

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX101618\
 Data File : VX005364.D
 Acq On : 16 Oct 2018 12:20
 Operator : JC/MD
 Sample : J5519-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BS-DRUM

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	1.642	67	80	84	rBV4	14981	42495	3.16%	0.208%
2	2.617	234	240	250	rVB	42038	77431	5.76%	0.379%
3	3.019	299	306	311	rBV2	10275	23232	1.73%	0.114%
4	4.769	581	593	603	rBV2	6708	22805	1.70%	0.112%
5	5.501	700	713	728	rBV	158956	450572	33.51%	2.204%
6	5.659	728	739	759	rBV	280977	817888	60.82%	4.001%
7	6.068	793	806	821	rBV	177198	457517	34.02%	2.238%
8	6.860	925	936	956	rBV	415694	991430	73.72%	4.850%
9	8.713	1233	1240	1265	rBV	833305	1344782	100.00%	6.578%
10	10.116	1463	1470	1489	rBV	799534	1126947	83.80%	5.513%
11	11.140	1630	1638	1653	rBV	686843	893172	66.42%	4.369%
12	11.433	1682	1686	1693	rBV	43701	68688	5.11%	0.336%
13	11.506	1693	1698	1707	rVB	59038	75029	5.58%	0.367%
14	11.670	1720	1725	1731	rVB	37349	49595	3.69%	0.243%
15	11.810	1743	1748	1754	rVB	68710	88950	6.61%	0.435%
16	12.024	1779	1783	1787	rBV	20350	25532	1.90%	0.125%
17	12.079	1787	1792	1798	rVV	957429	1154028	85.82%	5.645%
18	12.140	1798	1802	1808	rVB	135357	172844	12.85%	0.845%
19	12.298	1820	1828	1830	rBV3	90080	127591	9.49%	0.624%
20	12.323	1830	1832	1836	rVV	95660	105345	7.83%	0.515%
21	12.371	1836	1840	1848	rVB2	205266	269295	20.03%	1.317%
22	12.463	1848	1855	1860	rBV	18524	25439	1.89%	0.124%
23	12.518	1860	1864	1870	rVB	53786	64909	4.83%	0.318%
24	12.585	1870	1875	1877	rBV	111273	136374	10.14%	0.667%
25	12.609	1877	1879	1883	rVB	92071	99841	7.42%	0.488%
26	12.670	1883	1889	1892	rBV	240261	288122	21.43%	1.409%
27	12.701	1892	1894	1898	rVB	51062	53222	3.96%	0.260%
28	12.755	1898	1903	1911	rBV2	166251	272500	20.26%	1.333%
29	12.853	1915	1919	1922	rVB2	17144	19233	1.43%	0.094%
30	12.902	1922	1927	1932	rVB2	103842	135120	10.05%	0.661%
31	12.975	1932	1939	1943	rBV	173872	232197	17.27%	1.136%
32	13.018	1943	1946	1952	rVB	335296	389277	28.95%	1.904%
33	13.085	1952	1957	1958	rBV	45067	61311	4.56%	0.300%
34	13.103	1958	1960	1962	rVV2	46090	48865	3.63%	0.239%

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Integration Parameters: RTEINT.P

Integrator: RTE
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35	13.127	1962	1964	1966	rVB	29573	26114	1.94%	0.128%
36	13.194	1971	1975	1976	rVV	50190	55273	4.11%	0.270%
37	13.225	1976	1980	1984	rVV	196650	265863	19.77%	1.301%
38	13.268	1984	1987	1990	rVB3	33076	41129	3.06%	0.201%
39	13.310	1990	1994	1996	rBV2	60049	82058	6.10%	0.401%
40	13.347	1996	2000	2006	rVV2	740201	1045735	77.76%	5.115%
41	13.426	2010	2013	2017	rVV	51830	65735	4.89%	0.322%
42	13.481	2017	2022	2030	rVB	469941	584846	43.49%	2.861%
43	13.560	2030	2035	2038	rBV2	68467	99039	7.36%	0.484%
44	13.603	2038	2042	2045	rVV	165056	236530	17.59%	1.157%
45	13.639	2045	2048	2053	rVV3	211243	390016	29.00%	1.908%
46	13.700	2054	2058	2059	rVV	113104	158225	11.77%	0.774%
47	13.719	2059	2061	2067	rVB2	226410	318414	23.68%	1.558%
48	13.810	2073	2076	2079	rVB	15594	16339	1.21%	0.080%
49	13.853	2079	2083	2088	rBV2	30876	42448	3.16%	0.208%
50	13.914	2088	2093	2098	rVB	231561	286957	21.34%	1.404%
51	13.987	2098	2105	2109	rBV2	318545	446344	33.19%	2.183%
52	14.030	2109	2112	2115	rVV	114135	132819	9.88%	0.650%
53	14.066	2115	2118	2124	rVB4	23346	36707	2.73%	0.180%
54	14.133	2124	2129	2133	rBV	221133	269204	20.02%	1.317%
55	14.286	2147	2154	2162	rVV	680615	1128961	83.95%	5.522%
56	14.347	2162	2164	2166	rVV	17241	17032	1.27%	0.083%
57	14.377	2166	2169	2174	rVV	109642	141856	10.55%	0.694%
58	14.432	2174	2178	2182	rVV2	235871	291760	21.70%	1.427%
59	14.475	2182	2185	2190	rVB2	33320	44961	3.34%	0.220%
60	14.542	2190	2196	2201	rBV2	586289	766104	56.97%	3.748%
61	14.584	2201	2203	2208	rVB3	22965	28806	2.14%	0.141%
62	14.676	2213	2218	2219	rBV	222271	260527	19.37%	1.274%
63	14.694	2219	2221	2224	rVV	214407	258935	19.25%	1.267%
64	14.731	2224	2227	2232	rVB	156689	195798	14.56%	0.958%
65	14.786	2232	2236	2241	rBV2	46063	60250	4.48%	0.295%
66	14.840	2241	2245	2248	rBV	110213	143419	10.66%	0.702%
67	14.877	2248	2251	2254	rVV	98731	143457	10.67%	0.702%
68	14.908	2254	2256	2261	rVB2	85809	94656	7.04%	0.463%
69	14.962	2261	2265	2272	rBV2	100974	192087	14.28%	0.940%
70	15.115	2283	2290	2295	rBV	39680	62555	4.65%	0.306%
71	15.249	2307	2312	2320	rBV2	230059	418110	31.09%	2.045%

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72	15.322	2320	2324	2332	rVB4	25778	49466	3.68%	0.242%
73	15.395	2332	2336	2339	rVB4	10073	15667	1.17%	0.077%
74	15.444	2339	2344	2348	rBV2	57931	92800	6.90%	0.454%
75	15.548	2354	2361	2366	rBV2	179720	302362	22.48%	1.479%
76	15.596	2366	2369	2374	rVV	66075	118796	8.83%	0.581%
77	15.688	2378	2384	2388	rVV	224526	383048	28.48%	1.874%
78	15.724	2388	2390	2398	rVB	127663	184996	13.76%	0.905%
79	15.920	2418	2422	2427	rVB3	40679	75185	5.59%	0.368%
80	16.072	2441	2447	2451	rBV	20839	36317	2.70%	0.178%
81	16.169	2458	2463	2467	rBV	19706	33388	2.48%	0.163%
82	16.389	2489	2499	2502	rBV10	8479	25454	1.89%	0.125%
83	16.444	2505	2508	2515	rVB	11824	20090	1.49%	0.098%
84	16.694	2544	2549	2554	rVB2	13798	23531	1.75%	0.115%
85	16.749	2554	2558	2564	rVB3	11529	21301	1.58%	0.104%

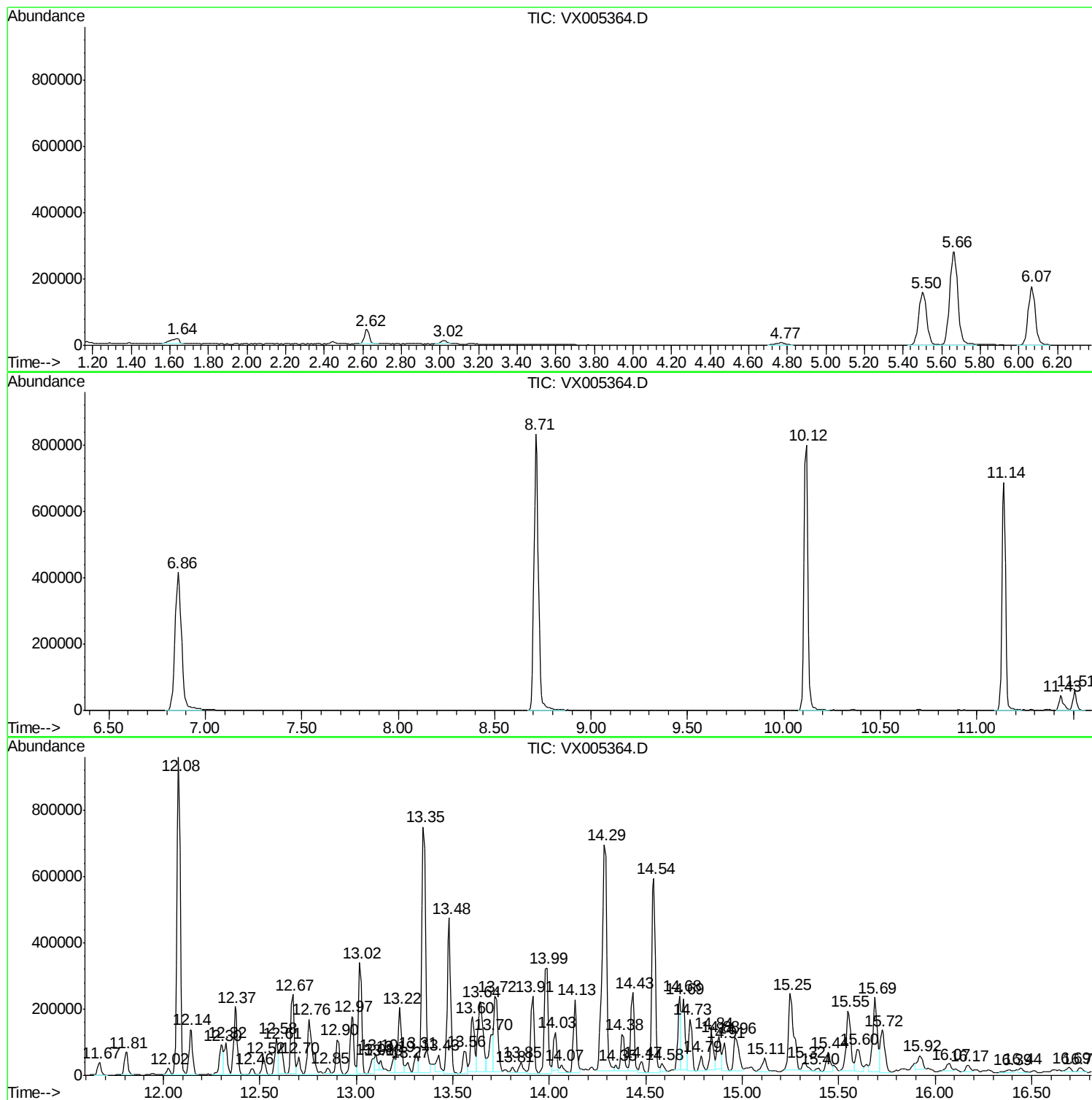
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.M
Quant Title : SW846 8260

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



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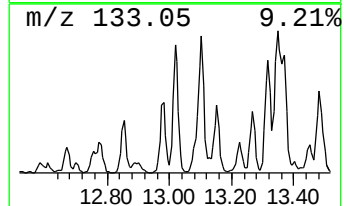
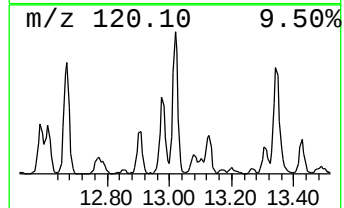
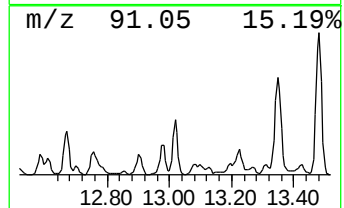
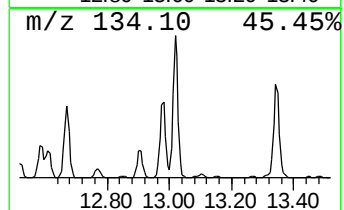
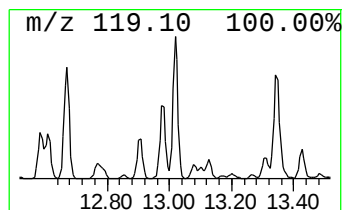
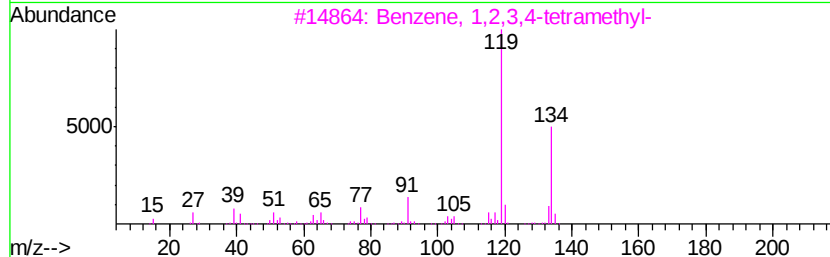
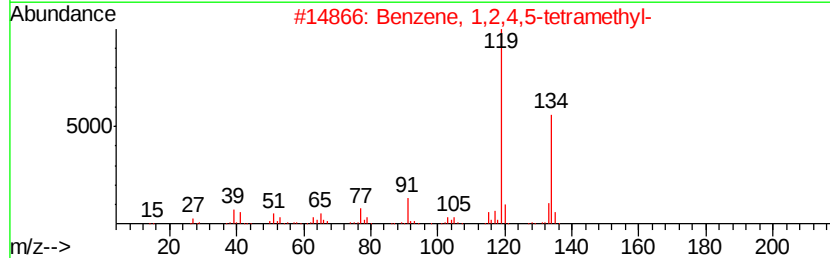
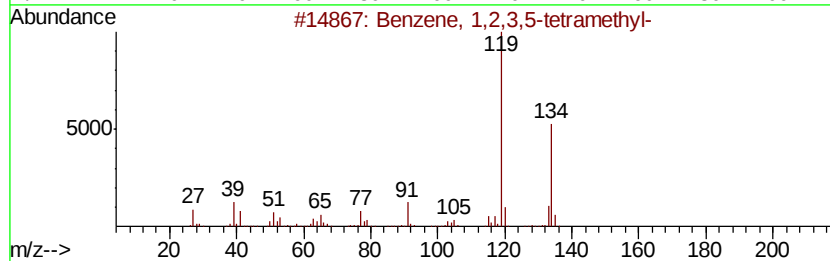
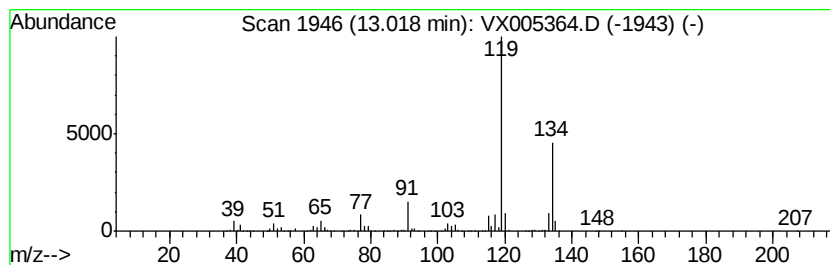
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	16.87 ug/l	389277	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



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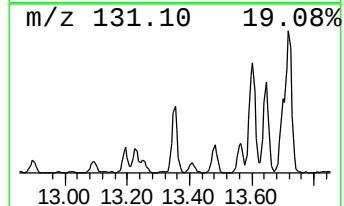
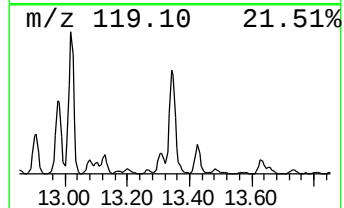
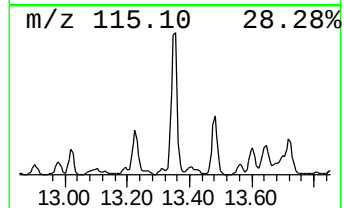
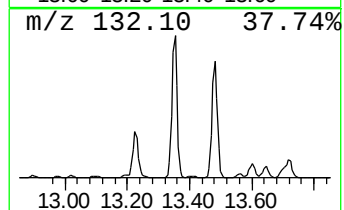
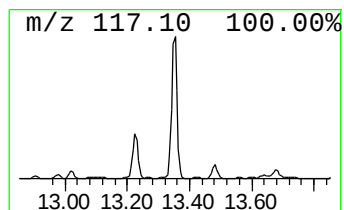
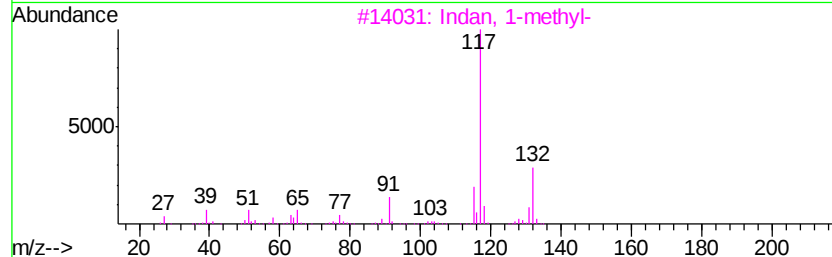
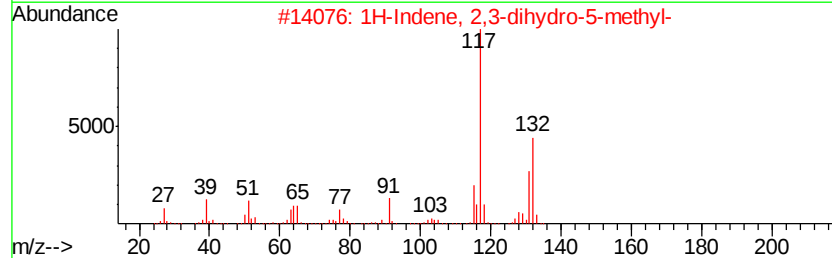
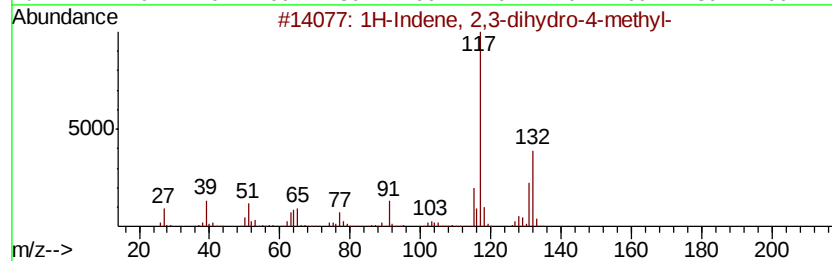
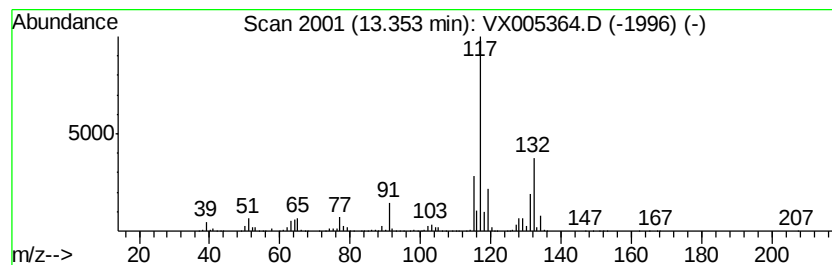
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	45.31 ug/l	1045740	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	92
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	86
3		Indan, 1-methyl-	132	C10H12	000767-58-8	70
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	64



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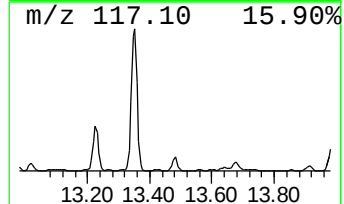
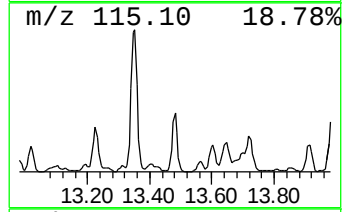
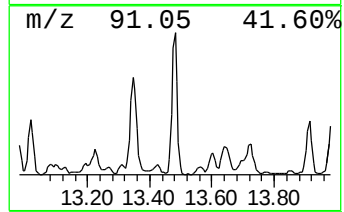
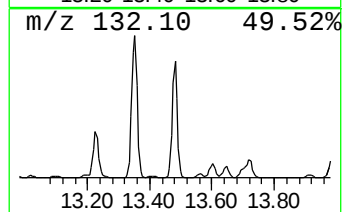
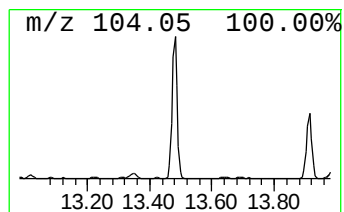
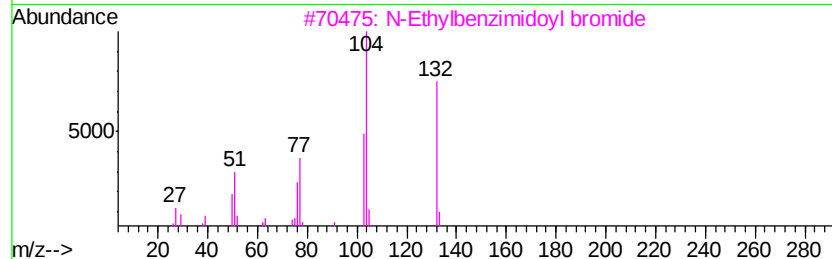
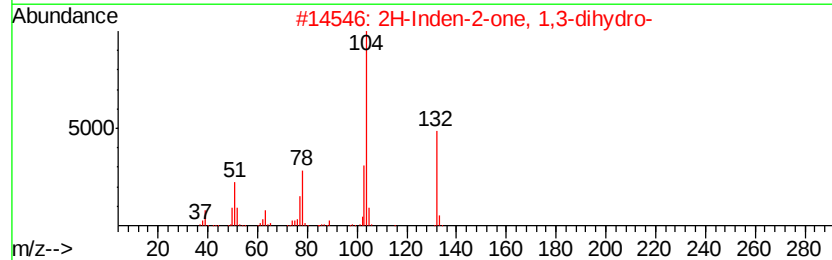
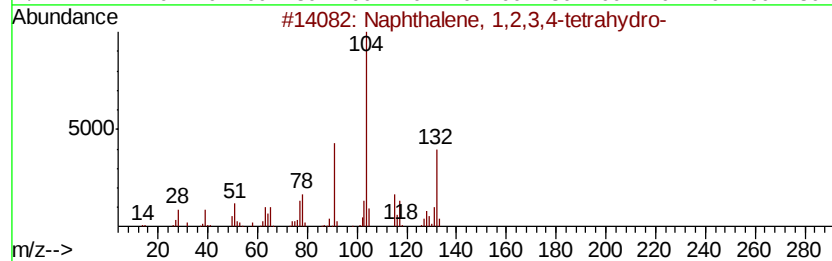
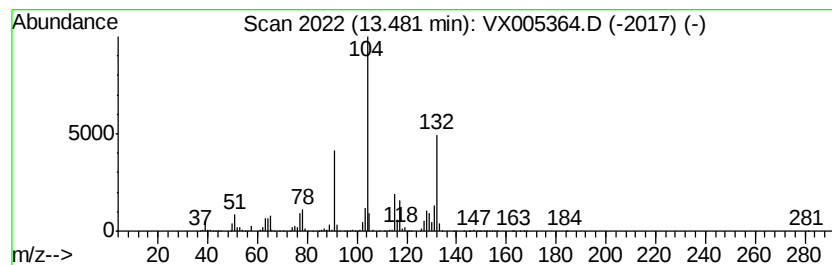
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.48	25.34 ug/l	584846	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2		2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	50
3		N-Ethylbenzimidoyl bromide	211	C9H10BrN	041182-87-0	40
4		2-Pheylethylamine, N,N-di(trifluoromethyl)-	313	C12H9F6N2	1000328-32-5	35
5		3(2H)-Furanone, dihydro-2,2-dimethyl-	190	C12H14O2	063678-00-2	35



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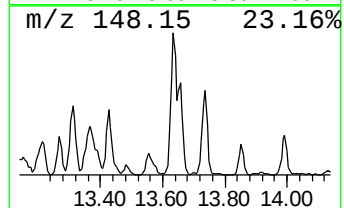
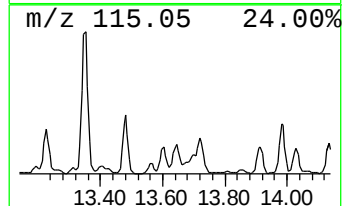
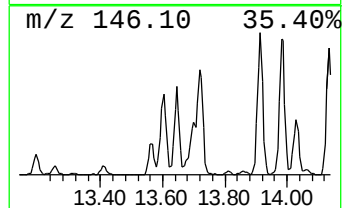
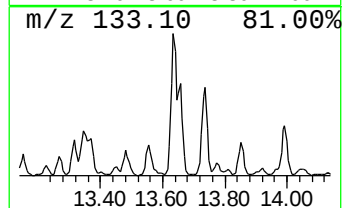
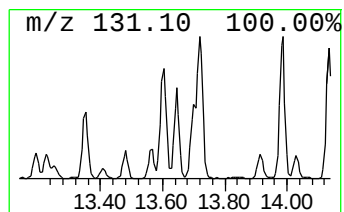
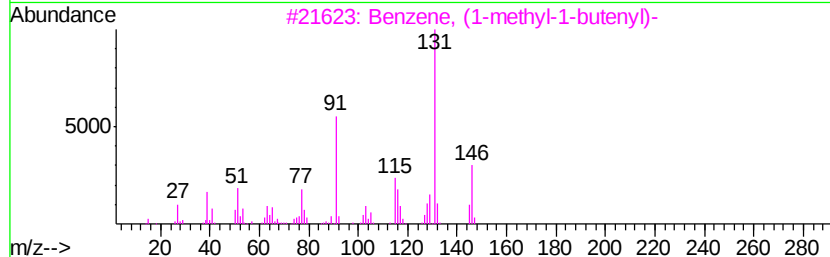
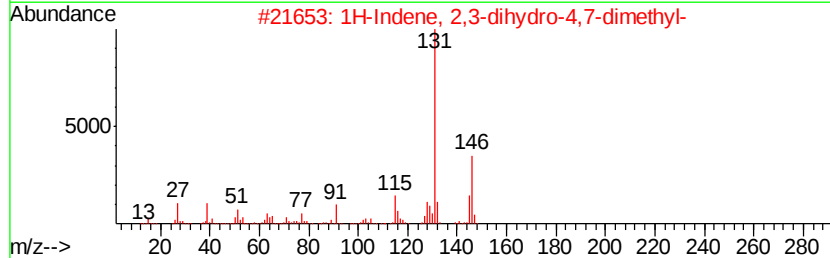
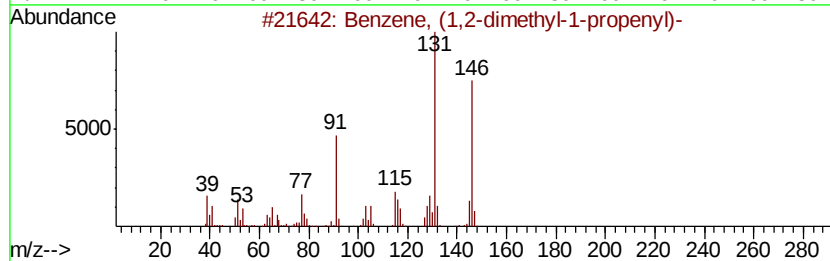
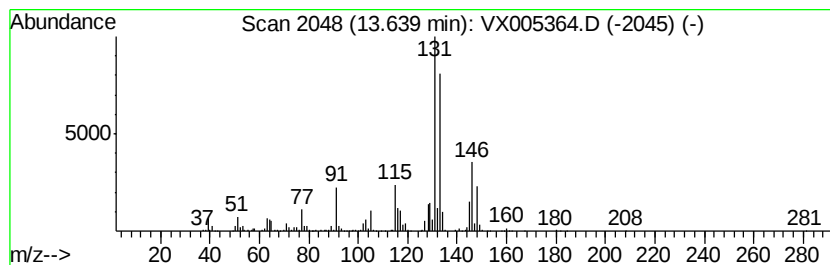
Instrument :
MSVOA_X
ClientSampleId :
BS-DRUM

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

Peak Number 4 Benzene, (1,2-dimethyl-1-propenyl)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.64	16.90 ug/l	390016	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 90
2		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 86
3		Benzene, (1-methyl-1-butenyl)-	146 C11H14	053172-84-2 70
4		1H-Indene, 2,3-dihydro-1,2-dimet...	146 C11H14	017057-82-8 70
5		Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3 70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX101618\
Data File : VX005364.D
Acq On : 16 Oct 2018 12:20
Operator : JC/MD
Sample : J5519-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 11 Sample Multiplier: 1

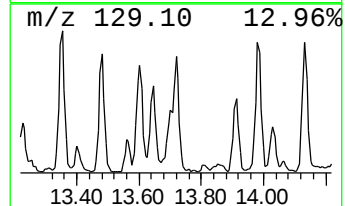
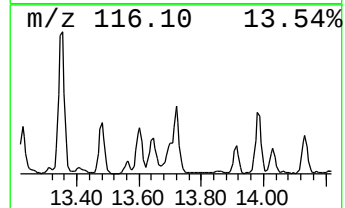
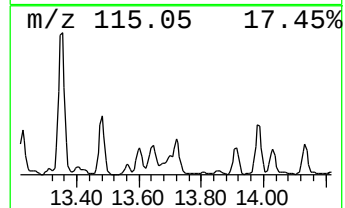
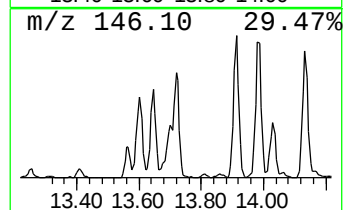
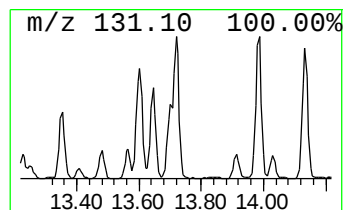
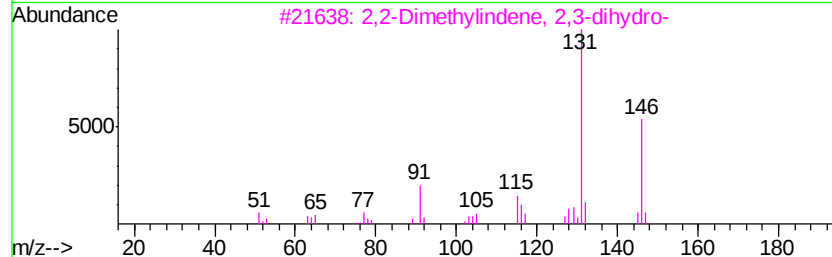
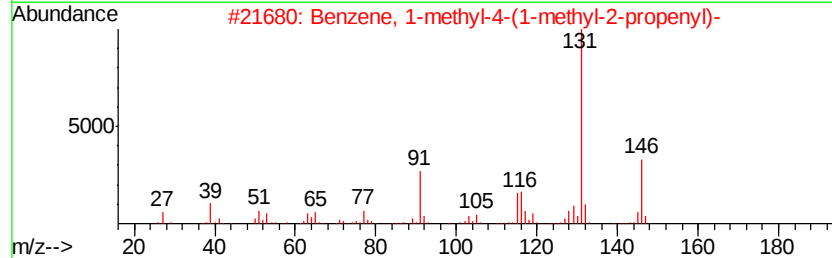
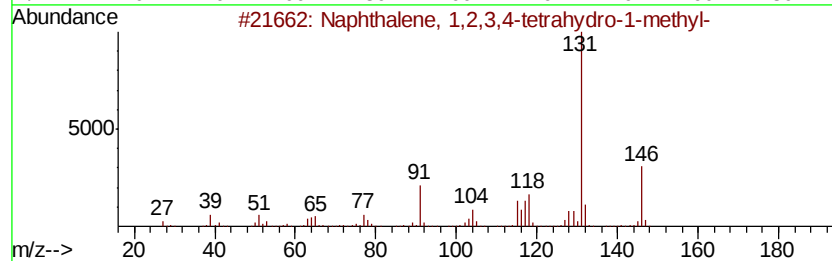
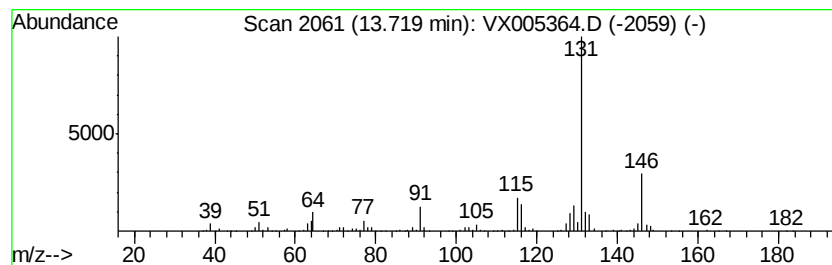
Instrument :
MSVOA_X
ClientSampleId :
BS-DRUM

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

Peak Number 5 Naphthalene, 1,2,3,4-tetra... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	13.80 ug/l	318414	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001559-81-5 90
2		Benzene, 1-methyl-4-(1-methyl-2-...	146 C11H14	097664-18-1 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		1H-Indene, 2,3-dihydro-1,2-dimet...	146 C11H14	017057-82-8 90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX101618\
Data File : VX005364.D
Acq On : 16 Oct 2018 12:20
Operator : JC/MD
Sample : J5519-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BS-DRUM

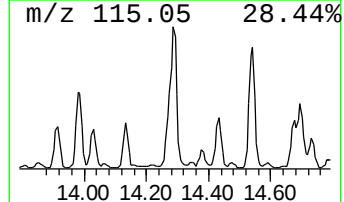
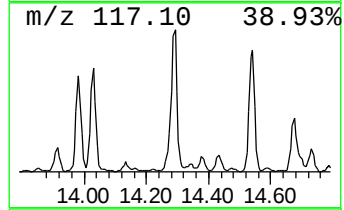
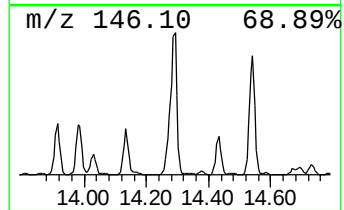
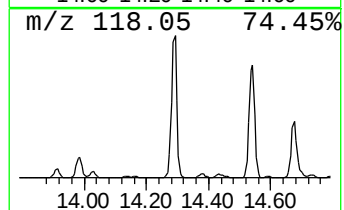
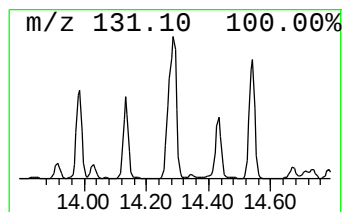
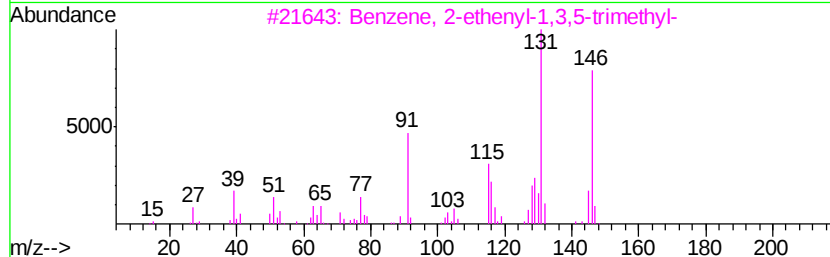
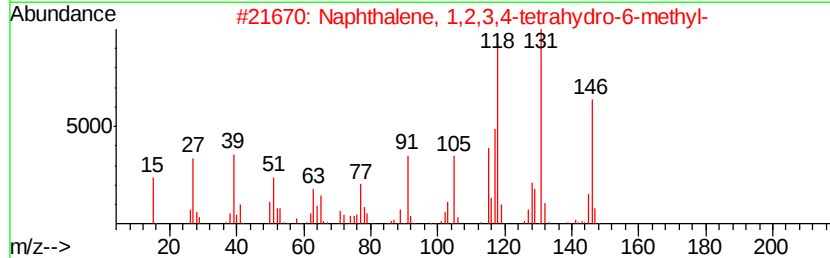
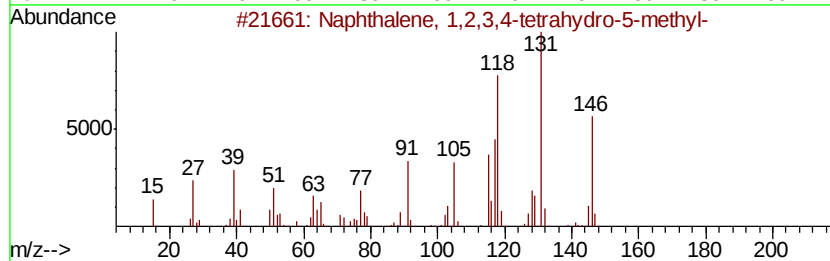
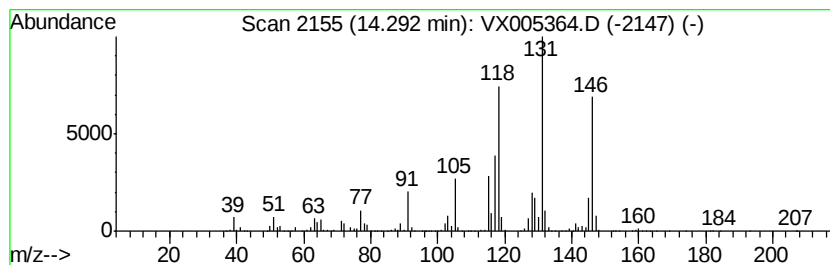
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.29	48.91 ug/l	1128960	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	97
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	96
3		Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	86
4		1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	81
5		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX101618\
Data File : VX005364.D
Acq On : 16 Oct 2018 12:20
Operator : JC/MD
Sample : J5519-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BS-DRUM

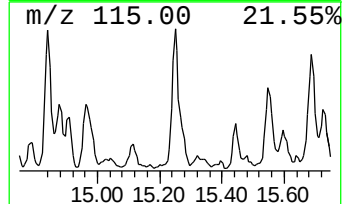
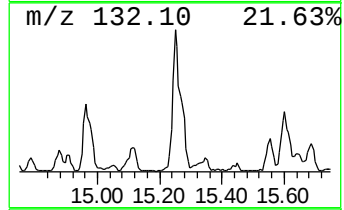
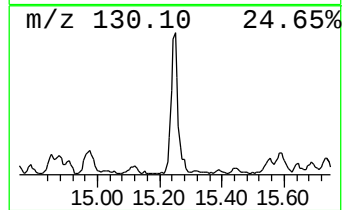
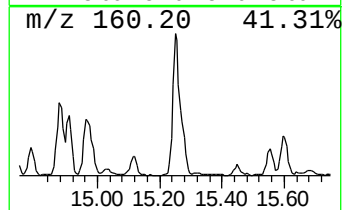
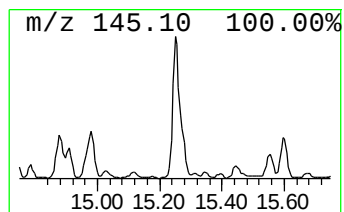
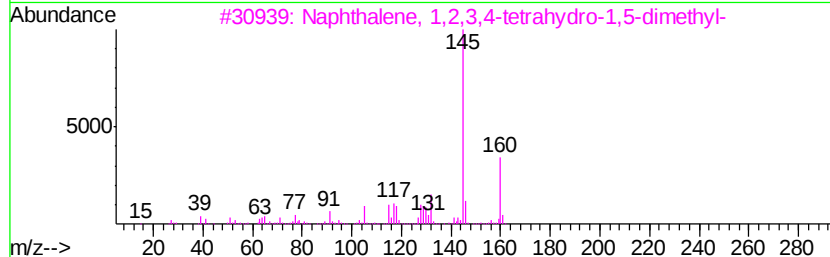
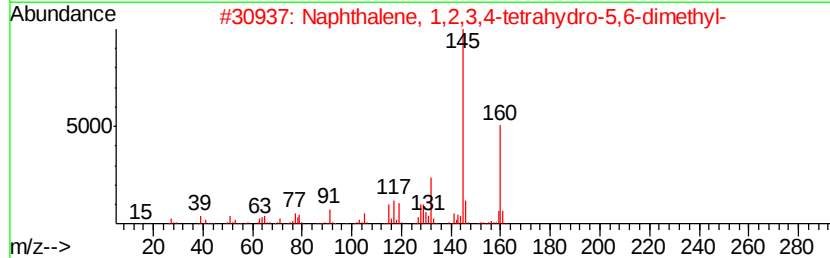
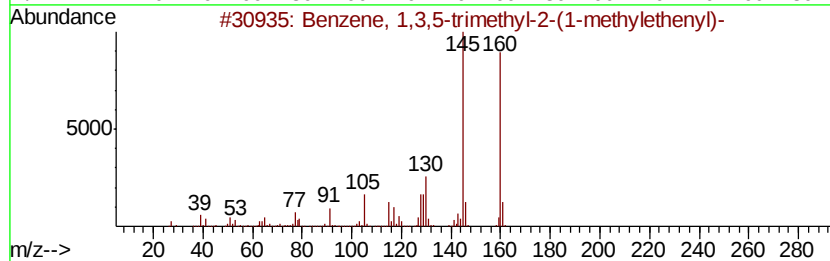
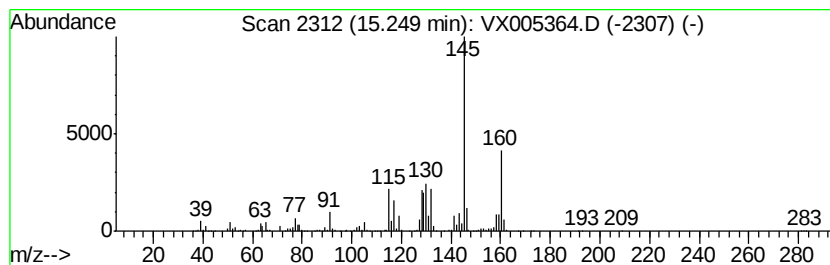
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,3,5-trimethyl-2-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.25	18.12 ug/l	418110	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3,5-trimethyl-2-(1-methyl-2-propenyl)-	160	C12H16	014679-13-1	95
2		Naphthalene, 1,2,3,4-tetrahydro-5,6-dimethyl-	160	C12H16	020027-77-4	92
3		Naphthalene, 1,2,3,4-tetrahydro-5,6-dimethyl-	160	C12H16	021564-91-0	83
4		Benzene, 1-(2-butenyl)-2,3-dimethyl-	160	C12H16	054340-85-1	81
5		Benzene, 4-(2-butenyl)-1,2-dimethyl-	160	C12H16	054340-86-2	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX101618\
Data File : VX005364.D
Acq On : 16 Oct 2018 12:20
Operator : JC/MD
Sample : J5519-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BS-DRUM

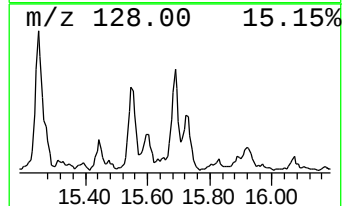
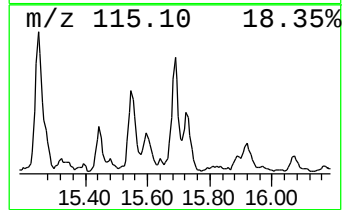
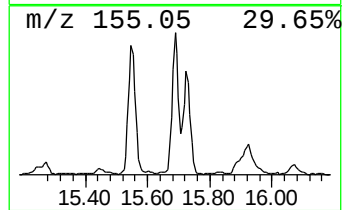
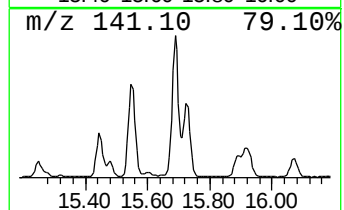
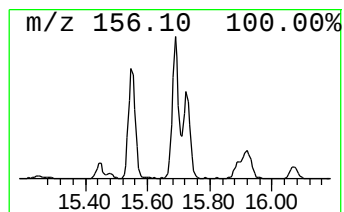
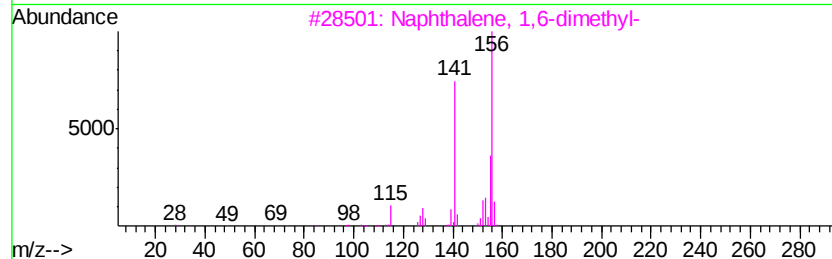
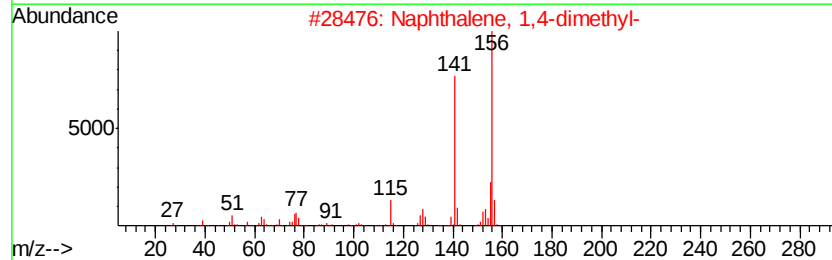
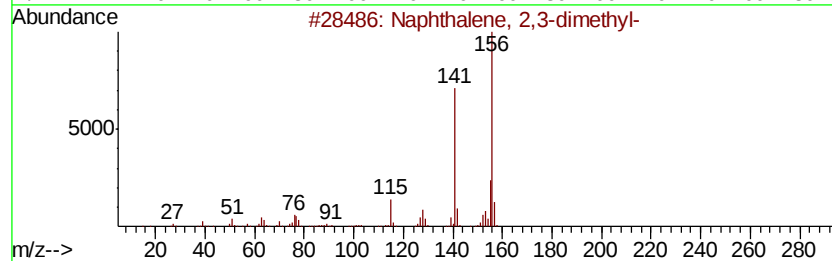
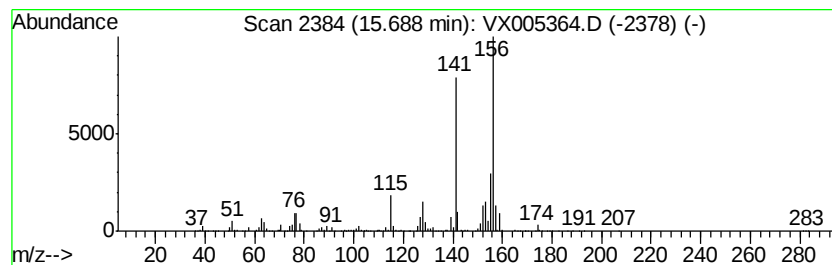
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 2,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.69	16.60 ug/l	383048	1,4-Dichlorobenzene-d4	12.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
4			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
5			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX101618\
Data File : VX005364.D
Acq On : 16 Oct 2018 12:20
Operator : JC/MD
Sample : J5519-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BS-DRUM

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1,2,3,5-...	13.02	16.9	ug/l	389277	4	12.08	1154030	50.0
1H-Indene, 2,3-di...	13.35	45.3	ug/l	1045740	4	12.08	1154030	50.0
Naphthalene, 1,2,...	13.48	25.3	ug/l	584846	4	12.08	1154030	50.0
Benzene, (1,2-dim...	13.64	16.9	ug/l	390016	4	12.08	1154030	50.0
Naphthalene, 1,2,...	13.72	13.8	ug/l	318414	4	12.08	1154030	50.0
Naphthalene, 1,2,...	14.29	48.9	ug/l	1128960	4	12.08	1154030	50.0
Benzene, 1,3,5-tr...	15.25	18.1	ug/l	418110	4	12.08	1154030	50.0
Naphthalene, 2,3-...	15.69	16.6	ug/l	383048	4	12.08	1154030	50.0