

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112420\
 Data File : VX019626.D
 Acq On : 25 Nov 2020 00:03
 Operator : JC/MD
 Sample : L4841-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 40741

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\82X112120W.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	1.471	59	63	68	rBV	212075	248884	1.81%	0.170%
2	2.118	162	169	175	rBV	95481	156706	1.14%	0.107%
3	2.422	213	219	234	rVB	1421927	2390046	17.39%	1.635%
4	2.599	241	248	269	rVB	595990	1302308	9.47%	0.891%
5	3.940	458	468	485	rBV	176005	521513	3.79%	0.357%
6	4.635	571	582	596	rBV	198042	577760	4.20%	0.395%
7	5.111	649	660	678	rBV2	57830	207065	1.51%	0.142%
8	5.458	707	717	730	rBV2	166806	485060	3.53%	0.332%
9	5.623	734	744	764	rVB	310267	881635	6.41%	0.603%
10	6.025	800	810	817	rBV	192230	500819	3.64%	0.343%
11	6.111	817	824	835	rVB	141194	377170	2.74%	0.258%
12	6.824	930	941	953	rBV	466420	1144863	8.33%	0.783%
13	7.513	1047	1054	1069	rVB	110202	219976	1.60%	0.150%
14	8.622	1229	1236	1242	rBV	80295	140348	1.02%	0.096%
15	8.696	1242	1248	1254	rVV	994719	1547198	11.25%	1.058%
16	8.763	1254	1259	1265	rVB	652767	981965	7.14%	0.672%
17	10.092	1472	1477	1492	rVB	946131	1334572	9.71%	0.913%
18	10.232	1495	1500	1506	rBV	976631	1292374	9.40%	0.884%
19	10.342	1511	1518	1524	rBV	4179847	5653758	41.13%	3.867%
20	10.543	1546	1551	1556	rBV	123587	154219	1.12%	0.105%
21	10.622	1559	1564	1568	rVV2	87668	138119	1.00%	0.094%
22	10.683	1568	1574	1578	rVV	2775690	3627692	26.39%	2.481%
23	10.720	1578	1580	1588	rVB	329209	369798	2.69%	0.253%
24	10.805	1588	1594	1599	rBV	218282	313850	2.28%	0.215%
25	10.915	1606	1612	1620	rVB2	85679	197051	1.43%	0.135%
26	11.000	1620	1626	1631	rBV	344832	430690	3.13%	0.295%
27	11.122	1640	1646	1652	rBV	819272	1100646	8.01%	0.753%
28	11.341	1674	1682	1688	rBV	1388164	2022496	14.71%	1.383%
29	11.421	1689	1695	1702	rVV	5441777	9464487	68.85%	6.473%
30	11.488	1702	1706	1720	rVV	2617070	3707528	26.97%	2.536%
31	11.598	1720	1724	1728	rVV2	116801	175026	1.27%	0.120%
32	11.652	1728	1733	1742	rVB	3142782	3807284	27.70%	2.604%
33	11.793	1751	1756	1761	rVB	11502768	13746937	100.00%	9.402%
34	11.878	1766	1770	1772	rBV2	131110	170719	1.24%	0.117%

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35	11.933	1775	1779	1784	rVB2	262934	366179	2.66%	0.250%
36	12.012	1784	1792	1795	rBV	384275	511934	3.72%	0.350%
37	12.061	1795	1800	1805	rVB2	1046980	1400008	10.18%	0.957%
38	12.128	1805	1811	1816	rBV	4619996	5483533	39.89%	3.750%
39	12.256	1827	1832	1835	rBV	8082551	9970531	72.53%	6.819%
40	12.280	1835	1836	1839	rVB	2395711	2079218	15.12%	1.422%
41	12.354	1844	1848	1855	rVB3	2062295	3141530	22.85%	2.149%
42	12.457	1858	1865	1868	rBV3	387044	670621	4.88%	0.459%
43	12.500	1868	1872	1878	rVB	795454	992570	7.22%	0.679%
44	12.567	1878	1883	1885	rBV	1377514	1755976	12.77%	1.201%
45	12.591	1885	1887	1892	rVV	1402565	1656481	12.05%	1.133%
46	12.652	1892	1897	1900	rVV	2417876	2760876	20.08%	1.888%
47	12.683	1900	1902	1906	rVB	380497	403256	2.93%	0.276%
48	12.738	1906	1911	1919	rBV2	1050766	1877867	13.66%	1.284%
49	12.835	1924	1927	1930	rBV	156015	184768	1.34%	0.126%
50	12.884	1932	1935	1940	rVB	734150	916964	6.67%	0.627%
51	12.963	1940	1948	1951	rBV	1613437	2138034	15.55%	1.462%
52	13.000	1951	1954	1961	rVB	2712642	3216190	23.40%	2.200%
53	13.061	1961	1964	1965	rBV	181042	184914	1.35%	0.126%
54	13.079	1965	1967	1970	rVB3	147420	161268	1.17%	0.110%
55	13.207	1980	1988	1992	rBV	2019368	2592487	18.86%	1.773%
56	13.250	1992	1995	1998	rVB2	225125	242497	1.76%	0.166%
57	13.329	1998	2008	2013	rBV2	4481253	6996755	50.90%	4.785%
58	13.384	2013	2017	2019	rVV3	318872	454280	3.30%	0.311%
59	13.408	2019	2021	2025	rVV2	184480	210571	1.53%	0.144%
60	13.463	2025	2030	2038	rVB	2791688	3643694	26.51%	2.492%
61	13.542	2038	2043	2046	rBV2	314270	441600	3.21%	0.302%
62	13.585	2046	2050	2053	rBV	498876	658296	4.79%	0.450%
63	13.622	2053	2056	2061	rVV3	616663	1096829	7.98%	0.750%
64	13.676	2063	2065	2067	rVV3	201830	268786	1.96%	0.184%
65	13.701	2067	2069	2074	rVB2	544727	764854	5.56%	0.523%
66	13.817	2083	2088	2094	rVB	4660767	5805720	42.23%	3.971%
67	13.890	2094	2100	2105	rVB	805924	1215253	8.84%	0.831%
68	13.963	2105	2112	2117	rBV3	874778	1317685	9.59%	0.901%
69	14.012	2117	2120	2123	rVB	261832	296520	2.16%	0.203%
70	14.073	2127	2130	2133	rVB3	372610	397476	2.89%	0.272%
71	14.115	2133	2137	2141	rBV	705635	840209	6.11%	0.575%

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72	14.219	2149	2154	2156	rBV2	260373	335679	2.44%	0.230%
73	14.268	2156	2162	2166	rVV	2154336	3384323	24.62%	2.315%
74	14.298	2166	2167	2173	rVB	384613	390925	2.84%	0.267%
75	14.359	2173	2177	2182	rBV2	346925	471869	3.43%	0.323%
76	14.414	2182	2186	2190	rVB	885474	1045397	7.60%	0.715%
77	14.457	2190	2193	2198	rVB4	124695	209358	1.52%	0.143%
78	14.518	2198	2203	2209	rBV2	1501597	2088749	15.19%	1.429%
79	14.676	2221	2229	2233	rBV2	1885520	2986728	21.73%	2.043%
80	14.713	2233	2235	2239	rVB2	393700	442255	3.22%	0.302%
81	14.768	2240	2244	2245	rBV	144902	164994	1.20%	0.113%
82	14.792	2245	2248	2250	rVV	379103	420454	3.06%	0.288%
83	14.816	2250	2252	2256	rVV	908108	1104450	8.03%	0.755%
84	14.859	2256	2259	2261	rVV2	196343	265831	1.93%	0.182%
85	14.884	2261	2263	2266	rVV2	181702	212916	1.55%	0.146%
86	14.945	2269	2273	2281	rVB2	258199	415800	3.02%	0.284%
87	15.030	2283	2287	2291	rVB	173737	245004	1.78%	0.168%
88	15.091	2292	2297	2301	rVB2	97643	149636	1.09%	0.102%
89	15.146	2301	2306	2310	rBV2	244997	392968	2.86%	0.269%
90	15.188	2310	2313	2315	rVV	193407	235686	1.71%	0.161%
91	15.225	2315	2319	2321	rVV	543731	774660	5.64%	0.530%
92	15.249	2321	2323	2327	rVB2	517036	594995	4.33%	0.407%
93	15.359	2338	2341	2346	rBV3	104448	169621	1.23%	0.116%
94	15.420	2346	2351	2355	rBV3	225348	360889	2.63%	0.247%
95	15.530	2364	2369	2374	rBV2	585609	993237	7.23%	0.679%
96	15.664	2386	2391	2394	rBV	392596	615982	4.48%	0.421%
97	15.700	2394	2397	2404	rVB	276991	444487	3.23%	0.304%
98	15.896	2424	2429	2434	rBV3	88628	183749	1.34%	0.126%
99	16.139	2465	2469	2479	rBV2	118358	276782	2.01%	0.189%
100	16.432	2511	2517	2522	rVB2	175702	333110	2.42%	0.228%

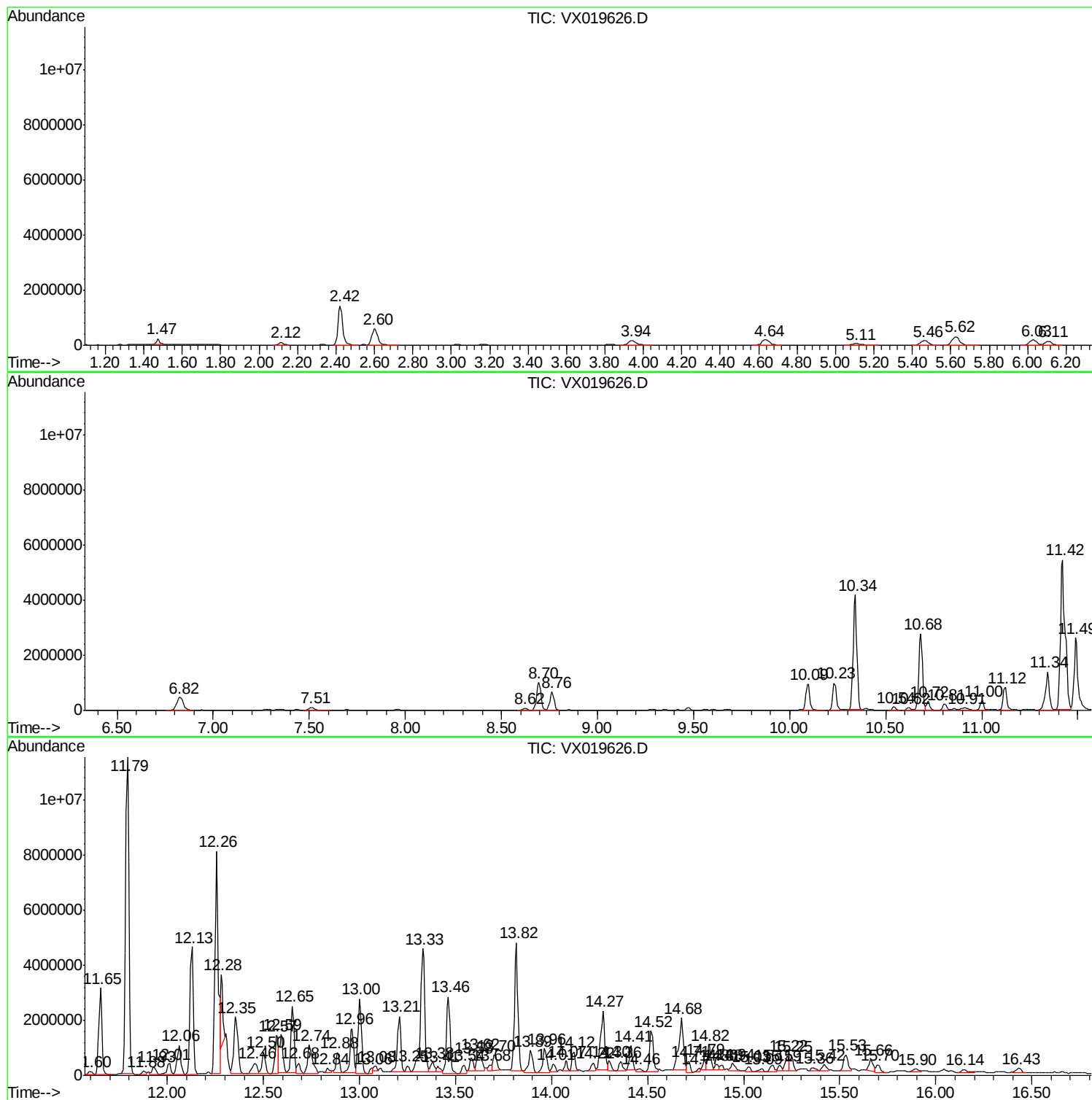
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112120W.M
Quant Title : SW846 8260

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



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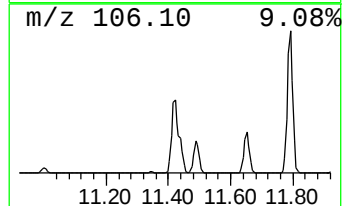
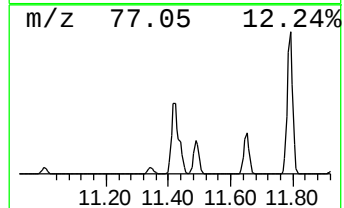
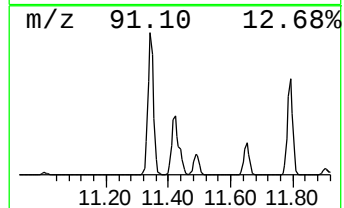
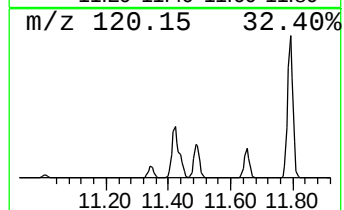
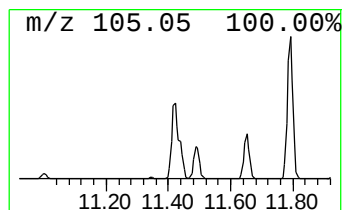
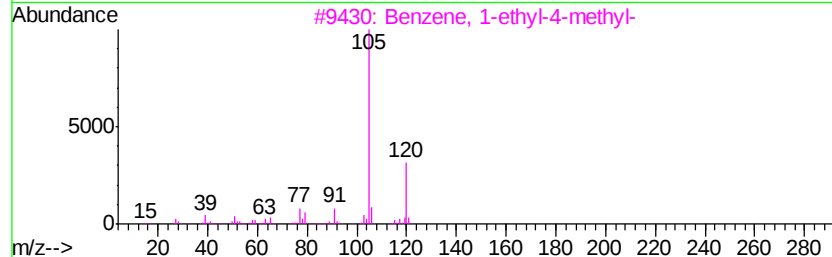
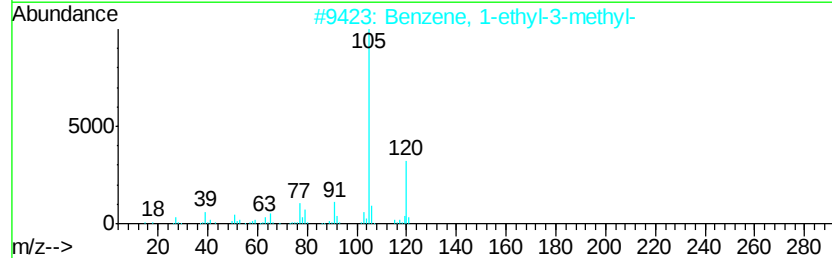
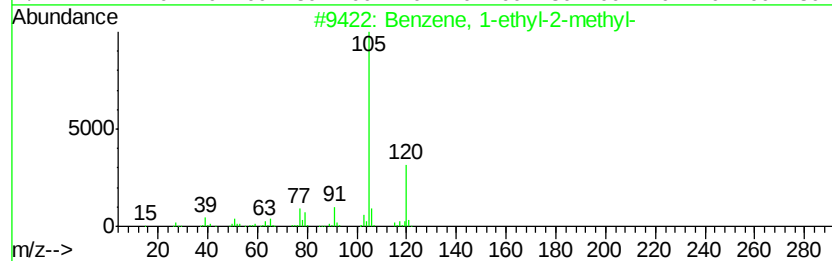
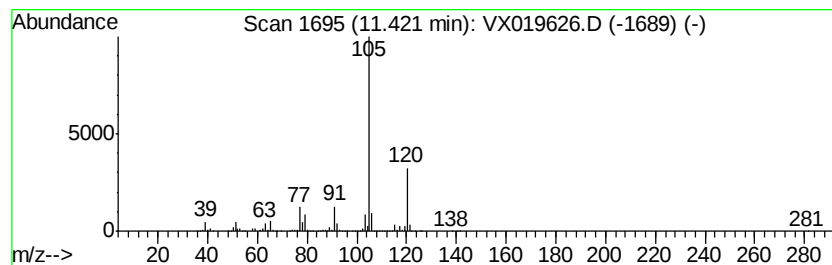
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112120W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.42	338.02 ug/l	9464490	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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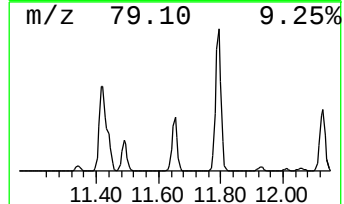
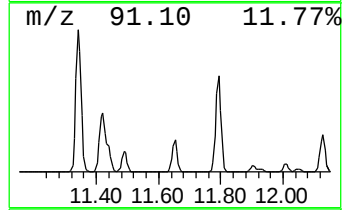
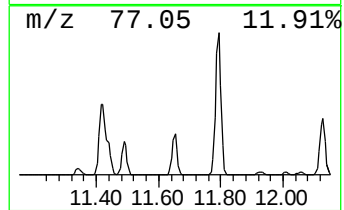
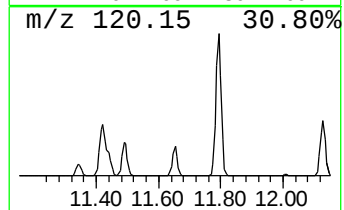
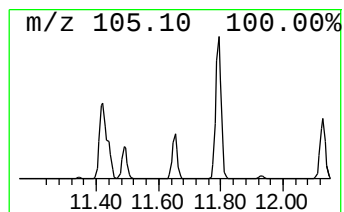
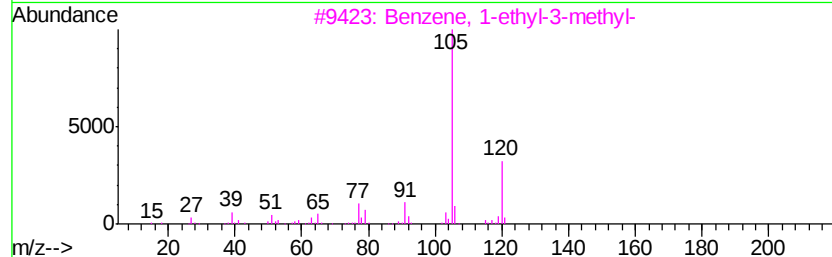
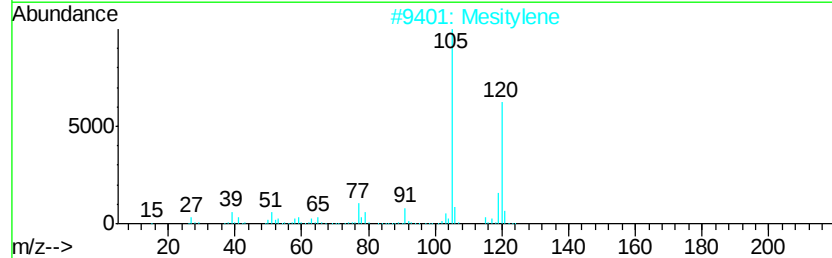
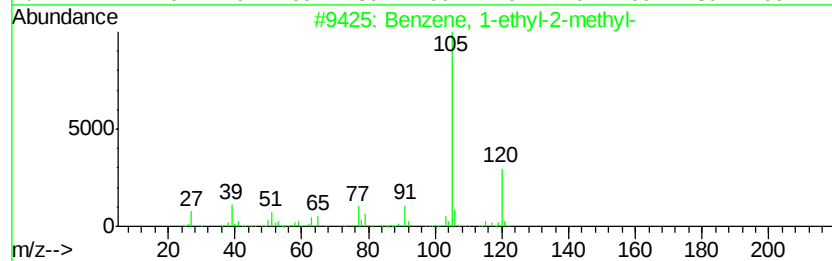
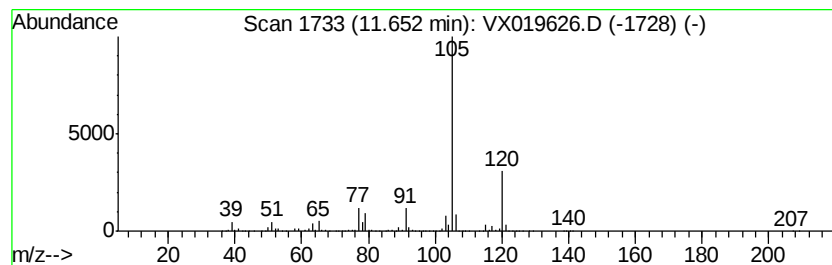
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Mesitylene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.65	135.97 ug/l	3807280	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2		Mesitylene	120	C9H12	000108-67-8	91
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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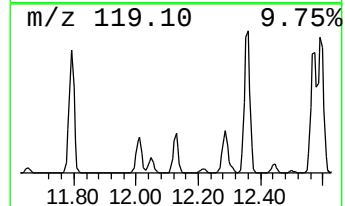
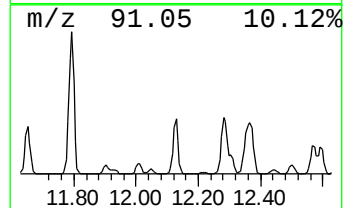
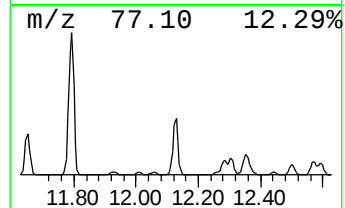
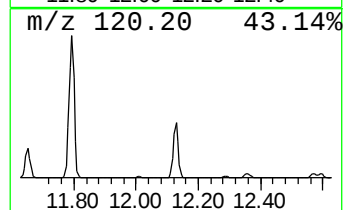
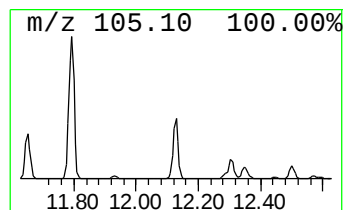
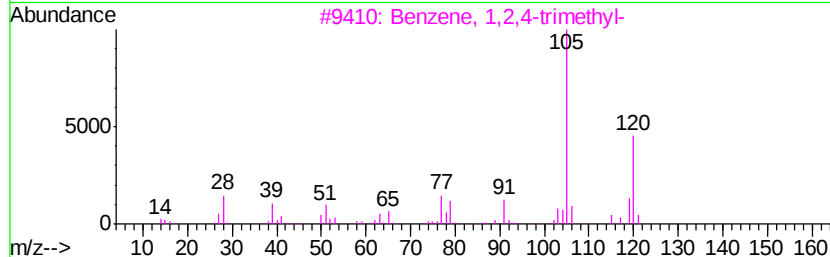
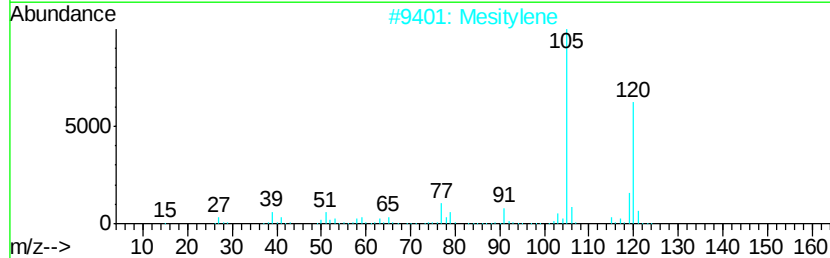
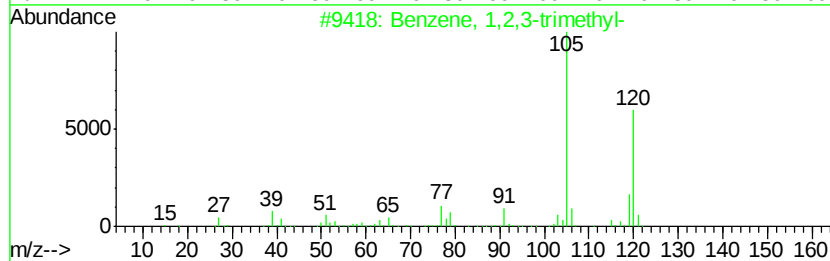
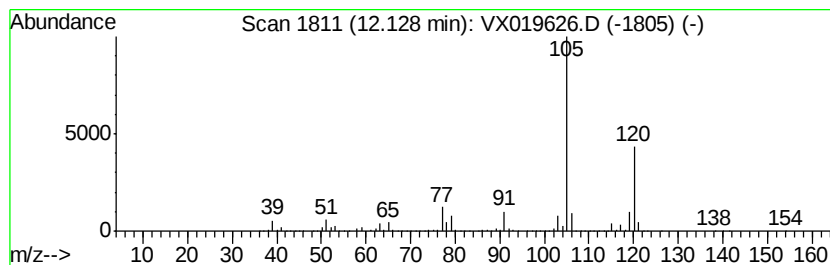
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Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1,2,3-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.13	195.84 ug/l	5483530	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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	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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112420\
Data File : VX019626.D
Acq On : 25 Nov 2020 00:03
Operator : JC/MD
Sample : L4841-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
40741

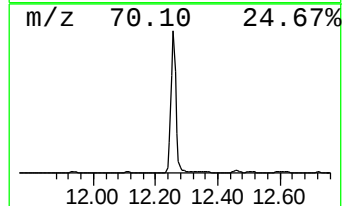
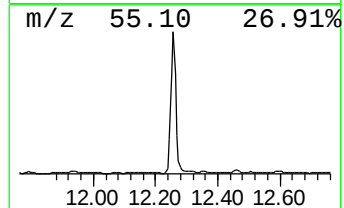
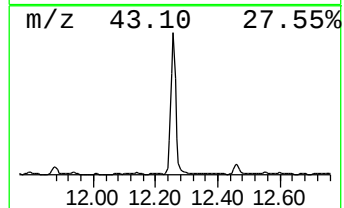
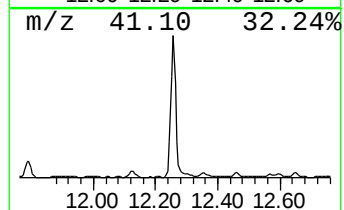
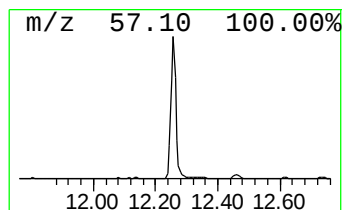
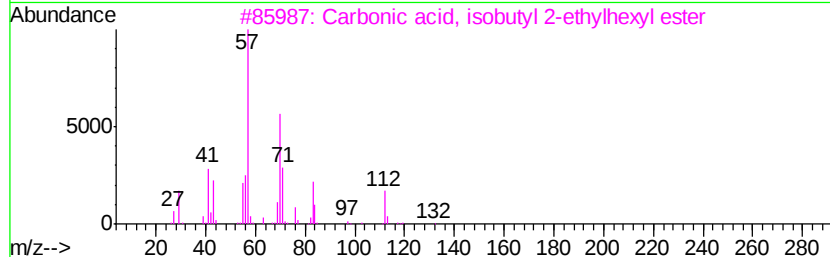
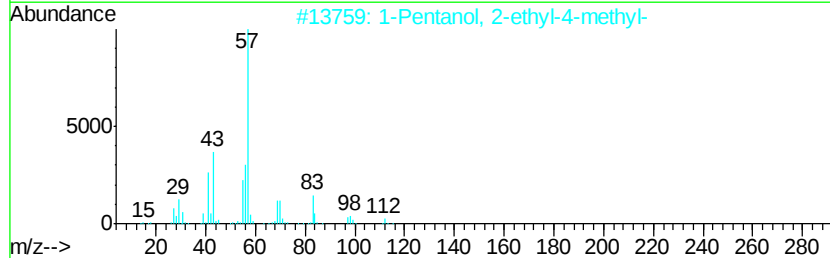
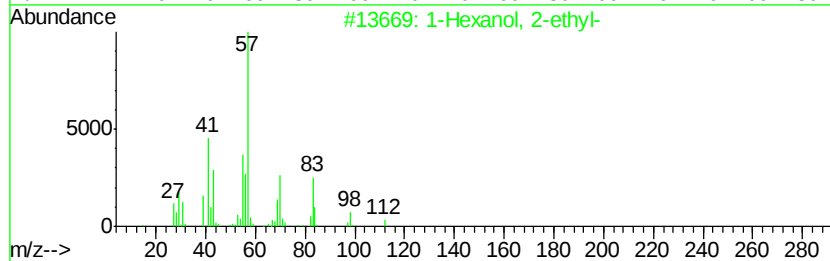
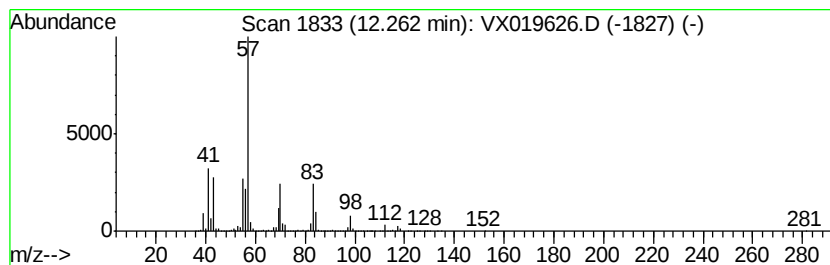
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Quant Title : SW846 8260

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

Peak Number 4 1-Hexanol, 2-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.26	356.09 ug/l	9970530	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	90
2		1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	56
3		Carbonic acid, isobutyl 2-ethylh...	230	C13H26O3	1000357-82-9	53
4		Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	50
5		2-Propyl-1-pentanol	130	C8H18O	058175-57-8	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112420\
Data File : VX019626.D
Acq On : 25 Nov 2020 00:03
Operator : JC/MD
Sample : L4841-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
40741

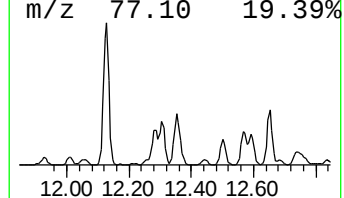
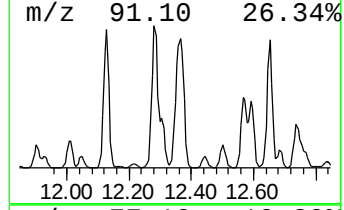
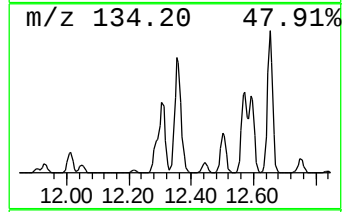
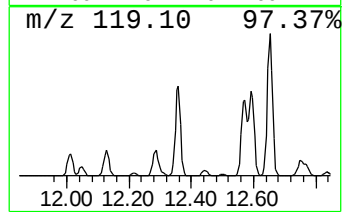
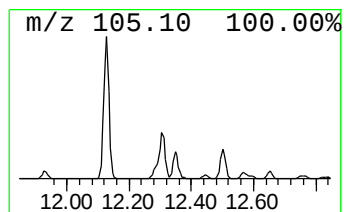
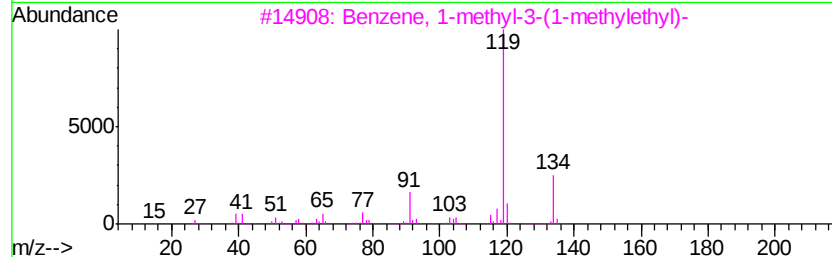
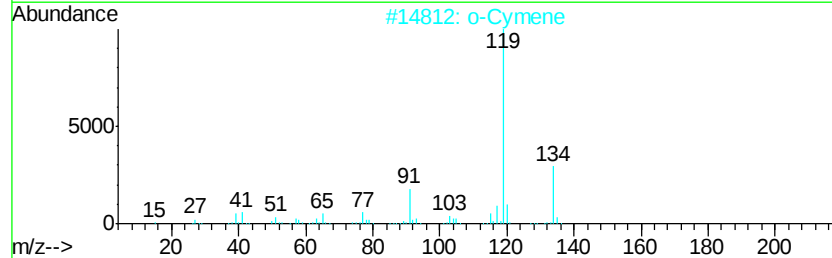
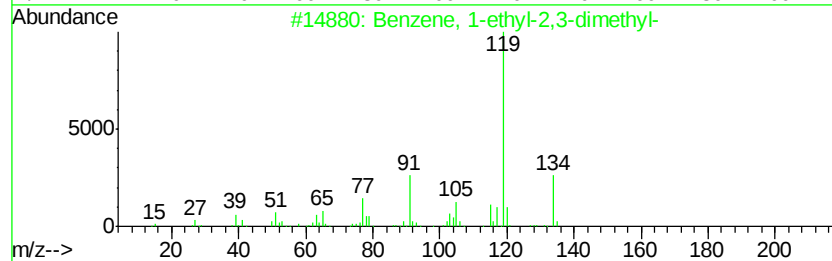
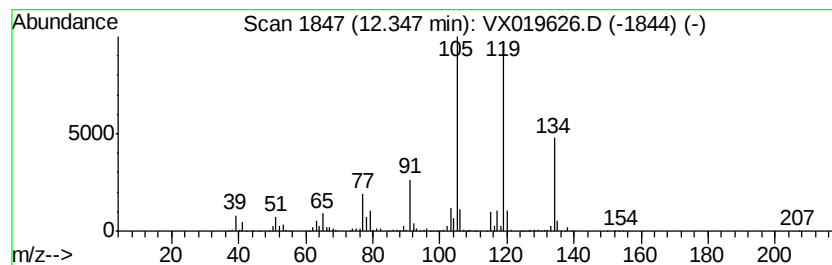
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.35	112.20 ug/l	3141530	1,4-Dichlorobenzene-d4	12.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
2		o-Cymene	134	C10H14	000527-84-4	94
3		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	94
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5		p-Cymene	134	C10H14	000099-87-6	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112420\
Data File : VX019626.D
Acq On : 25 Nov 2020 00:03
Operator : JC/MD
Sample : L4841-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
40741

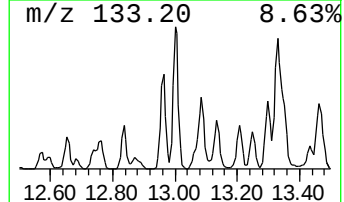
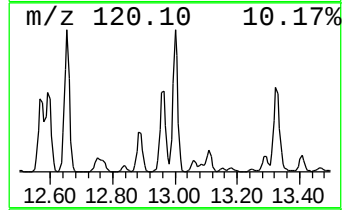
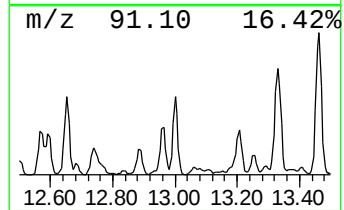
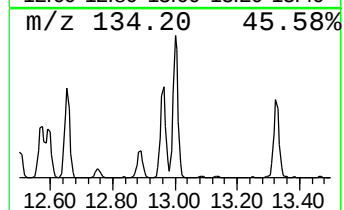
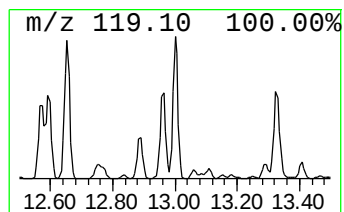
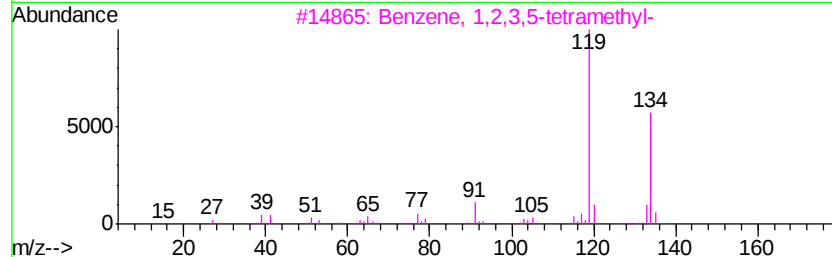
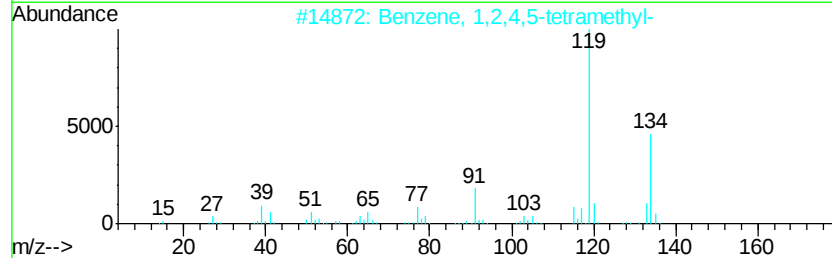
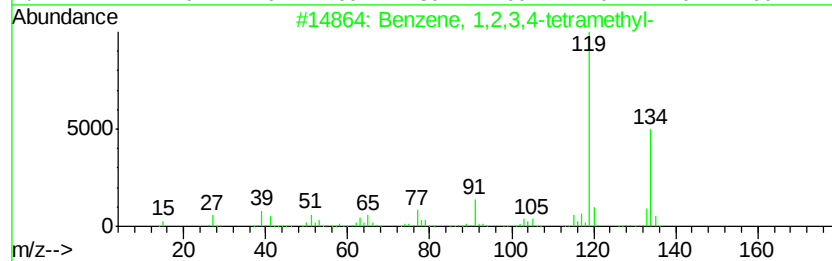
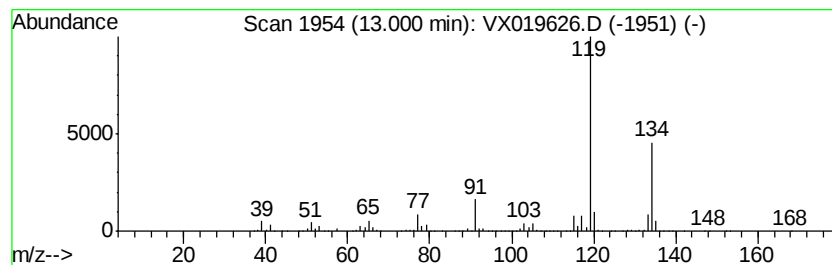
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.00	114.86 ug/l	3216190	1,4-Dichlorobenzene-d4	12.06

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112420\
Data File : VX019626.D
Acq On : 25 Nov 2020 00:03
Operator : JC/MD
Sample : L4841-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
40741

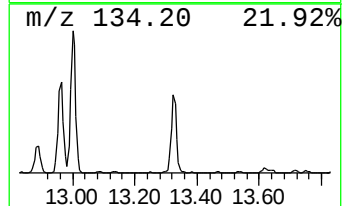
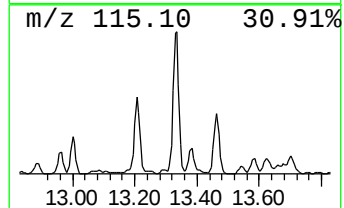
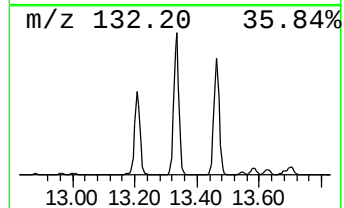
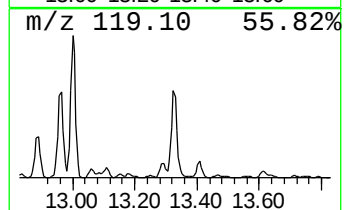
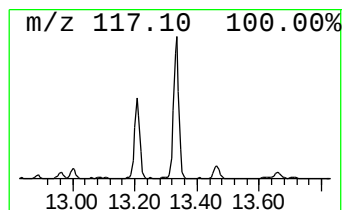
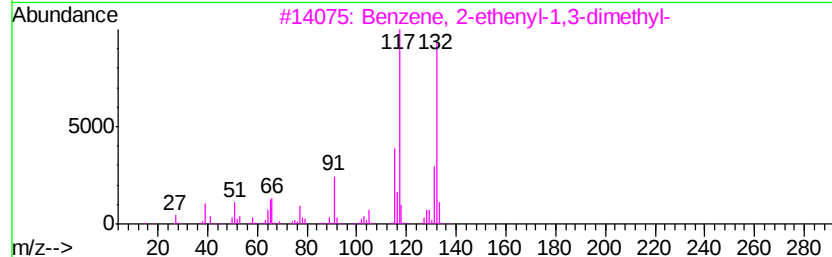
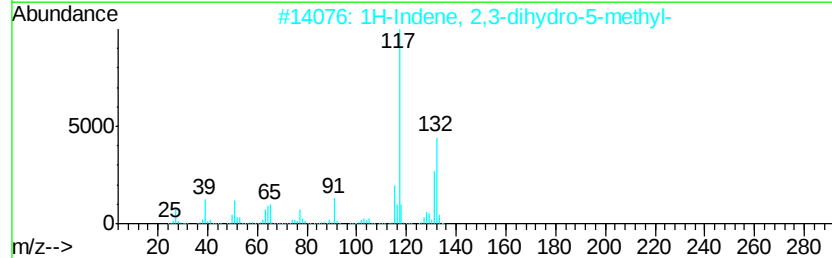
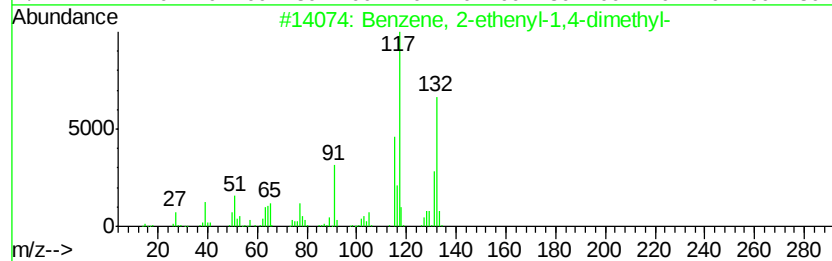
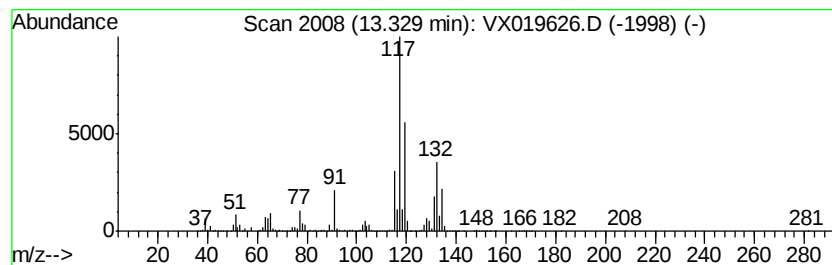
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.33	249.88 ug/l	6996760	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
3		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	70
4		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	70
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112420\
Data File : VX019626.D
Acq On : 25 Nov 2020 00:03
Operator : JC/MD
Sample : L4841-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
40741

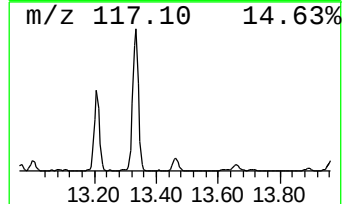
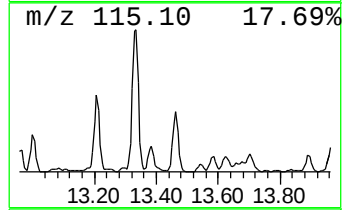
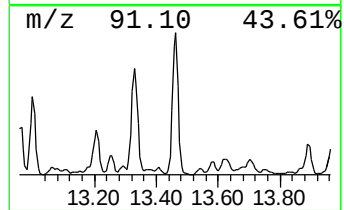
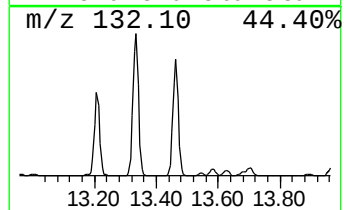
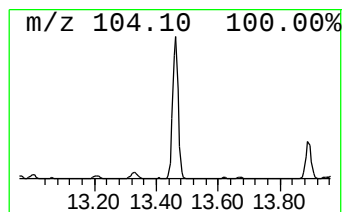
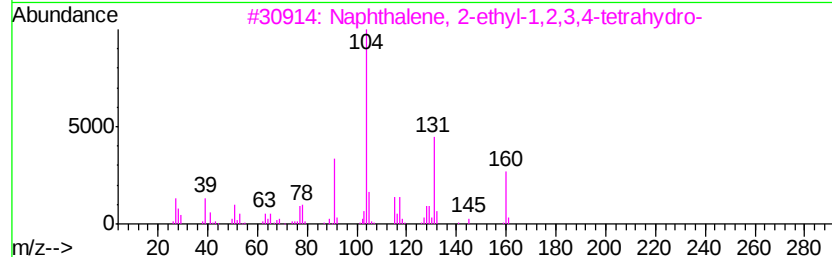
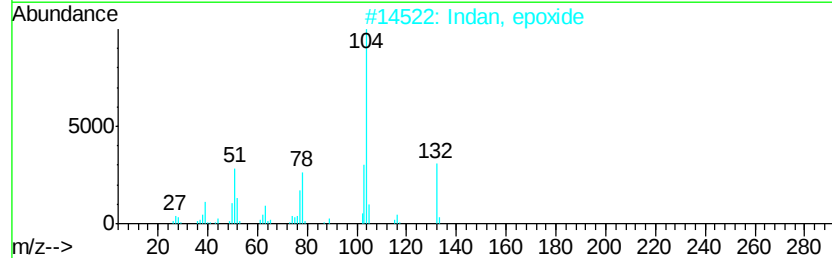
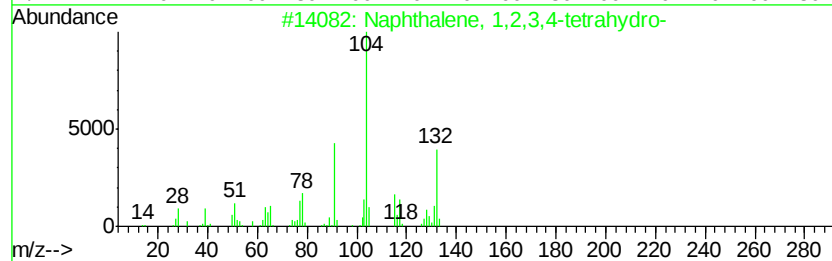
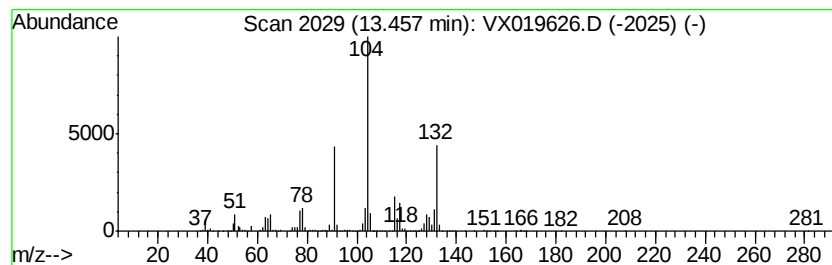
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Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.46	130.13 ug/l	3643690	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	93
2		Indan, epoxide	132	C9H8O	000768-22-9	50
3		Naphthalene, 2-ethyl-1,2,3,4-tetrahydro-	160	C12H16	032367-54-7	40
4		Benzene, 1,1'-(1,2-cyclobutanediyl)-	208	C16H16	007694-30-6	38
5		2-Furanone, 4-phenyltetrahydro-	162	C10H10O2	1000197-38-4	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112420\
Data File : VX019626.D
Acq On : 25 Nov 2020 00:03
Operator : JC/MD
Sample : L4841-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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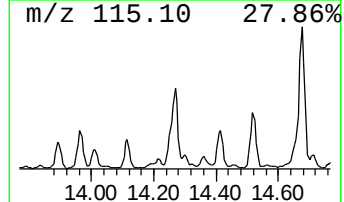
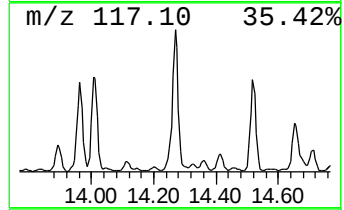
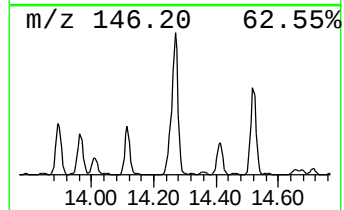
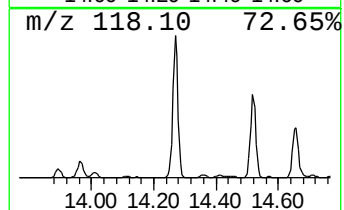
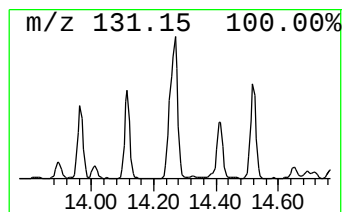
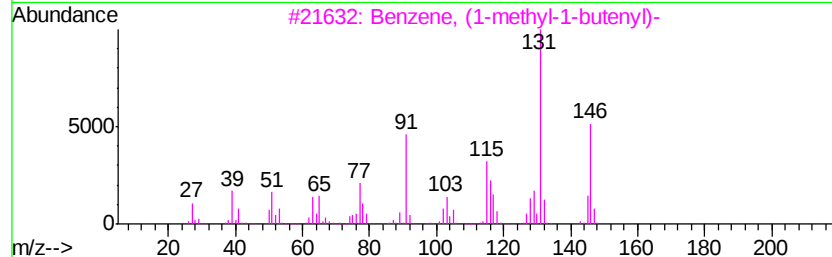
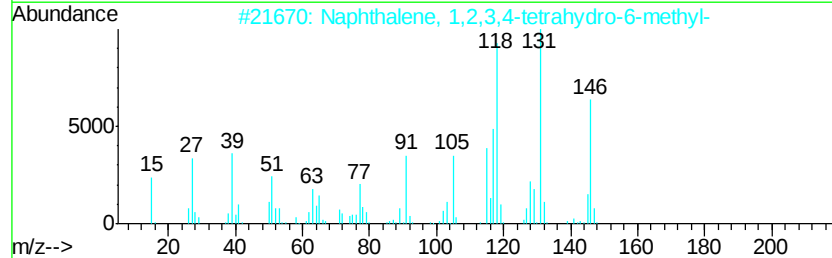
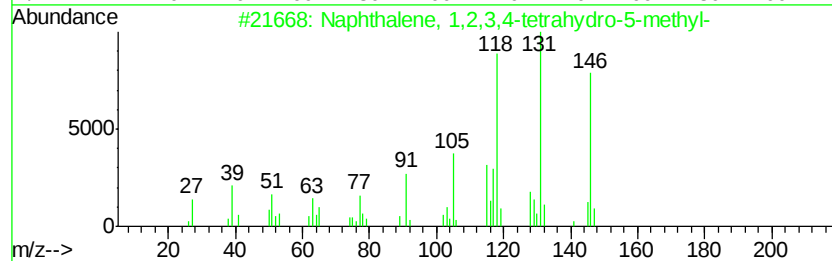
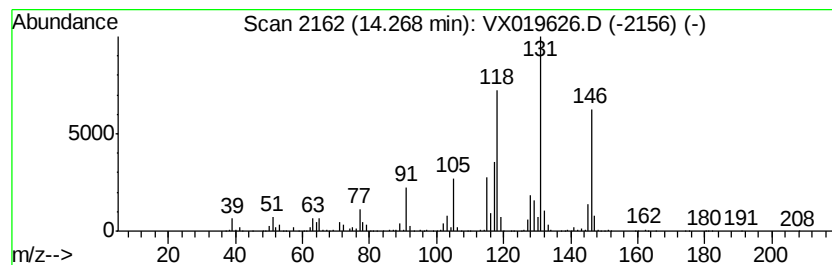
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetra... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.27	120.87 ug/l	3384320	1,4-Dichlorobenzene-d4	12.06

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	97
3		Benzene, (1-methyl-1-butenvl)-	146	C11H14	053172-84-2	76
4		Benzene, (3-methyl-2-butenvl)-	146	C11H14	004489-84-3	70
5		Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112420\
Data File : VX019626.D
Acq On : 25 Nov 2020 00:03
Operator : JC/MD
Sample : L4841-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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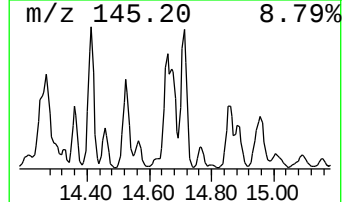
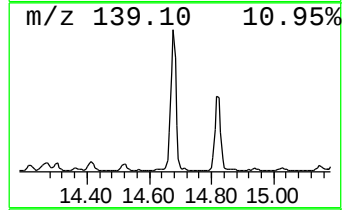
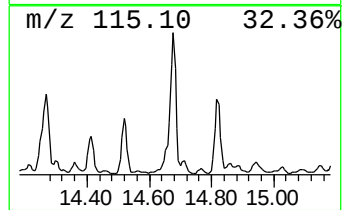
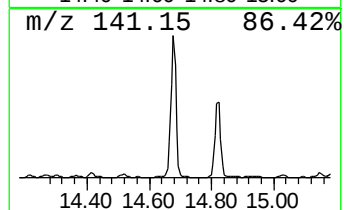
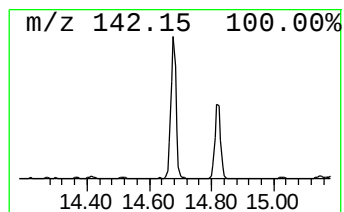
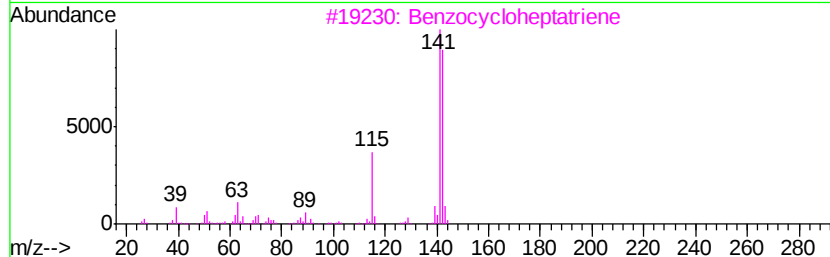
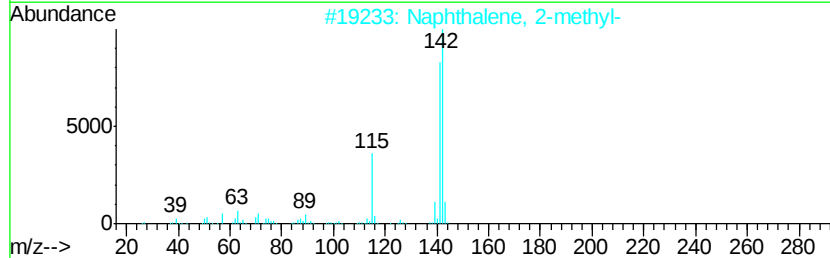
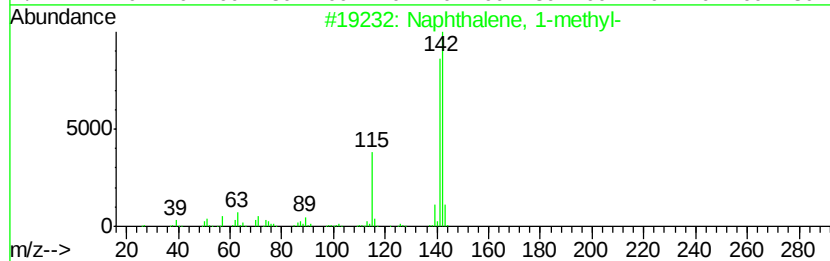
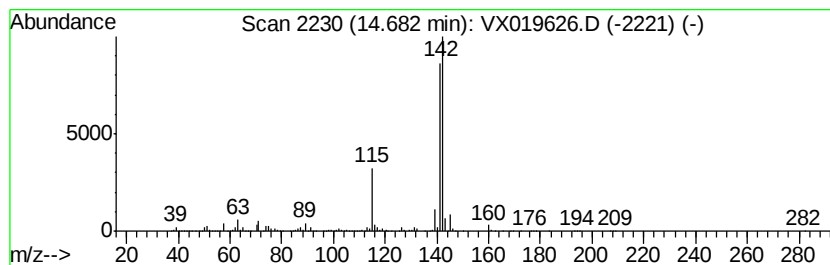
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112120W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.68	106.67 ug/l	2986730	1,4-Dichlorobenzene-d4	12.06

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	94
3		Benzocycloheptatriene	142	C11H10	000264-09-5	90
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	87
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX112420\
Data File : VX019626.D
Acq On : 25 Nov 2020 00:03
Operator : JC/MD
Sample : L4841-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
40741

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112120W.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-ethyl-...	11.42	338.0	ug/l	9464490	4	12.06	1400010	50.0
Mesitylene	11.65	136.0	ug/l	3807280	4	12.06	1400010	50.0
Benzene, 1,2,3-tr...	12.13	195.8	ug/l	5483530	4	12.06	1400010	50.0
1-Hexanol, 2-ethyl-	12.26	356.1	ug/l	9970530	4	12.06	1400010	50.0
Benzene, 1-ethyl-...	12.35	112.2	ug/l	3141530	4	12.06	1400010	50.0
Benzene, 1,2,3,4-...	13.00	114.9	ug/l	3216190	4	12.06	1400010	50.0
Benzene, 2-etheny...	13.33	249.9	ug/l	6996760	4	12.06	1400010	50.0
Naphthalene, 1,2,...	13.46	130.1	ug/l	3643690	4	12.06	1400010	50.0
Naphthalene, 1,2,...	14.27	120.9	ug/l	3384320	4	12.06	1400010	50.0
Naphthalene, 1-me...	14.68	106.7	ug/l	2986730	4	12.06	1400010	50.0