

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
 Data File : VX006007.D
 Acq On : 15 Nov 2018 06:46
 Operator : JC/MD
 Sample : J5915-02
 Misc : 5.00g/10ml/100ul/5.00mL/MSVOA_X/MEOH
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 T-2ME

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\82X110718W.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	4.745	577	589	601	rVB2	2755	9985	1.62%	0.093%
2	5.214	653	666	676	rBV5	3390	10628	1.72%	0.099%
3	5.501	701	713	727	rBV2	69372	215261	34.83%	2.006%
4	5.653	728	738	756	rVB	130555	377736	61.13%	3.520%
5	6.068	795	806	826	rBV3	77096	213858	34.61%	1.993%
6	6.324	839	848	860	rVB2	3360	10324	1.67%	0.096%
7	6.860	925	936	957	rBV	201667	475613	76.97%	4.432%
8	8.049	1124	1131	1141	rVB2	4108	8602	1.39%	0.080%
9	8.336	1173	1178	1182	rBV2	3938	6489	1.05%	0.060%
10	8.494	1200	1204	1214	rVB	4302	8027	1.30%	0.075%
11	8.713	1234	1240	1254	rVB	357122	576157	93.24%	5.369%
12	8.970	1276	1282	1290	rBV2	6186	10473	1.69%	0.098%
13	9.555	1371	1378	1385	rBV2	5419	11171	1.81%	0.104%
14	9.957	1440	1444	1452	rVB	23922	34250	5.54%	0.319%
15	10.061	1455	1461	1465	rBV	19585	27287	4.42%	0.254%
16	10.116	1465	1470	1481	rVB	383557	553306	89.54%	5.156%
17	10.256	1487	1493	1497	rVB	22454	30543	4.94%	0.285%
18	10.360	1497	1510	1515	rVV	65480	99924	16.17%	0.931%
19	10.414	1515	1519	1529	rVB2	28289	49253	7.97%	0.459%
20	10.640	1552	1556	1560	rVV	12266	16741	2.71%	0.156%
21	10.701	1560	1566	1568	rVV	12369	17250	2.79%	0.161%
22	10.725	1568	1570	1573	rVB	7471	8065	1.31%	0.075%
23	10.835	1580	1588	1592	rBV2	13925	25032	4.05%	0.233%
24	10.914	1592	1601	1604	rBV2	9506	17623	2.85%	0.164%
25	11.018	1612	1618	1623	rBV	13201	21132	3.42%	0.197%
26	11.085	1625	1629	1632	rBV	12485	17811	2.88%	0.166%
27	11.140	1632	1638	1648	rVB3	329817	548478	88.76%	5.111%
28	11.244	1648	1655	1659	rBV	33972	42980	6.96%	0.401%
29	11.359	1671	1674	1681	rVB	69809	92783	15.01%	0.865%
30	11.439	1681	1687	1694	rBV	193871	371817	60.17%	3.465%
31	11.506	1694	1698	1710	rVB2	134670	276969	44.82%	2.581%
32	11.603	1710	1714	1717	rBV	7407	9062	1.47%	0.084%
33	11.670	1720	1725	1732	rBV	85815	121780	19.71%	1.135%
34	11.737	1732	1736	1738	rBV2	12368	14975	2.42%	0.140%

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35	11.768	1738	1741	1743	rBV	13041	12102	1.96%	0.113%
36	11.810	1743	1748	1757	rVB	482400	617952	100.00%	5.759%
37	11.896	1757	1762	1764	rBV	21195	31666	5.12%	0.295%
38	11.920	1764	1766	1768	rBV	7522	6837	1.11%	0.064%
39	11.945	1768	1770	1774	rVB	20924	22664	3.67%	0.211%
40	11.987	1774	1777	1780	rBV	30218	38377	6.21%	0.358%
41	12.024	1780	1783	1787	rVB	30573	34675	5.61%	0.323%
42	12.079	1787	1792	1797	rBV	425484	570477	92.32%	5.316%
43	12.146	1797	1803	1807	rBV3	101176	163498	26.46%	1.524%
44	12.182	1807	1809	1810	rVV	13243	12558	2.03%	0.117%
45	12.201	1810	1812	1814	rVV	21480	22410	3.63%	0.209%
46	12.225	1814	1816	1824	rVB2	38117	48867	7.91%	0.455%
47	12.323	1824	1832	1836	rBV2	205616	391217	63.31%	3.646%
48	12.371	1836	1840	1847	rVB4	297072	480011	77.68%	4.473%
49	12.438	1847	1851	1852	rBV2	10958	14174	2.29%	0.132%
50	12.475	1852	1857	1861	rVB2	80804	96730	15.65%	0.901%
51	12.524	1861	1865	1871	rVB2	81231	122193	19.77%	1.139%
52	12.591	1871	1876	1877	rBV	141033	171510	27.75%	1.598%
53	12.609	1877	1879	1885	rVB	125254	143153	23.17%	1.334%
54	12.670	1885	1889	1893	rBV	254786	280389	45.37%	2.613%
55	12.774	1899	1906	1912	rVB4	54158	146537	23.71%	1.366%
56	12.853	1913	1919	1924	rBV4	21263	42040	6.80%	0.392%
57	12.908	1924	1928	1931	rVV	46374	58910	9.53%	0.549%
58	12.981	1931	1940	1943	rVV	141189	234753	37.99%	2.188%
59	13.018	1943	1946	1951	rVB	204968	275960	44.66%	2.572%
60	13.079	1951	1956	1957	rBV	52135	56598	9.16%	0.527%
61	13.103	1957	1960	1962	rVV2	85663	121268	19.62%	1.130%
62	13.127	1962	1964	1968	rVV	59980	60829	9.84%	0.567%
63	13.170	1968	1971	1973	rVV	18058	19102	3.09%	0.178%
64	13.219	1973	1979	1984	rVB2	88833	148410	24.02%	1.383%
65	13.268	1984	1987	1990	rBV2	45412	50631	8.19%	0.472%
66	13.316	1990	1995	1998	rVV3	103497	169908	27.50%	1.583%
67	13.353	1998	2001	2010	rVV5	160580	254720	41.22%	2.374%
68	13.426	2010	2013	2018	rVB2	69294	87587	14.17%	0.816%
69	13.481	2018	2022	2031	rVB6	23243	51567	8.34%	0.481%
70	13.560	2031	2035	2038	rBV3	31433	49244	7.97%	0.459%
71	13.603	2038	2042	2044	rBV	38395	45186	7.31%	0.421%

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72	13.633	2044	2047	2055	rVB2	89817	165449	26.77%	1.542%
73	13.731	2059	2063	2068	rVB3	53877	79806	12.91%	0.744%
74	13.774	2068	2070	2076	rVB3	14390	21888	3.54%	0.204%
75	13.828	2076	2079	2086	rVB3	20084	36648	5.93%	0.342%
76	13.889	2086	2089	2092	rBV4	16228	29289	4.74%	0.273%
77	13.987	2100	2105	2109	rVB4	35820	57706	9.34%	0.538%
78	14.030	2109	2112	2115	rVV	20279	24501	3.96%	0.228%
79	14.091	2119	2122	2125	rVB3	49158	55178	8.93%	0.514%
80	14.133	2125	2129	2133	rBV	60324	70071	11.34%	0.653%
81	14.219	2141	2143	2148	rVB2	11891	13777	2.23%	0.128%
82	14.273	2148	2152	2165	rVB	58436	99966	16.18%	0.932%
83	14.377	2165	2169	2174	rBV	39321	49500	8.01%	0.461%
84	14.432	2174	2178	2182	rVB	35075	40018	6.48%	0.373%
85	14.475	2182	2185	2191	rVB5	11823	19808	3.21%	0.185%
86	14.548	2191	2197	2205	rBV4	12461	33002	5.34%	0.308%
87	14.621	2205	2209	2213	rBV5	5480	9827	1.59%	0.092%
88	14.694	2213	2221	2226	rBV	111161	154025	24.93%	1.435%
89	14.786	2231	2236	2238	rBV3	9100	14257	2.31%	0.133%
90	14.810	2238	2240	2242	rVV	21679	20869	3.38%	0.194%
91	14.840	2242	2245	2249	rVB	44136	50391	8.15%	0.470%
92	15.249	2307	2312	2314	rBV	7588	11509	1.86%	0.107%
93	15.273	2314	2316	2320	rVB2	7048	8790	1.42%	0.082%
94	15.444	2339	2344	2348	rBV	13487	20679	3.35%	0.193%
95	15.554	2356	2362	2372	rVB3	32639	58677	9.50%	0.547%
96	15.688	2379	2384	2388	rBV2	29538	48377	7.83%	0.451%
97	15.724	2388	2390	2397	rVB	16397	21527	3.48%	0.201%
98	15.913	2413	2421	2426	rBV	6297	13287	2.15%	0.124%
99	16.072	2442	2447	2451	rBV2	4653	7573	1.23%	0.071%
100	16.450	2504	2509	2515	rBV6	3371	8078	1.31%	0.075%

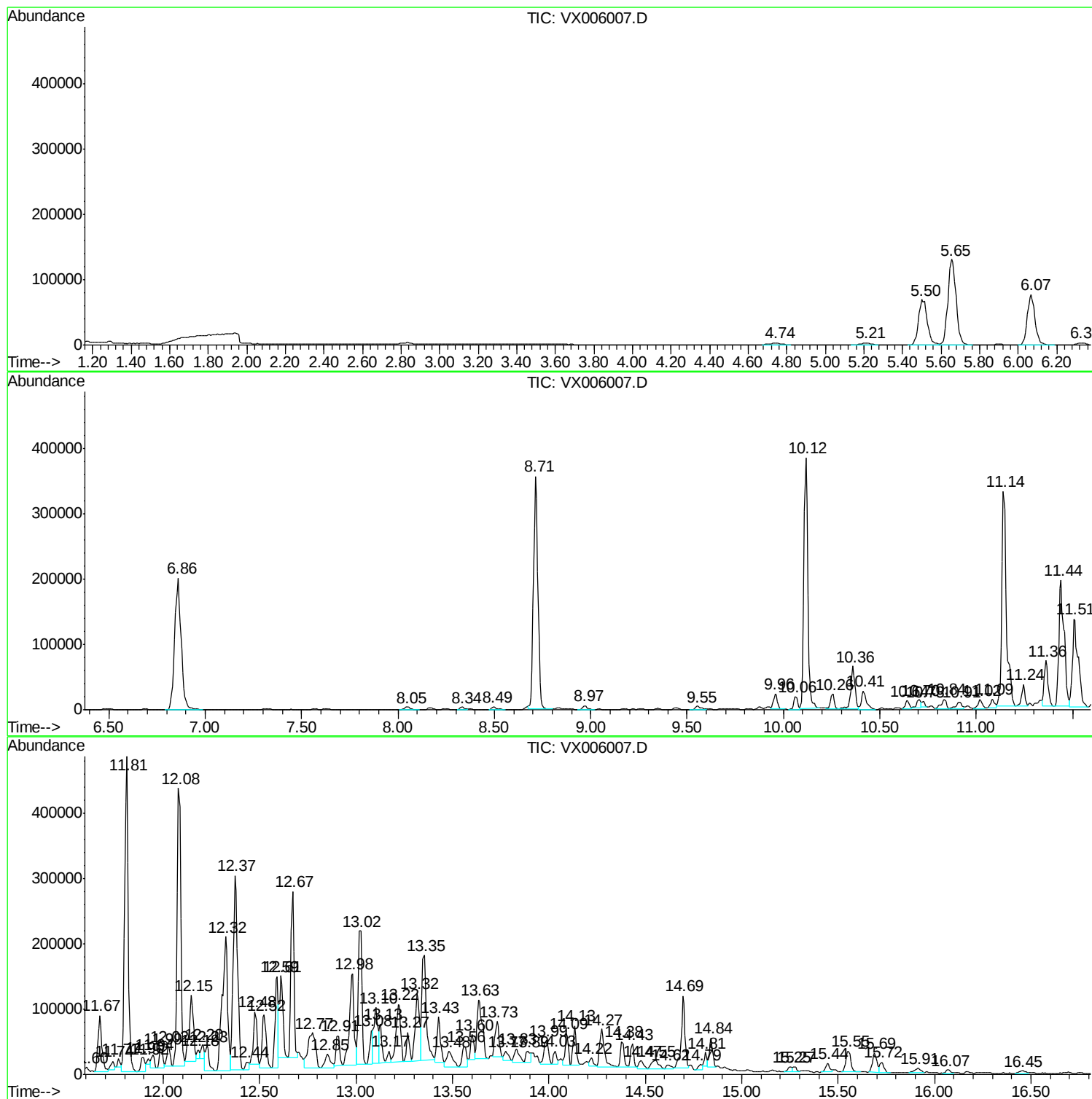
Sum of corrected areas: 10730568

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



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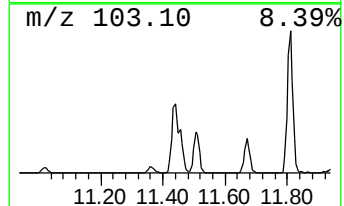
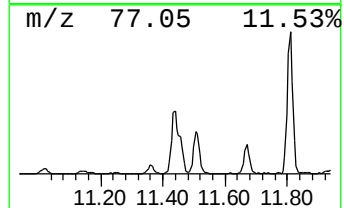
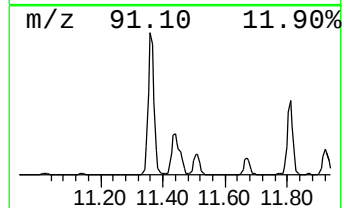
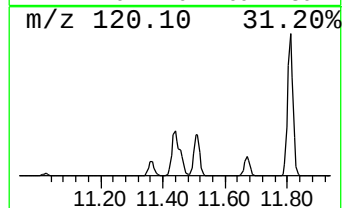
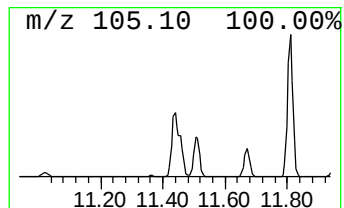
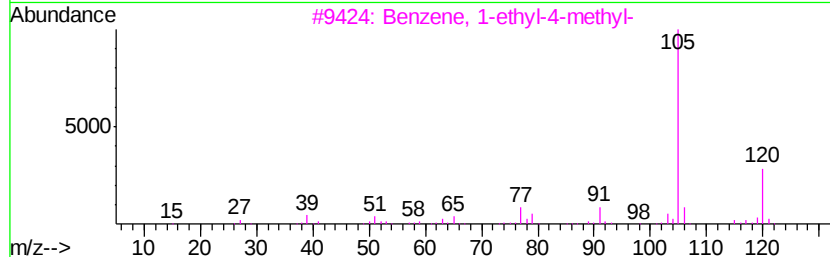
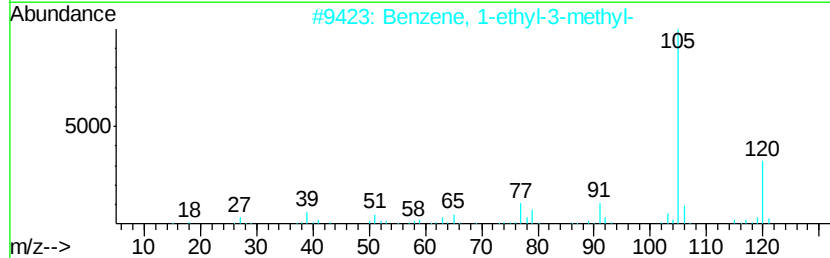
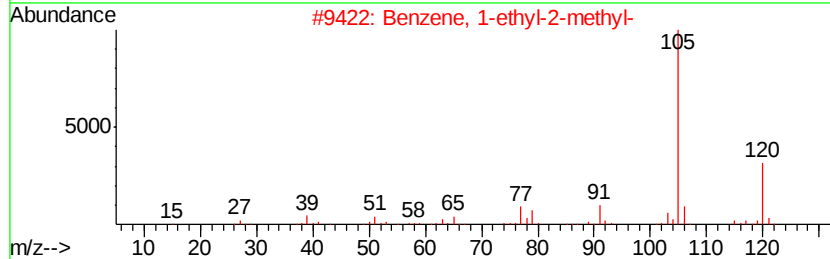
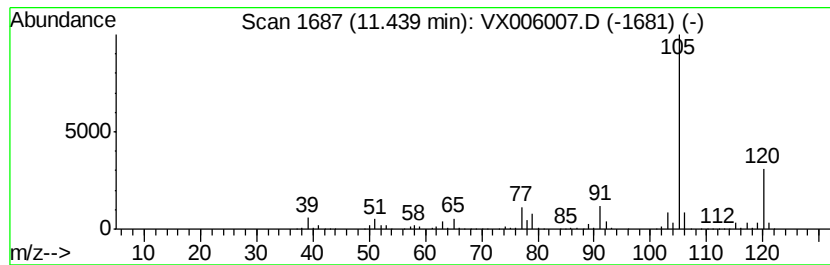
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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.
11.44	32.59 ug/l	371817	1,4-Dichlorobenzene-d4	12.08

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		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Mesitylene	120	C9H12	000108-67-8	91



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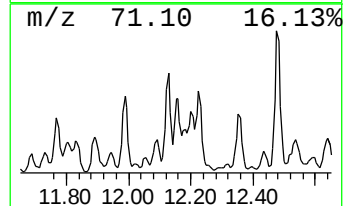
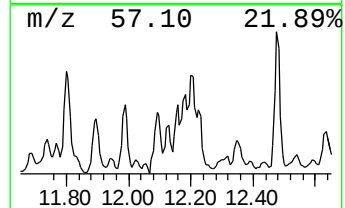
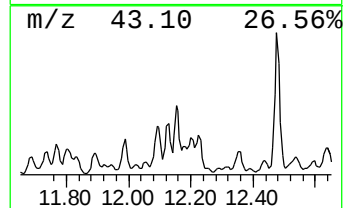
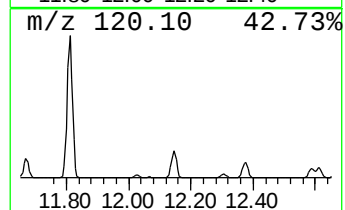
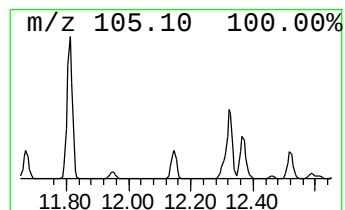
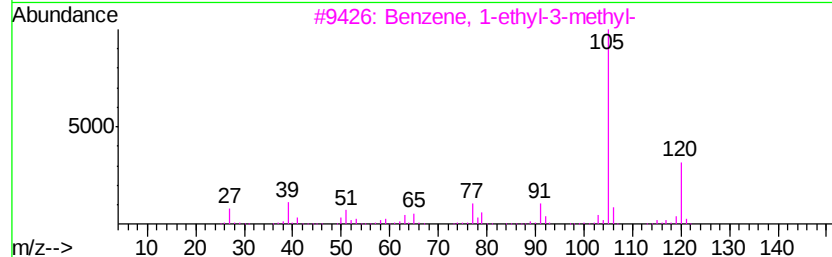
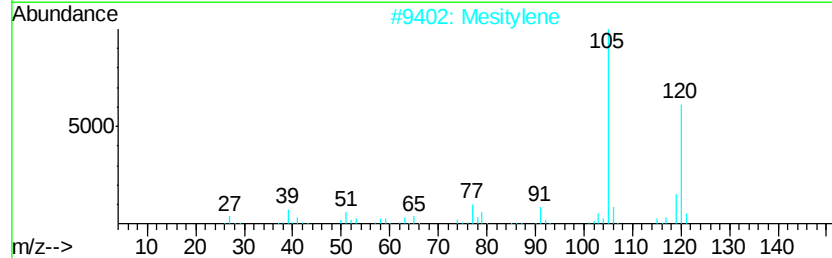
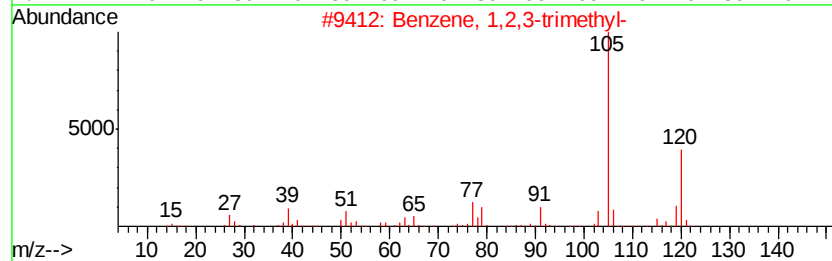
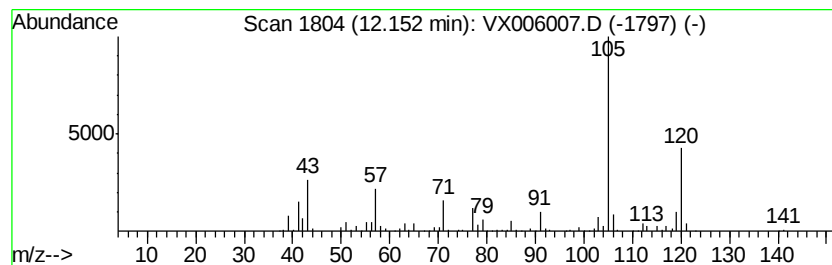
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TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1,2,3-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	14.33 ug/l	163498	1,4-Dichlorobenzene-d4	12.08

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-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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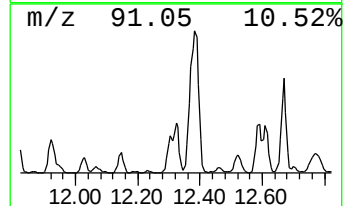
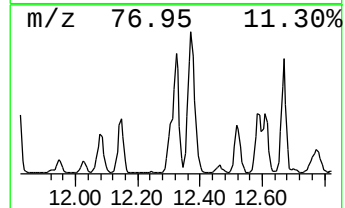
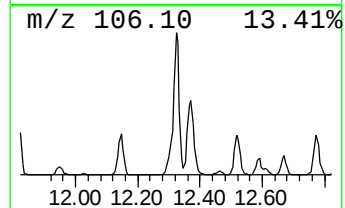
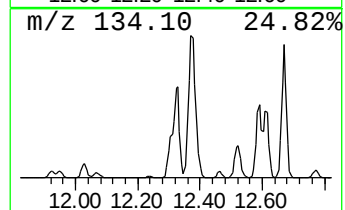
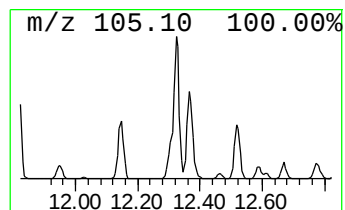
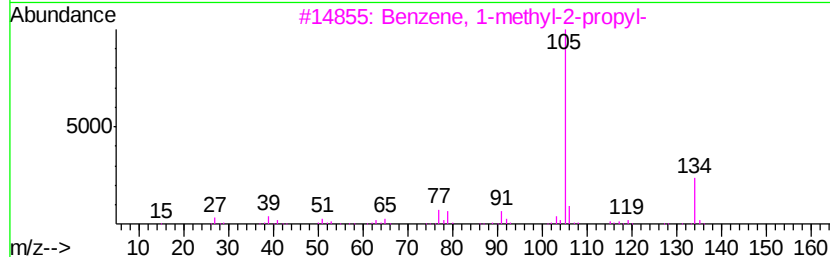
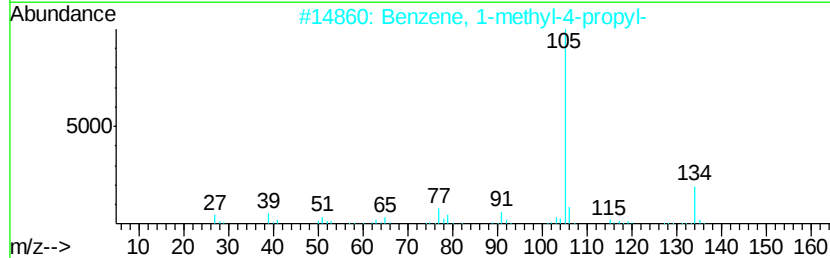
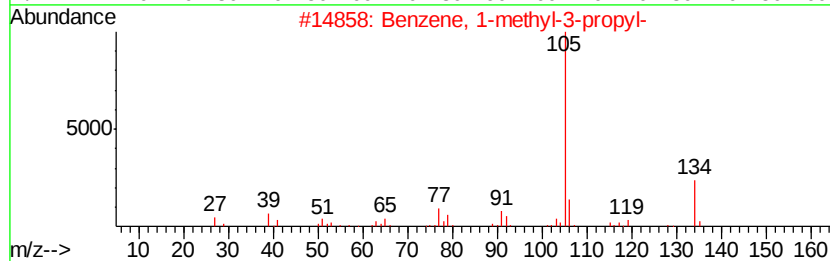
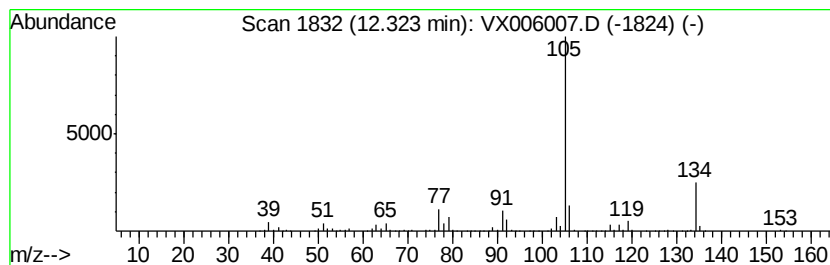
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-methyl-3-propyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.32	34.29 ug/l	391217	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	95
2		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
3		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91
4		Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	72
5		Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX006007.D
Acq On : 15 Nov 2018 06:46
Operator : JC/MD
Sample : J5915-02
Misc : 5.00g/10ml/100ul/5.00mL/MSVOA_X/MEOH
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T-2ME

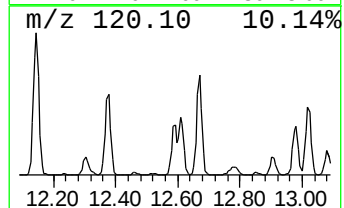
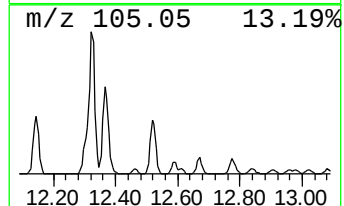
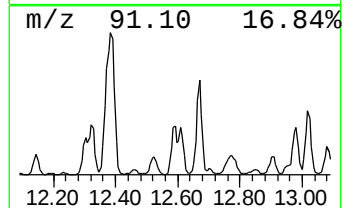
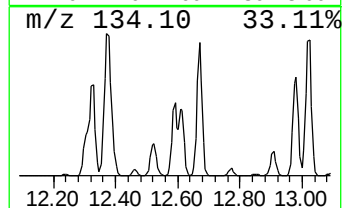
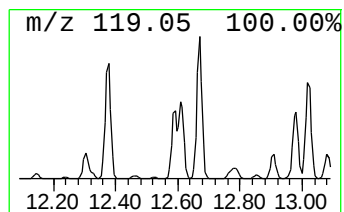
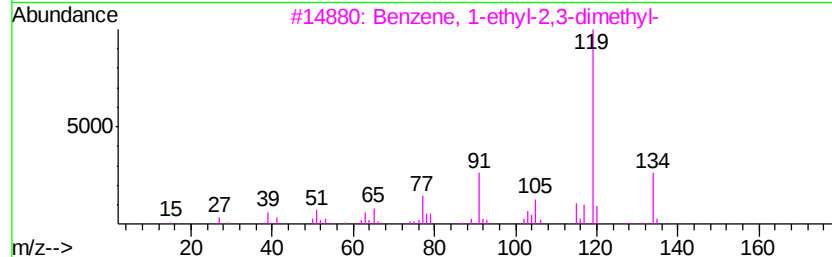
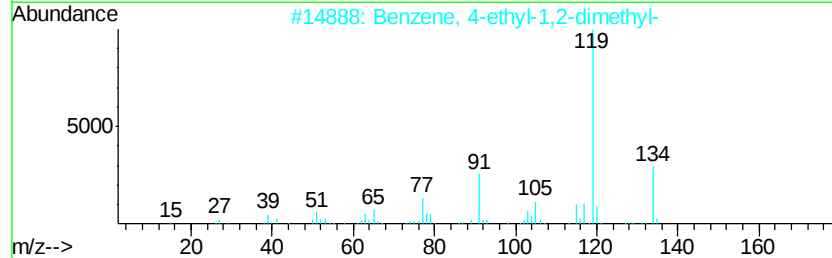
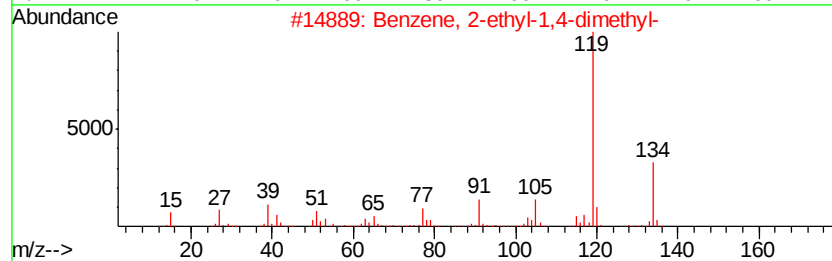
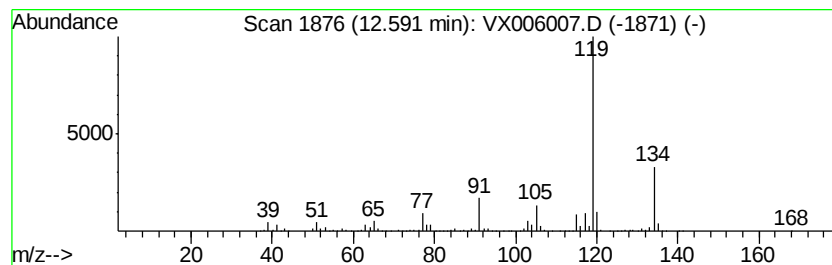
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.59	15.03 ug/l	171510	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4		o-Cymene	134	C10H14	000527-84-4	94
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX006007.D
Acq On : 15 Nov 2018 06:46
Operator : JC/MD
Sample : J5915-02
Misc : 5.00g/10ml/100ul/5.00mL/MSVOA_X/MEOH
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T-2ME

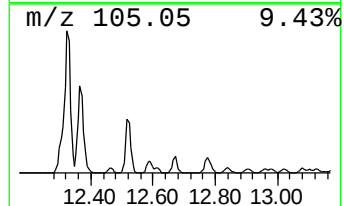
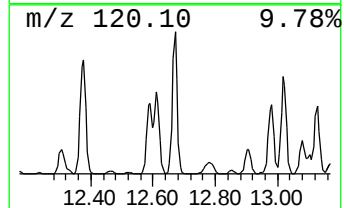
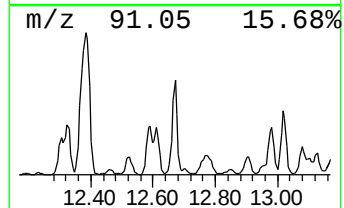
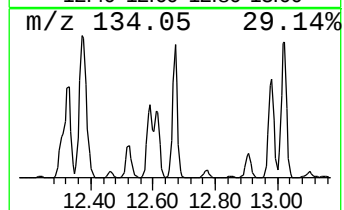
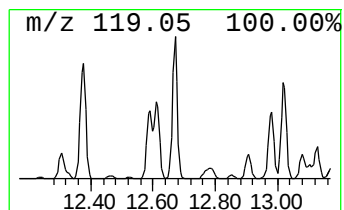
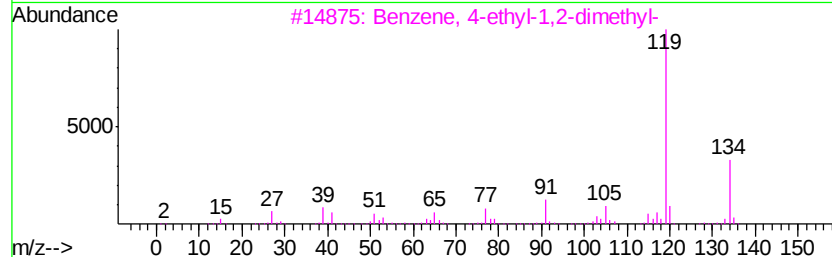
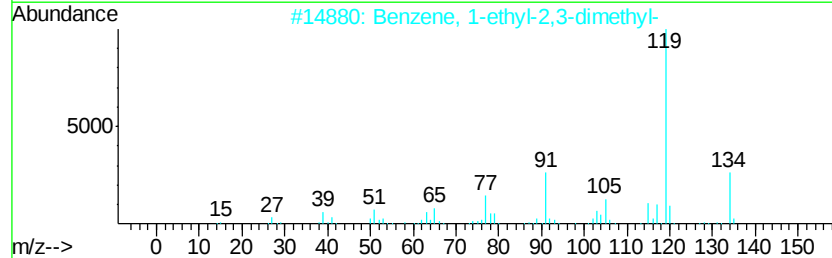
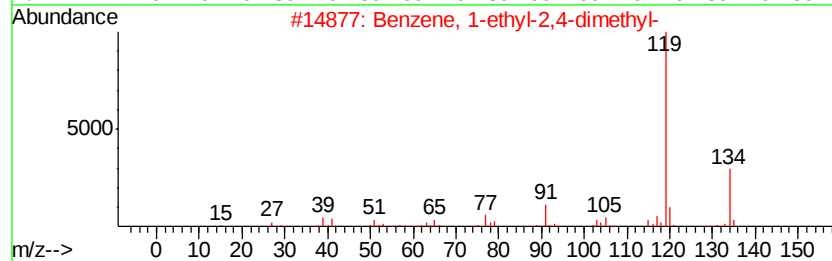
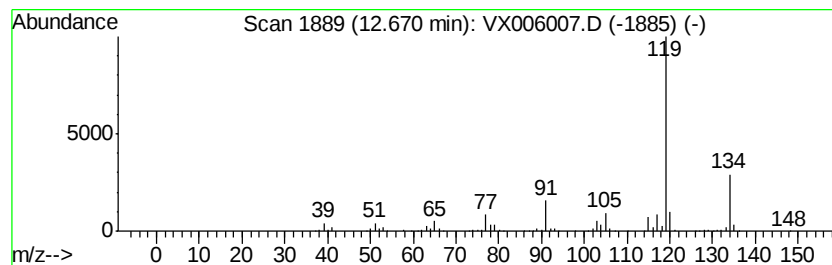
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	24.58 ug/l	280389	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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-1,4-dimethyl-	134	C10H14	001758-88-9	95
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX006007.D
Acq On : 15 Nov 2018 06:46
Operator : JC/MD
Sample : J5915-02
Misc : 5.00g/10ml/100ul/5.00mL/MSVOA_X/MEOH
ALS Vial : 44 Sample Multiplier: 1

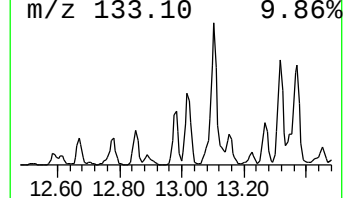
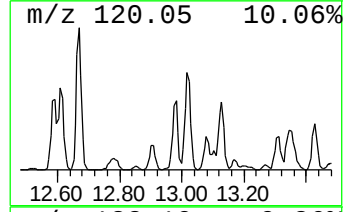
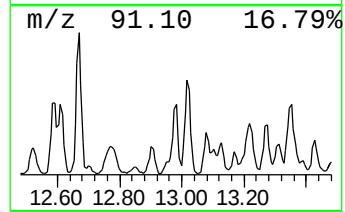
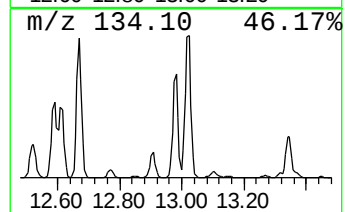
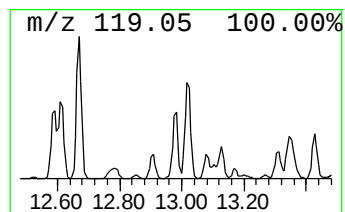
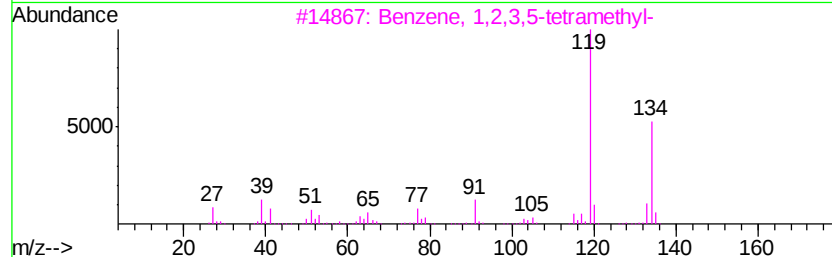
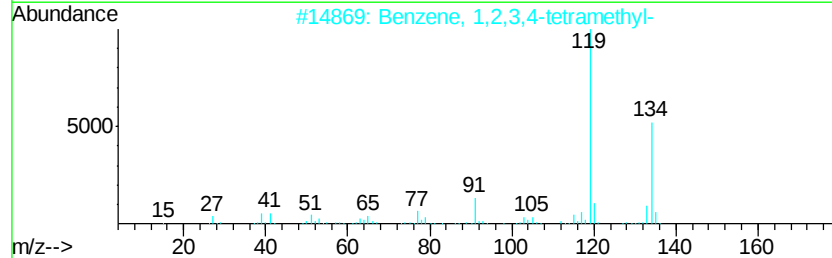
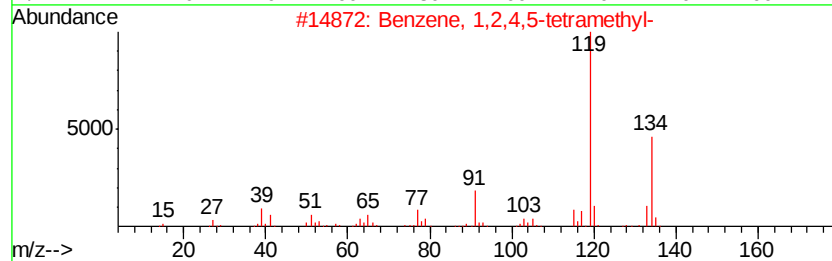
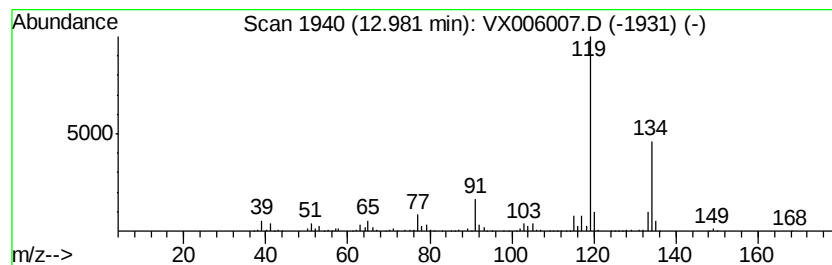
Instrument :
MSVOA_X
ClientSampleId :
T-2ME

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
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 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	20.58 ug/l	234753	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1,2,4,5-tetramethyl-	134 C10H14	000095-93-2	97
2	Benzene, 1,2,3,4-tetramethyl-	134 C10H14	000488-23-3	95
3	Benzene, 1,2,3,5-tetramethyl-	134 C10H14	000527-53-7	94
4	Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9	93
5	Benzene, 1-methyl-3-(1-methyleth...	134 C10H14	000535-77-3	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX006007.D
Acq On : 15 Nov 2018 06:46
Operator : JC/MD
Sample : J5915-02
Misc : 5.00g/10ml/100ul/5.00mL/MSVOA_X/MEOH
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T-2ME

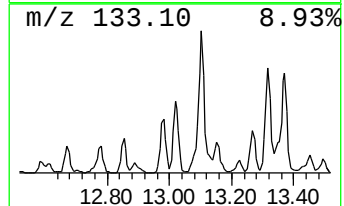
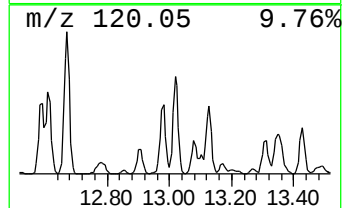
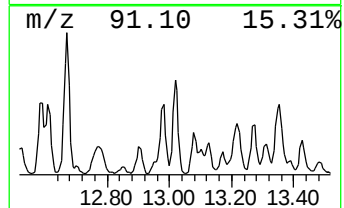
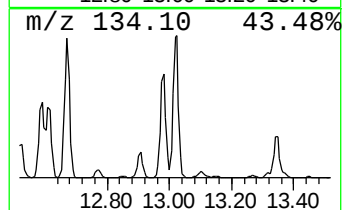
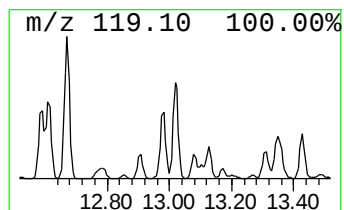
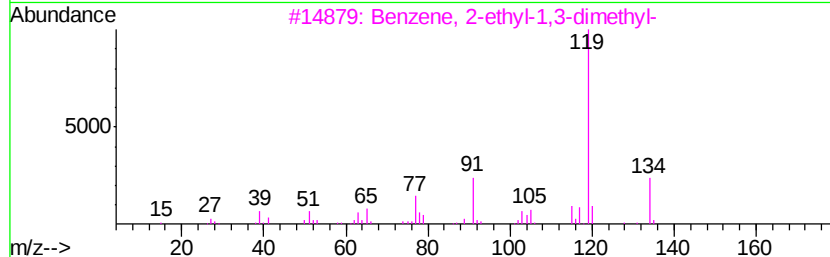
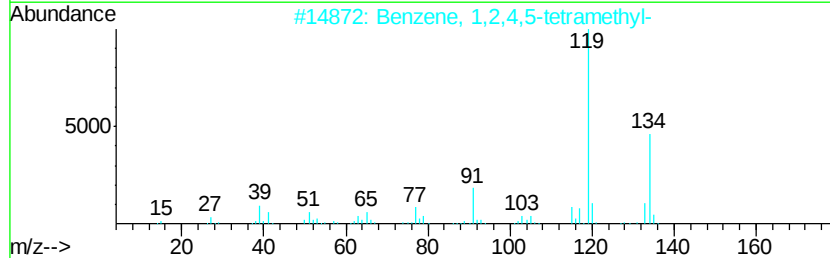
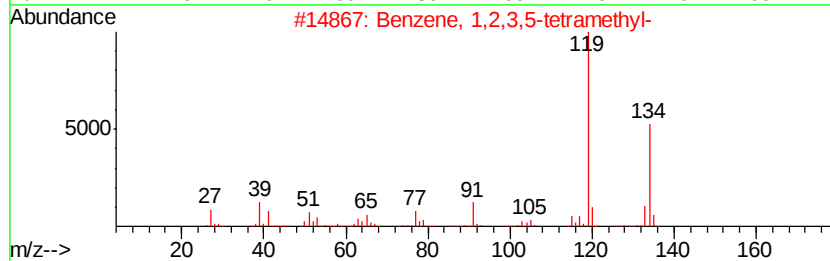
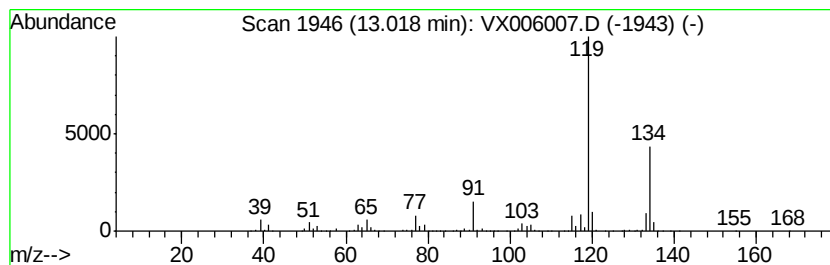
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	24.19 ug/l	275960	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	95
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
3		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX006007.D
Acq On : 15 Nov 2018 06:46
Operator : JC/MD
Sample : J5915-02
Misc : 5.00g/10ml/100ul/5.00mL/MSVOA_X/MEOH
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T-2ME

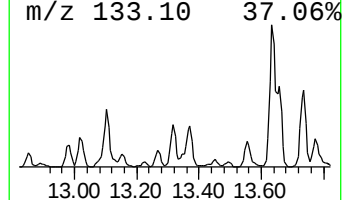
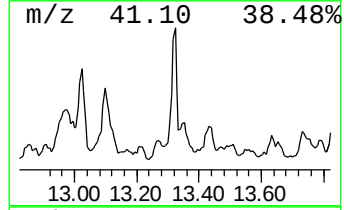
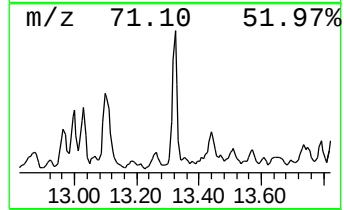
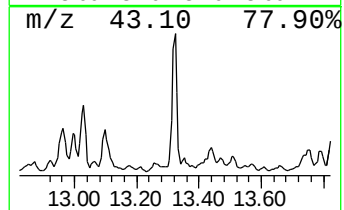
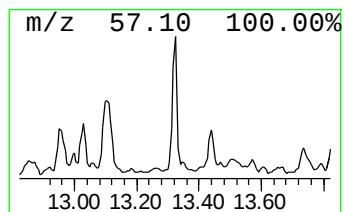
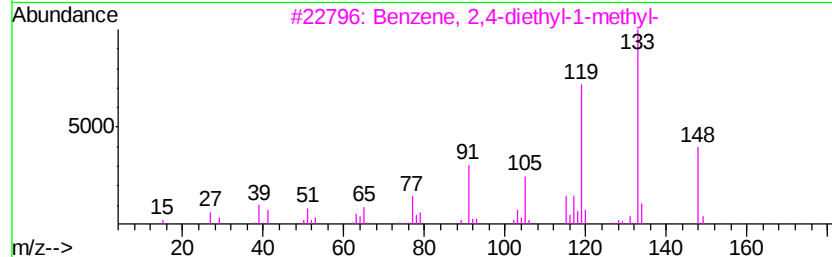
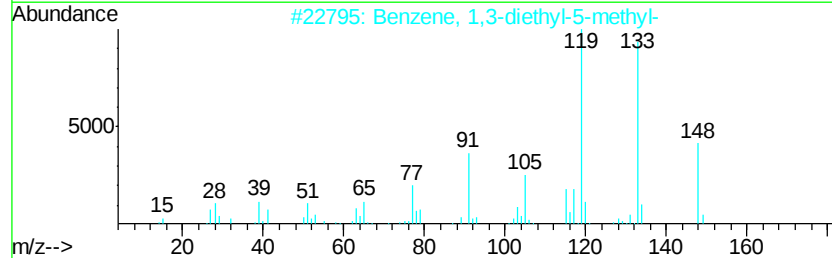
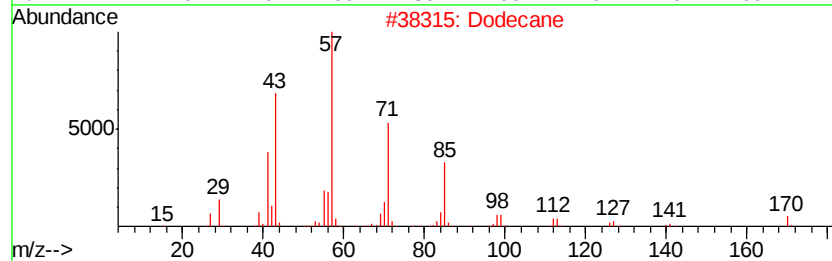
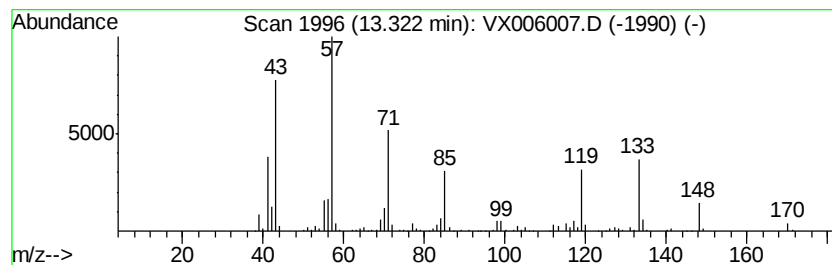
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Quant Title : SW846 8260

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

Peak Number 8 Dodecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.32	14.89 ug/l	169908	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane	170	C12H26	000112-40-3	91
2		Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	64
3		Benzene, 2,4-diethyl-1-methyl-	148	C11H16	001758-85-6	64
4		Disiloxane, 1,1,3,3-tetramethyl-	134	C4H14OSi2	003277-26-7	43
5		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	42



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX006007.D
Acq On : 15 Nov 2018 06:46
Operator : JC/MD
Sample : J5915-02
Misc : 5.00g/10ml/100ul/5.00mL/MSVOA_X/MEOH
ALS Vial : 44 Sample Multiplier: 1

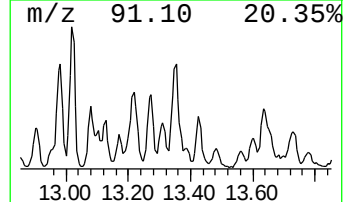
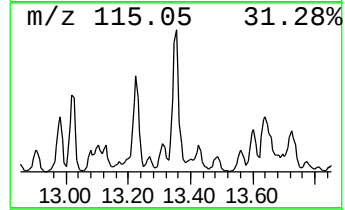
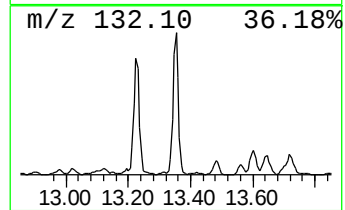
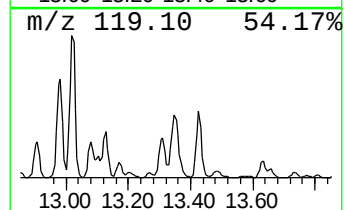
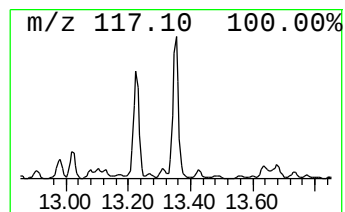
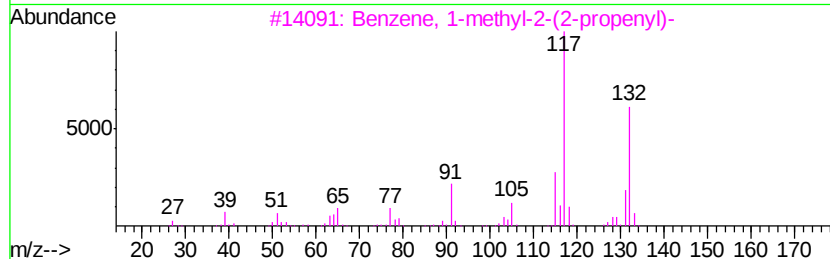
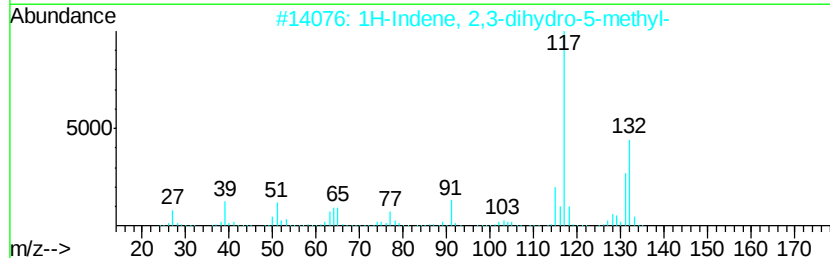
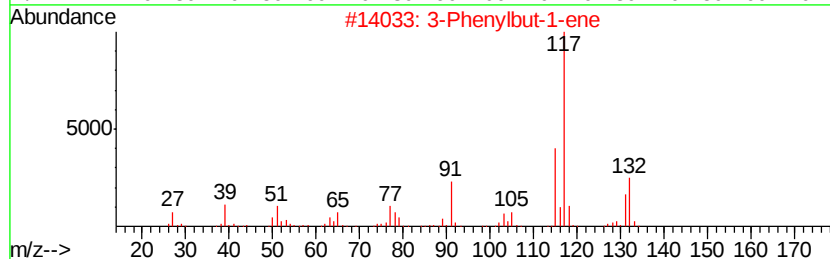
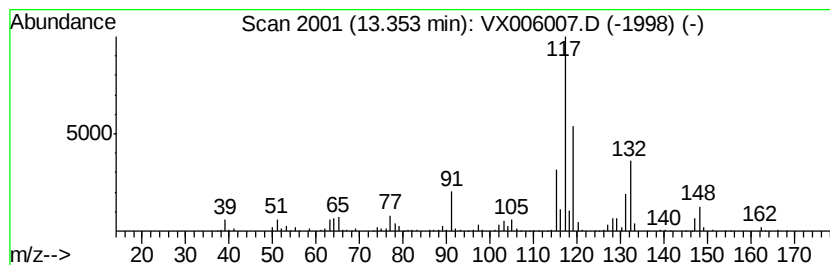
Instrument :
MSVOA_X
ClientSampleId :
T-2ME

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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

Peak Number 9 3-Phenylbut-1-ene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	22.33 ug/l	254720	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		3-Phenylbut-1-ene	132 C10H12	000934-10-1 76
2		1H-Indene, 2,3-dihydro-5-methyl-	132 C10H12	000874-35-1 70
3		Benzene, 1-methyl-2-(2-propenyl)-	132 C10H12	001587-04-8 70
4		1H-Indene, 2,3-dihydro-4-methyl-	132 C10H12	000824-22-6 70
5		Benzene, (1-methyl-1-propenyl)-,...	132 C10H12	000768-00-3 68



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX006007.D
Acq On : 15 Nov 2018 06:46
Operator : JC/MD
Sample : J5915-02
Misc : 5.00g/10ml/100ul/5.00mL/MSVOA_X/MEOH
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T-2ME

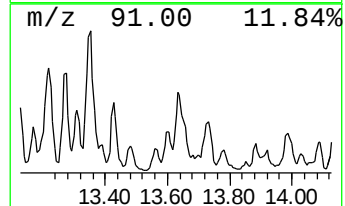
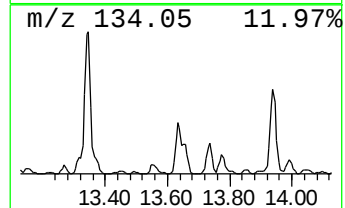
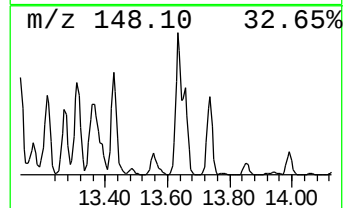
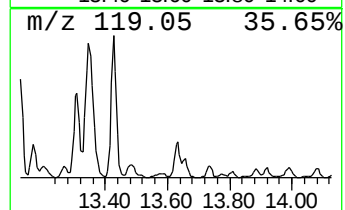
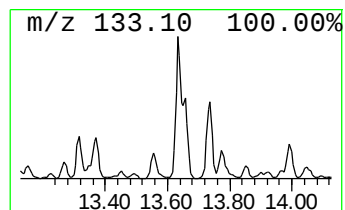
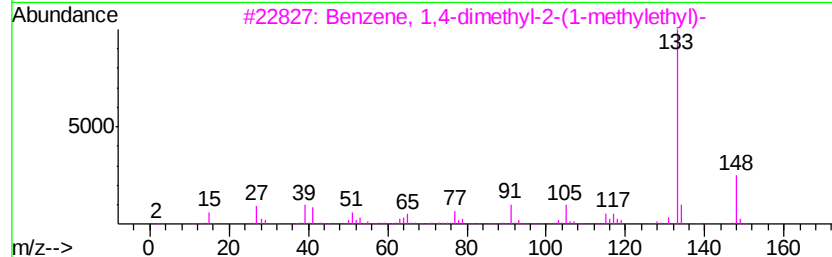
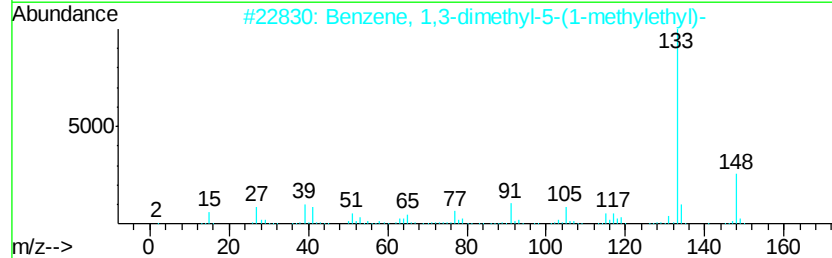
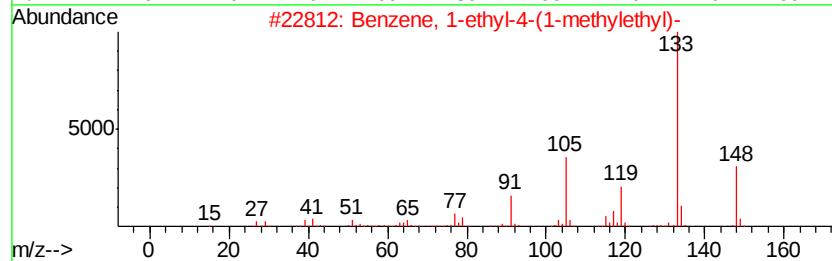
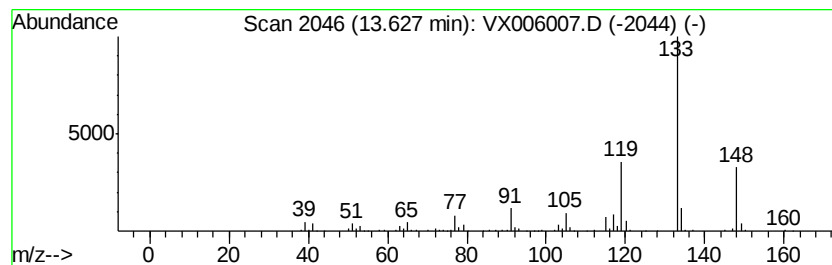
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1-ethyl-4-(1-methy... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.63	14.50 ug/l	165449	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	90
2		Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	87
3		Benzene, 1,4-dimethyl-2-(1-methy...	148	C11H16	004132-72-3	87
4		Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	87
5		Benzene, 2,4-dimethyl-1-(1-methy...	148	C11H16	004706-89-2	87



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX111418\
Data File : VX006007.D
Acq On : 15 Nov 2018 06:46
Operator : JC/MD
Sample : J5915-02
Misc : 5.00g/10ml/100ul/5.00mL/MSVOA_X/MEOH
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T-2ME

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.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-...	11.44	32.6	ug/l	371817	4	12.08	570477	50.0
Benzene, 1,2,3-tr...	12.15	14.3	ug/l	163498	4	12.08	570477	50.0
Benzene, 1-methyl...	12.32	34.3	ug/l	391217	4	12.08	570477	50.0
Benzene, 2-ethyl-...	12.59	15.0	ug/l	171510	4	12.08	570477	50.0
Benzene, 1-ethyl-...	12.67	24.6	ug/l	280389	4	12.08	570477	50.0
Benzene, 1,2,4,5-...	12.98	20.6	ug/l	234753	4	12.08	570477	50.0
Benzene, 1,2,3,5-...	13.02	24.2	ug/l	275960	4	12.08	570477	50.0
Dodecane	13.32	14.9	ug/l	169908	4	12.08	570477	50.0
3-Phenylbut-1-ene	13.35	22.3	ug/l	254720	4	12.08	570477	50.0
Benzene, 1-ethyl-...	13.63	14.5	ug/l	165449	4	12.08	570477	50.0