

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX101618\
 Data File : VX005361.D
 Acq On : 16 Oct 2018 11:00
 Operator : JC/MD
 Sample : J5521-13
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TWP-01

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\82X101218W.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	2.617	234	240	250	rBV	87426	161807	3.33%	0.296%
2	5.500	698	713	726	rBV	146923	422215	8.69%	0.772%
3	5.665	726	740	761	rVV	257401	782529	16.11%	1.431%
4	6.067	796	806	824	rBV	162633	427144	8.80%	0.781%
5	6.335	841	850	862	rVB2	38902	113020	2.33%	0.207%
6	6.860	924	936	960	rBV	367834	924585	19.04%	1.691%
7	7.061	962	969	978	rVB	35000	80234	1.65%	0.147%
8	8.055	1123	1132	1143	rVB	33037	70520	1.45%	0.129%
9	8.177	1144	1152	1161	rBV2	44626	93058	1.92%	0.170%
10	8.713	1225	1240	1256	rBV	757246	1300537	26.78%	2.378%
11	10.115	1464	1470	1479	rBV	760807	1056314	21.75%	1.932%
12	10.329	1501	1505	1509	rBV2	31579	49398	1.02%	0.090%
13	11.018	1613	1618	1623	rBV	54620	73860	1.52%	0.135%
14	11.139	1632	1638	1643	rBV	679328	855454	17.62%	1.564%
15	11.920	1757	1766	1767	rBV	51474	67889	1.40%	0.124%
16	11.944	1767	1770	1775	rVB	191307	243341	5.01%	0.445%
17	12.078	1787	1792	1804	rVB	910900	1126327	23.19%	2.060%
18	12.231	1813	1817	1821	rVB	51679	61105	1.26%	0.112%
19	12.304	1821	1829	1835	rBV	348011	449580	9.26%	0.822%
20	12.371	1835	1840	1847	rVB	122161	171897	3.54%	0.314%
21	12.462	1847	1855	1861	rBV	61182	89980	1.85%	0.165%
22	12.670	1870	1889	1892	rBV	565516	744356	15.33%	1.361%
23	12.700	1892	1894	1899	rVV	265323	299899	6.18%	0.548%
24	12.755	1899	1903	1911	rVB2	517742	847582	17.45%	1.550%
25	12.895	1919	1926	1931	rVB2	117165	193036	3.97%	0.353%
26	12.981	1931	1940	1945	rBV	658064	884176	18.21%	1.617%
27	13.078	1952	1956	1963	rVB2	180176	277990	5.72%	0.508%
28	13.151	1963	1968	1972	rBV2	118386	204848	4.22%	0.375%
29	13.194	1972	1975	1977	rBV2	208120	209546	4.31%	0.383%
30	13.224	1977	1980	1983	rVB	351416	365270	7.52%	0.668%
31	13.316	1990	1995	1997	rBV2	154468	214286	4.41%	0.392%
32	13.352	1997	2001	2007	rVV2	1385690	2013267	41.46%	3.682%
33	13.426	2007	2013	2019	rVV2	265952	498465	10.26%	0.912%
34	13.480	2019	2022	2026	rVB	59798	81732	1.68%	0.149%

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Integrator: RTE
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35	13.566	2031	2036	2038	rBV	324080	407692	8.40%	0.746%
36	13.602	2038	2042	2045	rVV	676599	914933	18.84%	1.673%
37	13.639	2045	2048	2052	rVV2	779112	1097985	22.61%	2.008%
38	13.724	2052	2062	2070	rVB3	924669	2157300	44.42%	3.945%
39	13.810	2070	2076	2079	rBV	135404	181226	3.73%	0.331%
40	13.852	2079	2083	2088	rBV2	161670	221447	4.56%	0.405%
41	13.913	2088	2093	2097	rVB	538833	720791	14.84%	1.318%
42	13.986	2097	2105	2109	rBV2	777487	1183463	24.37%	2.164%
43	14.029	2109	2112	2116	rBV	398483	467086	9.62%	0.854%
44	14.066	2116	2118	2121	rVB	94638	94510	1.95%	0.173%
45	14.133	2124	2129	2133	rBV	924040	1112592	22.91%	2.035%
46	14.218	2141	2143	2147	rVB	54180	53494	1.10%	0.098%
47	14.273	2147	2152	2161	rVV2	1047288	1951868	40.19%	3.570%
48	14.346	2161	2164	2166	rVV	59478	64975	1.34%	0.119%
49	14.377	2166	2169	2174	rVV	366135	458578	9.44%	0.839%
50	14.432	2174	2178	2182	rVB	887218	1069160	22.02%	1.955%
51	14.480	2182	2186	2191	rVB	161869	225765	4.65%	0.413%
52	14.541	2191	2196	2201	rBV2	812984	1130503	23.28%	2.067%
53	14.584	2201	2203	2209	rVB2	68229	84586	1.74%	0.155%
54	14.675	2209	2218	2219	rBV4	481404	692676	14.26%	1.267%
55	14.700	2219	2222	2225	rVV	964102	1282387	26.41%	2.345%
56	14.730	2225	2227	2232	rVB	314028	392662	8.09%	0.718%
57	14.785	2232	2236	2240	rBV2	166322	215725	4.44%	0.395%
58	14.840	2240	2245	2249	rVV	2947569	3712110	76.44%	6.789%
59	14.877	2249	2251	2254	rVV2	210163	272345	5.61%	0.498%
60	14.907	2254	2256	2260	rVV2	133303	158746	3.27%	0.290%
61	14.968	2260	2266	2272	rVV3	116026	257173	5.30%	0.470%
62	15.041	2273	2278	2282	rVB5	51860	117159	2.41%	0.214%
63	15.114	2283	2290	2294	rBV	72625	128168	2.64%	0.234%
64	15.248	2307	2312	2320	rVB	388252	671923	13.84%	1.229%
65	15.328	2320	2325	2332	rVB4	118176	216776	4.46%	0.396%
66	15.444	2336	2344	2347	rBV	706805	1074631	22.13%	1.965%
67	15.480	2347	2350	2355	rVB	451687	609296	12.55%	1.114%
68	15.553	2355	2362	2367	rBV	1934675	3134086	64.54%	5.732%
69	15.694	2374	2385	2388	rBV	2904296	4856260	100.00%	8.881%
70	15.730	2388	2391	2398	rVB	1587520	2254375	46.42%	4.123%
71	15.809	2399	2404	2410	rBV2	115744	224441	4.62%	0.410%

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72	15.919	2411	2422	2432	rVB	781006	2044917	42.11%	3.740%
73	16.072	2441	2447	2458	rVB	485334	823918	16.97%	1.507%
74	16.169	2458	2463	2469	rBV	223953	372867	7.68%	0.682%
75	16.285	2472	2482	2487	rVB	168820	359751	7.41%	0.658%
76	16.352	2487	2493	2495	rBV	58424	99434	2.05%	0.182%
77	16.389	2495	2499	2501	rVV2	40421	61403	1.26%	0.112%
78	16.443	2501	2508	2513	rVV2	182793	365536	7.53%	0.668%
79	16.510	2516	2519	2524	rVB2	68154	116627	2.40%	0.213%
80	16.565	2524	2528	2533	rBV	71548	122238	2.52%	0.224%
81	16.651	2533	2542	2544	rBV5	81956	186666	3.84%	0.341%
82	16.693	2544	2549	2554	rVB	239171	419815	8.64%	0.768%
83	16.754	2554	2559	2564	rBV	193838	383373	7.89%	0.701%

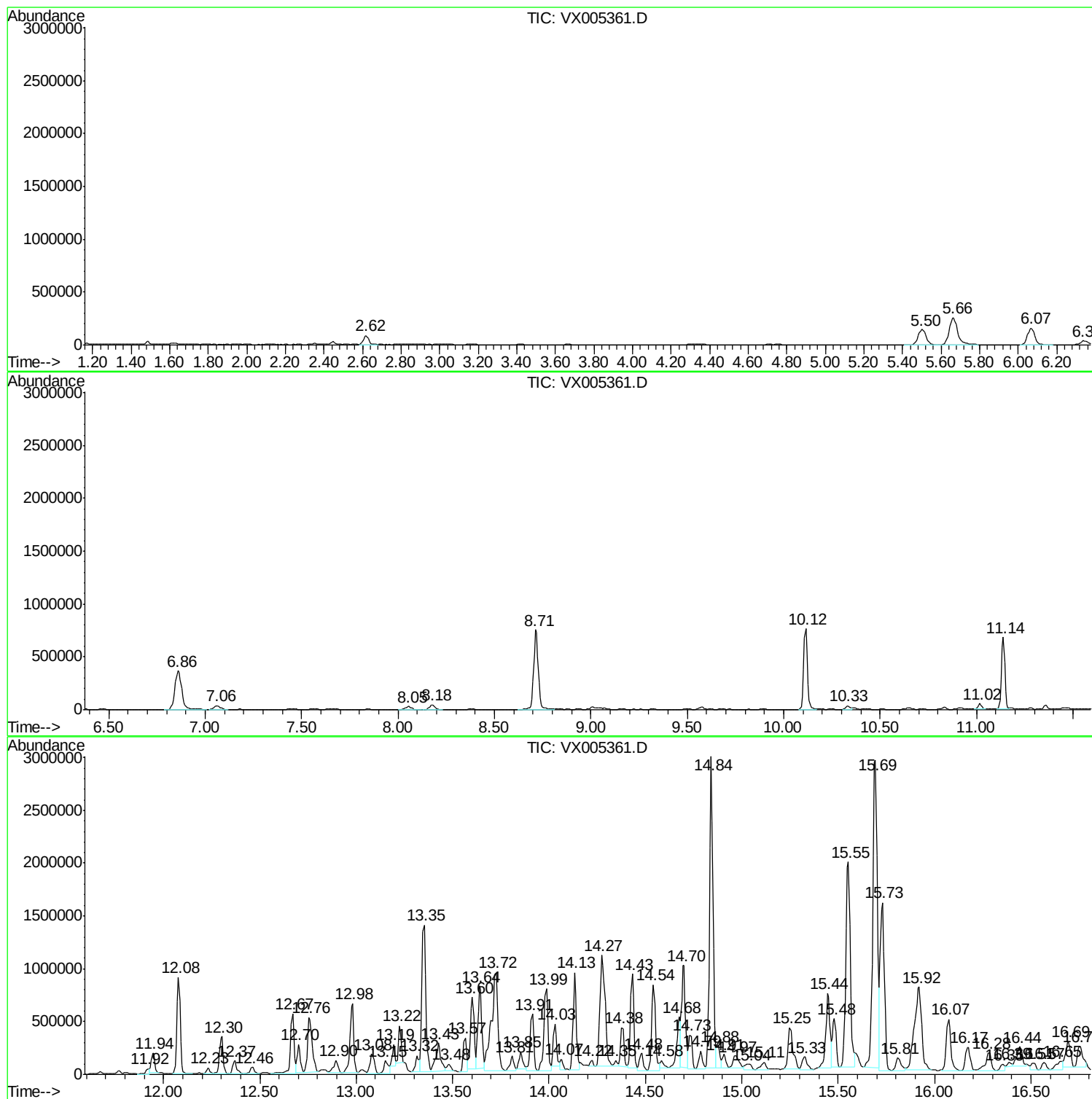
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Instrument :
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TWP-01

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.M
Quant Title : SW846 8260

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



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Instrument :
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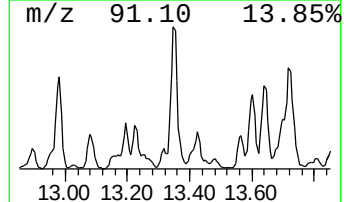
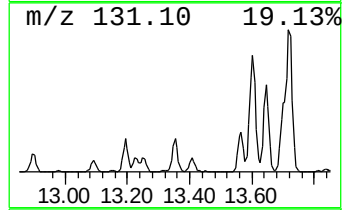
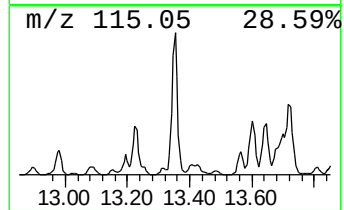
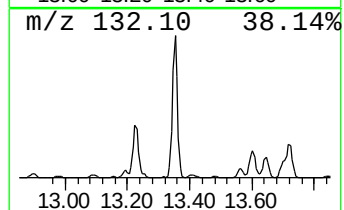
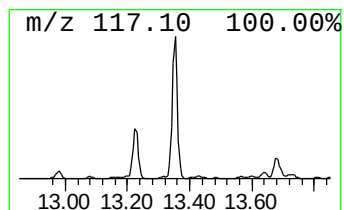
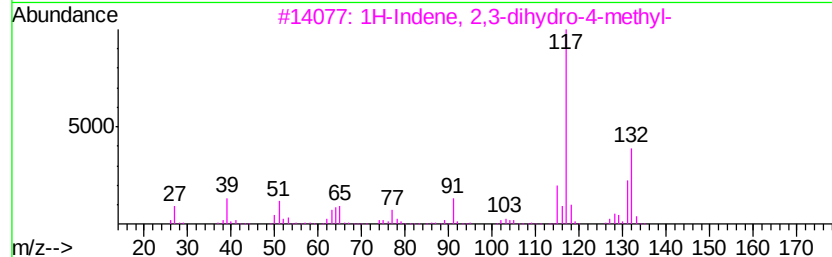
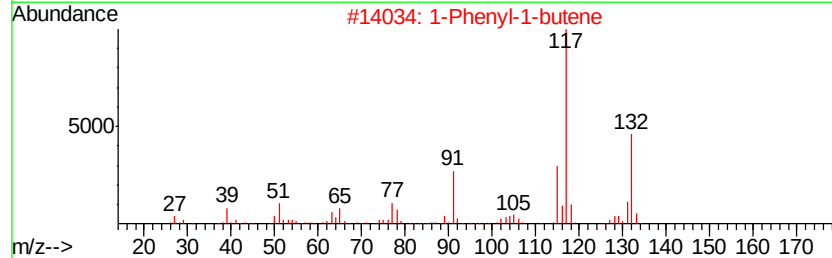
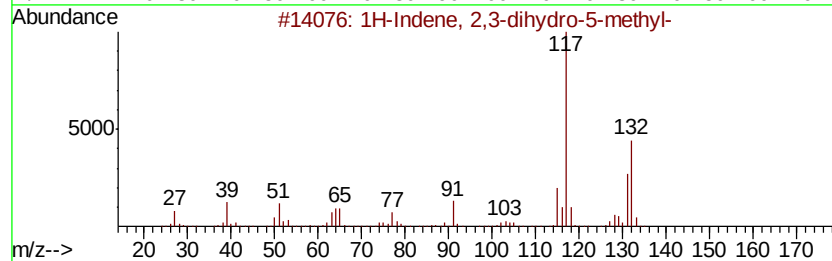
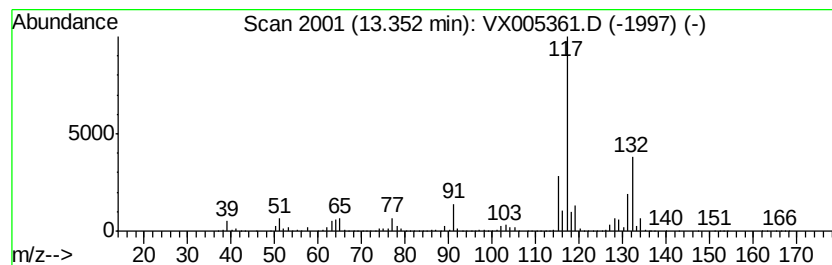
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.M
Quant Title : SW846 8260

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

Peak Number 1 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	89.37 ug/l	2013270	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
2		1-Phenyl-1-butene	132	C10H12	000824-90-8	87
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	86
5		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	83



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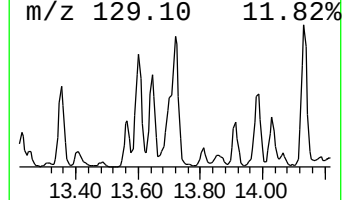
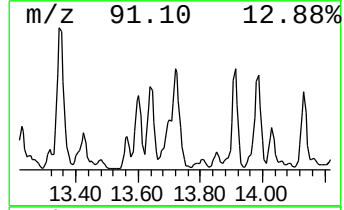
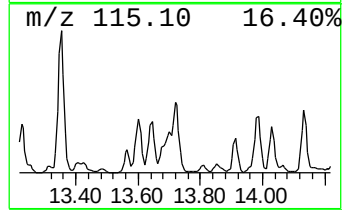
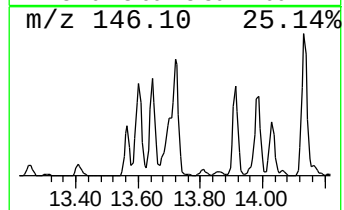
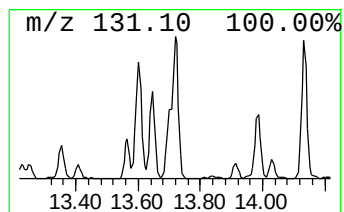
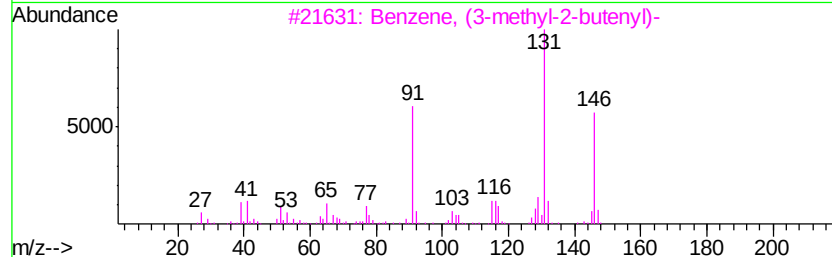
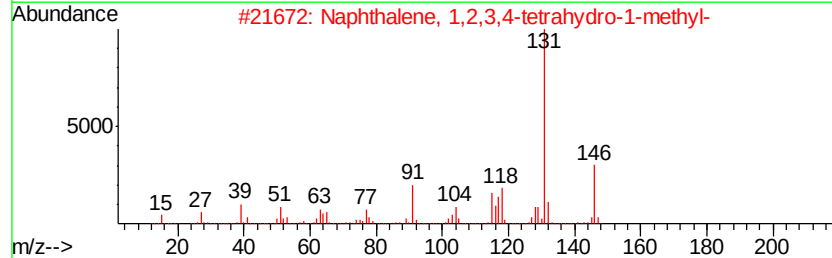
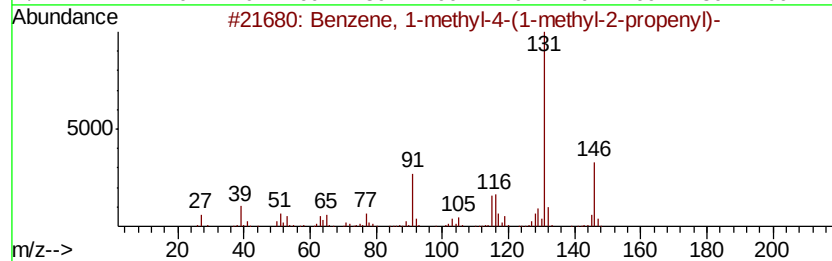
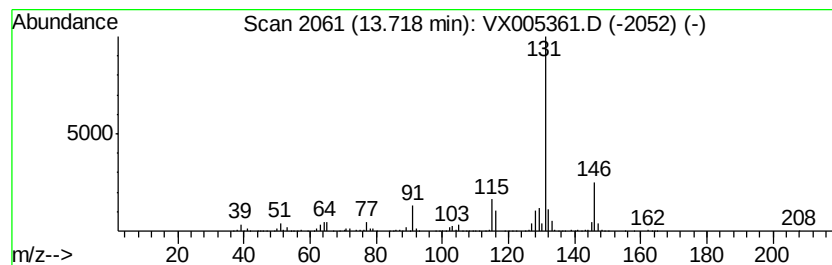
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1-methyl-4-(1-meth... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	95.77 ug/l	2157300	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-methyl-4-(1-methyl-2-...	146 C11H14	097664-18-1 90
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001559-81-5 90
3		Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3 90
4		2,2-Dimethylindene, 2,3-dihydro-	146 C11H14	020836-11-7 90
5		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 87



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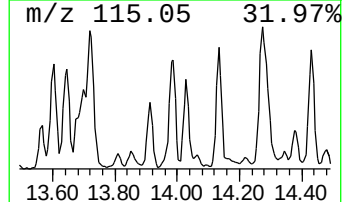
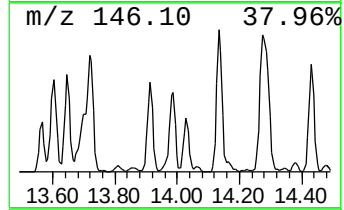
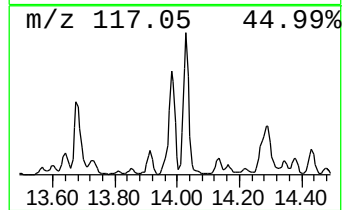
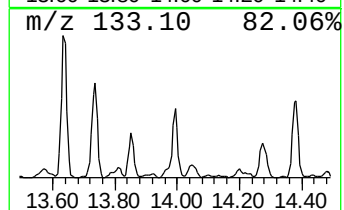
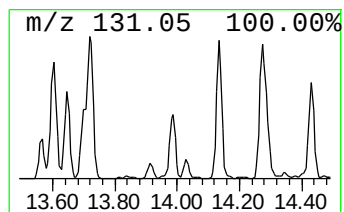
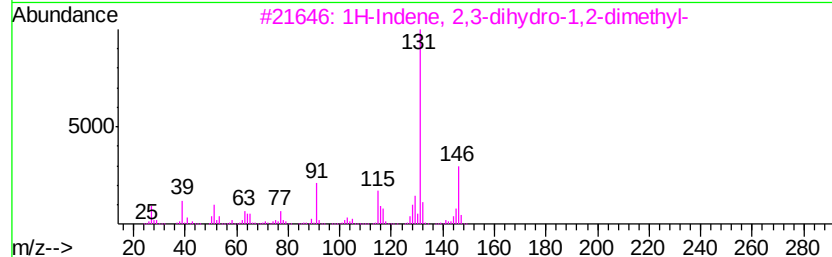
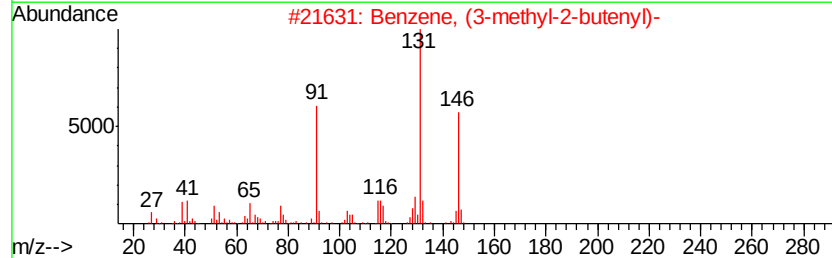
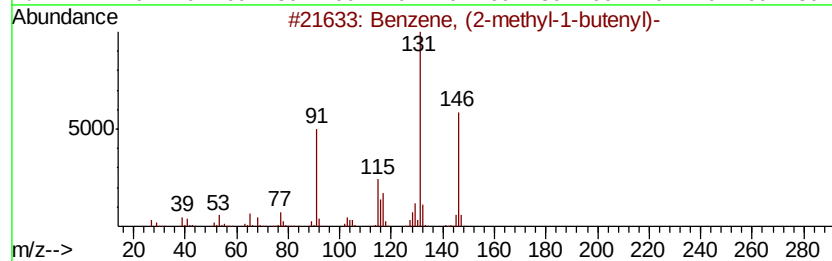
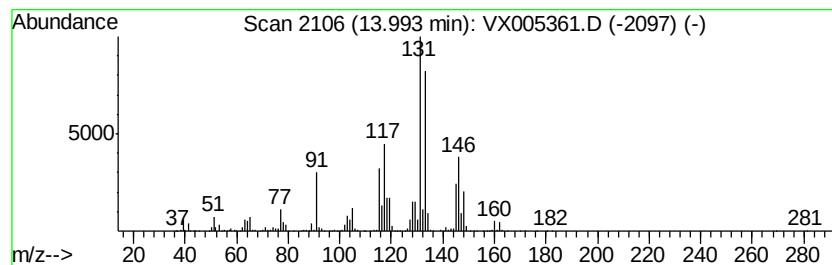
Instrument :
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ClientSampled :
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, (2-methyl-1-butenyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	52.54 ug/l	1183460	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (2-methyl-1-butenyl)-	146 C11H14	056253-64-6 90
2		Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3 60
3		1H-Indene, 2,3-dihydro-1,2-dimet...	146 C11H14	017057-82-8 60
4		Benzene, (1-methyl-1-butenyl)-	146 C11H14	053172-84-2 60
5		1H-Indene, 2,3-dihydro-4,6-dimet...	146 C11H14	001685-82-1 60



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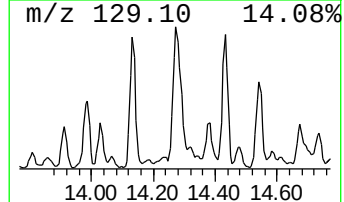
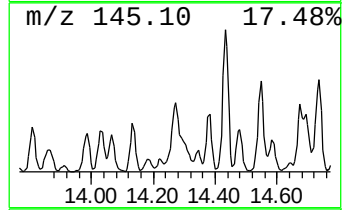
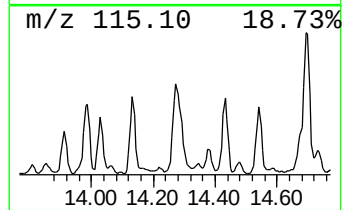
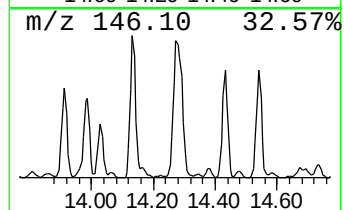
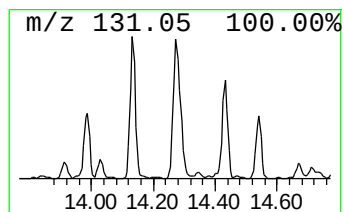
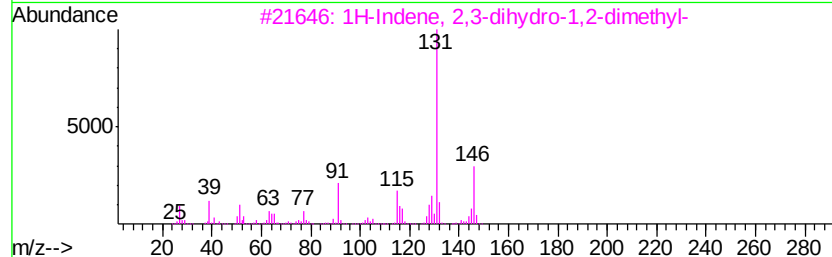
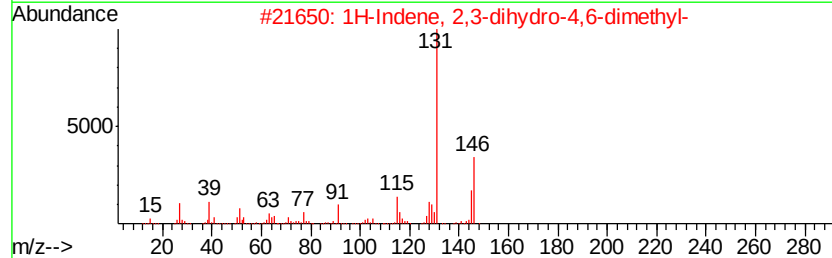
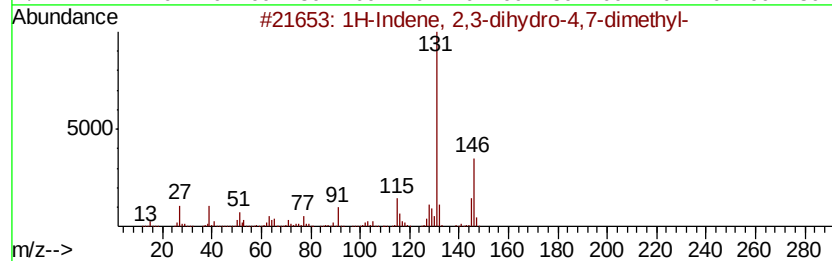
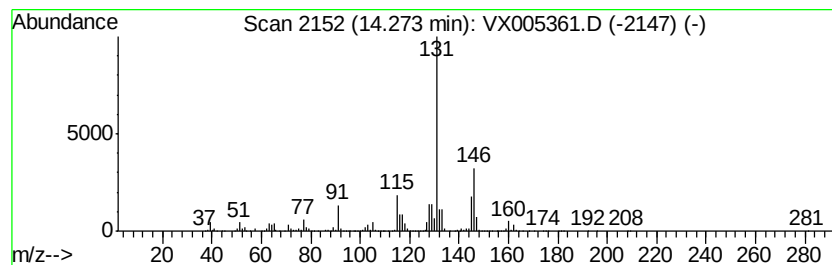
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Quant Title : SW846 8260

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

Peak Number 4 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.27	86.65 ug/l	1951870	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	95
2		1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	95
3		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91
4		1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	91
5		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	87



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX101618\
Data File : VX005361.D
Acq On : 16 Oct 2018 11:00
Operator : JC/MD
Sample : J5521-13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TWP-01

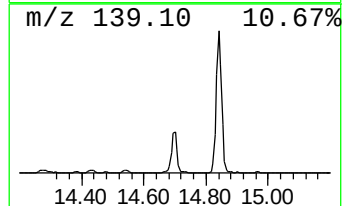
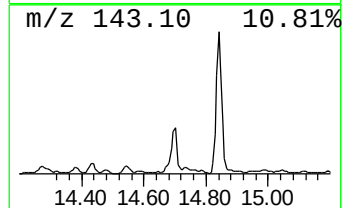
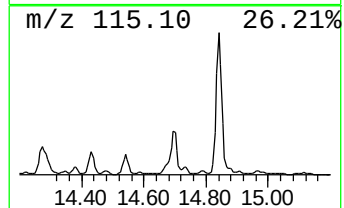
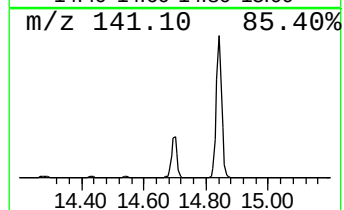
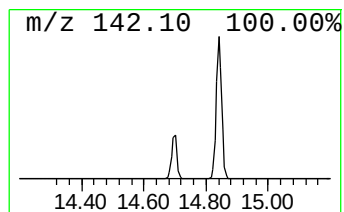
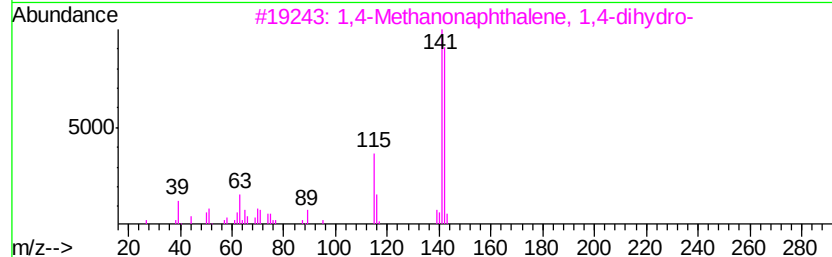
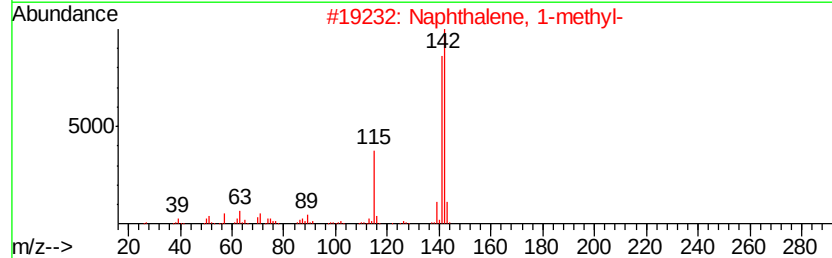
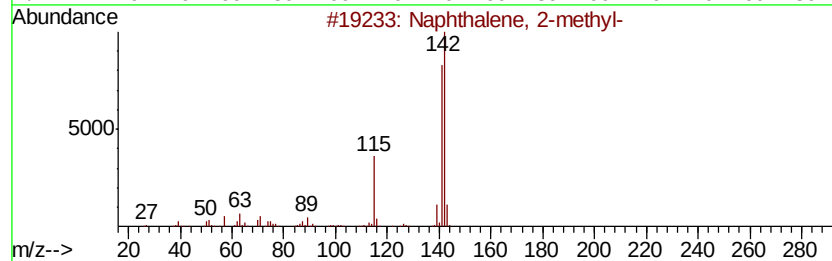
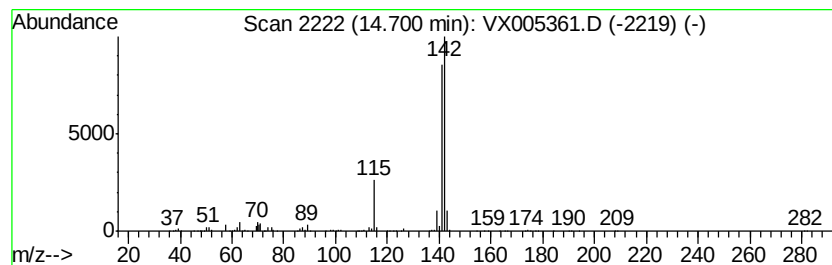
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Quant Title : SW846 8260

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

Peak Number 5 Naphthalene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.70	56.93 ug/l	1282390	1,4-Dichlorobenzene-d4	12.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
3		1,4-Methanonaphthalene, 1,4-dihv...	142	C11H10	004453-90-1	91
4		Benzocycloheptatriene	142	C11H10	000264-09-5	90
5		3,4-Difluorobenzaldehyde	142	C7H4F2O	034036-07-2	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX101618\
Data File : VX005361.D
Acq On : 16 Oct 2018 11:00
Operator : JC/MD
Sample : J5521-13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TWP-01

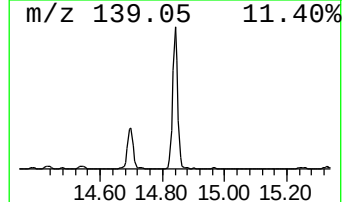
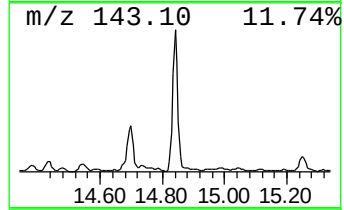
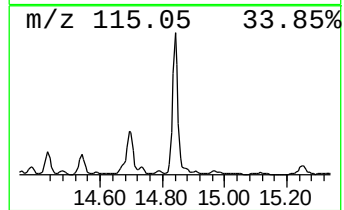
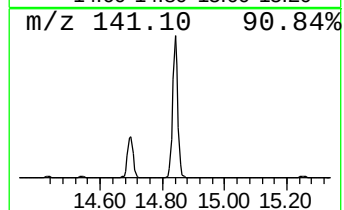
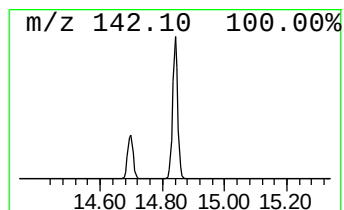
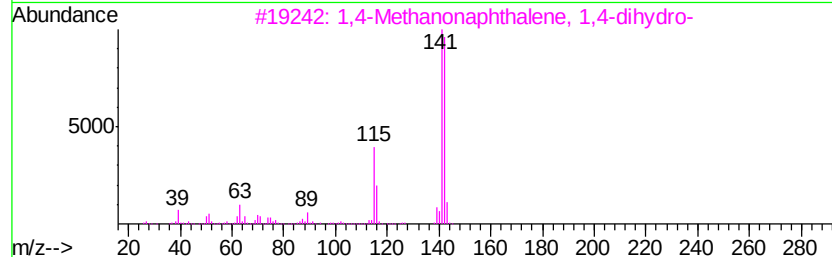
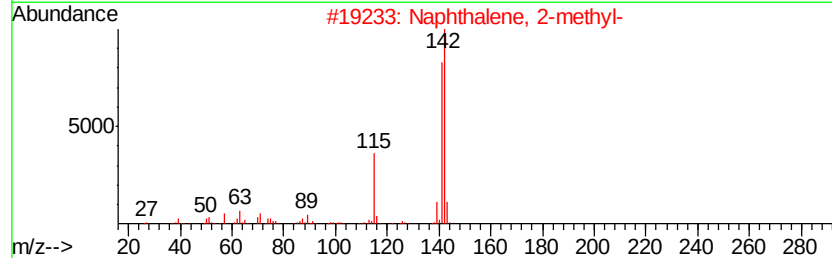
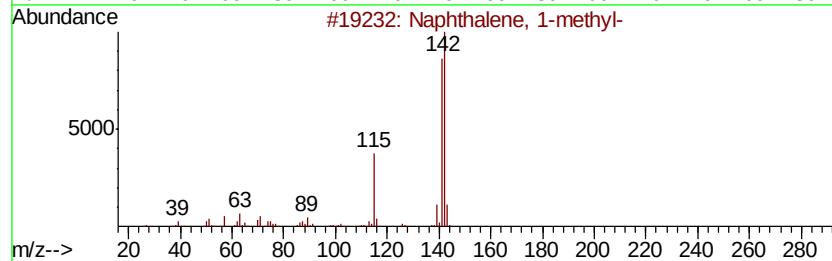
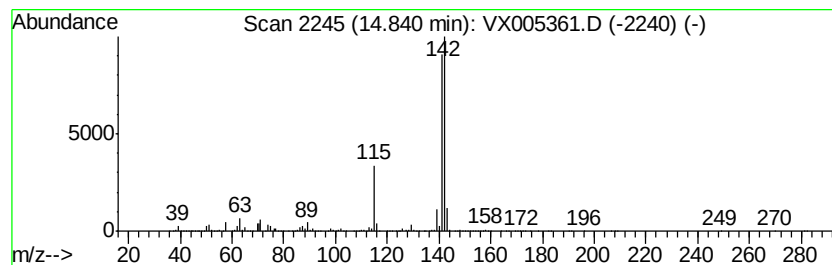
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Quant Title : SW846 8260

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

Peak Number 6 1,4-Methanonaphthalene, 1,4... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.84	164.79 ug/l	3712110	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	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	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX101618\
Data File : VX005361.D
Acq On : 16 Oct 2018 11:00
Operator : JC/MD
Sample : J5521-13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

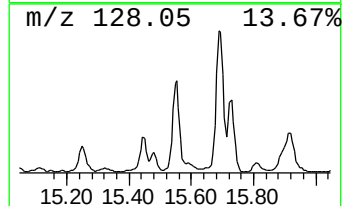
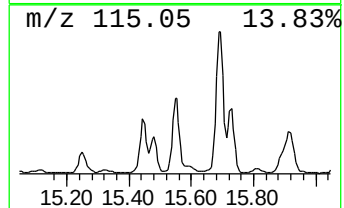
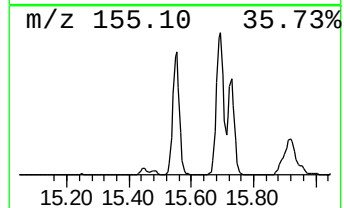
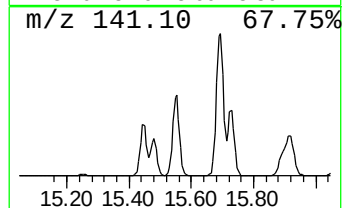
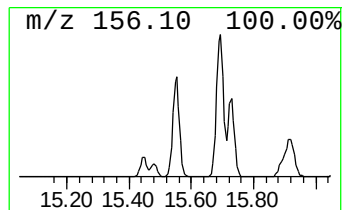
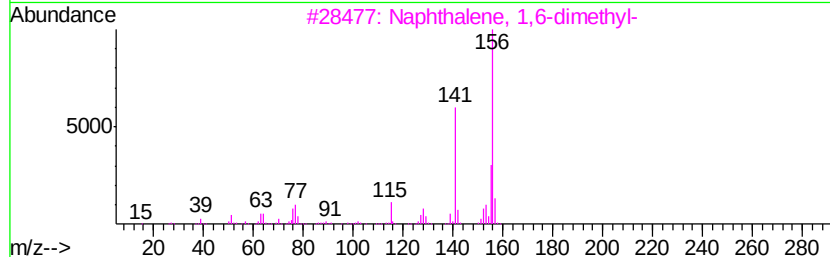
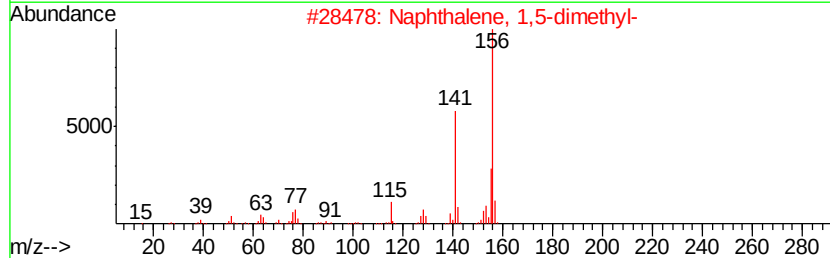
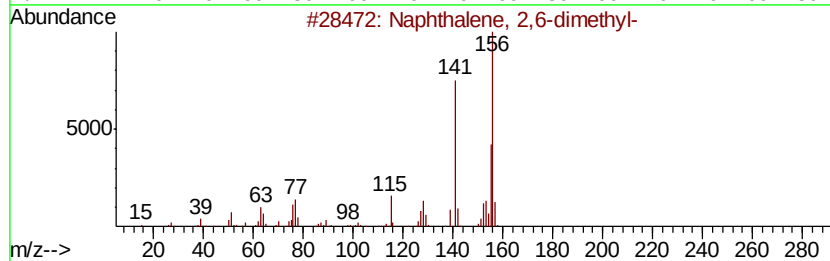
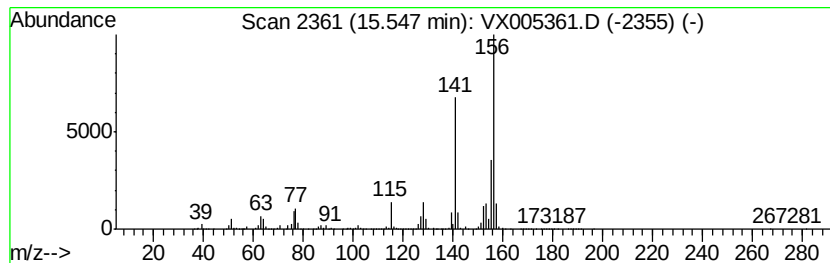
Instrument :
MSVOA_X
ClientSampled :
TWP-01

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.M
Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 2,6-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.55	139.13 ug/l	3134090	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 98
2		Naphthalene, 1,5-dimethyl-	156 C12H12	000571-61-9 97
3		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 97
4		Naphthalene, 1,7-dimethyl-	156 C12H12	000575-37-1 97
5		Naphthalene, 2,7-dimethyl-	156 C12H12	000582-16-1 97



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX101618\
Data File : VX005361.D
Acq On : 16 Oct 2018 11:00
Operator : JC/MD
Sample : J5521-13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TWP-01

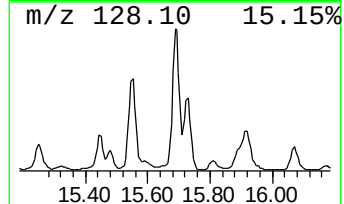
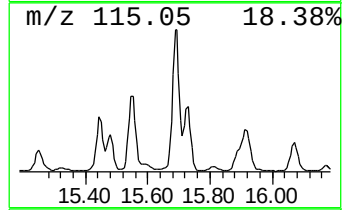
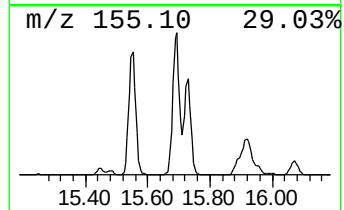
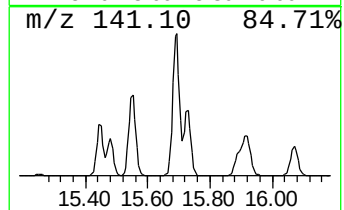
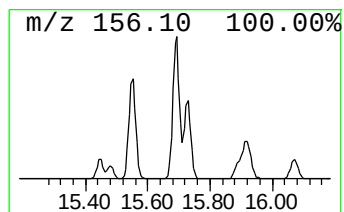
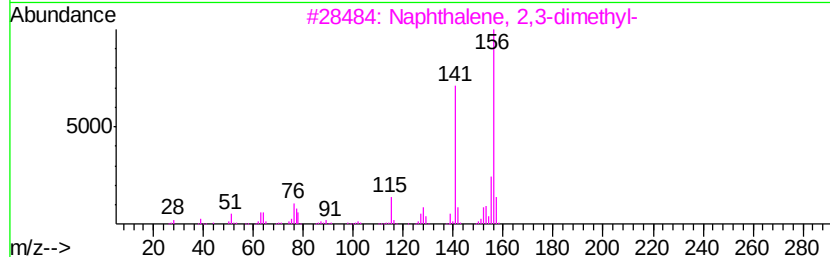
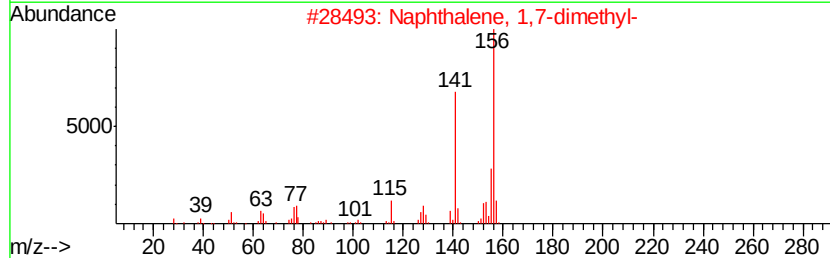
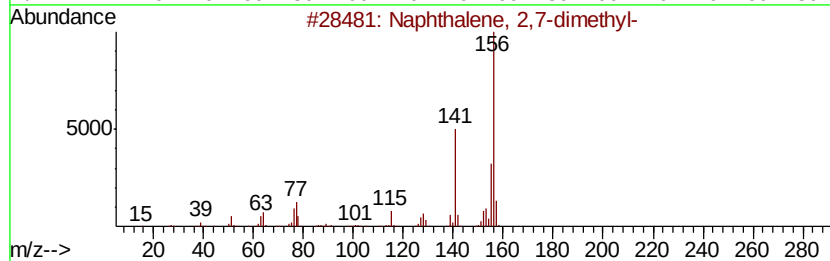
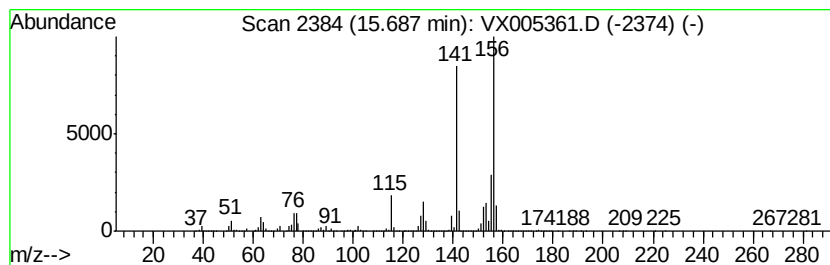
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Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 2,7-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.69	215.58 ug/l	4856260	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
2		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
3		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
5		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX101618\
Data File : VX005361.D
Acq On : 16 Oct 2018 11:00
Operator : JC/MD
Sample : J5521-13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TWP-01

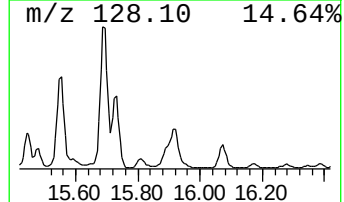
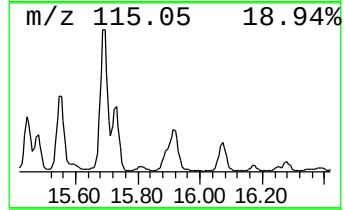
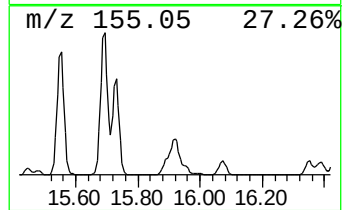
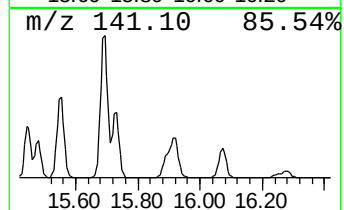
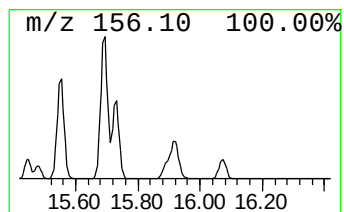
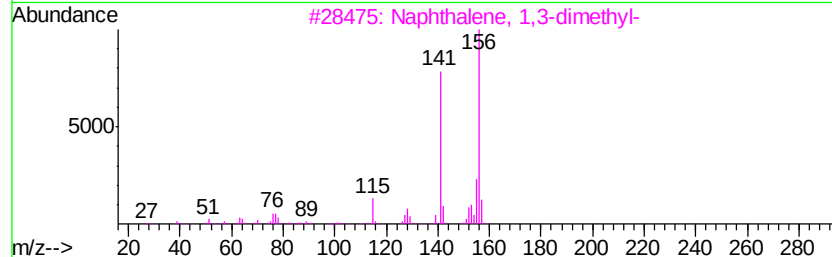
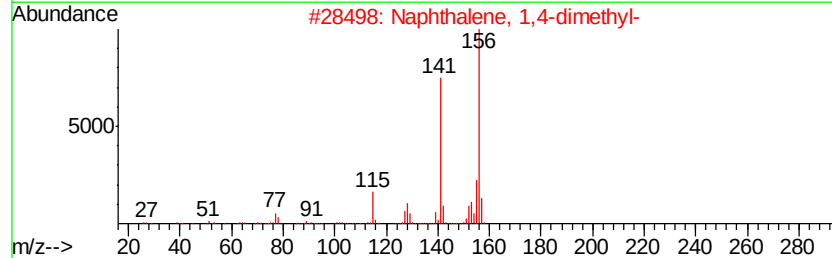
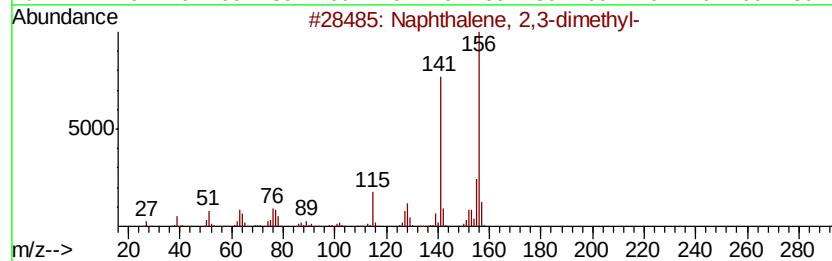
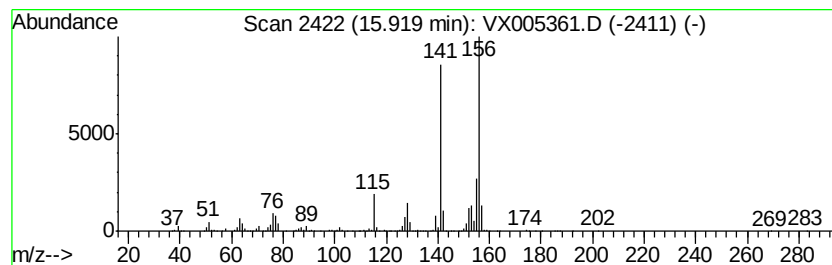
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 2,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.92	90.78 ug/l	2044920	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, 1,4-dimethyl-	156	C12H12	000571-58-4	97
3		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
4		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5		Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	96



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX101618\
Data File : VX005361.D
Acq On : 16 Oct 2018 11:00
Operator : JC/MD
Sample : J5521-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TWP-01

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.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
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Benzene, 1-methyl...	13.72	95.8	ug/l	2157300	4	12.08	1126330	50.0
Benzene, (2-methy...	13.99	52.5	ug/l	1183460	4	12.08	1126330	50.0
1H-Indene, 2,3-di...	14.27	86.7	ug/l	1951870	4	12.08	1126330	50.0
Naphthalene, 1-me...	14.70	56.9	ug/l	1282390	4	12.08	1126330	50.0
1,4-Methanonaphth...	14.84	164.8	ug/l	3712110	4	12.08	1126330	50.0
Naphthalene, 2,6-...	15.55	139.1	ug/l	3134090	4	12.08	1126330	50.0
Naphthalene, 2,7-...	15.69	215.6	ug/l	4856260	4	12.08	1126330	50.0
Naphthalene, 2,3-...	15.92	90.8	ug/l	2044920	4	12.08	1126330	50.0