

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX071819\
 Data File : VX010953.D
 Acq On : 18 Jul 2019 13:02
 Operator : JC/SP
 Sample : K3627-03
 Misc : 3.54g/5mL/100uL/5mL/MSVOA_X/MEOH
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TR-04-071519

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.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	1.648	66	81	82	rBV8	10577	31671	3.18%	0.282%
2	2.861	277	280	288	rVB2	5983	10431	1.05%	0.093%
3	3.019	300	306	313	rBV	7513	19362	1.95%	0.172%
4	5.501	702	713	726	rBV	102672	314432	31.59%	2.801%
5	5.653	727	738	757	rVB	201601	576937	57.97%	5.139%
6	6.061	794	805	821	rBV	110513	313393	31.49%	2.791%
7	6.854	926	935	952	rBV	292566	719966	72.34%	6.413%
8	8.713	1232	1240	1254	rBV	626130	995241	100.00%	8.865%
9	9.603	1380	1386	1392	rVB5	5310	11594	1.16%	0.103%
10	10.116	1464	1470	1482	rVB	568349	845931	85.00%	7.535%
11	10.359	1501	1510	1515	rBV6	5267	11088	1.11%	0.099%
12	10.414	1515	1519	1524	rVB	9555	13121	1.32%	0.117%
13	10.829	1580	1587	1591	rBV6	6568	11818	1.19%	0.105%
14	10.914	1596	1601	1604	rBV5	6222	10026	1.01%	0.089%
15	11.140	1632	1638	1648	rBV	521710	704675	70.80%	6.277%
16	11.243	1651	1655	1657	rVV2	8526	11927	1.20%	0.106%
17	11.438	1682	1687	1691	rBV	18267	34465	3.46%	0.307%
18	11.481	1691	1694	1696	rVV3	12366	18749	1.88%	0.167%
19	11.530	1696	1702	1708	rVB2	48328	78290	7.87%	0.697%
20	11.670	1714	1725	1732	rBV4	20006	46641	4.69%	0.415%
21	11.768	1738	1741	1744	rBV	11267	13884	1.40%	0.124%
22	11.810	1744	1748	1757	rVB	45102	71303	7.16%	0.635%
23	11.945	1764	1770	1775	rVB6	13273	20789	2.09%	0.185%
24	11.999	1775	1779	1781	rBV3	20473	26561	2.67%	0.237%
25	12.024	1781	1783	1786	rVB2	13346	14280	1.43%	0.127%
26	12.079	1786	1792	1797	rBV	721361	887643	89.19%	7.906%
27	12.146	1797	1803	1808	rVV3	29047	51333	5.16%	0.457%
28	12.219	1812	1815	1823	rVB4	16186	32343	3.25%	0.288%
29	12.304	1823	1829	1830	rBV	27175	39328	3.95%	0.350%
30	12.323	1830	1832	1836	rVB	29433	30946	3.11%	0.276%
31	12.371	1836	1840	1846	rVB3	64951	114011	11.46%	1.015%
32	12.475	1848	1857	1861	rBV	101927	138928	13.96%	1.237%
33	12.518	1861	1864	1870	rVB2	35738	50783	5.10%	0.452%
34	12.585	1871	1875	1877	rBV	28251	38841	3.90%	0.346%

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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
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35	12.609	1877	1879	1884	rVV	32823	42478	4.27%	0.378%
36	12.664	1885	1888	1892	rVB	29422	35456	3.56%	0.316%
37	12.780	1897	1907	1911	rBV4	34952	99567	10.00%	0.887%
38	12.871	1919	1922	1926	rVB2	31813	33990	3.42%	0.303%
39	12.938	1929	1933	1937	rBV5	22163	37516	3.77%	0.334%
40	12.975	1937	1939	1943	rVV2	21548	21203	2.13%	0.189%
41	13.018	1943	1946	1953	rVB2	53624	102781	10.33%	0.915%
42	13.097	1953	1959	1962	rBV3	42495	74635	7.50%	0.665%
43	13.127	1962	1964	1966	rVB	13365	11027	1.11%	0.098%
44	13.219	1972	1979	1984	rBV5	50989	104995	10.55%	0.935%
45	13.267	1984	1987	1990	rVV	36815	40647	4.08%	0.362%
46	13.316	1990	1995	1997	rVV3	103059	136648	13.73%	1.217%
47	13.347	1997	2000	2005	rVV3	145472	205760	20.67%	1.833%
48	13.389	2005	2007	2010	rVV	13989	14689	1.48%	0.131%
49	13.426	2010	2013	2017	rVV2	16720	21762	2.19%	0.194%
50	13.481	2017	2022	2030	rVB4	99282	193759	19.47%	1.726%
51	13.560	2030	2035	2038	rBV2	36937	58090	5.84%	0.517%
52	13.597	2038	2041	2044	rBV	31478	36984	3.72%	0.329%
53	13.639	2044	2048	2052	rBV3	52940	93645	9.41%	0.834%
54	13.725	2058	2062	2067	rBV4	38360	59877	6.02%	0.533%
55	13.853	2080	2083	2086	rBV4	28585	36068	3.62%	0.321%
56	13.908	2086	2092	2097	rVB4	90345	176213	17.71%	1.570%
57	13.981	2097	2104	2109	rBV4	91408	190307	19.12%	1.695%
58	14.030	2109	2112	2114	rVV	22690	26440	2.66%	0.236%
59	14.060	2114	2117	2119	rVV2	32214	45175	4.54%	0.402%
60	14.090	2119	2122	2125	rVV3	100069	116705	11.73%	1.039%
61	14.133	2125	2129	2132	rVV2	71144	95577	9.60%	0.851%
62	14.164	2132	2134	2137	rVV	21631	22085	2.22%	0.197%
63	14.219	2140	2143	2147	rVV	28362	30721	3.09%	0.274%
64	14.286	2147	2154	2160	rVB2	209585	359192	36.09%	3.199%
65	14.377	2165	2169	2174	rBV4	52301	75903	7.63%	0.676%
66	14.432	2174	2178	2181	rBV	74434	94278	9.47%	0.840%
67	14.481	2182	2186	2191	rVB7	43774	81650	8.20%	0.727%
68	14.536	2191	2195	2204	rVB2	141635	239220	24.04%	2.131%
69	14.609	2204	2207	2209	rBV3	29798	43267	4.35%	0.385%
70	14.670	2214	2217	2224	rVV2	212941	345795	34.74%	3.080%
71	14.731	2224	2227	2231	rVB2	107081	137209	13.79%	1.222%

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72	14.785	2231	2236	2237	rBV2	52781	67975	6.83%	0.605%
73	14.810	2237	2240	2243	rVV	102229	111638	11.22%	0.994%
74	14.840	2243	2245	2248	rVV3	34521	49329	4.96%	0.439%
75	14.877	2248	2251	2253	rVV2	56560	74195	7.45%	0.661%
76	14.901	2253	2255	2261	rVV4	69849	95308	9.58%	0.849%
77	14.962	2261	2265	2276	rVB3	52451	114382	11.49%	1.019%
78	15.109	2282	2289	2293	rBV4	42030	92193	9.26%	0.821%
79	15.243	2307	2311	2314	rBV	110675	165543	16.63%	1.474%
80	15.310	2319	2322	2324	rVV3	17900	23161	2.33%	0.206%
81	15.340	2324	2327	2331	rVB4	37714	49251	4.95%	0.439%
82	15.389	2331	2335	2339	rVB5	22470	38355	3.85%	0.342%
83	15.444	2339	2344	2348	rBV3	50723	87360	8.78%	0.778%
84	15.554	2357	2362	2367	rBV2	101600	149736	15.05%	1.334%
85	15.608	2367	2371	2373	rVV4	18844	27080	2.72%	0.241%
86	15.639	2374	2376	2379	rVV2	22395	24035	2.41%	0.214%
87	15.682	2379	2383	2387	rVB3	53426	76098	7.65%	0.678%
88	15.919	2418	2422	2425	rBV3	18333	31471	3.16%	0.280%
89	16.029	2436	2440	2444	rBV6	16978	31305	3.15%	0.279%
90	16.169	2458	2463	2464	rBV4	12255	19310	1.94%	0.172%
91	16.456	2505	2510	2516	rVB3	25552	47389	4.76%	0.422%

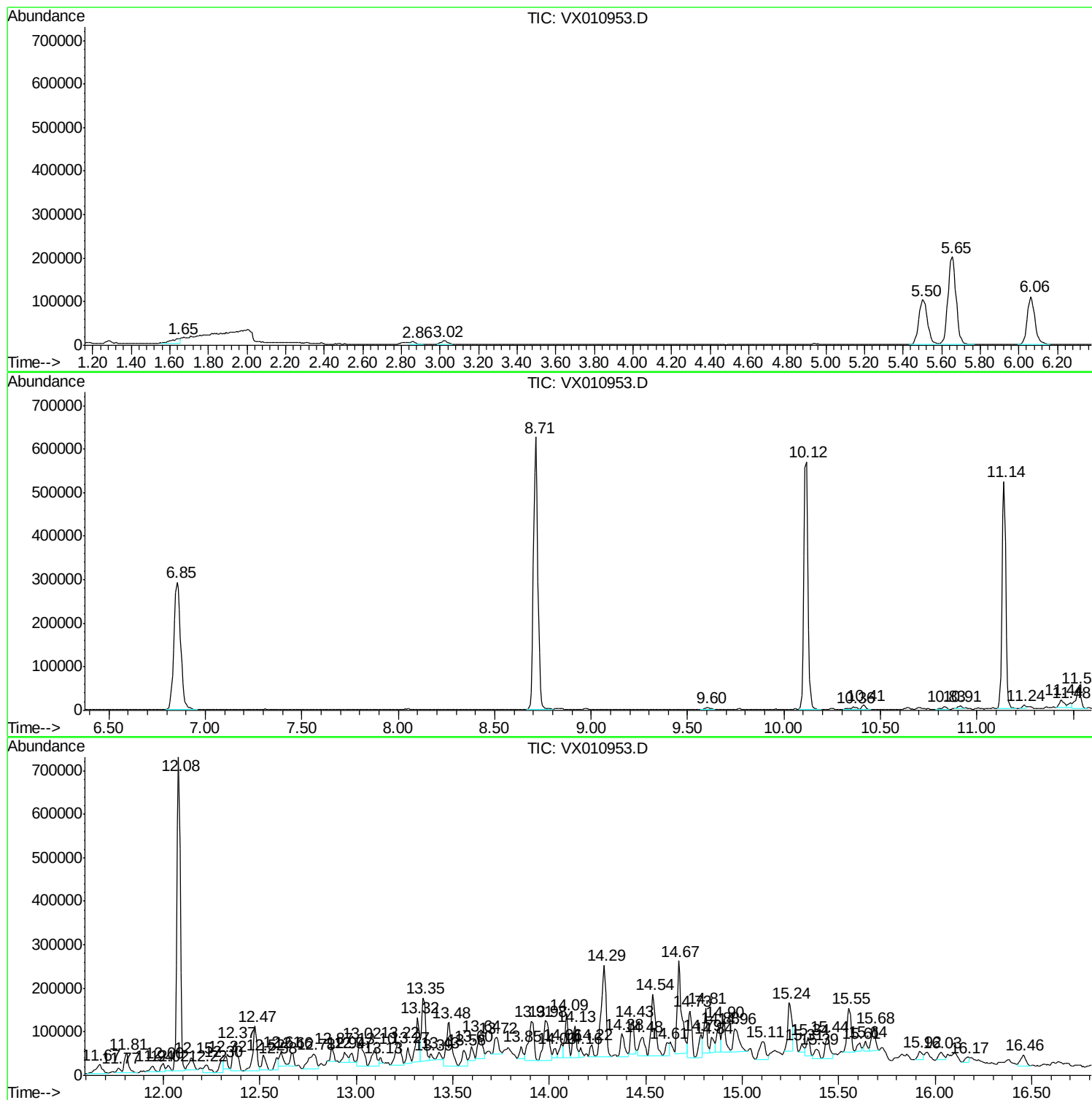
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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



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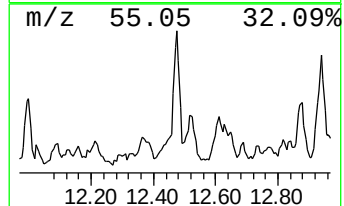
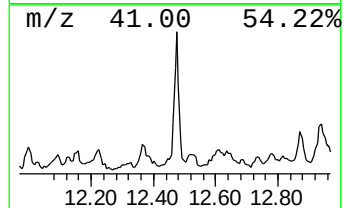
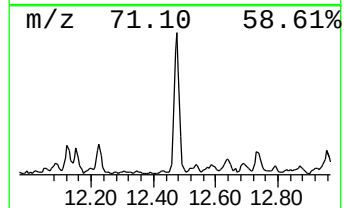
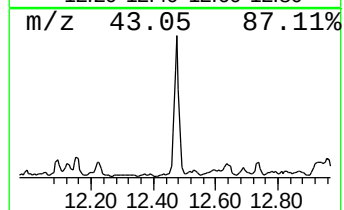
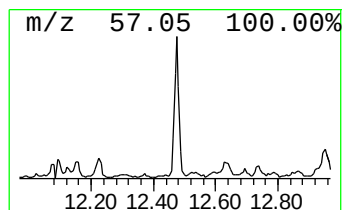
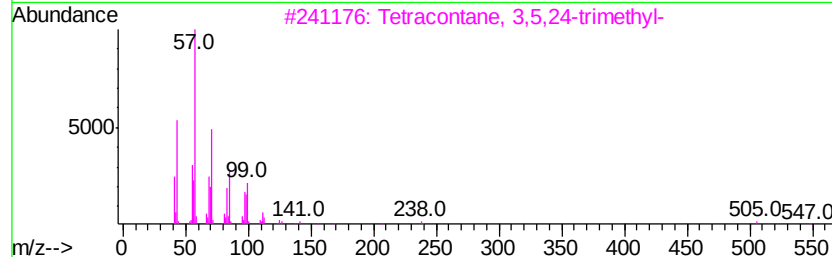
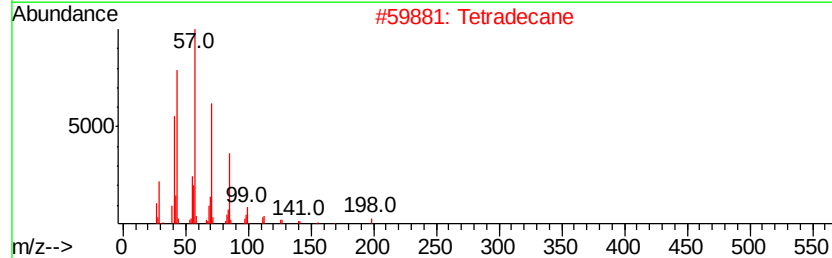
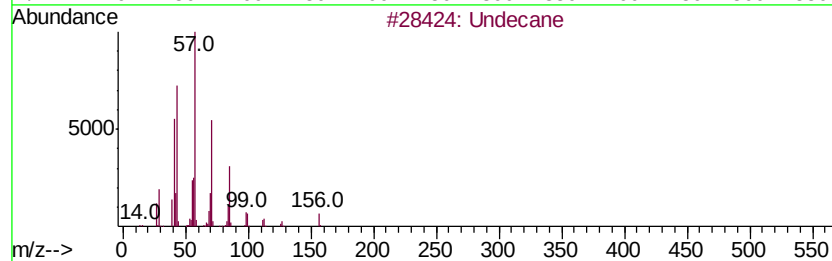
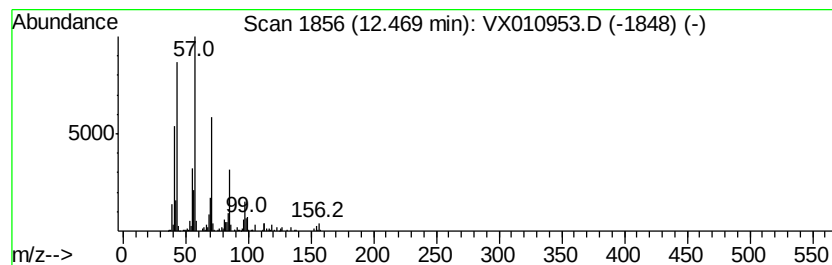
Instrument :
MSVOA_X
ClientSampleId :
TR-04-071519

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

Peak Number 1 Undecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.47	7.83 ug/l	138928	1,4-Dichlorobenzene-d4		12.08	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane		156	C11H24	001120-21-4	96
2	Tetradecane		198	C14H30	000629-59-4	72
3	Tetracontane, 3,5,24-trimethyl-		605	C43H88	055162-61-3	72
4	Hexadecane		226	C16H34	000544-76-3	64
5	Tridecane		184	C13H28	000629-50-5	59



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ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
TR-04-071519

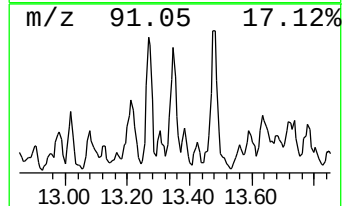
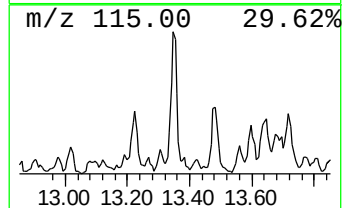
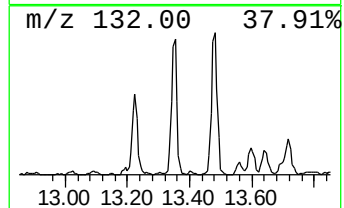
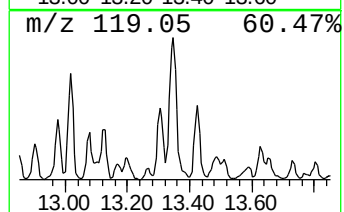
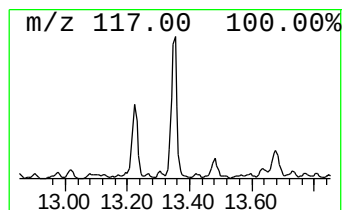
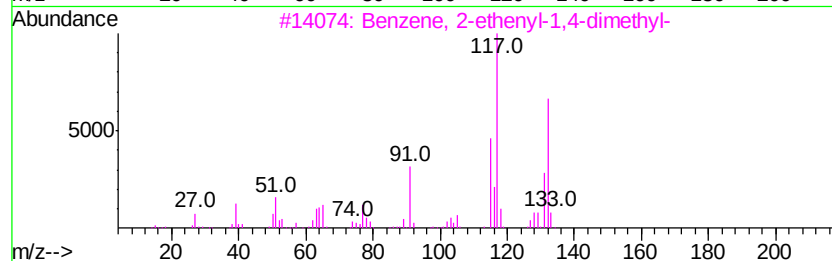
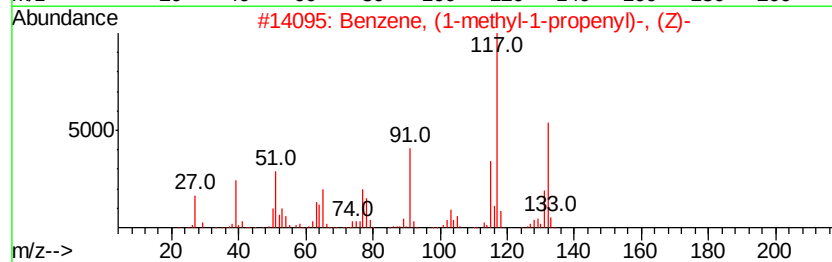
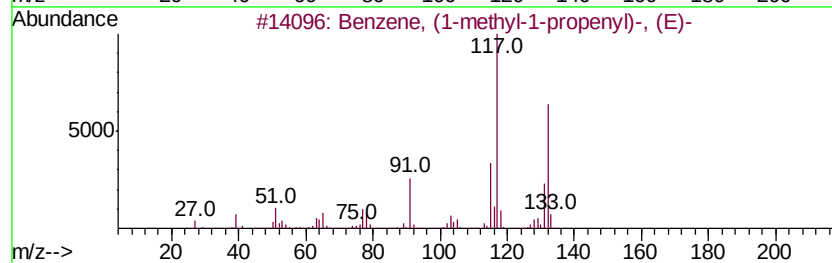
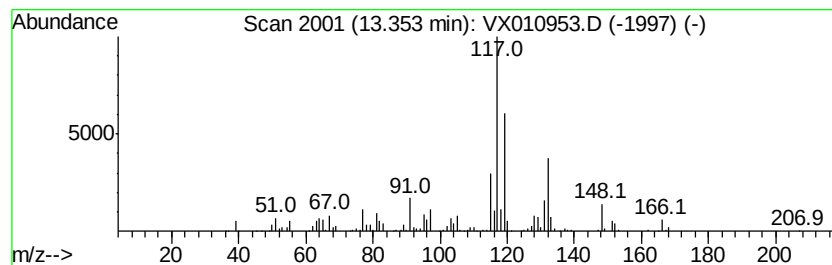
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, (1-methyl-1-propen... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	11.59 ug/l	205760	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90
2		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	89
3		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	74
4		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	64
5		2,4-Dimethylstyrene	132	C10H12	002234-20-0	59



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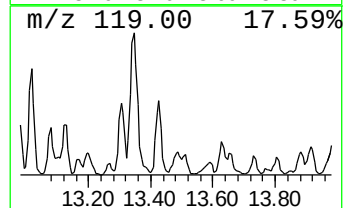
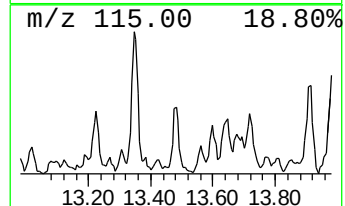
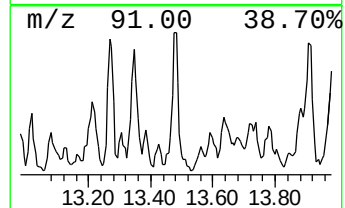
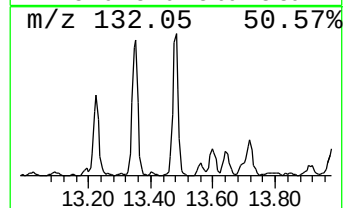
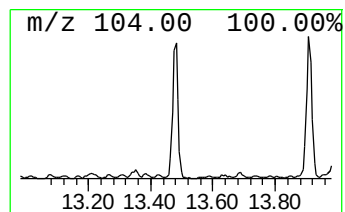
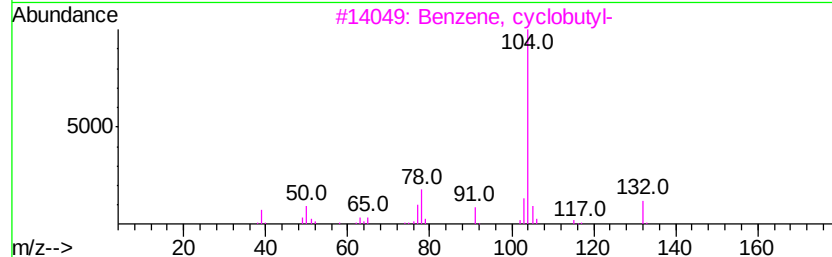
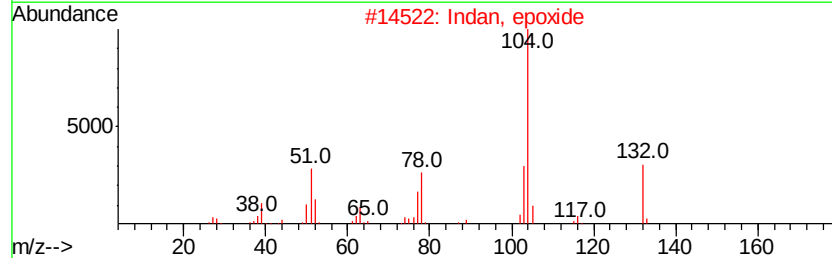
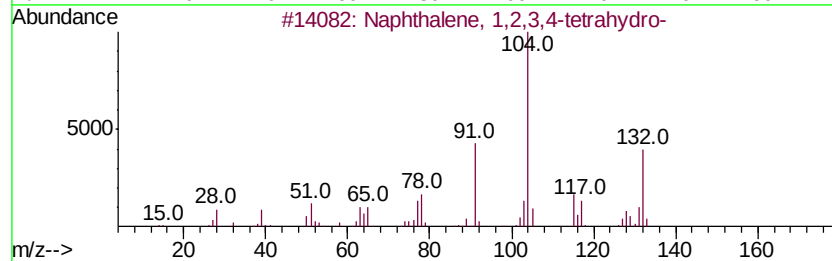
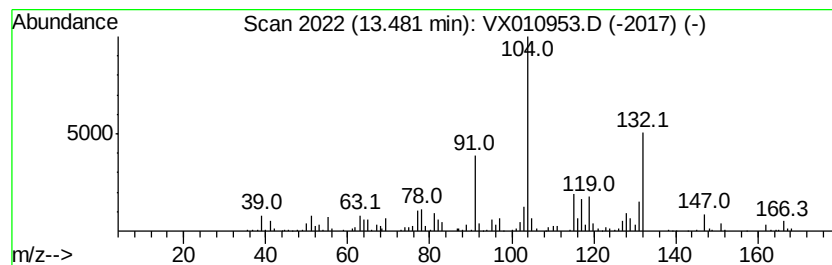
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.48	10.91 ug/l	193759	1,4-Dichlorobenzene-d4	12.08	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	90
2	Indan, epoxide	132	C9H8O	000768-22-9	58
3	Benzene, cyclobutyl-	132	C10H12	004392-30-7	53
4	2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	50
5	Naphthalene, 2-ethyl-1,2,3,4-tet...	160	C12H16	032367-54-7	47



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Misc : 3.54µL/100µL/5mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

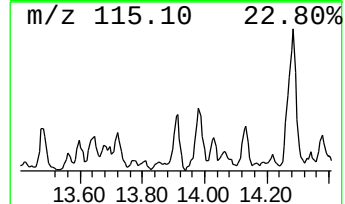
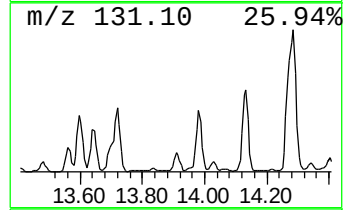
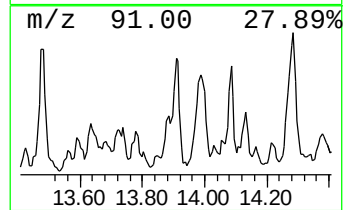
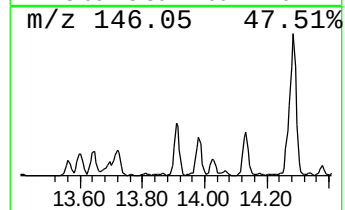
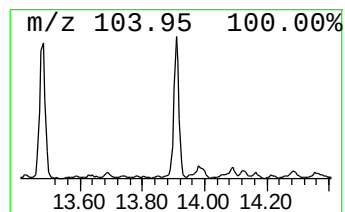
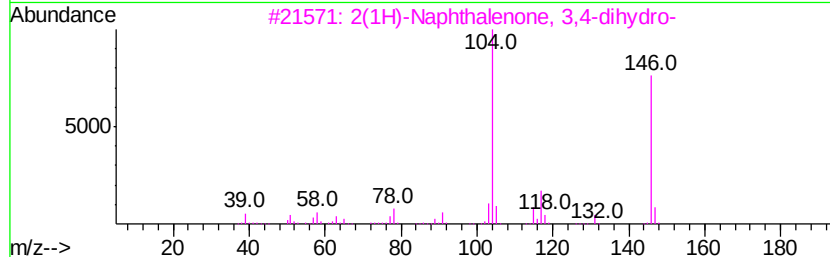
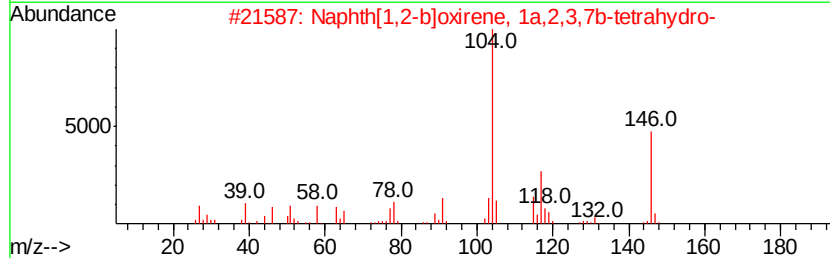
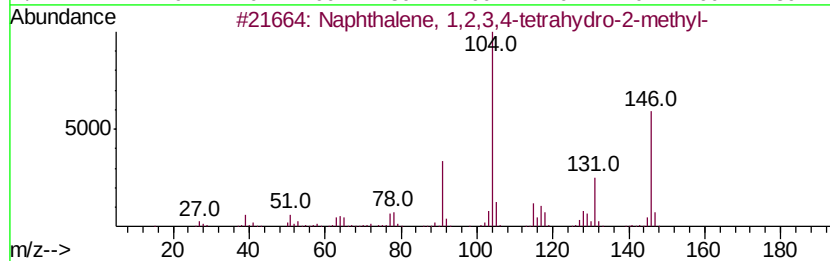
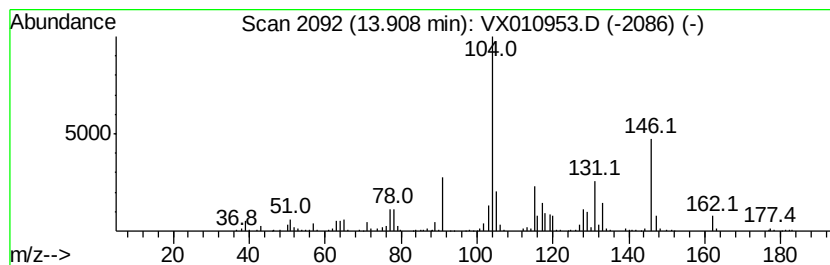
Instrument :
MSVOA_X
ClientSampleID :
TR-04-071519

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.91	9.93 ug/l	176213	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	003877-19-8 81
2		Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146 C10H10O	002461-34-9 64
3		2(1H)-Naphthalenone, 3,4-dihydro-	146 C10H10O	000530-93-8 50
4		Naphthalene, 2-ethyl-1,2,3,4-tet...	160 C12H16	032367-54-7 47
5		2-Methylbenzyl cyanide	131 C9H9N	022364-68-7 41



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX071819\
Data File : VX010953.D
Acq On : 18 Jul 2019 13:02
Operator : JC/SP
Sample : K3627-03
Misc : 3.54g/5mL/100uL/5mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

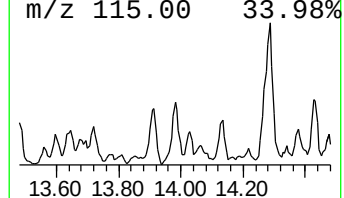
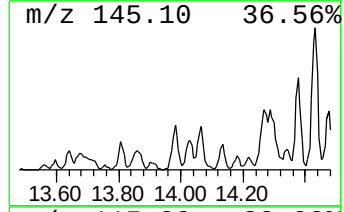
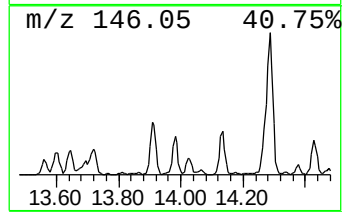
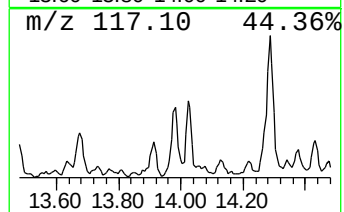
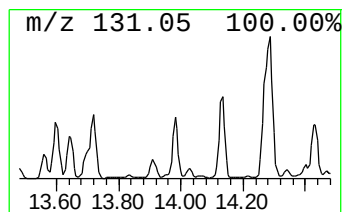
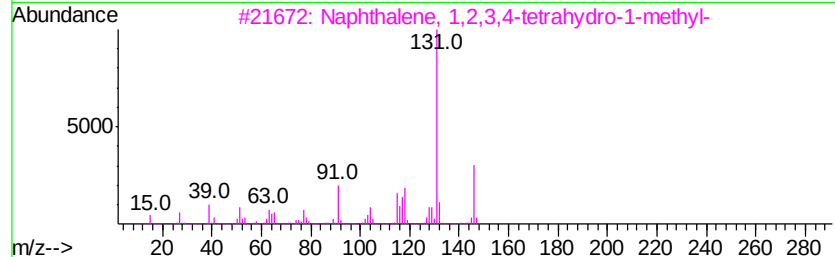
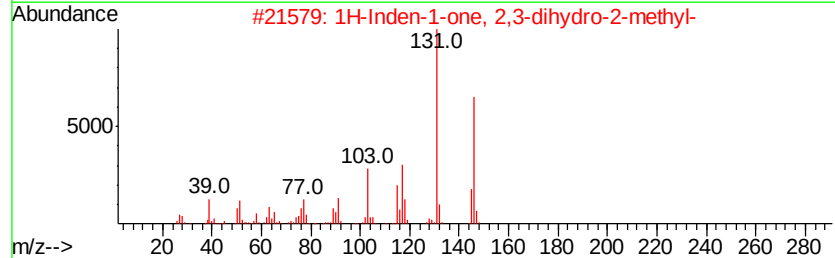
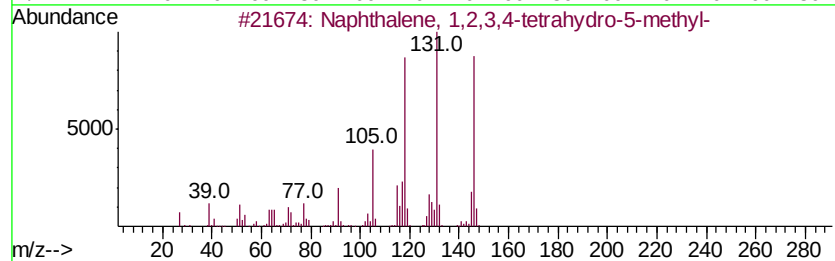
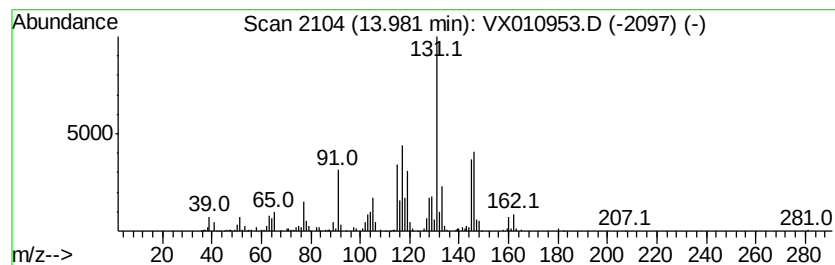
Instrument :
MSVOA_X
ClientSampled :
TR-04-071519

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

Peak Number 5 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.98	10.72 ug/l	190307	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	76
2	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	70
3	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	64
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	60
5	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX071819\
Data File : VX010953.D
Acq On : 18 Jul 2019 13:02
Operator : JC/SP
Sample : K3627-03
Misc : 3.54µ/5mL/100µL/5mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

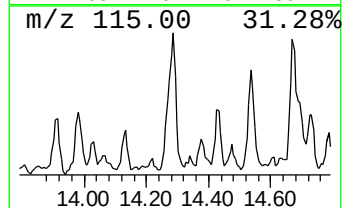
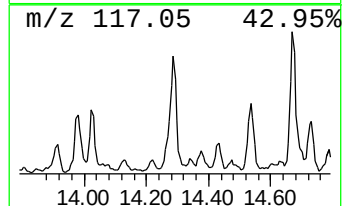
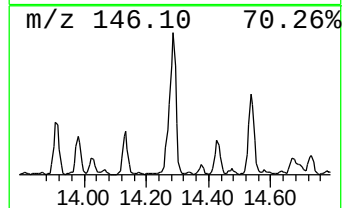
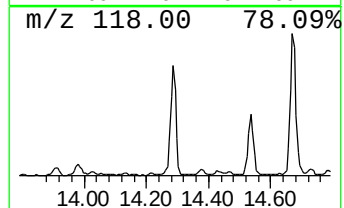
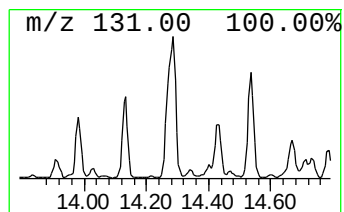
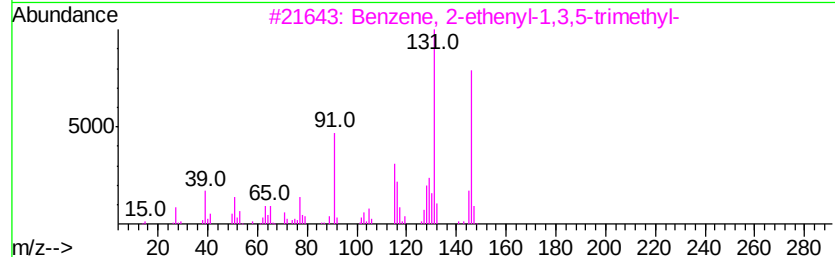
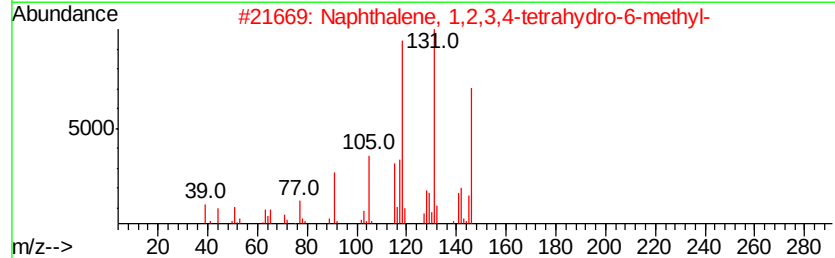
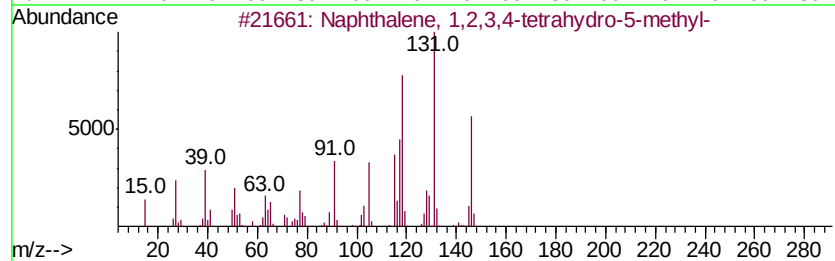
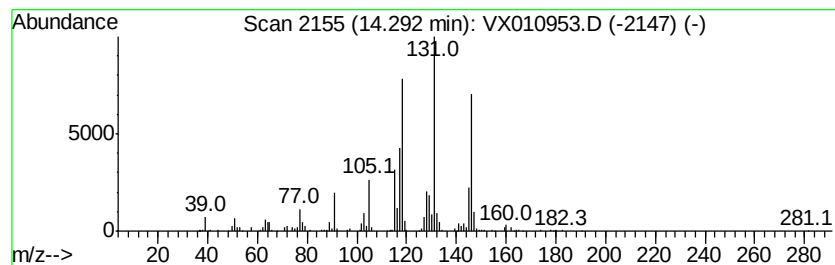
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ClientSampleId :
TR-04-071519

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.29	20.23 ug/l	359192	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 95
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 93
3		Benzene, 2-ethenyl-1,3,5-trimethyl-	146 C11H14	000769-25-5 86
4		Benzene, (1-methyl-1-butenyl)-	146 C11H14	053172-84-2 86
5		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX071819\
Data File : VX010953.D
Acq On : 18 Jul 2019 13:02
Operator : JC/SP
Sample : K3627-03
Misc : 3.54uL/5mL/100uL/5mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

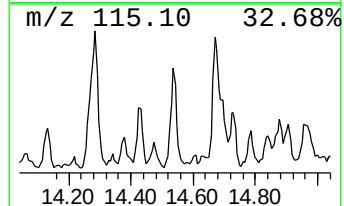
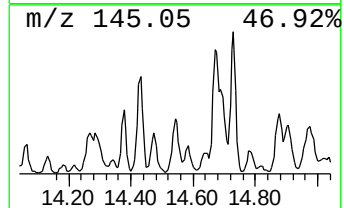
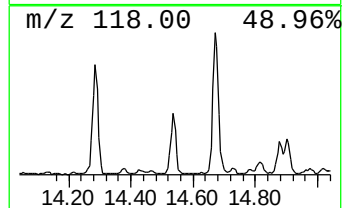
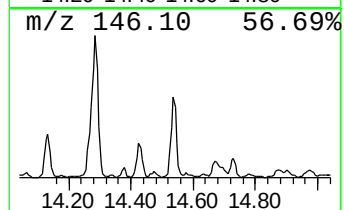
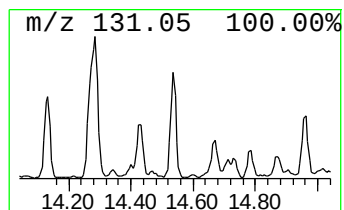
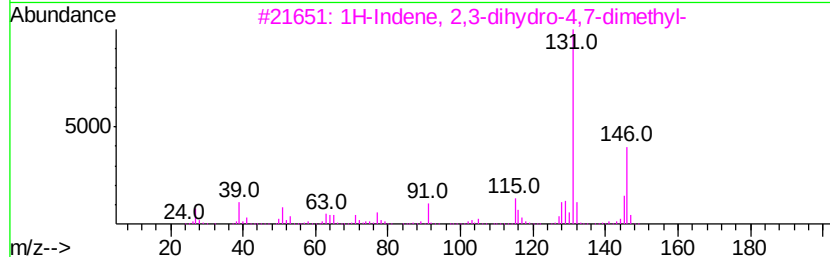
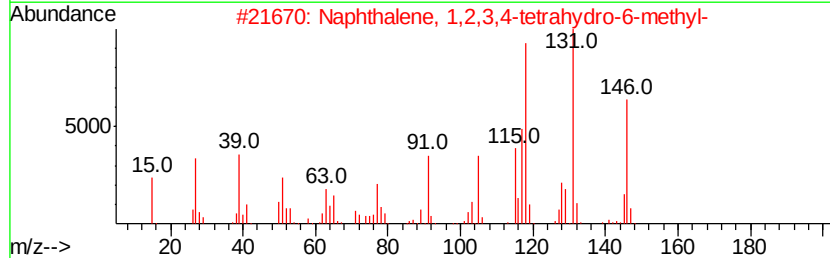
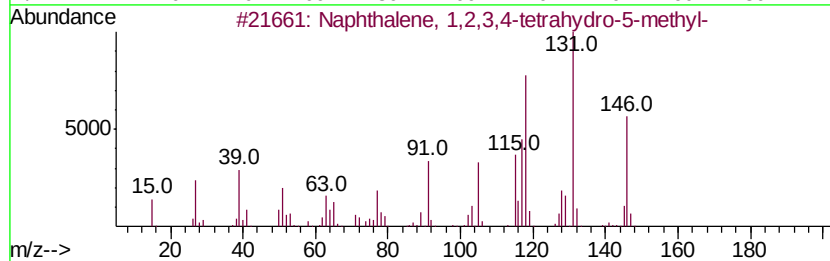
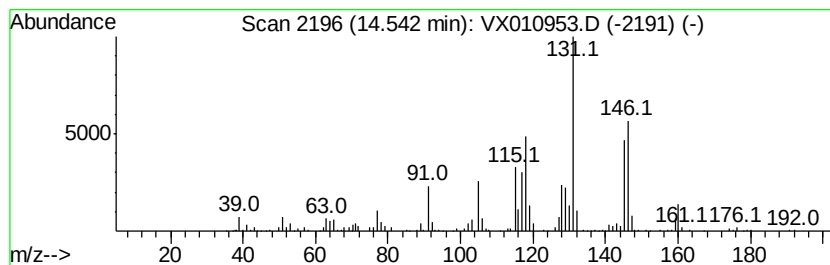
Instrument :
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ClientSampleID :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

Peak Number 7 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.54	13.48 ug/l	239220	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 93
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 86
3		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 64
4		Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3 64
5		Naphthalene, 6-ethyl-1,2,3,4-tet...	160 C12H16	022531-20-0 64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX071819\
Data File : VX010953.D
Acq On : 18 Jul 2019 13:02
Operator : JC/SP
Sample : K3627-03
Misc : 3.54µ/5mL/100µL/5mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
TR-04-071519

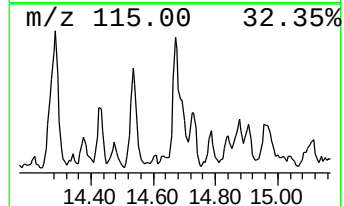
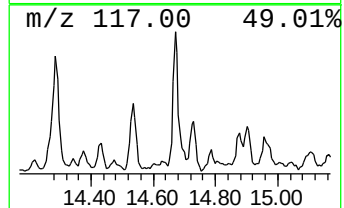
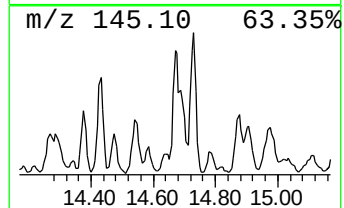
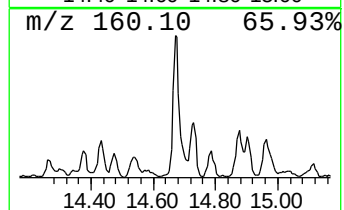
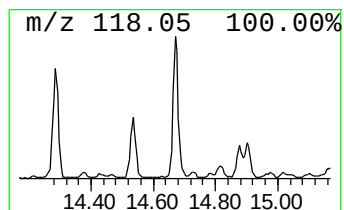
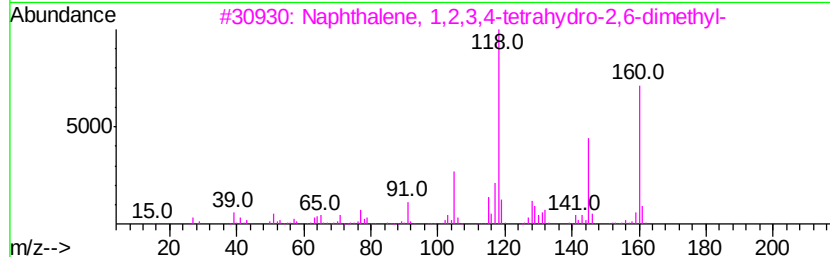
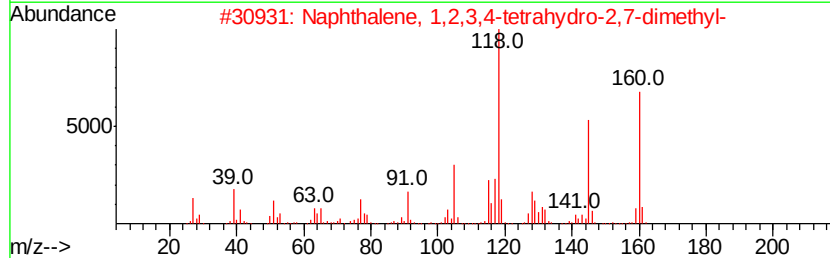
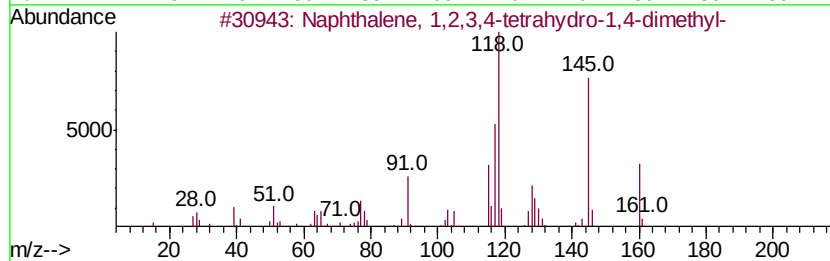
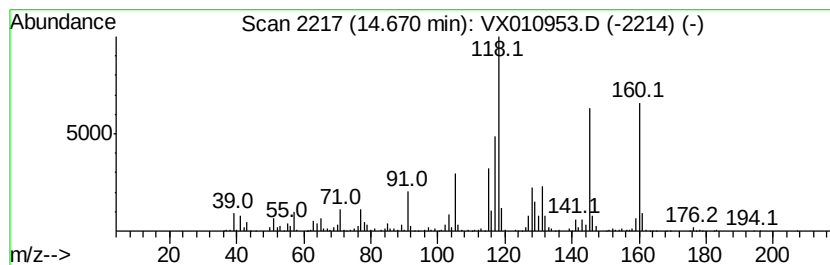
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.67	19.48 ug/l	345795	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	95
2		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	95
3		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	007524-63-2	76
4		1(2H)-Naphthalenone, 3,4-dihydro...	160	C11H12O	014944-23-1	58
5		1(2H)-Naphthalenone, 3,4-dihydro...	160	C11H12O	001590-08-5	52



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX071819\
Data File : VX010953.D
Acq On : 18 Jul 2019 13:02
Operator : JC/SP
Sample : K3627-03
Misc : 3.54g/5mL/100uL/5mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

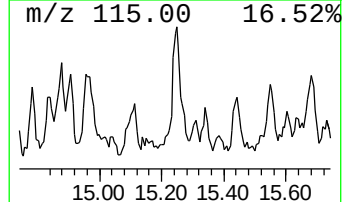
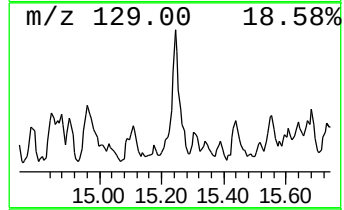
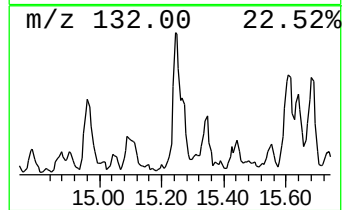
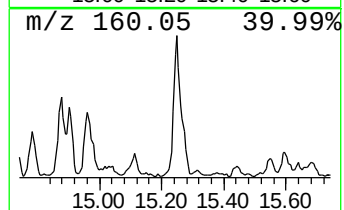
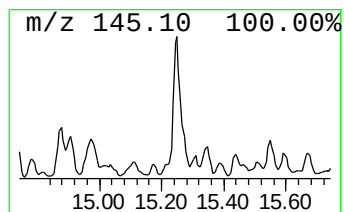
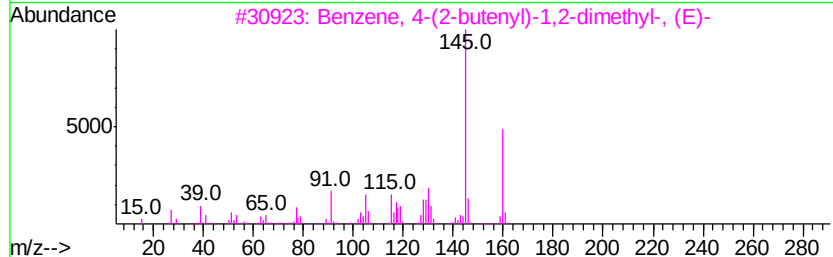
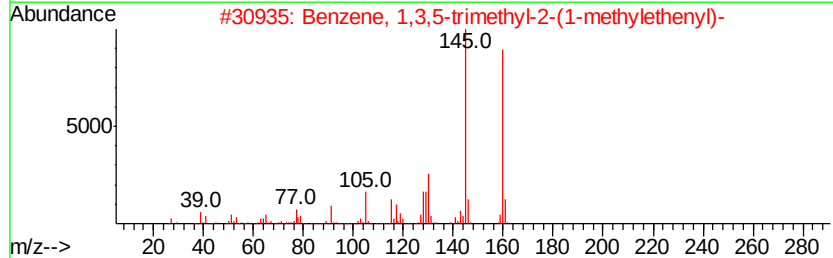
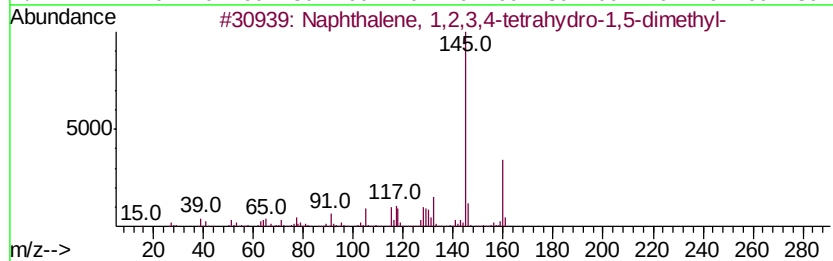
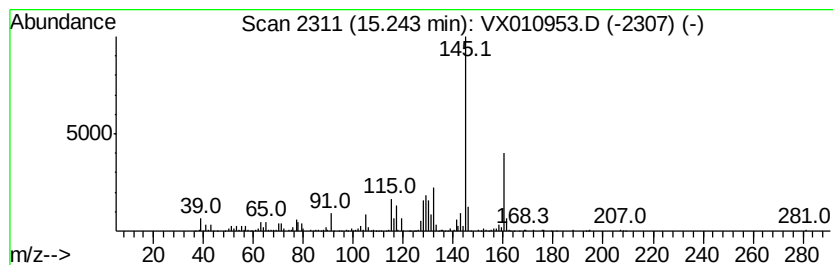
Instrument :
MSVOA_X
ClientSampleId :
TR-04-071519

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,3,5-trimethyl-2-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.24	9.32 ug/l	165543	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	021564-91-0 91
2		Benzene, 1,3,5-trimethyl-2-(1-me...	160 C12H16	014679-13-1 91
3		Benzene, 4-(2-butenyl)-1,2-dimet...	160 C12H16	054340-86-2 91
4		1H-Indene, 2,3-dihydro-1,1,5-tri...	160 C12H16	040650-41-7 87
5		Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	025419-33-4 83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX071819\
Data File : VX010953.D
Acq On : 18 Jul 2019 13:02
Operator : JC/SP
Sample : K3627-03
Misc : 3.54µ/5mL/100µL/5mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

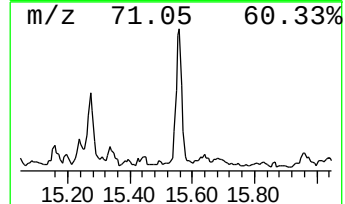
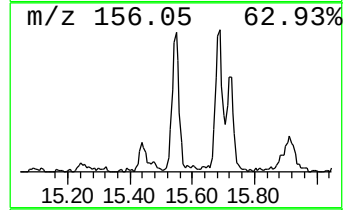
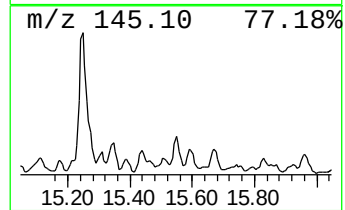
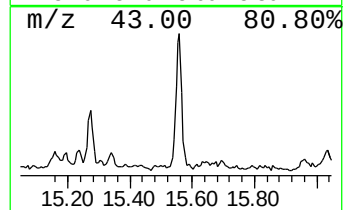
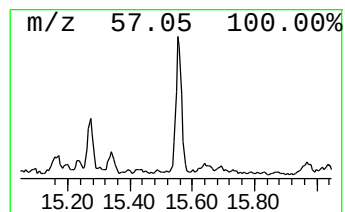
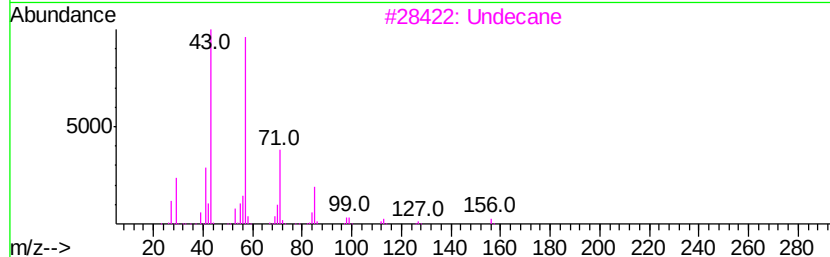
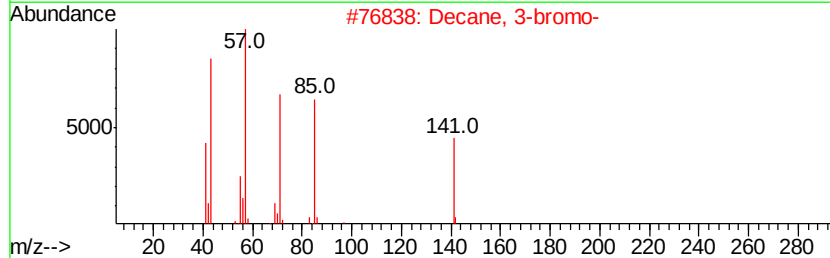
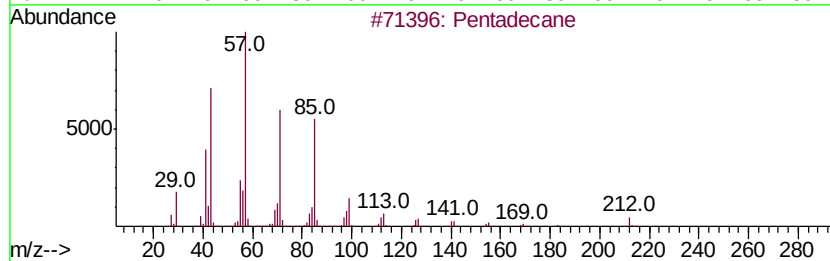
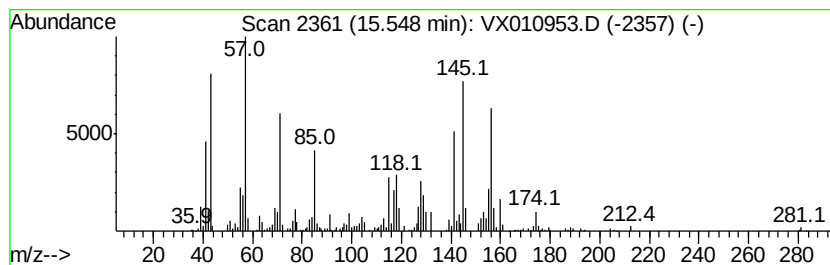
Instrument :
MSVOA_X
ClientSampleId :
TR-04-071519

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

Peak Number 10 Pentadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.55	8.43 ug/l	149736	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Pentadecane	212 C15H32	000629-62-9	60
2	Decane, 3-bromo-	220 C10H21Br	030571-71-2	43
3	Undecane	156 C11H24	001120-21-4	43
4	Dodecane	170 C12H26	000112-40-3	42
5	Hexadecane	226 C16H34	000544-76-3	42



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX071819\
Data File : VX010953.D
Acq On : 18 Jul 2019 13:02
Operator : JC/SP
Sample : K3627-03
Misc : 3.54g/5mL/100uL/5mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TR-04-071519

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.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
Undecane	12.47	7.8	ug/l	138928	4	12.08	887643	50.0
Benzene, (1-methy...	13.35	11.6	ug/l	205760	4	12.08	887643	50.0
Naphthalene, 1,2,...	13.48	10.9	ug/l	193759	4	12.08	887643	50.0
Naphthalene, 1,2,...	13.91	9.9	ug/l	176213	4	12.08	887643	50.0
Naphthalene, 1,2,...	13.98	10.7	ug/l	190307	4	12.08	887643	50.0
Naphthalene, 1,2,...	14.29	20.2	ug/l	359192	4	12.08	887643	50.0
1H-Indene, 2,3-di...	14.54	13.5	ug/l	239220	4	12.08	887643	50.0
Naphthalene, 1,2,...	14.67	19.5	ug/l	345795	4	12.08	887643	50.0
Benzene, 1,3,5-tr...	15.24	9.3	ug/l	165543	4	12.08	887643	50.0
Pentadecane	15.55	8.4	ug/l	149736	4	12.08	887643	50.0