

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091318\
 Data File : VX004528.D
 Acq On : 13 Sep 2018 20:35
 Operator : JC/MD
 Sample : J4892-09
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PR-MW-02-091218

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\82X091118W.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	1.788	99	104	112	rVB	57140	79900	4.66%	0.350%
2	2.001	134	139	149	rVB2	16275	29267	1.71%	0.128%
3	2.263	177	182	191	rVB	22751	37103	2.16%	0.163%
4	2.617	234	240	251	rVB	45237	82673	4.82%	0.363%
5	2.842	269	277	280	rBV	32919	72643	4.23%	0.319%
6	2.885	280	284	289	rVV	60841	132328	7.71%	0.580%
7	2.928	289	291	306	rVB4	31024	67120	3.91%	0.294%
8	3.166	321	330	342	rVB2	60067	135772	7.91%	0.595%
9	3.513	378	387	400	rVB4	11183	32848	1.91%	0.144%
10	3.812	425	436	446	rBV2	28736	72296	4.21%	0.317%
11	3.915	446	453	460	rBV3	18590	48468	2.82%	0.213%
12	4.190	487	498	507	rBV2	17064	47986	2.80%	0.210%
13	4.397	522	532	544	rVB2	97037	282544	16.46%	1.239%
14	4.519	544	552	566	rVB3	12579	38376	2.24%	0.168%
15	5.287	666	678	679	rBV4	24968	45143	2.63%	0.198%
16	5.507	698	714	722	rBV2	184204	557500	32.49%	2.445%
17	5.580	722	726	731	rVV4	28047	77091	4.49%	0.338%
18	5.665	731	740	768	rVB	264030	831876	48.47%	3.648%
19	5.921	772	782	796	rVB5	15852	49985	2.91%	0.219%
20	6.067	796	806	816	rBV2	193029	501577	29.23%	2.200%
21	6.263	829	838	847	rBV6	19331	83206	4.85%	0.365%
22	6.354	847	853	862	rVV4	33104	89580	5.22%	0.393%
23	6.458	862	870	881	rVB4	25756	74113	4.32%	0.325%
24	6.695	902	909	919	rVB2	8289	21968	1.28%	0.096%
25	6.860	926	936	944	rBV	353796	874792	50.97%	3.836%
26	6.933	946	948	961	rVV4	53173	125747	7.33%	0.551%
27	7.061	962	969	977	rVV3	10317	25518	1.49%	0.112%
28	7.147	977	983	992	rVB2	12388	27009	1.57%	0.118%
29	7.256	992	1001	1011	rBV2	44470	103530	6.03%	0.454%
30	7.457	1022	1034	1046	rBV4	52928	160096	9.33%	0.702%
31	7.732	1073	1079	1092	rVB2	11654	25981	1.51%	0.114%
32	7.927	1104	1111	1113	rBV2	14076	24561	1.43%	0.108%
33	7.957	1113	1116	1123	rVV3	17843	36380	2.12%	0.160%
34	8.055	1123	1132	1144	rVB4	12891	41654	2.43%	0.183%

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Integration Parameters: RTEINT.P

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35	8.177	1144	1152	1154	rBV2	23559	44212	2.58%	0.194%
36	8.213	1154	1158	1163	rVV	55377	100262	5.84%	0.440%
37	8.262	1163	1166	1176	rVB4	11890	23643	1.38%	0.104%
38	8.573	1211	1217	1226	rVB2	47884	87667	5.11%	0.384%
39	8.719	1234	1241	1248	rVV	927605	1389472	80.97%	6.093%
40	9.225	1319	1324	1334	rVB	47457	79104	4.61%	0.347%
41	9.518	1369	1372	1376	rVV	23533	36856	2.15%	0.162%
42	9.561	1376	1379	1386	rVB4	13917	28487	1.66%	0.125%
43	10.116	1464	1470	1476	rBV	744044	992891	57.86%	4.354%
44	10.256	1489	1493	1499	rVB	38436	51096	2.98%	0.224%
45	11.018	1613	1618	1627	rBV	1350977	1716138	100.00%	7.526%
46	11.140	1632	1638	1644	rBV	771531	969525	56.49%	4.252%
47	11.359	1669	1674	1685	rVB	533530	649783	37.86%	2.850%
48	11.670	1715	1725	1733	rBV	50016	71241	4.15%	0.312%
49	11.768	1733	1741	1745	rBV	18997	26350	1.54%	0.116%
50	11.920	1761	1766	1767	rBV	30812	33613	1.96%	0.147%
51	11.944	1767	1770	1776	rVB	92338	121455	7.08%	0.533%
52	12.079	1787	1792	1798	rBV	946277	1158886	67.53%	5.082%
53	12.146	1798	1803	1808	rVV	702292	846430	49.32%	3.712%
54	12.231	1813	1817	1822	rVB	32665	39800	2.32%	0.175%
55	12.298	1822	1828	1836	rBV	1036341	1319861	76.91%	5.788%
56	12.371	1836	1840	1841	rBV	69014	80543	4.69%	0.353%
57	12.463	1849	1855	1860	rBV	72056	91275	5.32%	0.400%
58	12.518	1860	1864	1871	rVB	18419	25280	1.47%	0.111%
59	12.670	1880	1889	1892	rBV	283132	353777	20.61%	1.551%
60	12.700	1892	1894	1899	rVV	150793	170238	9.92%	0.747%
61	12.755	1899	1903	1911	rVB2	419090	586507	34.18%	2.572%
62	12.902	1920	1927	1932	rVB2	106203	157644	9.19%	0.691%
63	12.981	1932	1940	1944	rBV	225586	299541	17.45%	1.314%
64	13.024	1944	1947	1952	rVB2	14052	21998	1.28%	0.096%
65	13.084	1952	1957	1964	rVB3	42837	66282	3.86%	0.291%
66	13.152	1964	1968	1972	rBV2	22349	37738	2.20%	0.165%
67	13.194	1972	1975	1977	rBV2	54064	59225	3.45%	0.260%
68	13.225	1977	1980	1990	rVB	241069	311930	18.18%	1.368%
69	13.347	1990	2000	2006	rBV2	1095148	1495306	87.13%	6.557%
70	13.402	2006	2009	2012	rBV2	27680	36468	2.13%	0.160%
71	13.481	2018	2022	2031	rVB	97069	128856	7.51%	0.565%

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72	13.566	2031	2036	2038	rBV2	46506	64955	3.78%	0.285%
73	13.603	2038	2042	2045	rVV	94970	136086	7.93%	0.597%
74	13.645	2045	2049	2052	rVV2	129028	197895	11.53%	0.868%
75	13.700	2052	2058	2059	rVV2	89505	149296	8.70%	0.655%
76	13.719	2059	2061	2070	rVB2	139739	190541	11.10%	0.836%
77	13.810	2070	2076	2078	rBV2	30358	40710	2.37%	0.179%
78	13.853	2078	2083	2088	rVB2	54167	91208	5.31%	0.400%
79	13.914	2088	2093	2097	rVB	99625	122265	7.12%	0.536%
80	13.987	2097	2105	2109	rBV2	136632	202205	11.78%	0.887%
81	14.029	2109	2112	2116	rVV	47720	55019	3.21%	0.241%
82	14.133	2124	2129	2132	rBV	56774	74051	4.31%	0.325%
83	14.164	2132	2134	2141	rVB3	29264	33218	1.94%	0.146%
84	14.285	2147	2154	2162	rVB	196469	345134	20.11%	1.514%
85	14.377	2165	2169	2173	rVV	47362	58553	3.41%	0.257%
86	14.432	2174	2178	2182	rVB	61550	75841	4.42%	0.333%
87	14.481	2182	2186	2190	rVB	18736	26312	1.53%	0.115%
88	14.542	2190	2196	2201	rBV2	190855	256884	14.97%	1.127%
89	14.694	2214	2221	2225	rBV	423490	570891	33.27%	2.504%
90	14.731	2225	2227	2232	rVB2	46594	52726	3.07%	0.231%
91	14.840	2240	2245	2250	rBV	595205	758328	44.19%	3.326%
92	14.907	2254	2256	2260	rVV3	22476	25933	1.51%	0.114%
93	14.980	2260	2268	2273	rVB6	10979	27810	1.62%	0.122%
94	15.249	2304	2312	2320	rBV2	26719	48526	2.83%	0.213%
95	15.444	2339	2344	2347	rBV	33463	47942	2.79%	0.210%
96	15.480	2347	2350	2354	rVB	19401	25096	1.46%	0.110%
97	15.547	2354	2361	2367	rBV	79045	132638	7.73%	0.582%
98	15.688	2376	2384	2388	rBV	114784	195946	11.42%	0.859%
99	15.724	2388	2390	2398	rVB	53179	73020	4.25%	0.320%
100	15.919	2412	2422	2431	rVB2	27617	78279	4.56%	0.343%

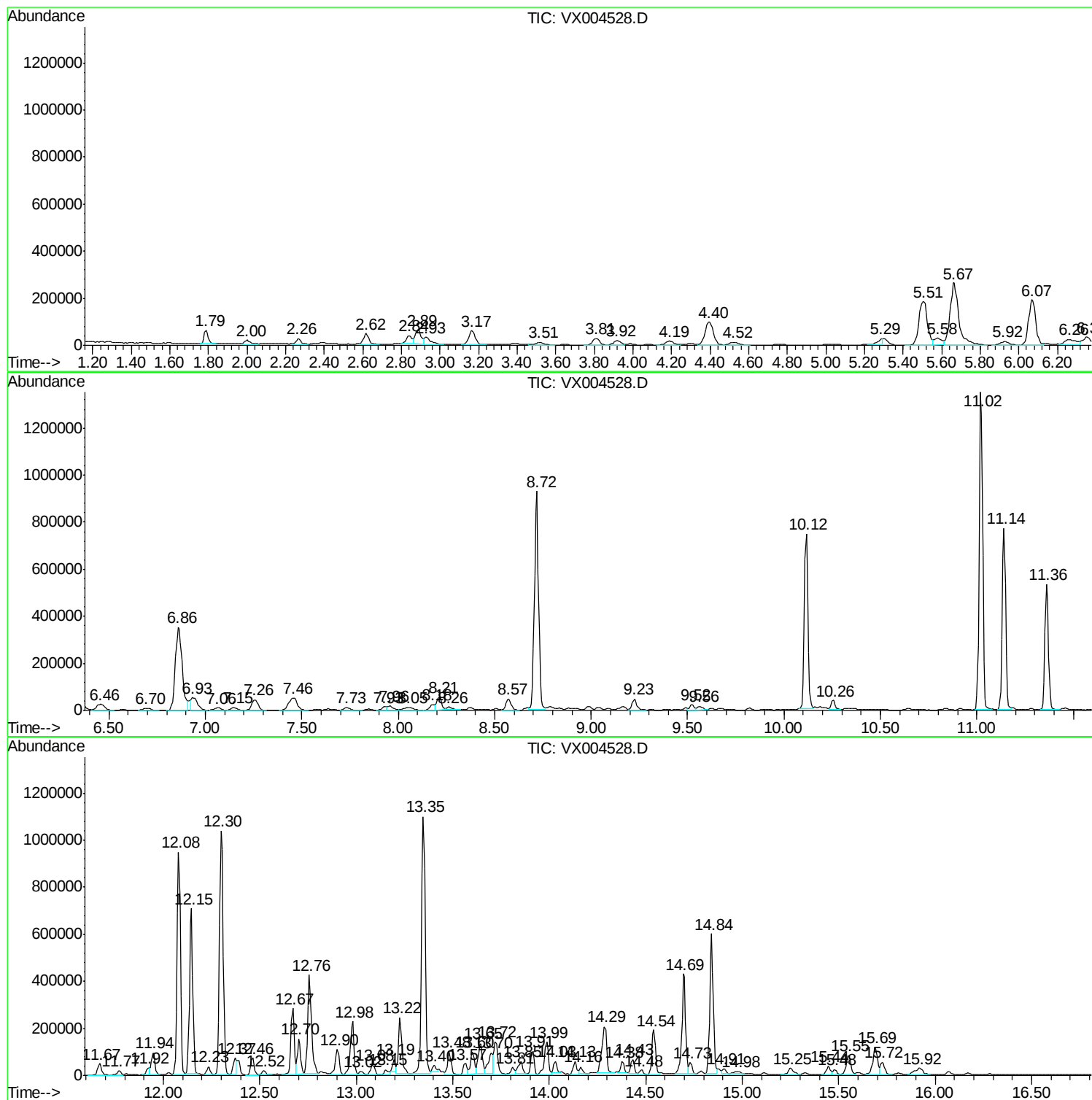
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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



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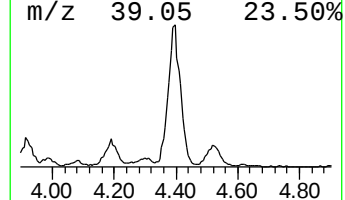
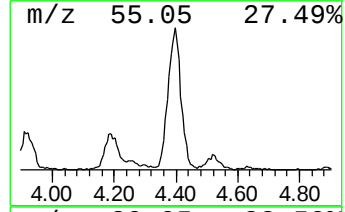
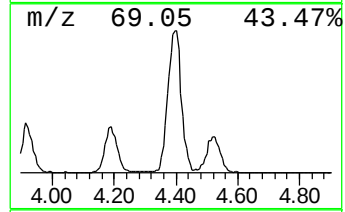
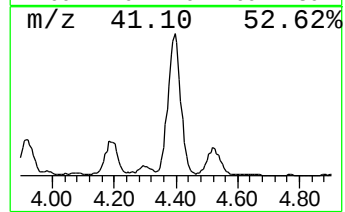
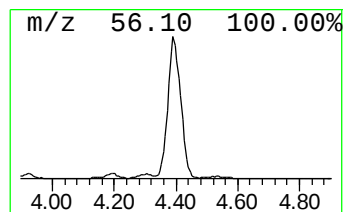
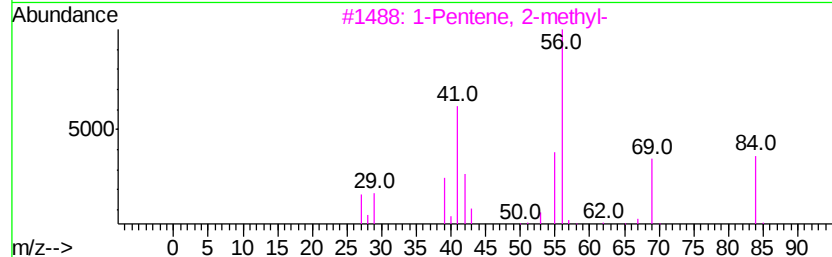
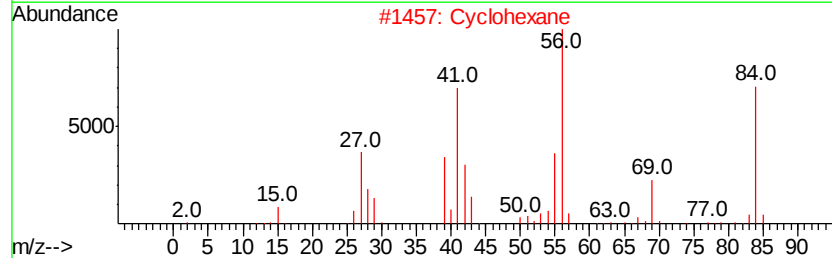
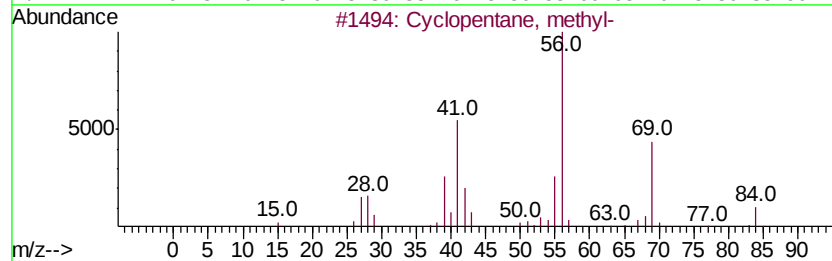
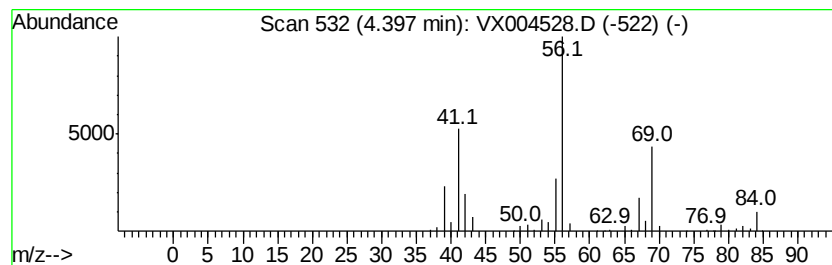
Instrument :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclopentane, methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
4.40	16.98 ug/l	282544	Pentafluorobenzene	5.66		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cvclopentane, methyl-	84	C6H12	000096-37-7	91
2		Cyclohexane	84	C6H12	000110-82-7	86
3		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	80
4		Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
5		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	50



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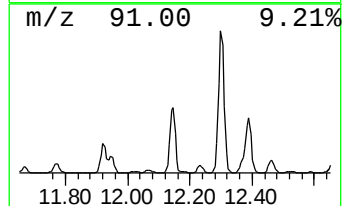
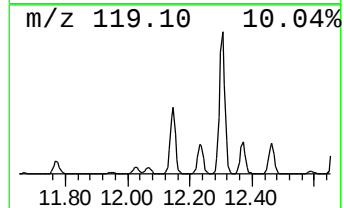
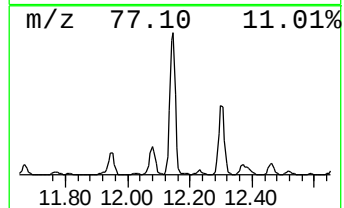
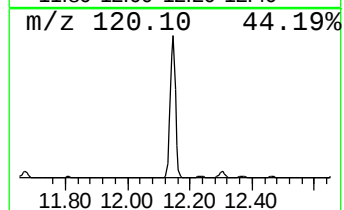
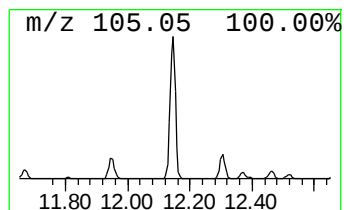
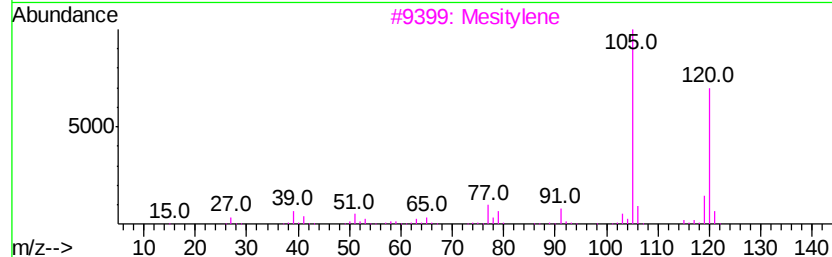
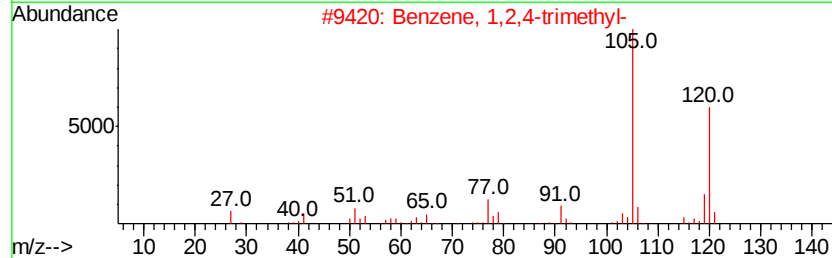
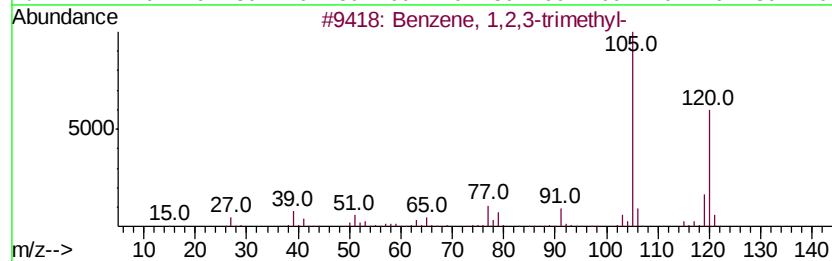
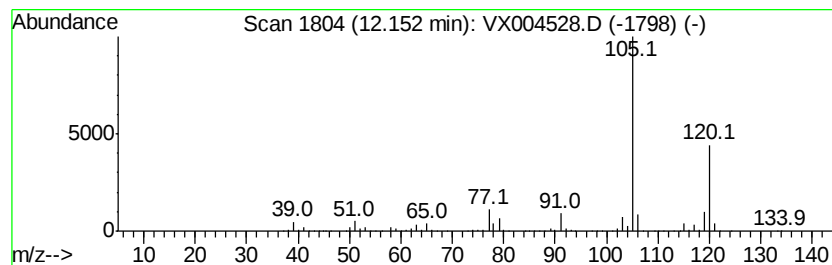
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1,2,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	36.52 ug/l	846430	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2,3-trimethyl-	120 C9H12	000526-73-8 95
2		Benzene, 1,2,4-trimethyl-	120 C9H12	000095-63-6 94
3		Mesitylene	120 C9H12	000108-67-8 91
4		Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4 91
5		Benzene, 1-ethyl-4-methyl-	120 C9H12	000622-96-8 90



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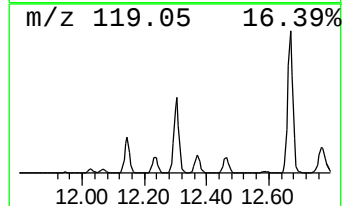
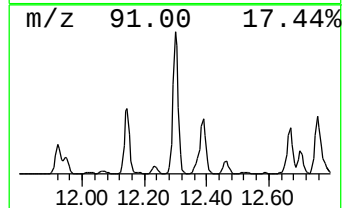
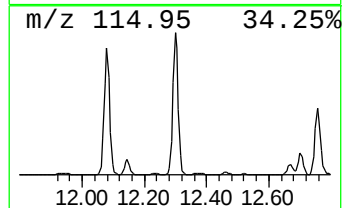
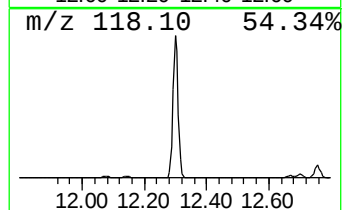
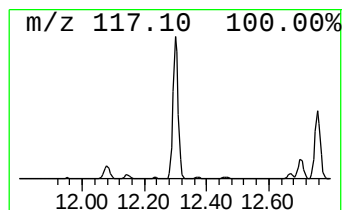
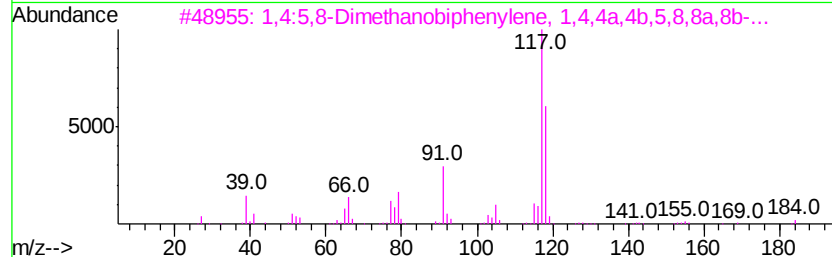
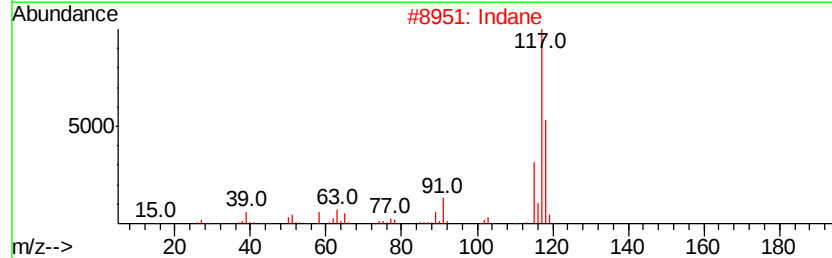
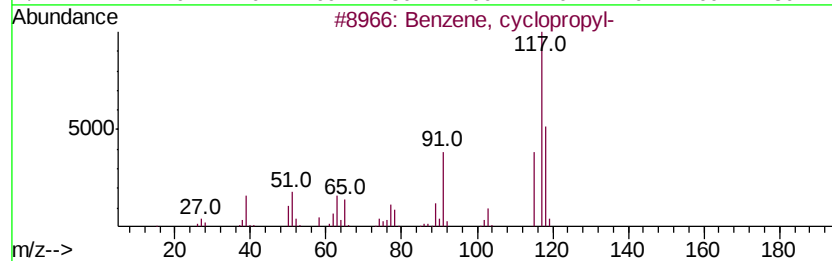
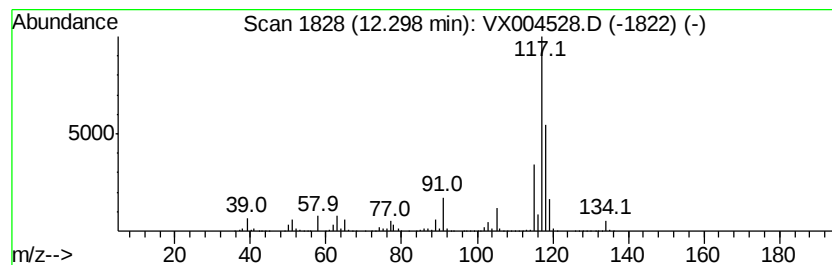
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, cyclopropyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	56.95 ug/l	1319860	1,4-Dichlorobenzene-d4	12.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, cvclopropyl-	118	C9H10	000873-49-4	76
2		Indane	118	C9H10	000496-11-7	76
3		1,4:5,8-Dimethanobiphenylene, 1,...	184	C14H16	001624-13-1	72
4		Benzene, 1-propenyl-	118	C9H10	000637-50-3	68
5		Deltacyclene	118	C9H10	007785-10-6	58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091318\
Data File : VX004528.D
Acq On : 13 Sep 2018 20:35
Operator : JC/MD
Sample : J4892-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

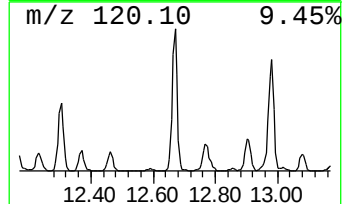
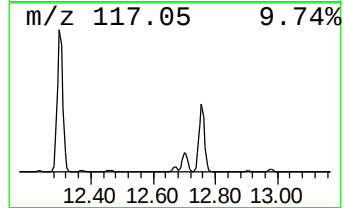
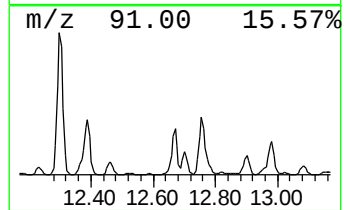
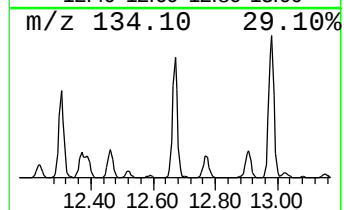
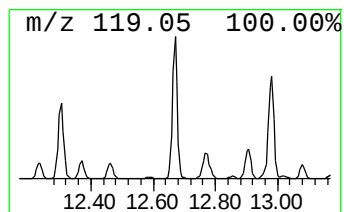
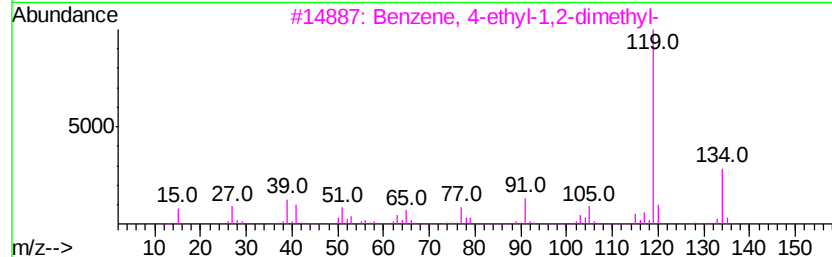
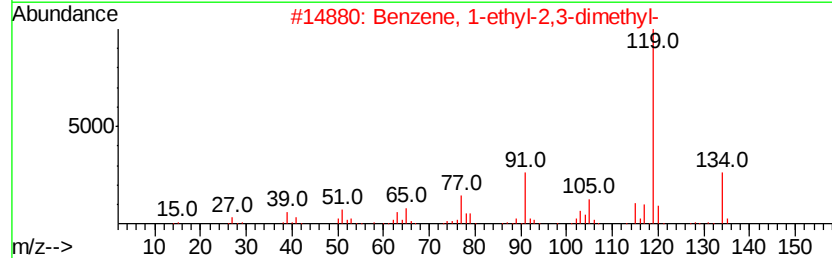
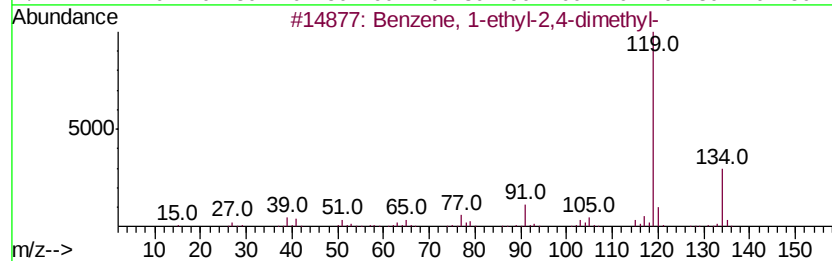
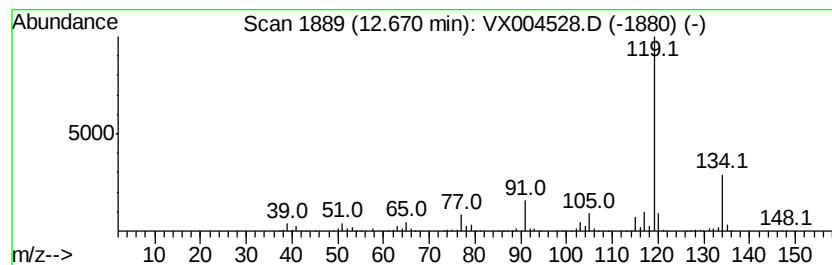
Instrument :
MSVOA_X
ClientSampleId :
PR-MW-02-091218

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	15.26 ug/l	353777	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9	97
2	Benzene, 1-ethyl-2,3-dimethyl-	134 C10H14	000933-98-2	96
3	Benzene, 4-ethyl-1,2-dimethyl-	134 C10H14	000934-80-5	95
4	Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4	95
5	Benzene, 1-methyl-3-(1-methyleth...	134 C10H14	000535-77-3	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091318\
Data File : VX004528.D
Acq On : 13 Sep 2018 20:35
Operator : JC/MD
Sample : J4892-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PR-MW-02-091218

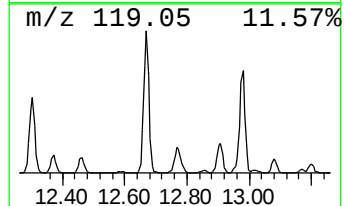
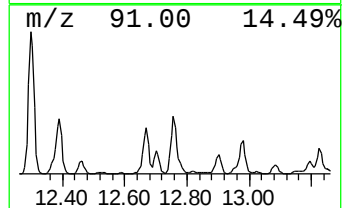
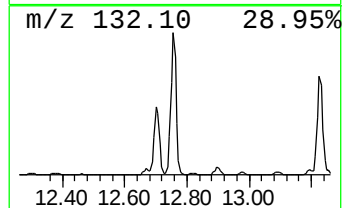
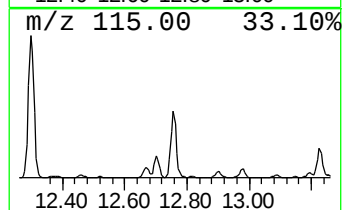
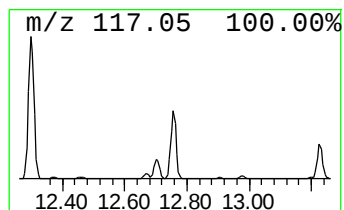
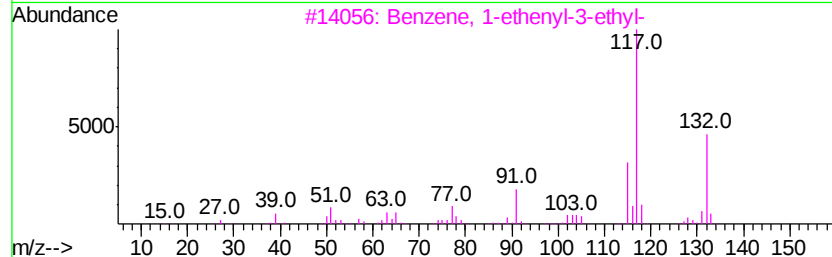
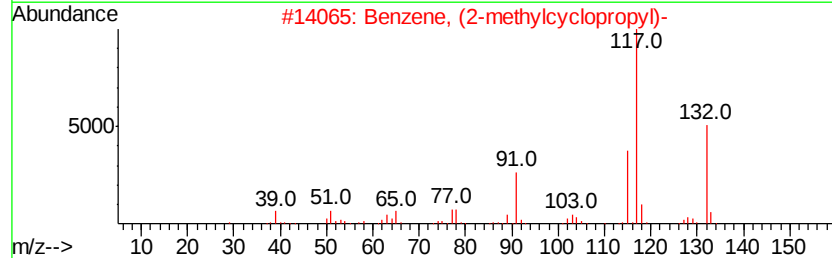
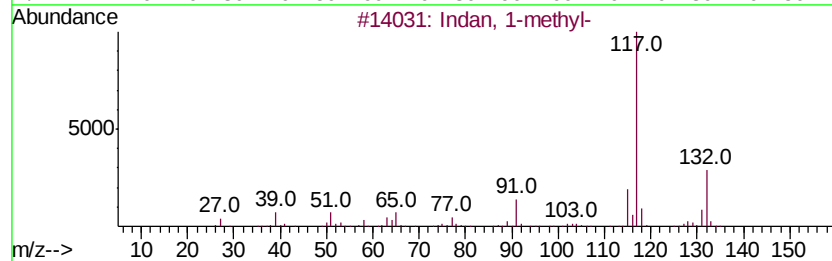
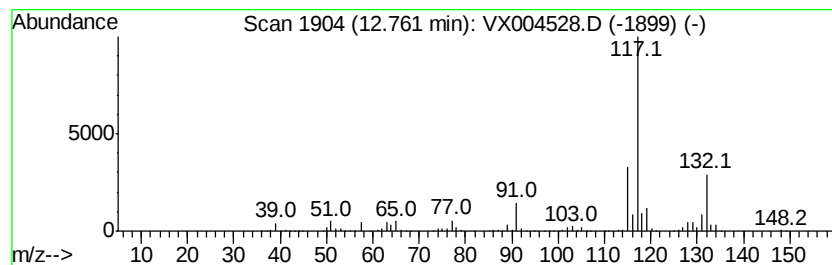
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	25.30 ug/l	586507	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indan, 1-methyl-	132	C10H12	000767-58-8	87
2		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	87
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
4		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	80
5		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	80



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091318\
Data File : VX004528.D
Acq On : 13 Sep 2018 20:35
Operator : JC/MD
Sample : J4892-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PR-MW-02-091218

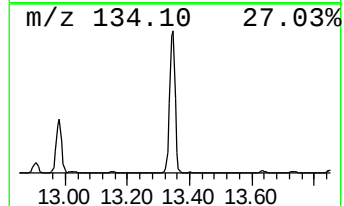
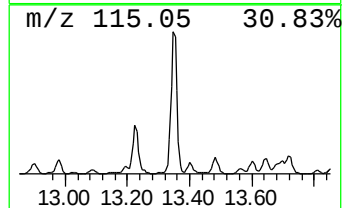
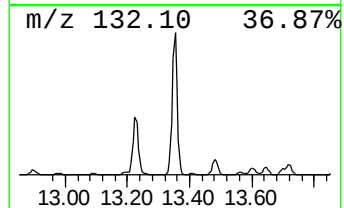
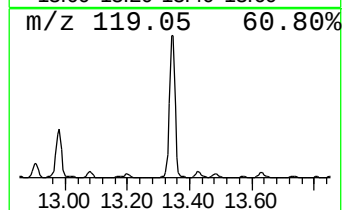
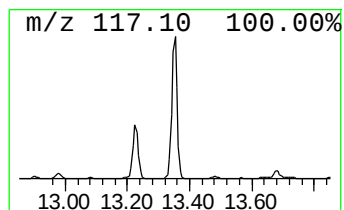
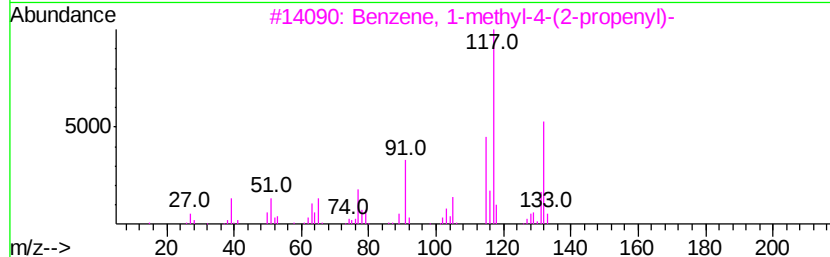
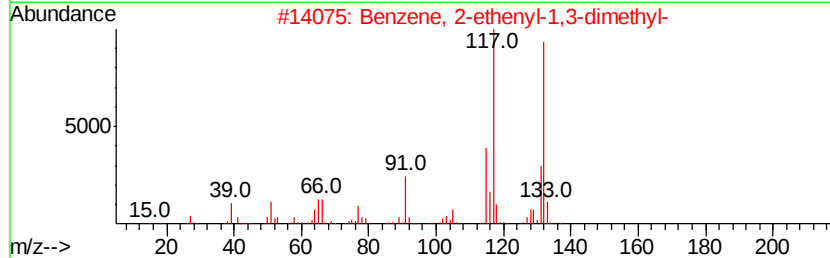
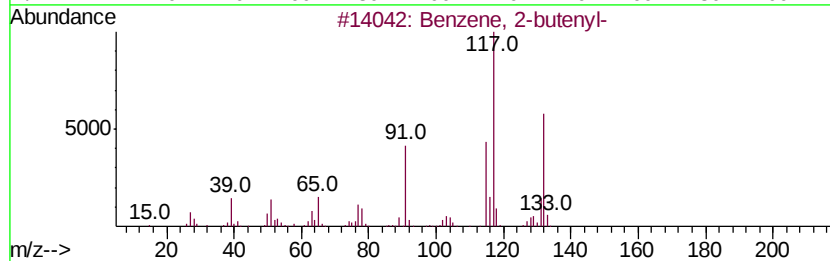
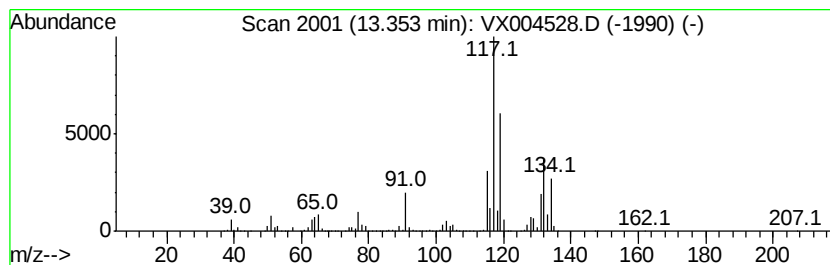
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 2-butenyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	64.51 ug/l	1495310	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	64
2		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	64
3		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	64
4		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	64
5		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	55



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091318\
Data File : VX004528.D
Acq On : 13 Sep 2018 20:35
Operator : JC/MD
Sample : J4892-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

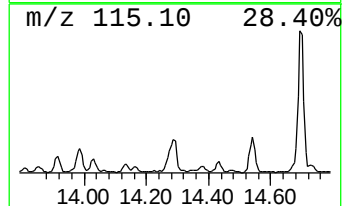
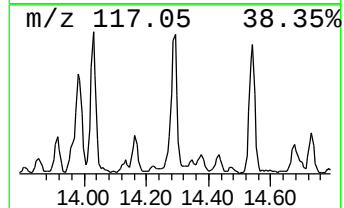
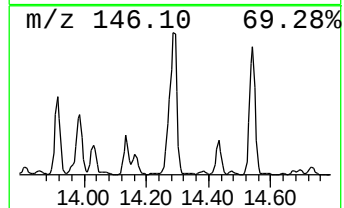
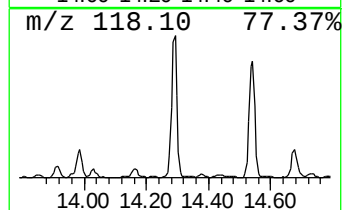
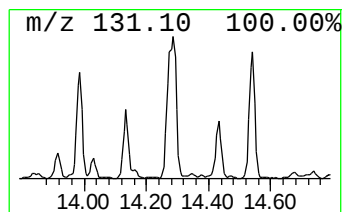
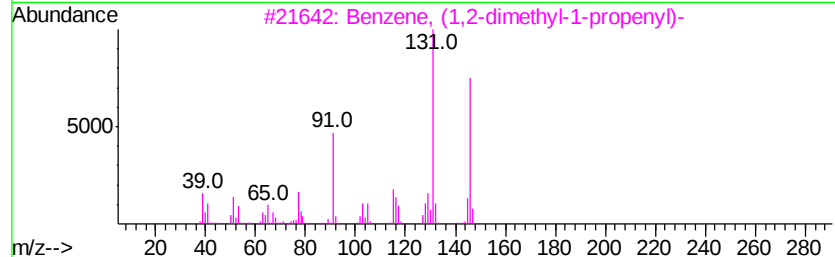
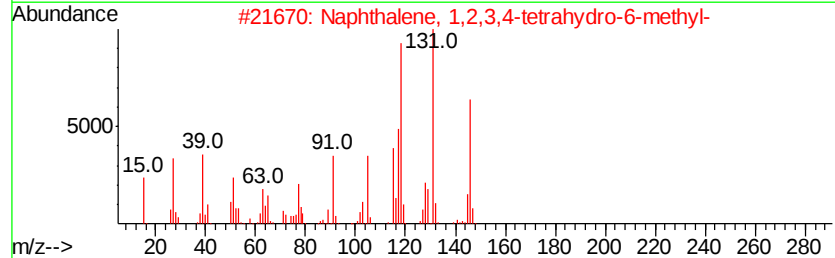
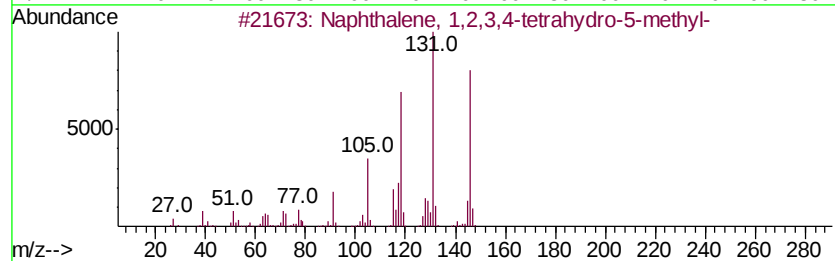
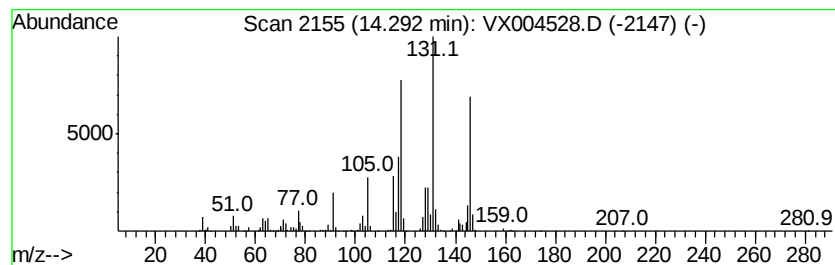
Instrument :
MSVOA_X
ClientSampleId :
PR-MW-02-091218

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.29	14.89 ug/l	345134	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 94
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 86
3		Benzene, (1,2-dimethyl-1-propenyl)-	146 C11H14	000769-57-3 68
4		Benzene, 2-ethenyl-1,3,5-trimethyl-	146 C11H14	000769-25-5 62
5		Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3 58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091318\
Data File : VX004528.D
Acq On : 13 Sep 2018 20:35
Operator : JC/MD
Sample : J4892-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PR-MW-02-091218

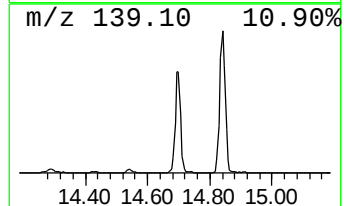
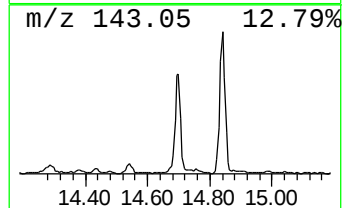
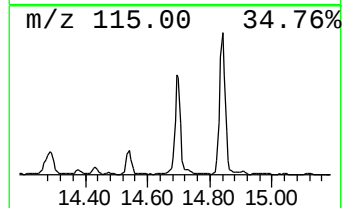
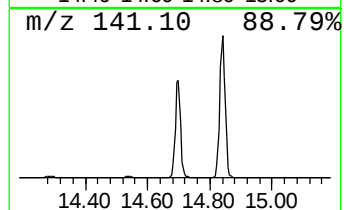
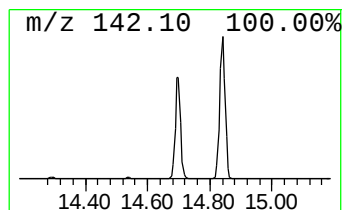
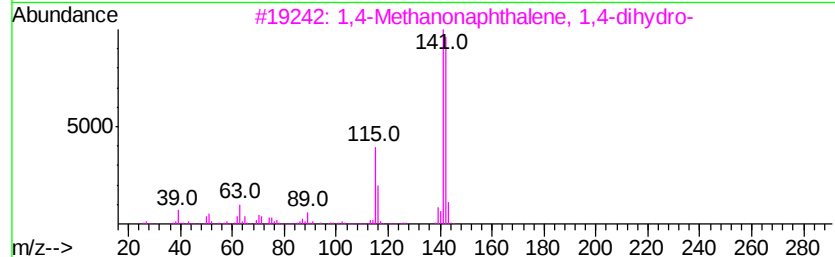
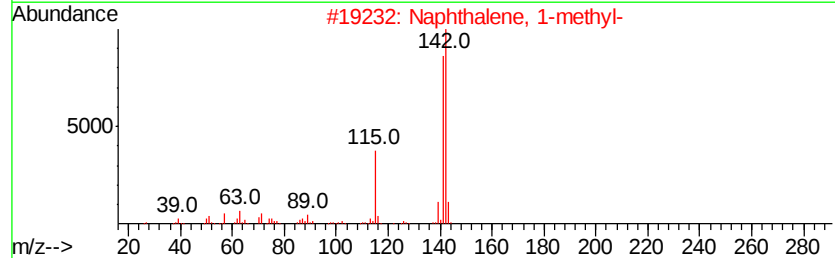
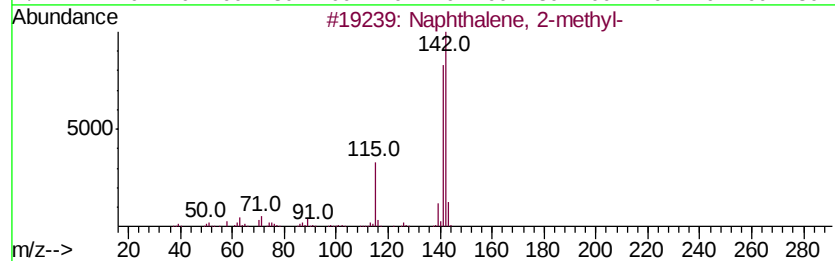
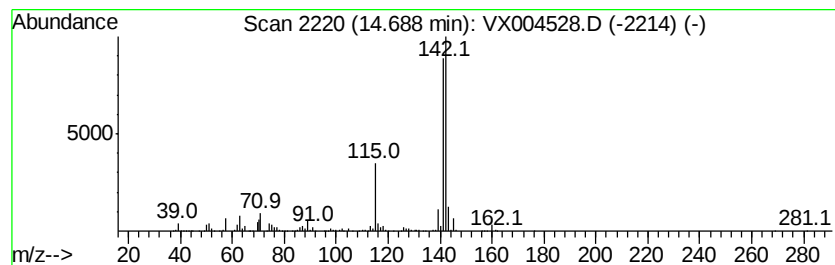
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	24.63 ug/l	570891	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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	90
4		Benzocycloheptatriene	142	C11H10	000264-09-5	90
5		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	59



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091318\
Data File : VX004528.D
Acq On : 13 Sep 2018 20:35
Operator : JC/MD
Sample : J4892-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PR-MW-02-091218

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.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
Cyclopentane, met...	4.40	17.0	ug/l	282544	1	5.66	831876	50.0
Benzene, 1,2,3-tr...	12.15	36.5	ug/l	846430	4	12.08	1158890	50.0
Benzene, cyclopro...	12.30	57.0	ug/l	1319860	4	12.08	1158890	50.0
Benzene, 1-ethyl-...	12.67	15.3	ug/l	353777	4	12.08	1158890	50.0
Indan, 1-methyl-	12.76	25.3	ug/l	586507	4	12.08	1158890	50.0
Benzene, 2-butenyl-	13.35	64.5	ug/l	1495310	4	12.08	1158890	50.0
Naphthalene, 1,2,...	14.29	14.9	ug/l	345134	4	12.08	1158890	50.0
Naphthalene, 1-me...	14.69	24.6	ug/l	570891	4	12.08	1158890	50.0