

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX061618\
 Data File : VX002658.D
 Acq On : 16 Jun 2018 20:18
 Operator : JC/MD
 Sample : J3467-07
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PR-MW-02-061118

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\82X061518W.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	1.075	76	80	94	rBV	1211371	1466220	100.00%	8.214%
2	1.270	108	112	117	rBV	15482	25809	1.76%	0.145%
3	1.337	117	123	126	rBV2	39038	85732	5.85%	0.480%
4	1.794	193	198	207	rVB2	38002	59552	4.06%	0.334%
5	2.002	228	232	242	rVB3	12272	20504	1.40%	0.115%
6	2.270	272	276	284	rVB2	10686	18340	1.25%	0.103%
7	2.624	328	334	342	rBV	18736	34664	2.36%	0.194%
8	2.843	364	370	373	rBV2	20576	40706	2.78%	0.228%
9	2.892	373	378	383	rVB2	29839	58431	3.99%	0.327%
10	3.166	414	423	432	rVB	37951	91640	6.25%	0.513%
11	3.812	521	529	538	rBV2	14000	35452	2.42%	0.199%
12	3.922	540	547	554	rBV4	8829	22415	1.53%	0.126%
13	4.190	583	591	601	rBV5	7918	20977	1.43%	0.118%
14	4.398	615	625	638	rVB2	52870	155199	10.58%	0.869%
15	4.532	638	647	656	rBV4	5993	19570	1.33%	0.110%
16	5.306	762	774	784	rVB4	11504	46052	3.14%	0.258%
17	5.507	795	807	816	rBV	134239	398312	27.17%	2.231%
18	5.672	825	834	858	rVB	208821	656159	44.75%	3.676%
19	5.928	863	876	887	rBV5	12035	37790	2.58%	0.212%
20	6.074	889	900	909	rBV	150240	383882	26.18%	2.151%
21	6.263	920	931	938	rBV4	14821	51570	3.52%	0.289%
22	6.355	940	946	956	rVV3	27445	80861	5.51%	0.453%
23	6.464	956	964	974	rVB4	16327	46686	3.18%	0.262%
24	6.867	1019	1030	1038	rBV	299818	722807	49.30%	4.049%
25	7.263	1086	1095	1105	rBV2	24629	56638	3.86%	0.317%
26	7.464	1116	1128	1140	rBV4	40518	119395	8.14%	0.669%
27	7.738	1167	1173	1182	rVB3	9344	18740	1.28%	0.105%
28	7.958	1206	1209	1218	rVV7	11872	25683	1.75%	0.144%
29	8.055	1218	1225	1235	rVB2	14246	38008	2.59%	0.213%
30	8.184	1237	1246	1248	rBV2	23375	48502	3.31%	0.272%
31	8.214	1248	1251	1256	rVB	33138	50437	3.44%	0.283%
32	8.574	1305	1310	1319	rVB4	29449	58411	3.98%	0.327%
33	8.671	1319	1326	1328	rBV3	12172	22218	1.52%	0.124%
34	8.720	1328	1334	1341	rVB	747415	1137461	77.58%	6.372%

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35	9.165	1403	1407	1412	rVB	11195	17411	1.19%	0.098%
36	9.226	1412	1417	1422	rBV	26846	39541	2.70%	0.222%
37	9.525	1462	1466	1469	rBV	18195	23367	1.59%	0.131%
38	9.574	1469	1474	1479	rVB7	9174	21102	1.44%	0.118%
39	10.073	1552	1556	1558	rBV	16090	21620	1.47%	0.121%
40	10.116	1558	1563	1569	rVB	630840	823835	56.19%	4.615%
41	10.256	1582	1586	1590	rVB	19855	25945	1.77%	0.145%
42	10.646	1642	1650	1657	rBV8	6887	17861	1.22%	0.100%
43	10.842	1670	1682	1685	rBV5	8951	17997	1.23%	0.101%
44	10.915	1690	1694	1705	rVB10	9363	21787	1.49%	0.122%
45	11.024	1706	1712	1718	rBV	795307	994223	67.81%	5.570%
46	11.140	1724	1731	1736	rBV	649004	809128	55.18%	4.533%
47	11.366	1759	1768	1776	rBV	300067	377983	25.78%	2.118%
48	11.671	1813	1818	1827	rVB	34641	50668	3.46%	0.284%
49	11.774	1827	1835	1838	rBV	16929	23678	1.61%	0.133%
50	11.927	1851	1860	1861	rBV	25518	34917	2.38%	0.196%
51	11.951	1861	1864	1870	rVB	73658	88934	6.07%	0.498%
52	12.079	1880	1885	1891	rBV	725926	925186	63.10%	5.183%
53	12.146	1891	1896	1902	rVB	368887	433915	29.59%	2.431%
54	12.238	1906	1911	1915	rVB	24743	30962	2.11%	0.173%
55	12.305	1915	1922	1929	rBV	740779	911737	62.18%	5.108%
56	12.372	1929	1933	1935	rBV	55484	76851	5.24%	0.431%
57	12.463	1943	1948	1953	rBV	52756	67079	4.57%	0.376%
58	12.524	1953	1958	1963	rVB	16228	22330	1.52%	0.125%
59	12.671	1974	1982	1985	rBV	276564	349391	23.83%	1.957%
60	12.707	1985	1988	1992	rVV	109237	132730	9.05%	0.744%
61	12.762	1992	1997	2004	rVV	326378	452468	30.86%	2.535%
62	12.902	2013	2020	2025	rVB3	66291	107203	7.31%	0.601%
63	12.981	2025	2033	2038	rBV	218076	276872	18.88%	1.551%
64	13.030	2038	2041	2046	rVB4	12514	18628	1.27%	0.104%
65	13.085	2046	2050	2057	rVB3	31067	48436	3.30%	0.271%
66	13.158	2057	2062	2065	rBV	17369	28947	1.97%	0.162%
67	13.201	2065	2069	2071	rBV2	34355	44040	3.00%	0.247%
68	13.231	2071	2074	2083	rVB	172151	217931	14.86%	1.221%
69	13.353	2083	2094	2100	rBV2	750797	1100626	75.07%	6.166%
70	13.408	2100	2103	2105	rBV2	15628	21042	1.44%	0.118%
71	13.487	2112	2116	2119	rVB2	33759	42693	2.91%	0.239%

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72	13.567	2125	2129	2132	rBV2	31643	43720	2.98%	0.245%
73	13.603	2132	2135	2138	rVV	58928	81933	5.59%	0.459%
74	13.646	2138	2142	2146	rVV3	89244	143129	9.76%	0.802%
75	13.701	2146	2151	2153	rVV2	60364	108192	7.38%	0.606%
76	13.725	2153	2155	2163	rVB2	94413	125727	8.57%	0.704%
77	13.811	2163	2169	2172	rBV	21626	30074	2.05%	0.168%
78	13.859	2172	2177	2181	rVB	33977	45497	3.10%	0.255%
79	13.914	2181	2186	2191	rVB	71913	90401	6.17%	0.506%
80	13.987	2191	2198	2202	rBV2	95995	144533	9.86%	0.810%
81	14.030	2202	2205	2210	rVB2	30555	37027	2.53%	0.207%
82	14.134	2218	2222	2225	rBV	37928	50671	3.46%	0.284%
83	14.170	2225	2228	2234	rVB2	18953	24215	1.65%	0.136%
84	14.292	2241	2248	2255	rVB2	134664	235139	16.04%	1.317%
85	14.384	2259	2263	2267	rVV	36467	46405	3.16%	0.260%
86	14.438	2267	2272	2276	rVB	41074	55854	3.81%	0.313%
87	14.481	2276	2279	2284	rVB2	16791	22256	1.52%	0.125%
88	14.542	2284	2289	2298	rBV2	130796	185389	12.64%	1.039%
89	14.701	2307	2315	2319	rBV	267654	368253	25.12%	2.063%
90	14.737	2319	2321	2325	rVB2	37485	39666	2.71%	0.222%
91	14.847	2333	2339	2344	rBV	408430	555159	37.86%	3.110%
92	14.908	2347	2349	2353	rVV3	18934	25287	1.72%	0.142%
93	14.981	2353	2361	2367	rVV7	9173	26042	1.78%	0.146%
94	15.255	2401	2406	2413	rVB3	22514	42254	2.88%	0.237%
95	15.451	2433	2438	2442	rBV	28548	47465	3.24%	0.266%
96	15.554	2448	2455	2460	rBV	64822	103056	7.03%	0.577%
97	15.694	2471	2478	2481	rBV	84576	136814	9.33%	0.766%
98	15.731	2481	2484	2492	rVB	50156	77234	5.27%	0.433%
99	15.920	2505	2515	2527	rVB3	22645	65313	4.45%	0.366%
100	16.078	2536	2541	2550	rVV3	10150	19354	1.32%	0.108%

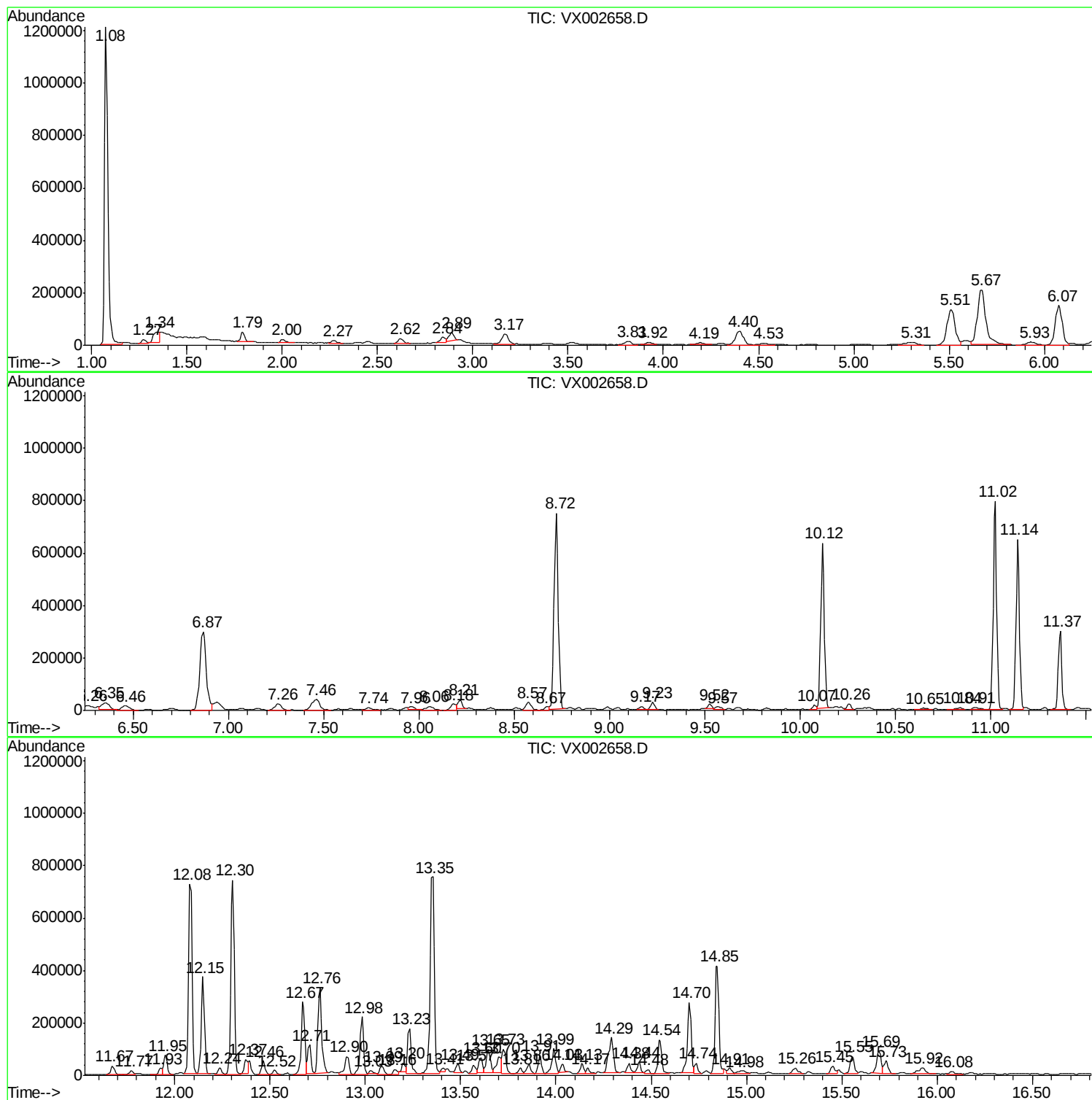
Sum of corrected areas: 17849946

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061518W.M
Quant Title : SW846 8260

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



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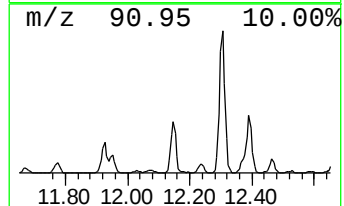
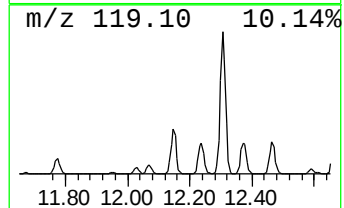
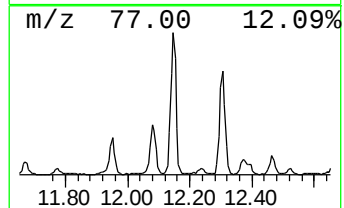
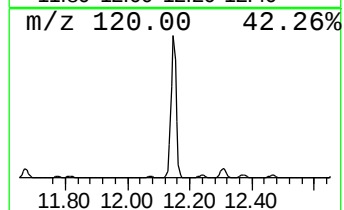
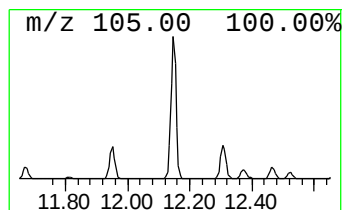
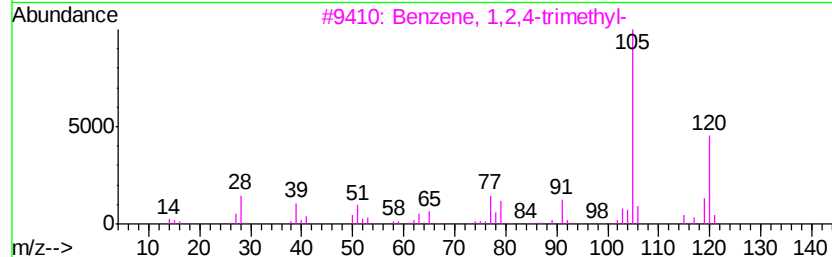
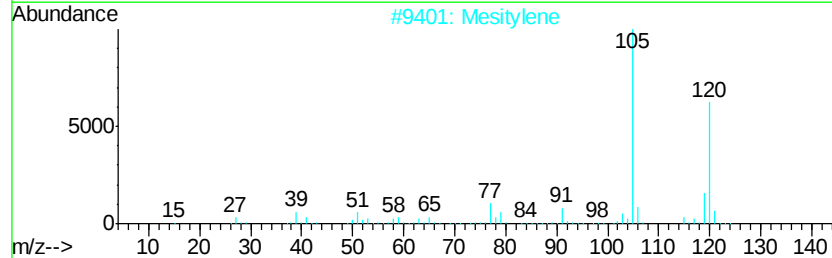
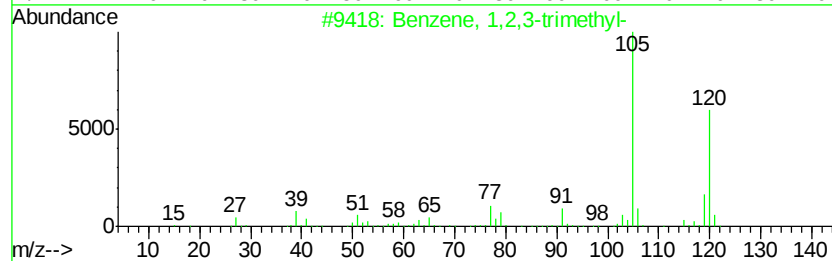
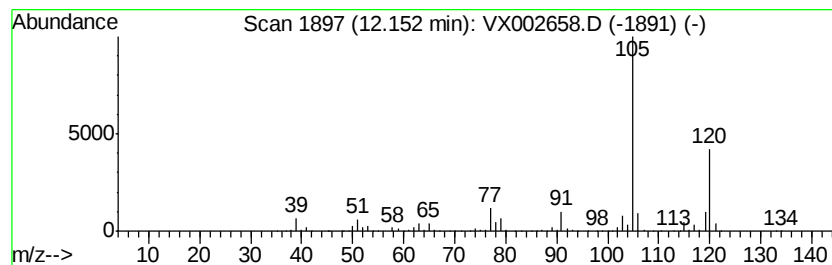
Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061518W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	23.45 ug/l	433915	1,4-Dichlorobenzene-d4	12.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2,3-trimethyl-	120 C9H12	000526-73-8 95
2		Mesitylene	120 C9H12	000108-67-8 94
3		Benzene, 1,2,4-trimethyl-	120 C9H12	000095-63-6 94
4		Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4 90
5		Benzene, 1-ethyl-2-methyl-	120 C9H12	000611-14-3 87



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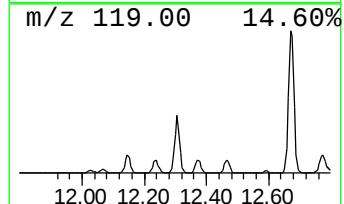
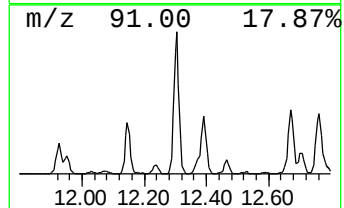
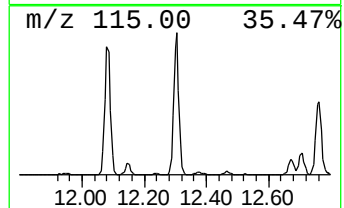
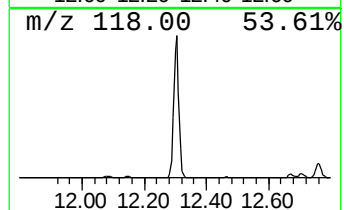
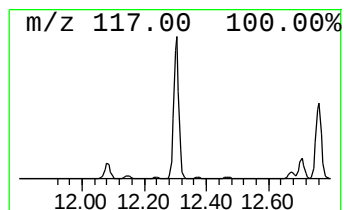
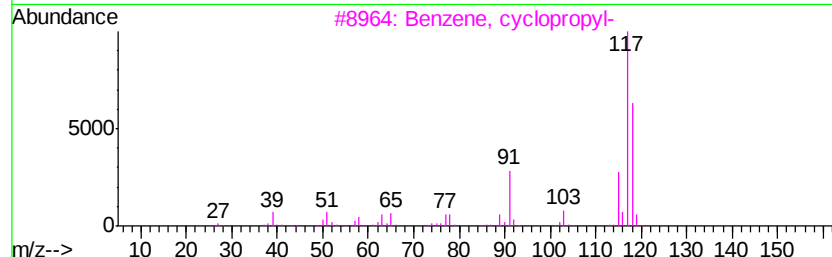
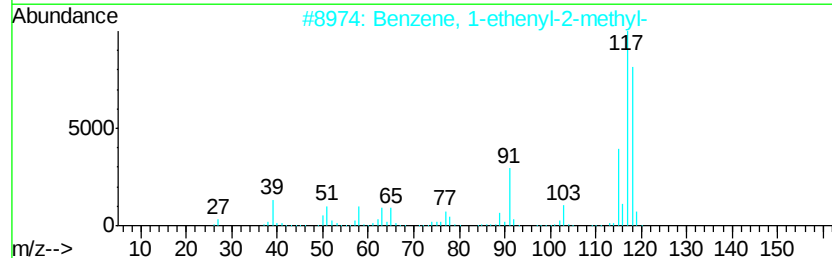
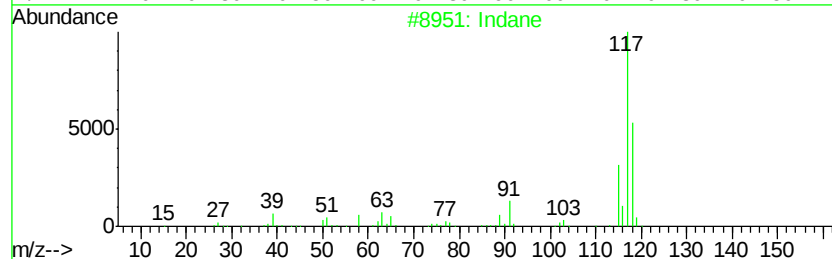
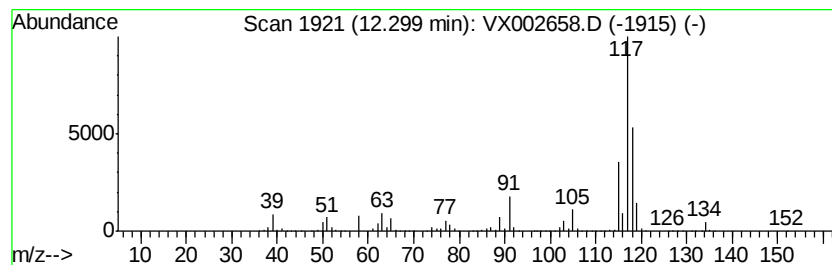
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	49.27 ug/l	911737	1,4-Dichlorobenzene-d4	12.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indane		118 C9H10	000496-11-7 92
2	Benzene, 1-ethenyl-2-methyl-		118 C9H10	000611-15-4 87
3	Benzene, cyclopropyl-		118 C9H10	000873-49-4 76
4	Benzene, 2-propenyl-		118 C9H10	000300-57-2 76
5	(Z)-1-Phenylpropene		118 C9H10	000766-90-5 62



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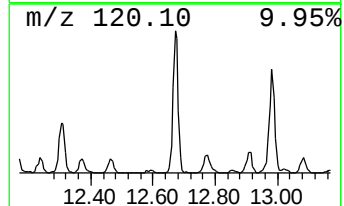
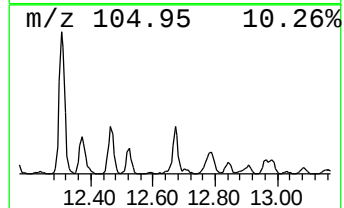
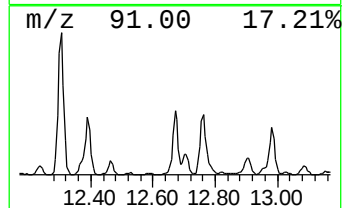
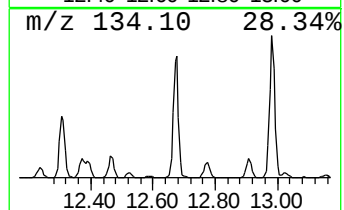
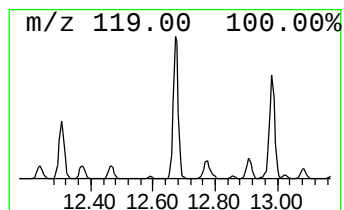
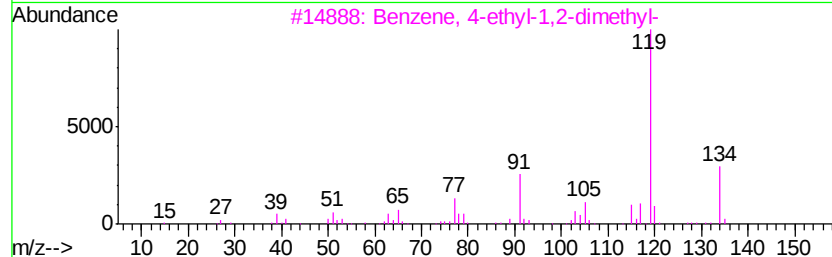
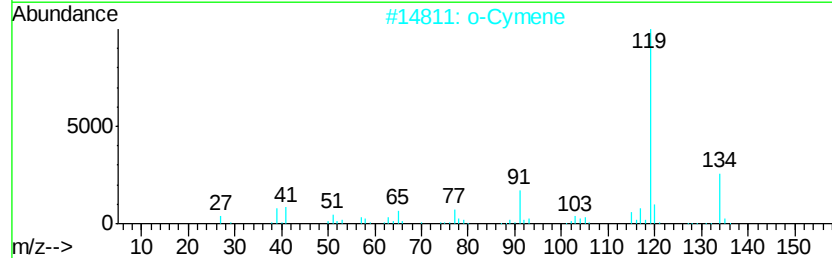
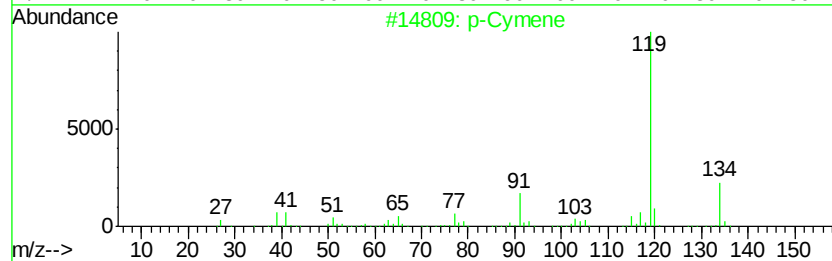
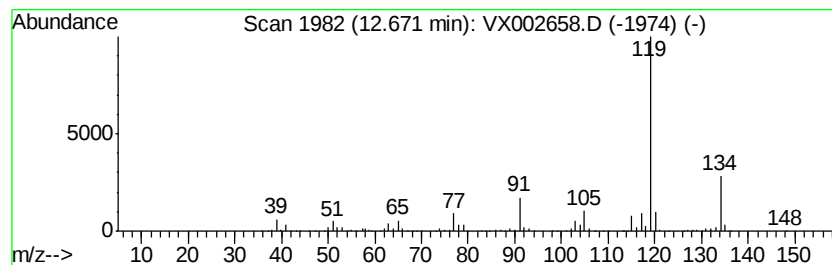
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TIC Integration Parameters: LSCINT.P

Peak Number 4 p-Cymene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	18.88 ug/l	349391	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	p-Cymene	134	C10H14	000099-87-6	95
2		o-Cymene	134	C10H14	000527-84-4	95
3		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061618\
Data File : VX002658.D
Acq On : 16 Jun 2018 20:18
Operator : JC/MD
Sample : J3467-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PR-MW-02-061118

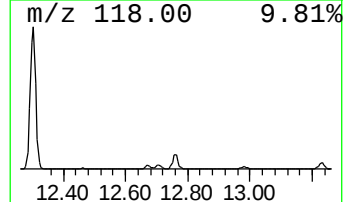
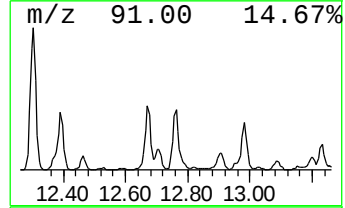
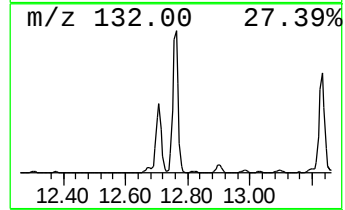
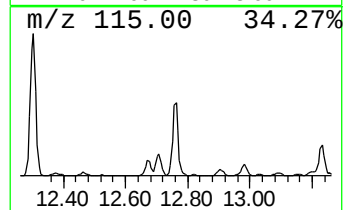
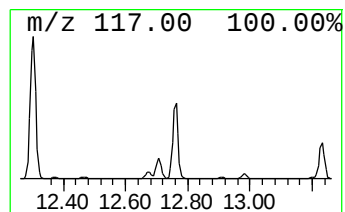
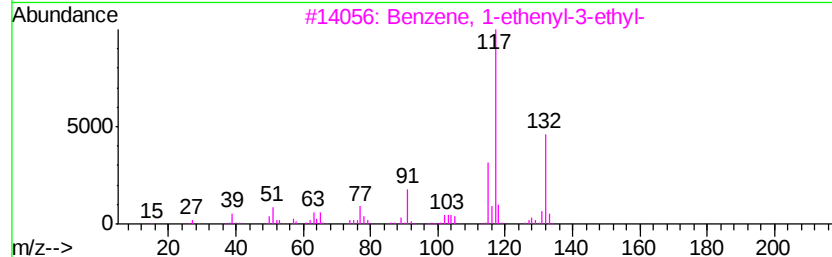
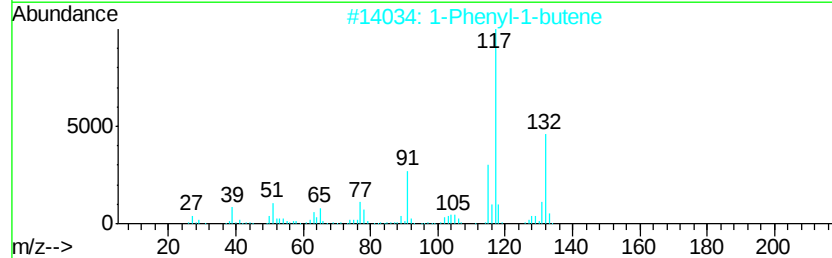
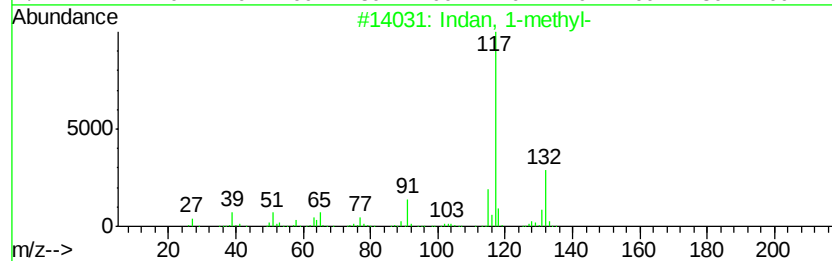
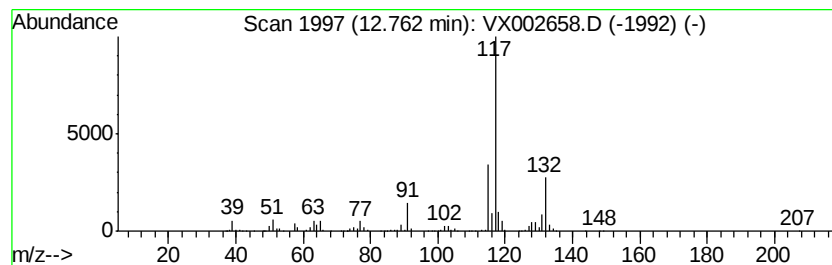
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061518W.M
Quant Title : SW846 8260

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	24.45 ug/l	452468	1,4-Dichlorobenzene-d4	12.09

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	87
2	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
4	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	86
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	80



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061618\
Data File : VX002658.D
Acq On : 16 Jun 2018 20:18
Operator : JC/MD
Sample : J3467-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PR-MW-02-061118

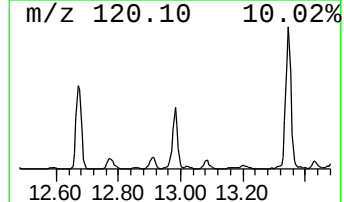
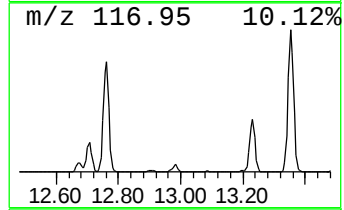
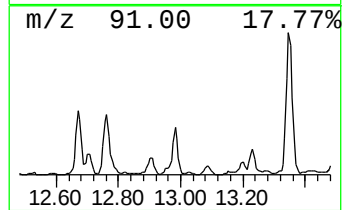
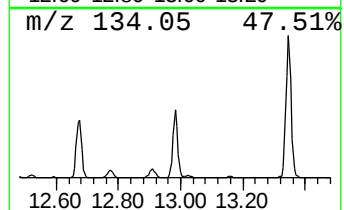
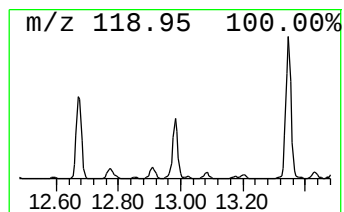
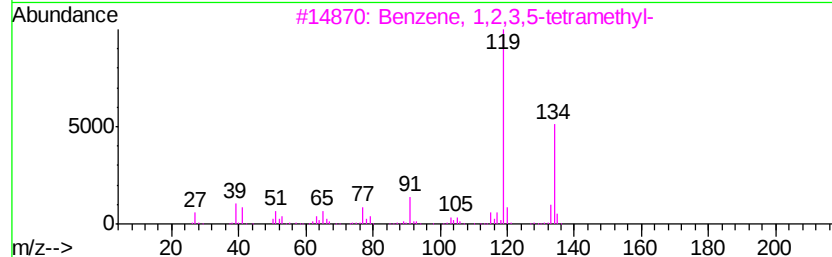
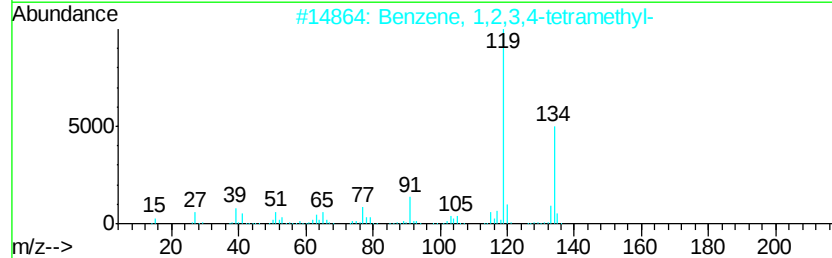
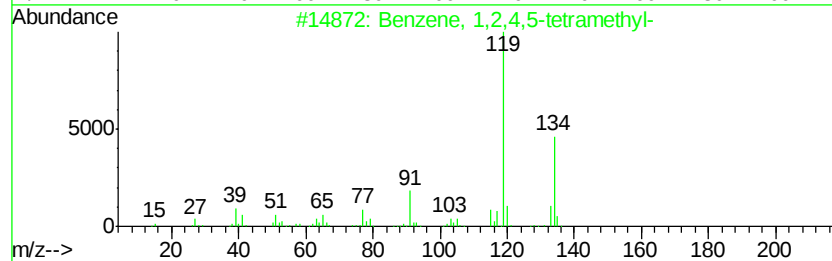
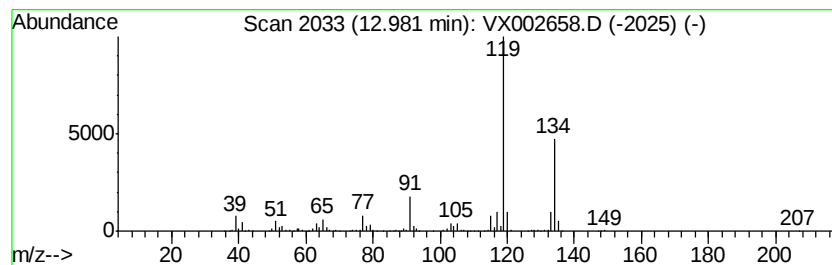
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	14.96 ug/l	276872	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4		o-Cymene	134	C10H14	000527-84-4	94
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061618\
Data File : VX002658.D
Acq On : 16 Jun 2018 20:18
Operator : JC/MD
Sample : J3467-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

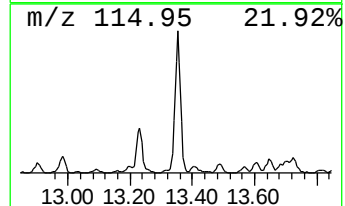
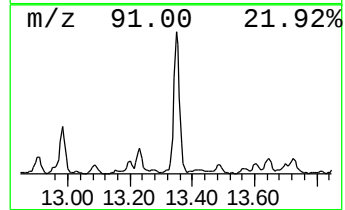
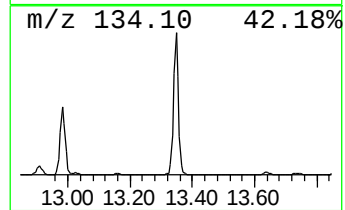
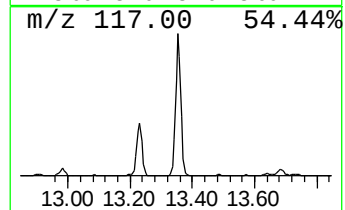
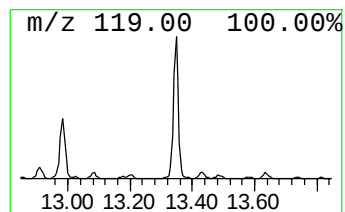
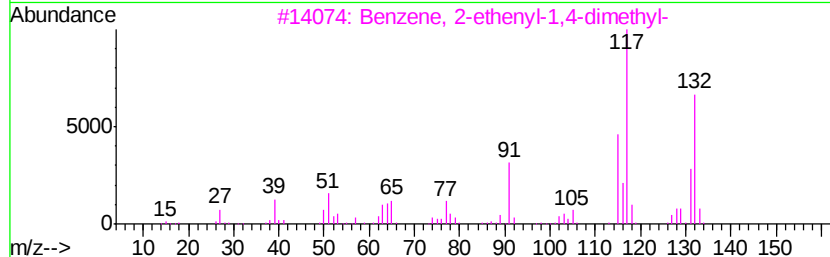
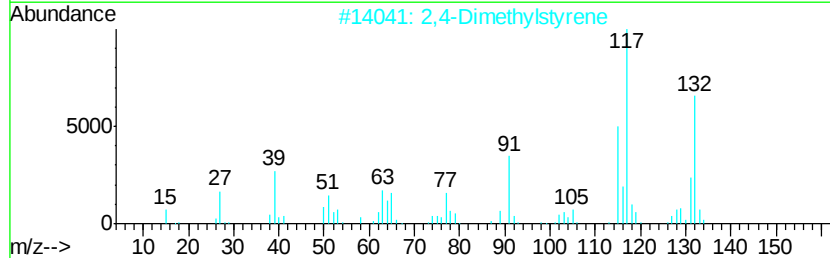
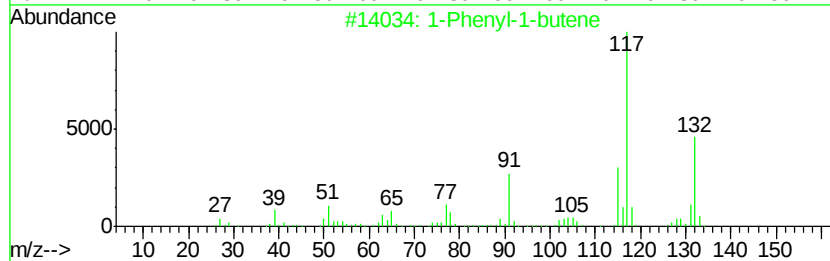
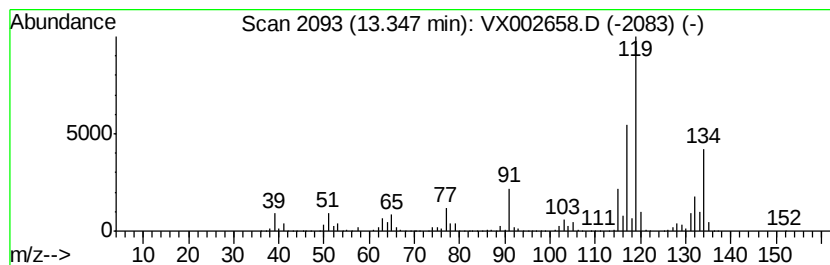
Instrument :
MSVOA_X
ClientSampleId :
PR-MW-02-061118

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061518W.M
Quant Title : SW846 8260

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

Peak Number 7 1-Phenyl-1-butene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	59.48 ug/l	1100630	1,4-Dichlorobenzene-d4	12.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Phenyl-1-butene	132 C10H12	000824-90-8	91
2	2,4-Dimethylstyrene	132 C10H12	002234-20-0	86
3	Benzene, 2-ethenyl-1,4-dimethyl-	132 C10H12	002039-89-6	86
4	3-Phenylbut-1-ene	132 C10H12	000934-10-1	55
5	Benzene, (1-methyl-1-propenyl)-,...	132 C10H12	000767-99-7	55



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061618\
Data File : VX002658.D
Acq On : 16 Jun 2018 20:18
Operator : JC/MD
Sample : J3467-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PR-MW-02-061118

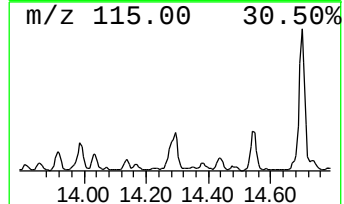
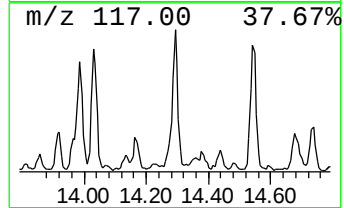
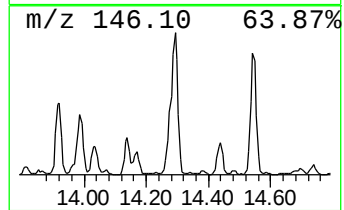
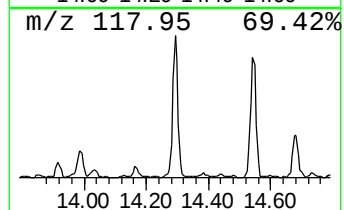
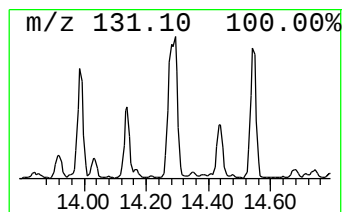
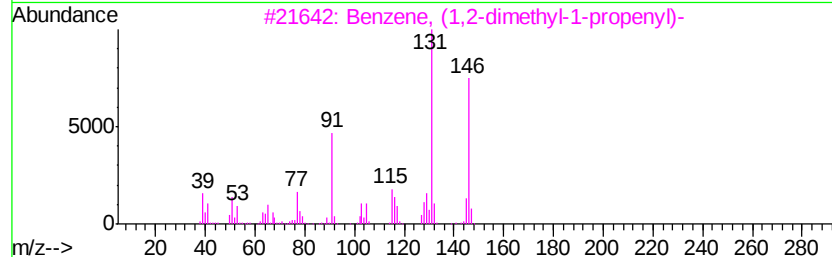
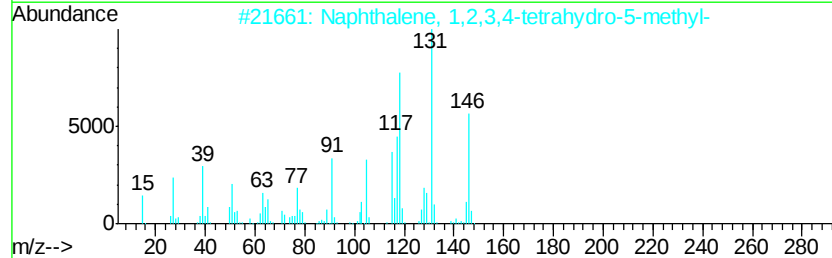
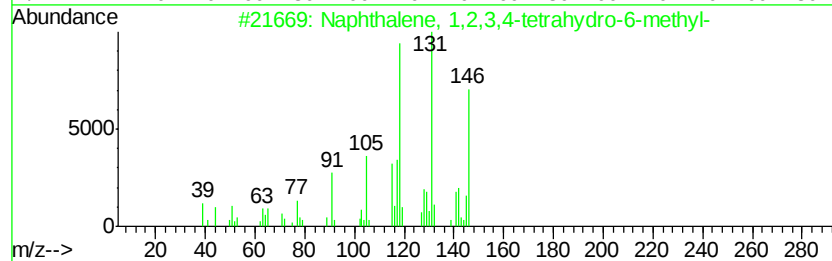
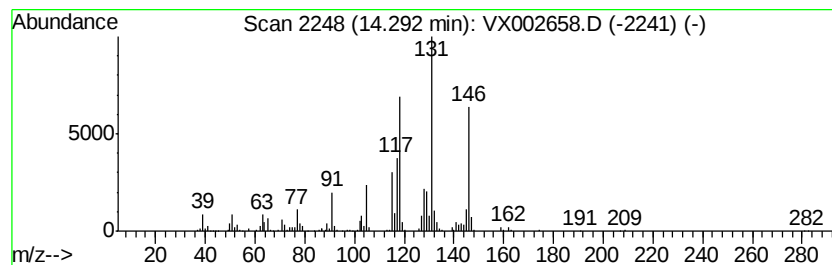
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Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.29	12.71 ug/l	235139	1,4-Dichlorobenzene-d4	12.09

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	93
3		Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	76
4		Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	70
5		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX061618\
Data File : VX002658.D
Acq On : 16 Jun 2018 20:18
Operator : JC/MD
Sample : J3467-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PR-MW-02-061118

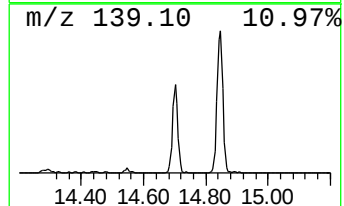
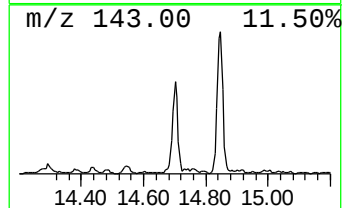
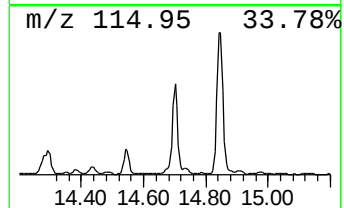
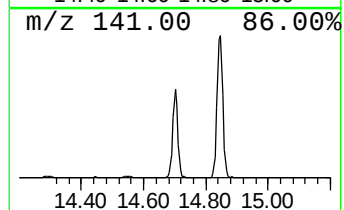
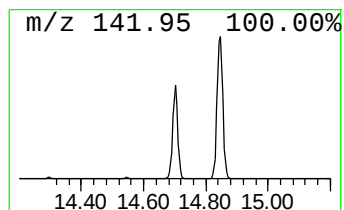
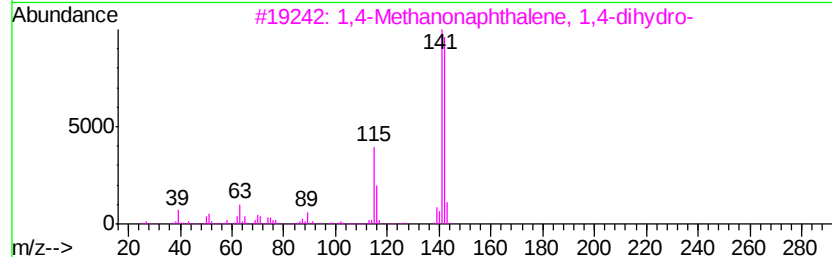
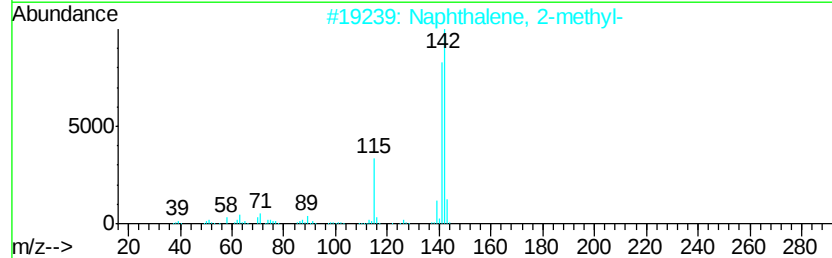
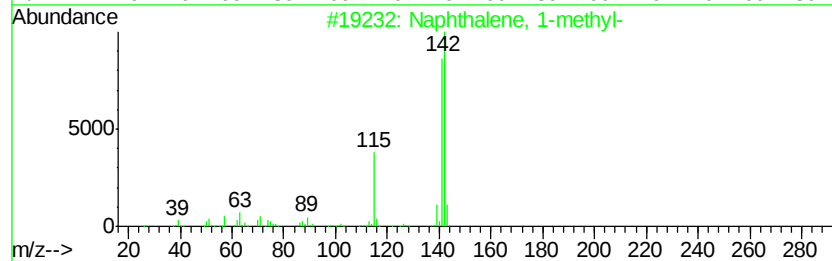
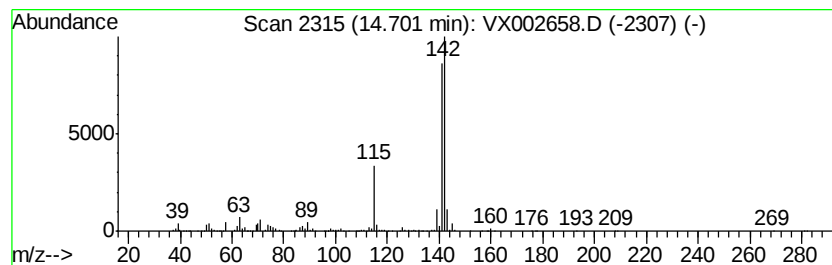
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061518W.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.70	19.90 ug/l	368253	1,4-Dichlorobenzene-d4	12.09

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		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX061618\
Data File : VX002658.D
Acq On : 16 Jun 2018 20:18
Operator : JC/MD
Sample : J3467-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PR-MW-02-061118

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061518W.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-tr...	12.15	23.4	ug/l	433915	4	12.09	925186	50.0
Indane	12.30	49.3	ug/l	911737	4	12.09	925186	50.0
p-Cymene	12.67	18.9	ug/l	349391	4	12.09	925186	50.0
Indan, 1-methyl-	12.76	24.4	ug/l	452468	4	12.09	925186	50.0
Benzene, 1,2,4,5-...	12.98	15.0	ug/l	276872	4	12.09	925186	50.0
1-Phenyl-1-butene	13.35	59.5	ug/l	1100630	4	12.09	925186	50.0
Naphthalene, 1,2,...	14.29	12.7	ug/l	235139	4	12.09	925186	50.0
Naphthalene, 1-me...	14.70	19.9	ug/l	368253	4	12.09	925186	50.0