	1798	rBV	891346	1084787	84.86%	8.069%
19	12.140	1798	1802	1808	rVB	17825	23765	1.86%	0.177%
20	12.231	1813	1817	1821	rVV	14226	18281	1.43%	0.136%
21	12.298	1821	1828	1835	rVV	335644	422022	33.01%	3.139%
22	12.365	1835	1839	1849	rVB2	40916	74919	5.86%	0.557%
23	12.457	1849	1854	1861	rBV	33696	45416	3.55%	0.338%
24	12.664	1882	1888	1891	rBV	115521	146876	11.49%	1.093%
25	12.700	1891	1894	1898	rVV	64960	78793	6.16%	0.586%
26	12.755	1898	1903	1910	rVB2	209836	286273	22.39%	2.129%
27	12.896	1919	1926	1931	rVB3	23699	39432	3.08%	0.293%
28	12.975	1931	1939	1943	rBV	104563	143870	11.25%	1.070%
29	13.017	1943	1946	1952	rVB2	10742	13623	1.07%	0.101%
30	13.084	1952	1957	1962	rVB3	24783	39562	3.09%	0.294%
31	13.152	1962	1968	1971	rBV2	14336	23659	1.85%	0.176%
32	13.194	1971	1975	1977	rBV2	32008	38175	2.99%	0.284%
33	13.225	1977	1980	1990	rVB	133369	180885	14.15%	1.345%
34	13.310	1990	1994	1995	rBV2	15665	17013	1.33%	0.127%

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35	13.347	1995	2000	2006	rVB2	409142	535942	41.93%	3.987%
36	13.402	2006	2009	2011	rBV4	20072	24018	1.88%	0.179%
37	13.481	2018	2022	2027	rVB	48783	62980	4.93%	0.468%
38	13.560	2031	2035	2038	rVV2	23958	33524	2.62%	0.249%
39	13.597	2038	2041	2044	rVV	66985	94485	7.39%	0.703%
40	13.639	2044	2048	2051	rVV2	82284	122549	9.59%	0.912%
41	13.676	2051	2054	2055	rVV	45461	52217	4.08%	0.388%
42	13.719	2058	2061	2070	rVB2	125809	184341	14.42%	1.371%
43	13.810	2070	2076	2077	rBV	15672	19784	1.55%	0.147%
44	13.834	2077	2080	2088	rVV3	27246	62019	4.85%	0.461%
45	13.908	2088	2092	2097	rVB	88357	112100	8.77%	0.834%
46	13.981	2097	2104	2108	rBV2	133032	179395	14.03%	1.334%
47	14.029	2108	2112	2115	rVV	46033	59140	4.63%	0.440%
48	14.060	2115	2117	2124	rVB2	9944	12890	1.01%	0.096%
49	14.133	2124	2129	2132	rBV	84718	109574	8.57%	0.815%
50	14.279	2146	2153	2161	rBV4	111865	249603	19.53%	1.857%
51	14.377	2165	2169	2174	rBV	47426	58700	4.59%	0.437%
52	14.432	2174	2178	2182	rVB	99025	124995	9.78%	0.930%
53	14.474	2182	2185	2190	rVB	15579	21211	1.66%	0.158%
54	14.535	2190	2195	2200	rBV2	139217	193787	15.16%	1.441%
55	14.584	2200	2203	2208	rVB2	13428	18818	1.47%	0.140%
56	14.694	2208	2221	2225	rBV	376098	542880	42.47%	4.038%
57	14.731	2225	2227	2231	rVB2	44161	49659	3.88%	0.369%
58	14.785	2231	2236	2239	rBV2	22094	27447	2.15%	0.204%
59	14.840	2239	2245	2249	rBV	478699	641383	50.17%	4.771%
60	14.907	2254	2256	2260	rVV2	20254	22212	1.74%	0.165%
61	14.962	2260	2265	2271	rVV3	22827	45449	3.56%	0.338%
62	15.109	2283	2289	2294	rBV3	9636	13925	1.09%	0.104%
63	15.249	2304	2312	2319	rBV	51526	90989	7.12%	0.677%
64	15.322	2319	2324	2331	rVB2	18649	30294	2.37%	0.225%
65	15.444	2337	2344	2347	rBV2	45269	70981	5.55%	0.528%
66	15.474	2347	2349	2354	rVV	34811	43645	3.41%	0.325%
67	15.547	2354	2361	2365	rVV	97011	154931	12.12%	1.152%
68	15.590	2365	2368	2374	rVV6	17829	33956	2.66%	0.253%
69	15.688	2374	2384	2387	rVV	147441	236975	18.54%	1.763%
70	15.724	2387	2390	2398	rVB	63616	97140	7.60%	0.723%
71	15.810	2398	2404	2410	rBV4	6543	15780	1.23%	0.117%

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72	15.919	2410	2422	2431	rVB2	35907	113935	8.91%	0.847%
73	16.072	2440	2447	2458	rVB	26450	50713	3.97%	0.377%
74	16.419	2499	2504	2516	rVB3	6580	19883	1.56%	0.148%

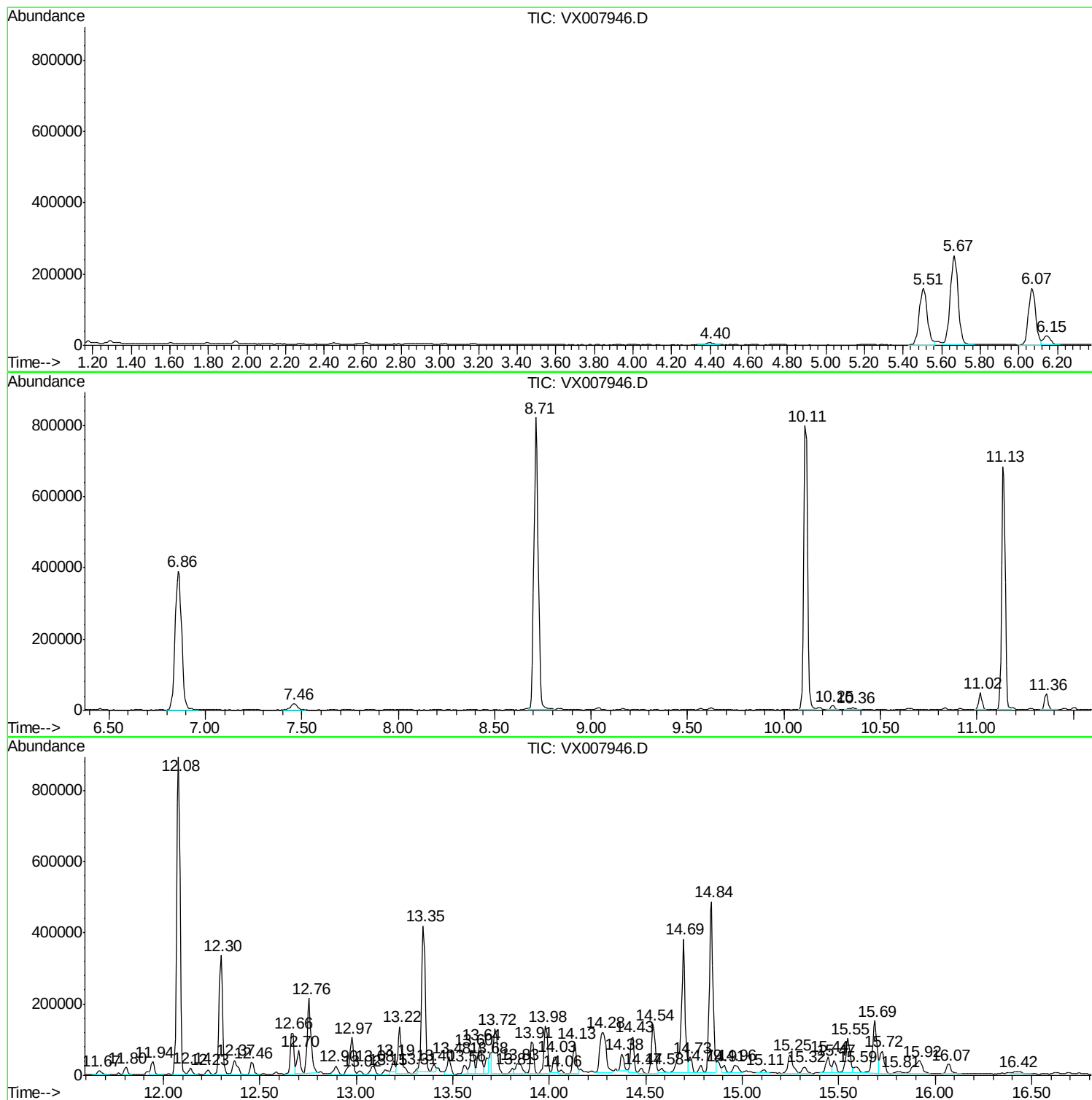
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X030619W.M
Quant Title : SW846 8260

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



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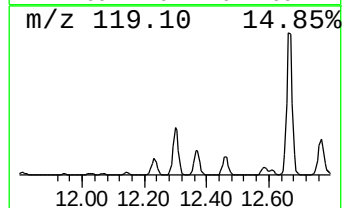
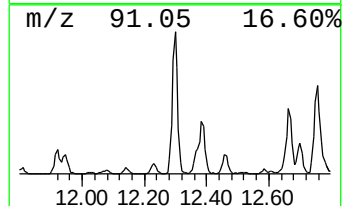
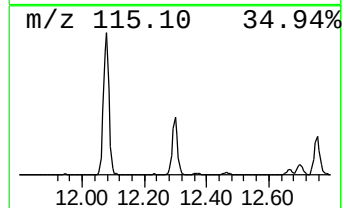
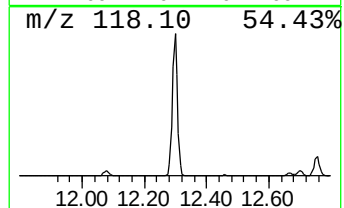
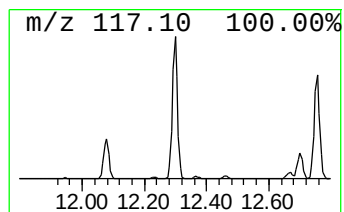
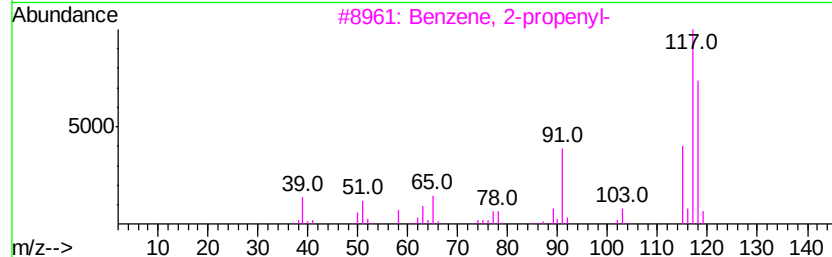
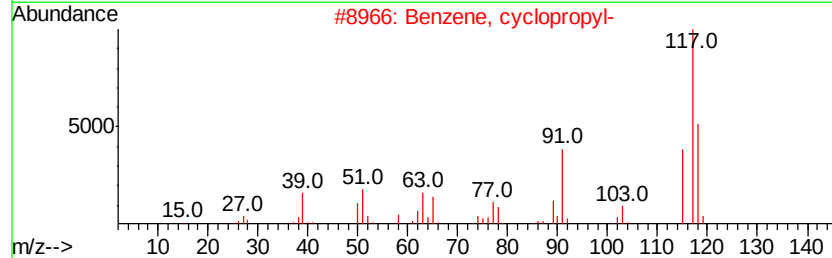
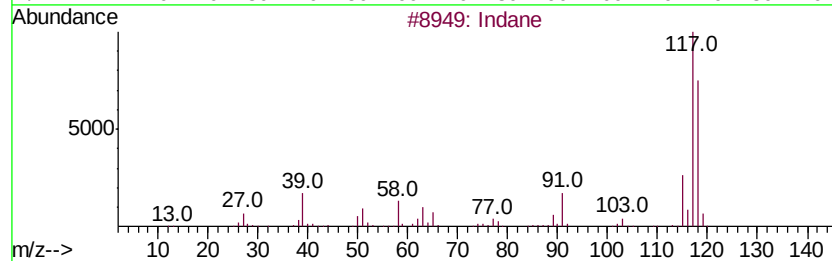
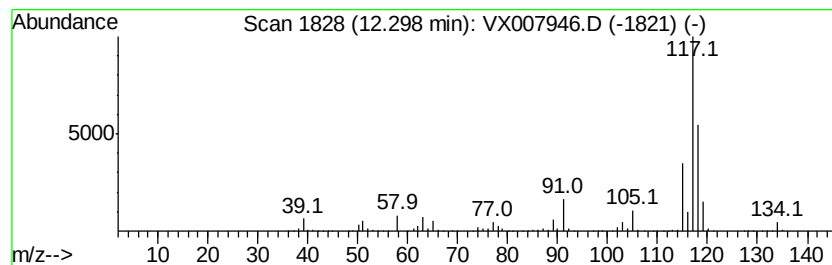
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X030619W.M
Quant Title : SW846 8260

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

Peak Number 1 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	19.45 ug/l	422022	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indane		118 C9H10	000496-11-7 81
2	Benzene, cyclopropyl-		118 C9H10	000873-49-4 76
3	Benzene, 2-propenyl-		118 C9H10	000300-57-2 74
4	Benzene, 1-propenyl-		118 C9H10	000637-50-3 74
5	Deltacyclene		118 C9H10	007785-10-6 64



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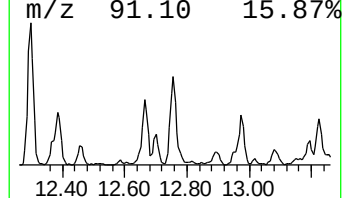
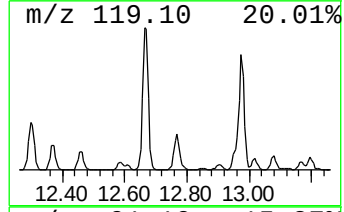
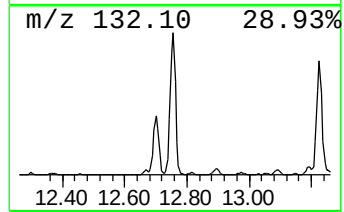
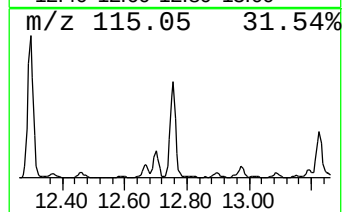
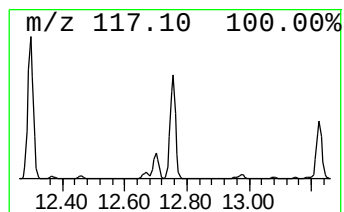
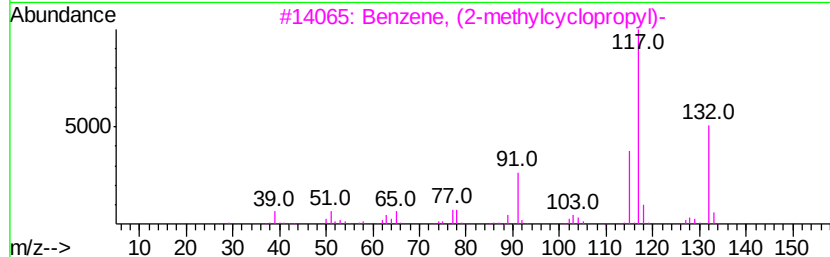
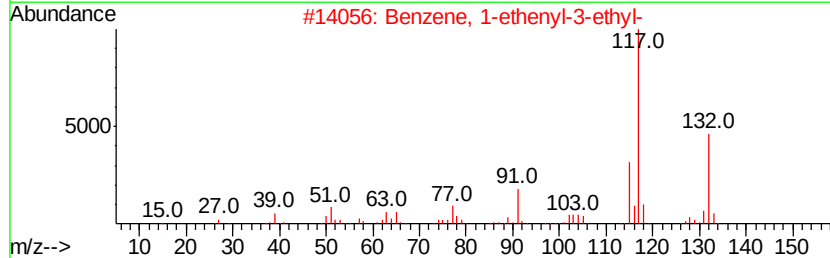
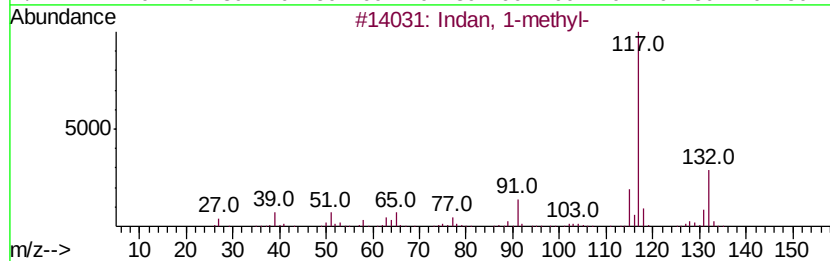
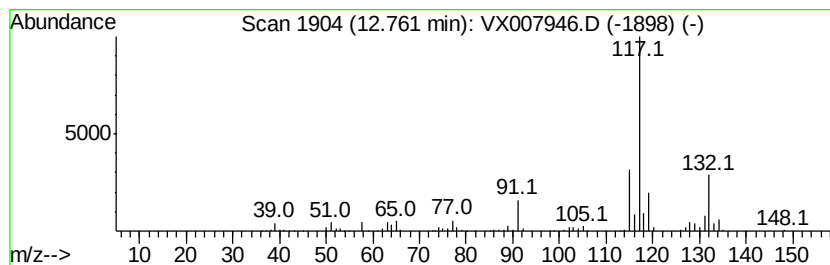
Instrument :
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ClientSampled :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X030619W.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Indan, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.76	13.19 ug/l	286273	1,4-Dichlorobenzene-d4	12.08	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	87
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
3	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	87
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	83
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	80



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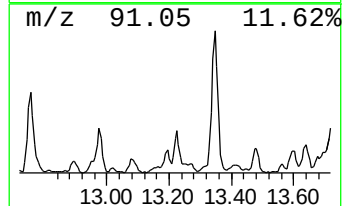
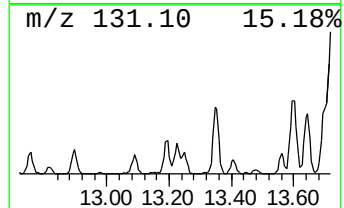
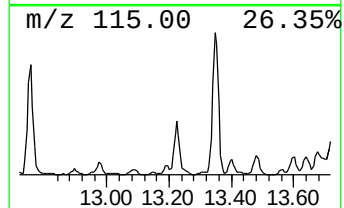
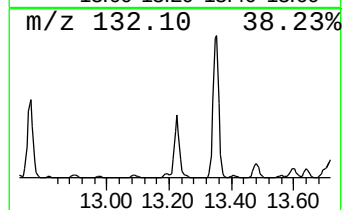
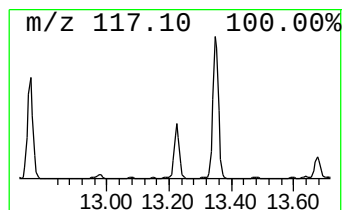
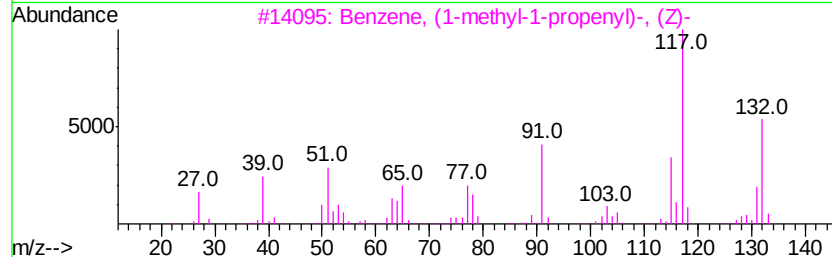
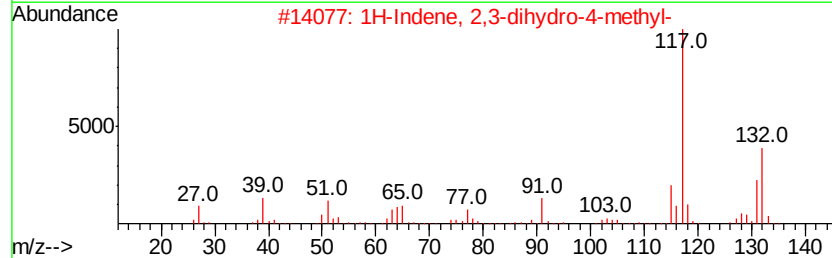
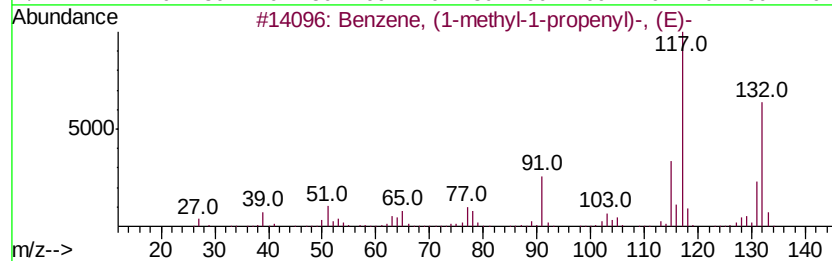
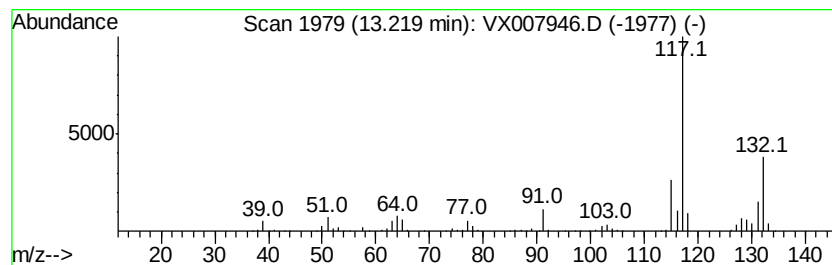
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, (1-methyl-1-propen... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	8.34 ug/l	180885	1,4-Dichlorobenzene-d4	12.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	91
2		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
3		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	90
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
5		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	87



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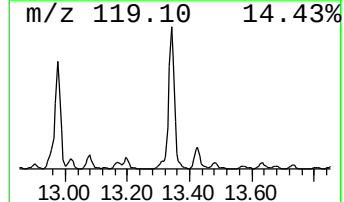
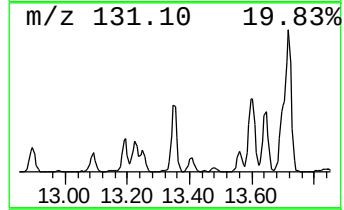
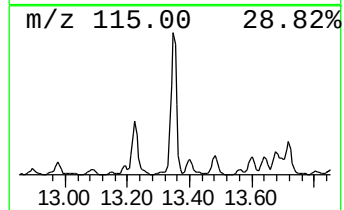
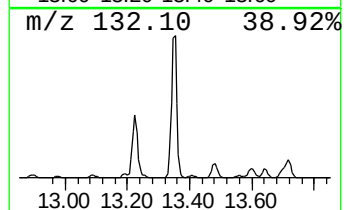
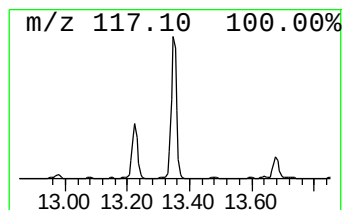
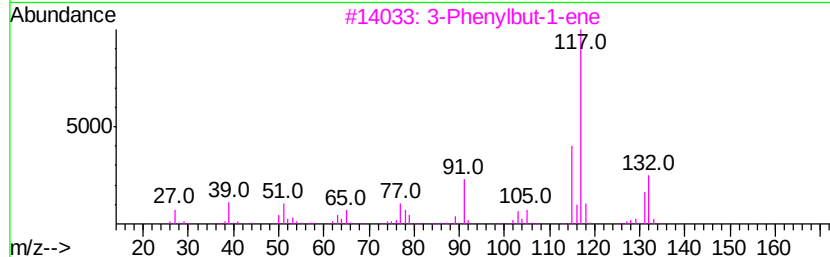
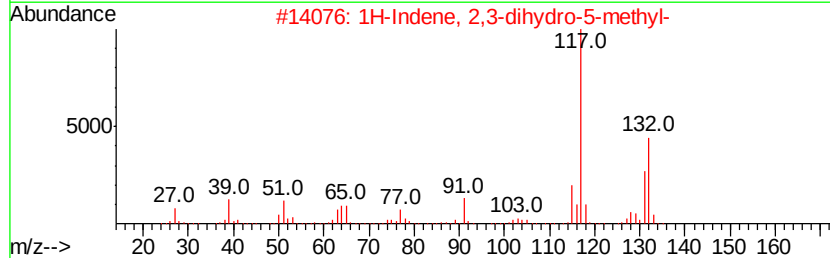
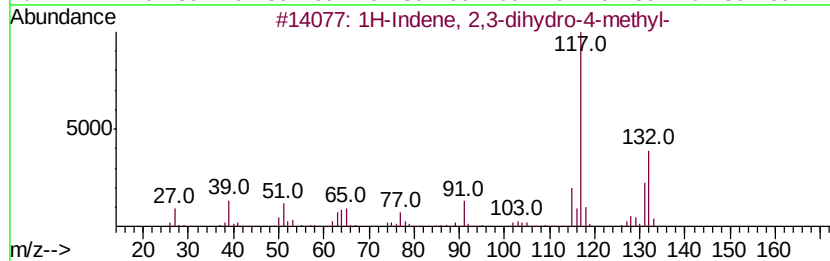
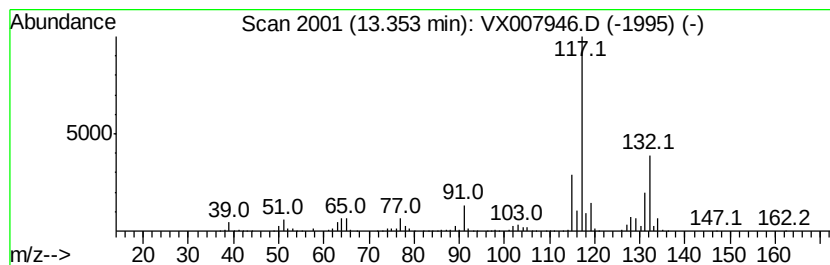
Instrument :
MSVOA_X
ClientSampleId :
ERM-33

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X030619W.M
Quant Title : SW846 8260

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

Peak Number 4 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	24.70 ug/l	535942	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132 C10H12	000824-22-6	93
2	1H-Indene, 2,3-dihydro-5-methyl-	132 C10H12	000874-35-1	87
3	3-Phenylbut-1-ene	132 C10H12	000934-10-1	87
4	Benzene, (1-methyl-1-propenyl)-,...	132 C10H12	000768-00-3	83
5	Indan, 1-methyl-	132 C10H12	000767-58-8	81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX030819\
Data File : VX007946.D
Acq On : 08 Mar 2019 22:15
Operator : JC/SP
Sample : K1800-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 28 Sample Multiplier: 1

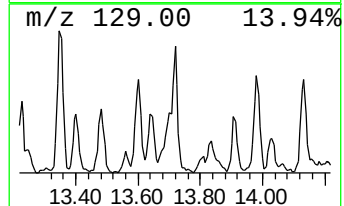
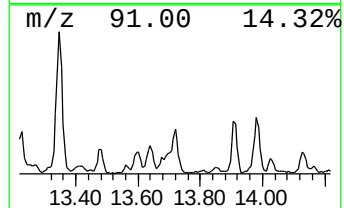
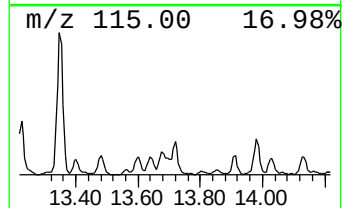
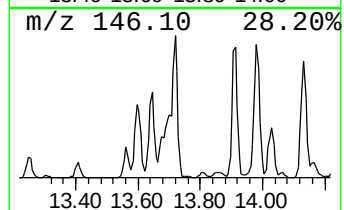
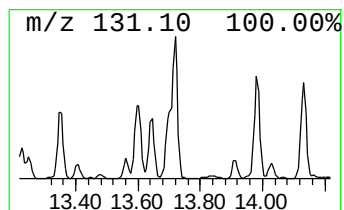
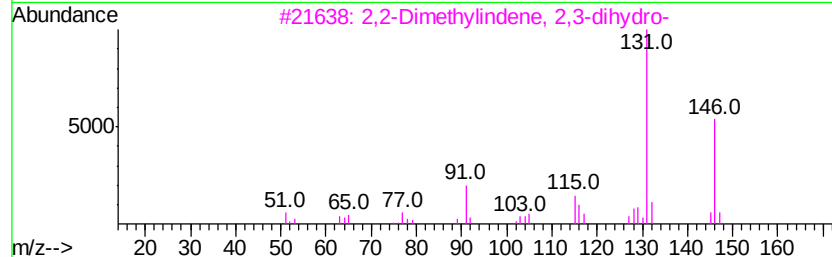
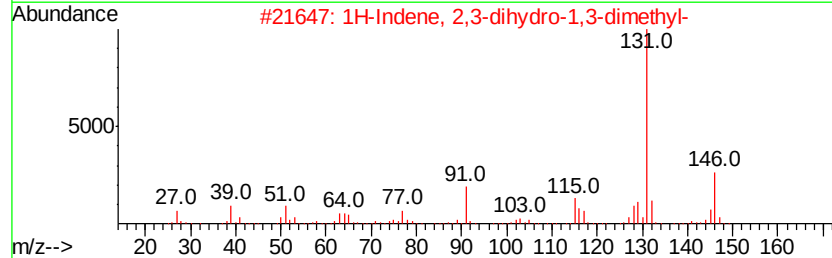
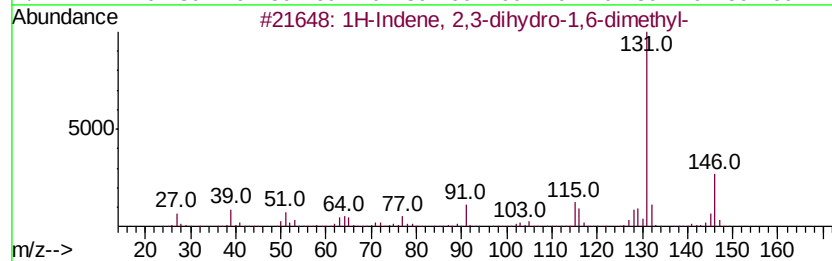
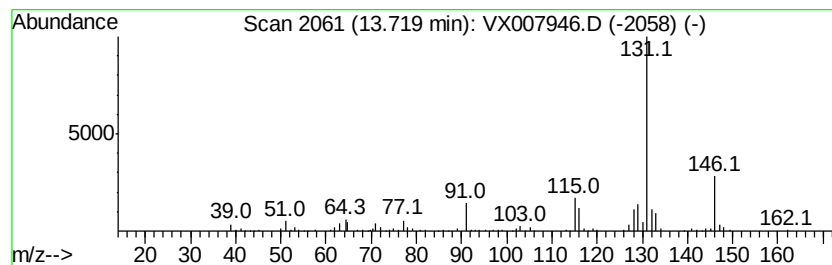
Instrument :
MSVOA_X
ClientSampleId :
ERM-33

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X030619W.M
Quant Title : SW846 8260

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

Peak Number 5 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	8.50 ug/l	184341	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1H-Indene, 2,3-dihydro-1,6-dimet...	146 C11H14	017059-48-2 93
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 90
4		Benzene, 1-methyl-2-(1-methyl-2-...	146 C11H14	097664-19-2 90
5		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001559-81-5 90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX030819\
Data File : VX007946.D
Acq On : 08 Mar 2019 22:15
Operator : JC/SP
Sample : K1800-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 28 Sample Multiplier: 1

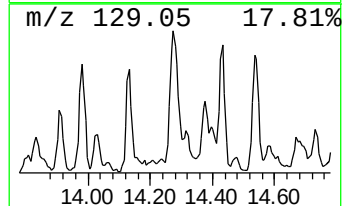
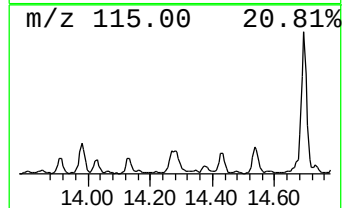
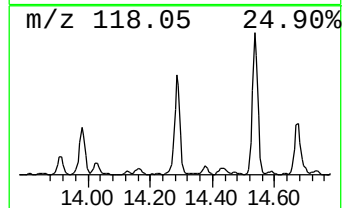
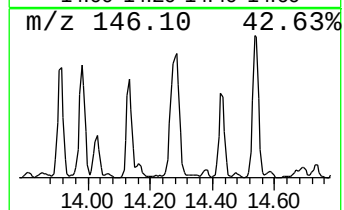
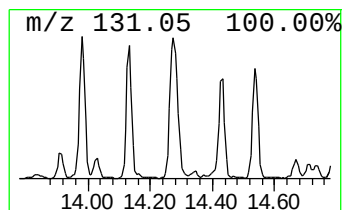
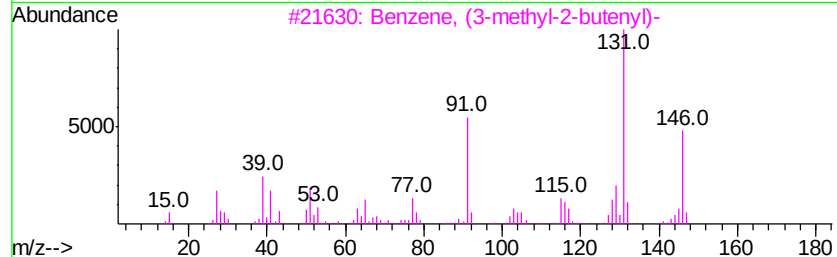
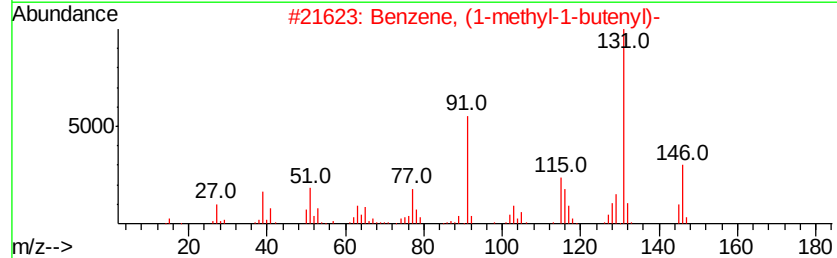
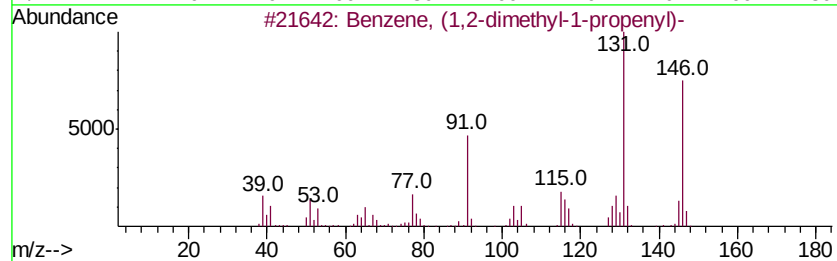
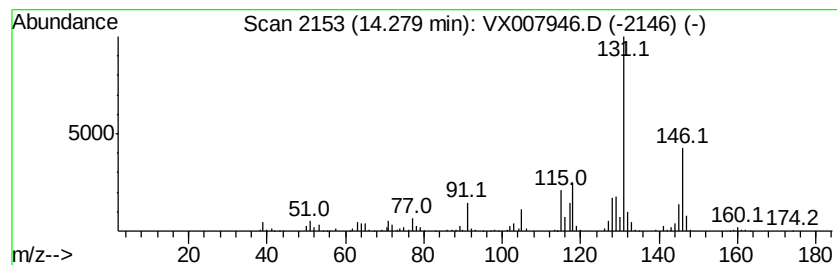
Instrument :
MSVOA_X
ClientSampleId :
ERM-33

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X030619W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.28	11.50 ug/l	249603	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (1,2-dimethyl-1-propenyl)-	146 C11H14	000769-57-3 93
2		Benzene, (1-methyl-1-butenyl)-	146 C11H14	053172-84-2 91
3		Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3 87
4		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 83
5		Benzene, (2-methyl-1-butenyl)-	146 C11H14	056253-64-6 83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX030819\
Data File : VX007946.D
Acq On : 08 Mar 2019 22:15
Operator : JC/SP
Sample : K1800-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
ERM-33

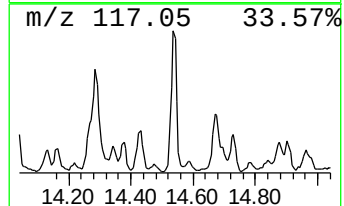
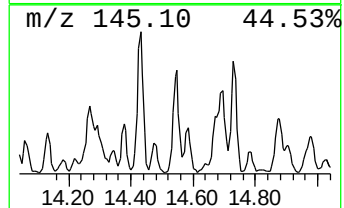
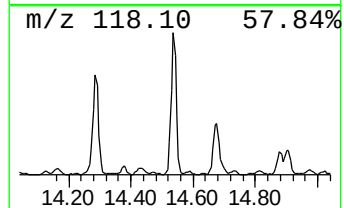
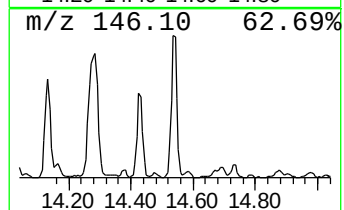
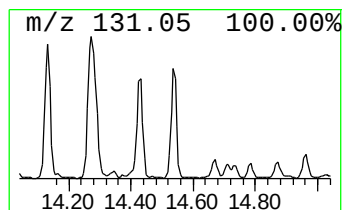
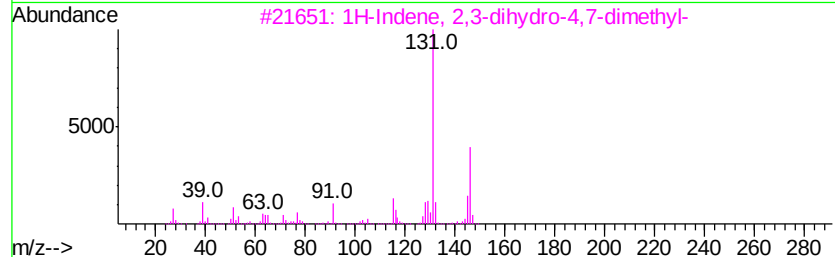
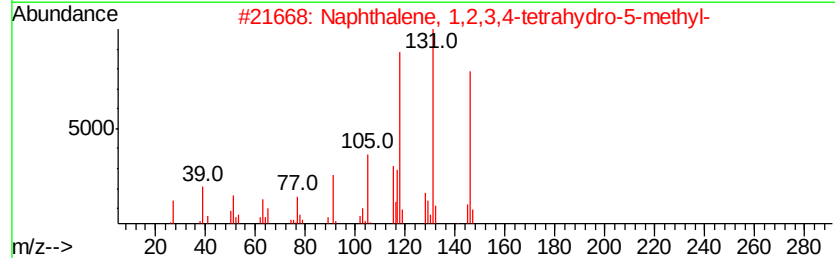
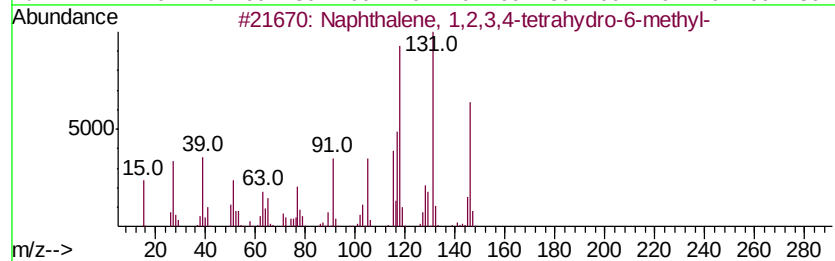
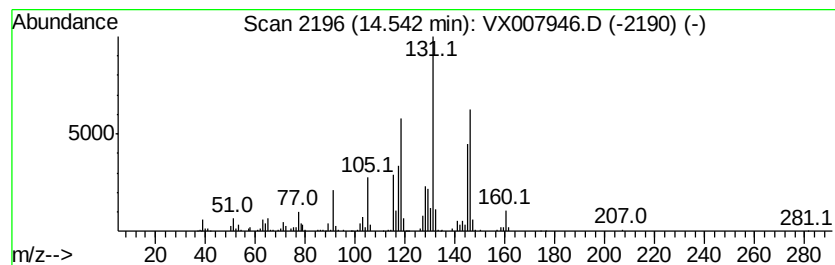
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Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.54	8.93 ug/l	193787	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	94
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	94
3		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70
4		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	64
5		Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX030819\
Data File : VX007946.D
Acq On : 08 Mar 2019 22:15
Operator : JC/SP
Sample : K1800-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 28 Sample Multiplier: 1

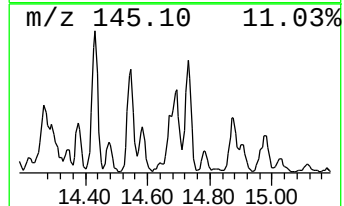
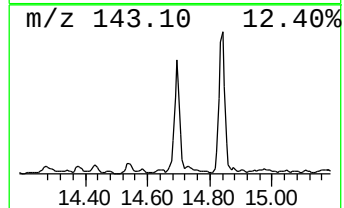
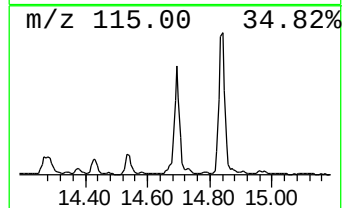
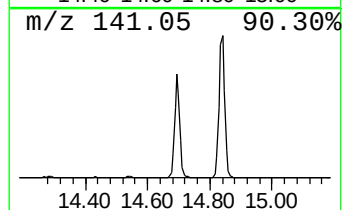
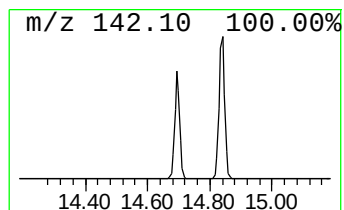
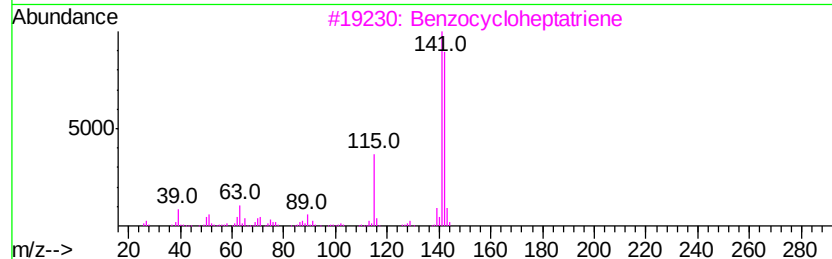
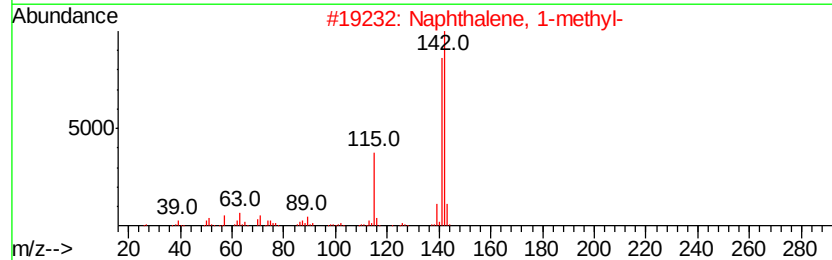
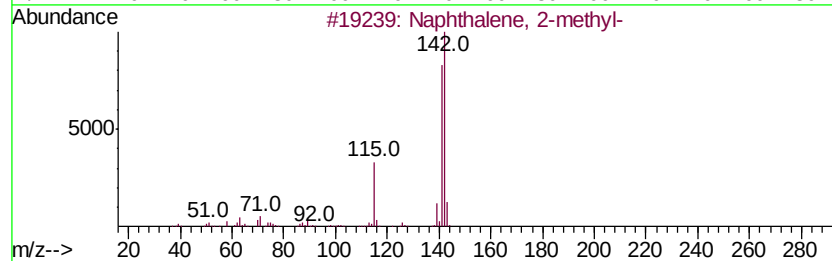
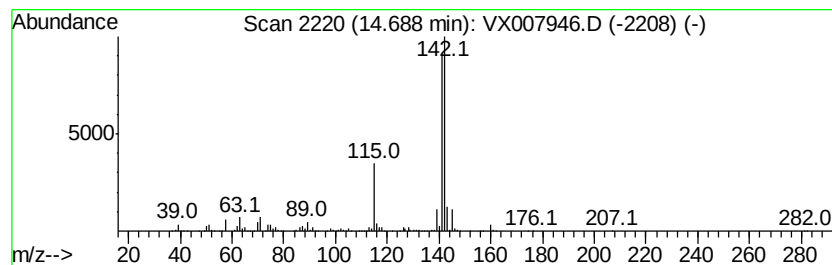
Instrument :
MSVOA_X
ClientSampleId :
ERM-33

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X030619W.M
Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	25.02 ug/l	542880	1,4-Dichlorobenzene-d4	12.08
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		Benzocycloheptatriene	142 C11H10	000264-09-5 90
4		1,4-Methanonaphthalene, 1,4-dihy...	142 C11H10	004453-90-1 90
5		1H-Indene, 1-ethylidene-	142 C11H10	002471-83-2 43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX030819\
Data File : VX007946.D
Acq On : 08 Mar 2019 22:15
Operator : JC/SP
Sample : K1800-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
ERM-33

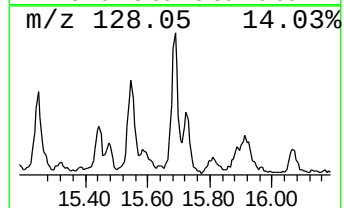
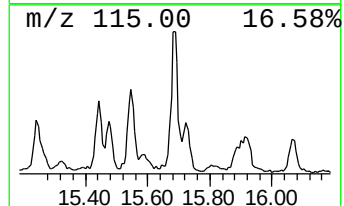
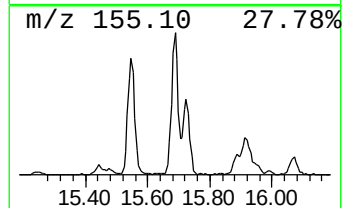
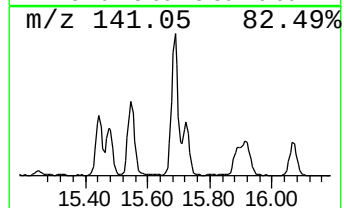
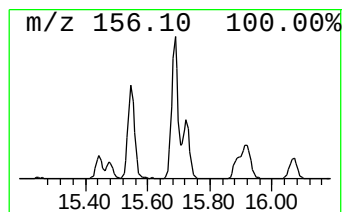
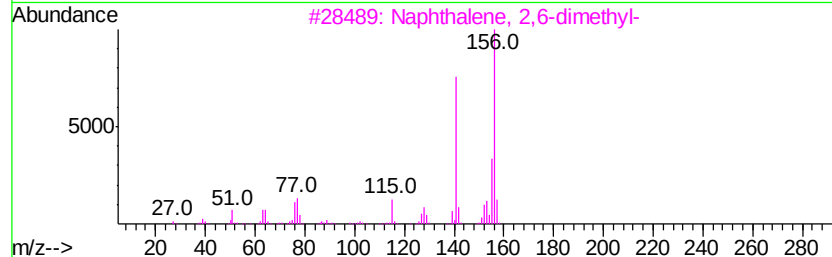
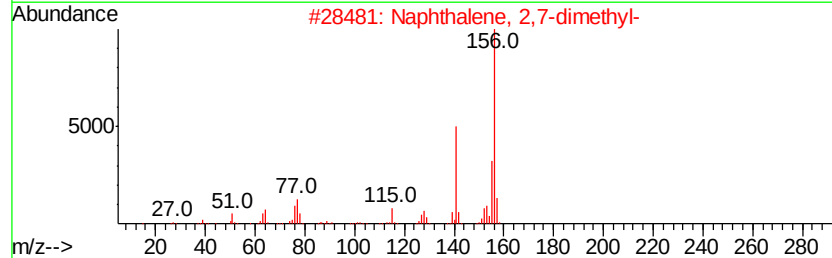
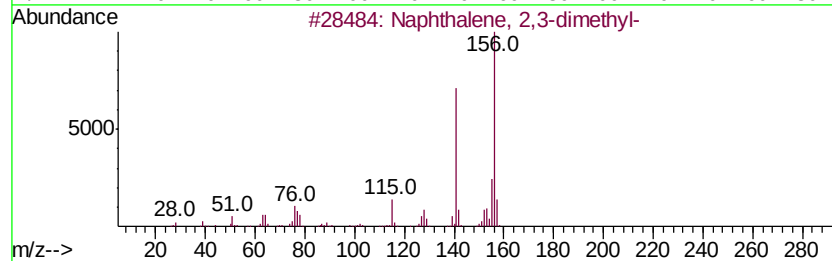
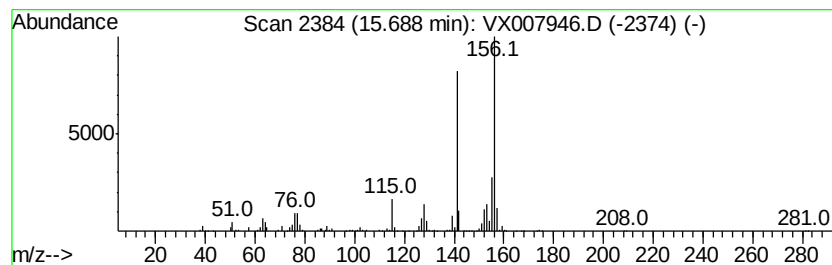
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X030619W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.69	10.92 ug/l	236975	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
4		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
5		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX030819\
Data File : VX007946.D
Acq On : 08 Mar 2019 22:15
Operator : JC/SP
Sample : K1800-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
ERM-33

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X030619W.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
Indane	12.30	19.4	ug/l	422022	4	12.08	1084790	50.0
Indan, 1-methyl-	12.76	13.2	ug/l	286273	4	12.08	1084790	50.0
Benzene, (1-methy...	13.22	8.3	ug/l	180885	4	12.08	1084790	50.0
1H-Indene, 2,3-di...	13.35	24.7	ug/l	535942	4	12.08	1084790	50.0
1H-Indene, 2,3-di...	13.72	8.5	ug/l	184341	4	12.08	1084790	50.0
Benzene, (1,2-dim...	14.28	11.5	ug/l	249603	4	12.08	1084790	50.0
Naphthalene, 1,2,...	14.54	8.9	ug/l	193787	4	12.08	1084790	50.0
Naphthalene, 1-me...	14.69	25.0	ug/l	542880	4	12.08	1084790	50.0
Naphthalene, 2,3-...	15.69	10.9	ug/l	236975	4	12.08	1084790	50.0