LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX072020\
Data File : VX017437.D
Acq On : 20 Jul 2020 16:28
Operator : JC/SP
Sample : L3329-13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : LT-4

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\82X071320W.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	2.416	211	218	227	rBV	21243	40211	2.59%	0.222%
2	2.575	235	244	253	rBV	42398	82949	5.34%	0.457%
3	5.464	700	718	730	rBV	171936	500375	32.24%	2.756%
4	5.629	736	745	759	rVB2	260674	724968	46.71%	3.994%
5	6.031	800	811	828	rBV	176749	468930	30.21%	2.583%
6	6.830	930	942	955	rBV	426006	1007373	64.91%	5.549%
7	7.440	1033	1042	1049	rBV3	16386	42277	2.72%	0.233%
8	8.695	1234	1248	1262	rBV	918087	1428858	92.06%	7.871%
9	9.019	1295	1301	1307	rVB7	7694	16052	1.03%	0.088%
10	10.098	1468	1478	1487	rBV	997816	1360706	87.67%	7.496%
11	10.323	1510	1515	1524	rVB9	6855	19417	1.25%	0.107%
12	11.122	1638	1646	1652	rBV	944122	1142166	73.59%	6.292%
13	11.652	1728	1733	1741	rVB4	12784	26986	1.74%	0.149%
14	11.750	1741	1749	1754	rBV	14831	27629	1.78%	0.152%
15	11.902	1770	1774	1778	rBV	15964	24552	1.58%	0.135%
16	12.061	1795	1800	1812	rVB	1265791	1552050	100.00%	8.550%
17	12.280	1829	1836	1841	rBV	491001	623282	40.16%	3.433%
18	12.445	1856	1863	1868	rBV	66711	85929	5.54%	0.473%
19	12.652	1887	1897	1900	rBV	178134	231589	14.92%	1.276%
20	12.683	1900	1902	1907	rVV	110654	135267	8.72%	0.745%
21	12.737	1907	1911	1918	rVB	446485	523379	33.72%	2.883%
22	12.878	1927	1934	1939	rVB	50655	75945	4.89%	0.418%
23	12.963	1940	1948	1952	rBV	351720	449013	28.93%	2.473%
24	13.006	1952	1955	1960	rVB2	17146	24155	1.56%	0.133%
25	13.073	1960	1966	1971	rBV	42026	56996	3.67%	0.314%
26	13.134	1972	1976	1979	rBV	30857	34312	2.21%	0.189%
27	13.176	1979	1983	1985	rBV	71111	83225	5.36%	0.458%
28	13.207	1985	1988	1992	rBV	208702	224815	14.49%	1.238%
29	13.298	1998	2003	2004	rBV3	29390	38438	2.48%	0.212%
30	13.329	2004	2008	2014	rVV2	659553	922268	59.42%	5.081%
31	13.390	2014	2018	2023	rVV2	38070	74191	4.78%	0.409%
32	13.463	2027	2030	2036	rVB2	31894	45024	2.90%	0.248%
33	13.548	2036	2044	2046	rBV2	61564	96965	6.25%	0.534%
34	13.585	2046	2050	2053	rVV	179586	251894	16.23%	1.388%

LSC Area Percent Report

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Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : LT-4

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\82X071320W.M
Title : SW846 8260

35	13.621	2053	2056	2060	rVV3	185038	261417	16.84%	1.440%
36	13.682	2060	2066	2067	rVV2	141287	225330	14.52%	1.241%
37	13.701	2067	2069	2079	rVB2	266090	408275	26.31%	2.249%
38	13.792	2079	2084	2087	rBV	37764	50215	3.24%	0.277%
39	13.835	2087	2091	2097	rVB2	43160	65830	4.24%	0.363%
40	13.896	2097	2101	2106	rBV2	77679	124984	8.05%	0.689%
41	13.963	2106	2112	2117	rBV2	158617	230096	14.83%	1.268%
42	14.012	2117	2120	2124	rVB	79495	100022	6.44%	0.551%
43	14.115	2132	2137	2140	rBV	179947	224940	14.49%	1.239%
44	14.146	2140	2142	2148	rVB2	30312	34530	2.22%	0.190%
45	14.255	2155	2160	2169	rBV	278654	428023	27.58%	2.358%
46	14.359	2173	2177	2182	rVV	124026	158637	10.22%	0.874%
47	14.414	2182	2186	2191	rVB2	319678	453892	29.24%	2.500%
48	14.463	2191	2194	2199	rVB2	28044	38887	2.51%	0.214%
49	14.524	2199	2204	2208	rBV2	202850	281764	18.15%	1.552%
50	14.566	2208	2211	2217	rVB3	18949	24473	1.58%	0.135%
51	14.639	2217	2223	2227	rBV	171547	272852	17.58%	1.503%
52	14.719	2232	2236	2240	rVB3	60397	85719	5.52%	0.472%
53	14.768	2240	2244	2248	rVB6	27721	41067	2.65%	0.226%
54	14.828	2249	2254	2262	rBV3	85828	187300	12.07%	1.032%
55	14.944	2267	2273	2280	rBV9	25014	61165	3.94%	0.337%
56	15.048	2286	2290	2294	rVB3	25298	37408	2.41%	0.206%
57	15.091	2294	2297	2302	rBV4	26432	32311	2.08%	0.178%
58	15.225	2313	2319	2327	rBV	71558	134802	8.69%	0.743%
59	15.304	2327	2332	2336	rVV7	24486	41676	2.69%	0.230%
60	15.456	2352	2357	2365	rVB	103989	155142	10.00%	0.855%
61	15.566	2370	2375	2382	rVB2	66057	124312	8.01%	0.685%
62	15.664	2382	2391	2397	rBV	373234	595626	38.38%	3.281%
63	15.786	2405	2411	2418	rBV2	41644	87306	5.63%	0.481%
64	15.865	2418	2424	2426	rVV2	88263	153768	9.91%	0.847%
65	15.895	2426	2429	2440	rVB	119617	205347	13.23%	1.131%
66	16.042	2448	2453	2465	rVB	105360	194177	12.51%	1.070%
67	16.389	2505	2510	2516	rVB2	41351	77198	4.97%	0.425%
68	16.542	2530	2535	2539	rBV7	10503	18436	1.19%	0.102%
69	16.596	2540	2544	2551	rVV4	13555	35049	2.26%	0.193%
70	16.657	2551	2554	2560	rVV6	16279	30948	1.99%	0.170%
71	16.724	2560	2565	2575	rVB5	10182	27951	1.80%	0.154%

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Integrator: RTE
Smoothing : ON Filtering: 5
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Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071320W.M
Title : SW846 8260

Sum of corrected areas: 18153061

LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX072020\
Data File : VX017437.D
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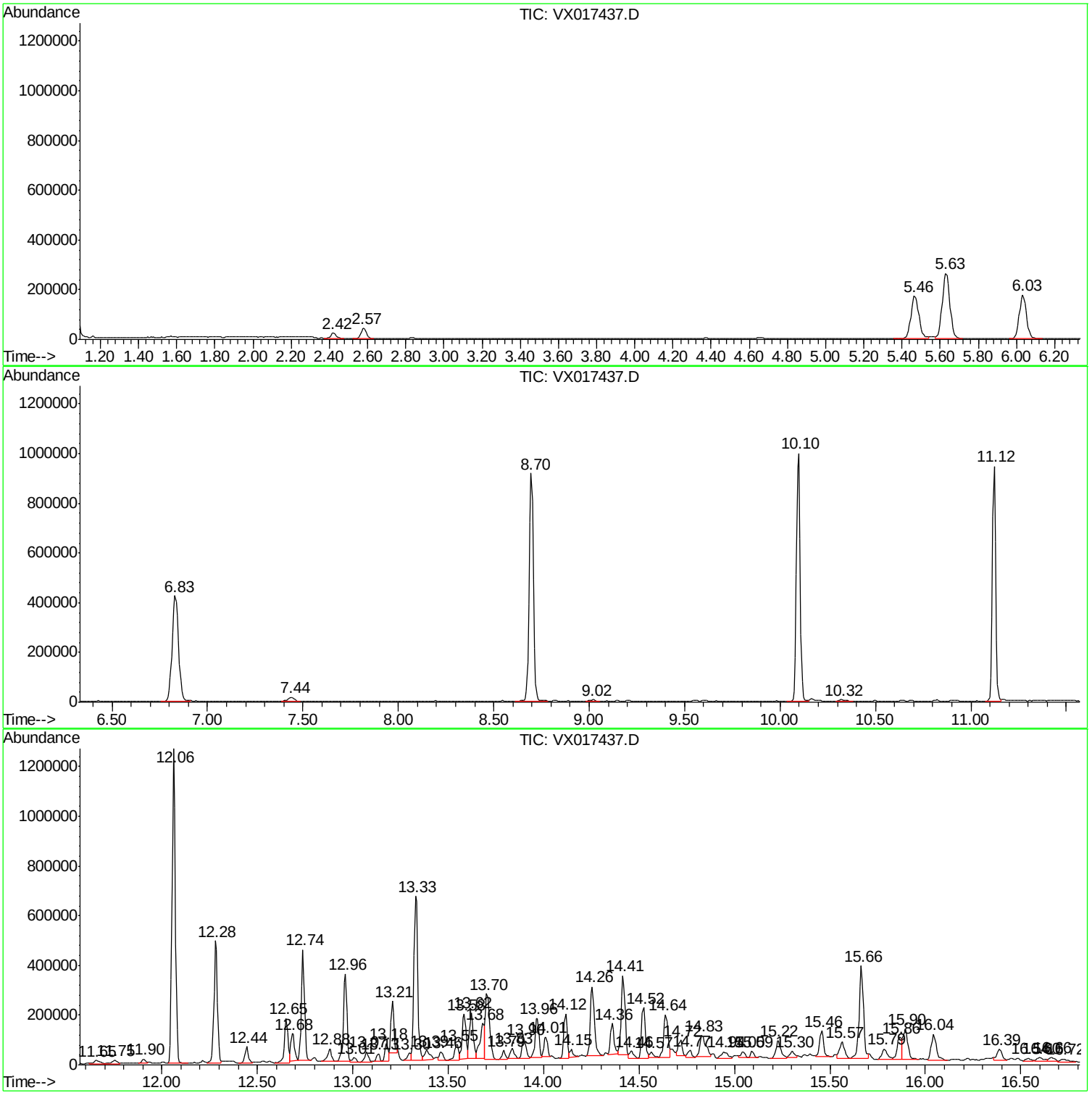
Operator : JC/SP
Sample : L3329-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : LT-4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071320W.M
Quant Title : SW846 8260

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



Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX072020\
Data File : VX017437.D
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Sample : L3329-13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : LT-4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071320W.M
Quant Title : SW846 8260

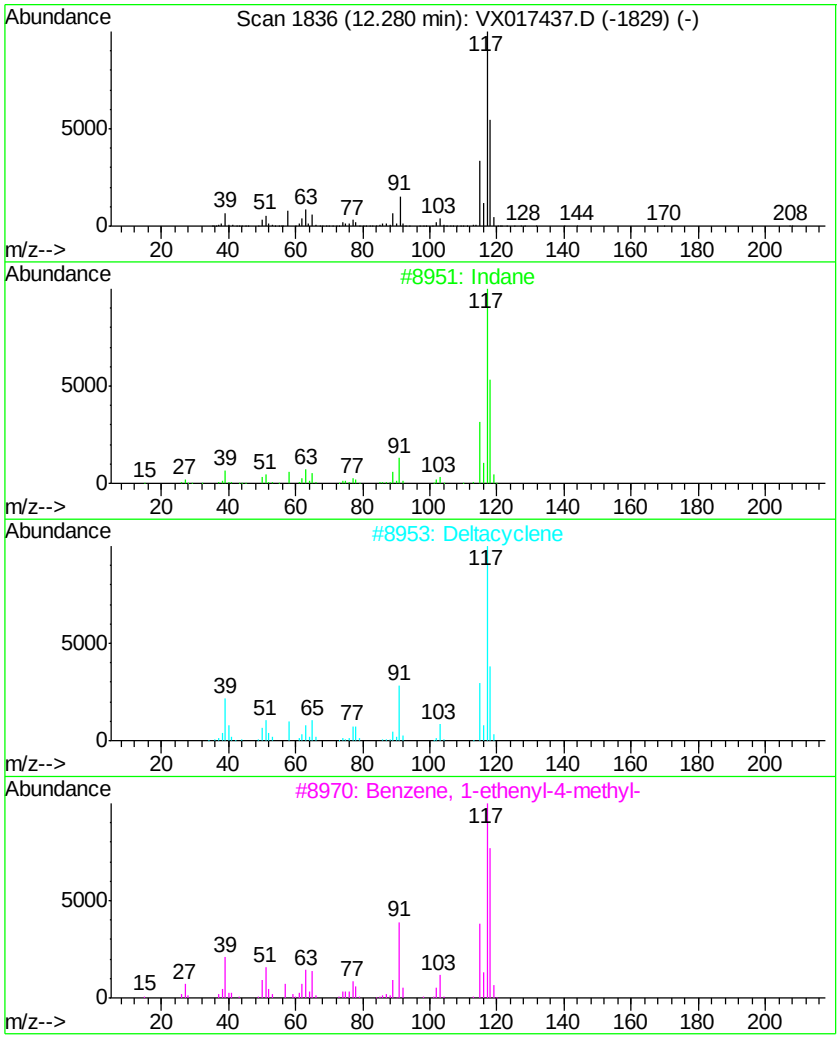
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

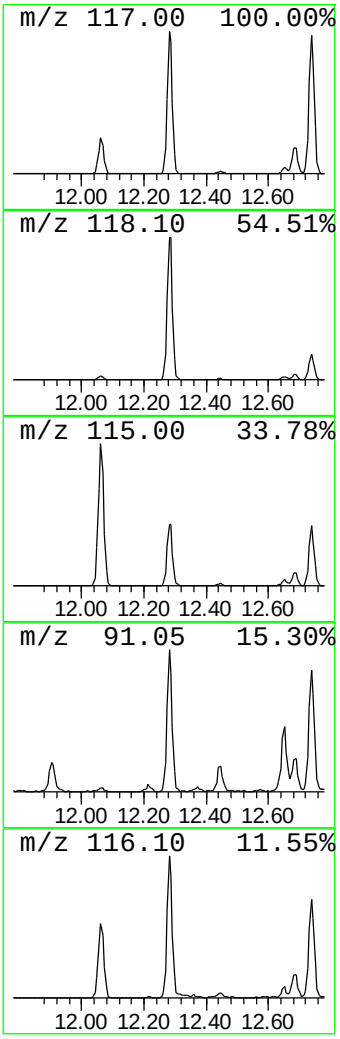
****

Peak Number 1 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.28	20.08 ug/l	623282	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	95
2	Deltacyclene	118	C9H10	007785-10-6	72
3	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	72
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	72
5	7-Methylenecycloocta-1,3,5-triene	118	C9H10	002570-13-0	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX072020\
Data File : VX017437.D
Acq On : 20 Jul 2020 16:28
Operator : JC/SP
Sample : L3329-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : LT-4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071320W.M
Quant Title : SW846 8260

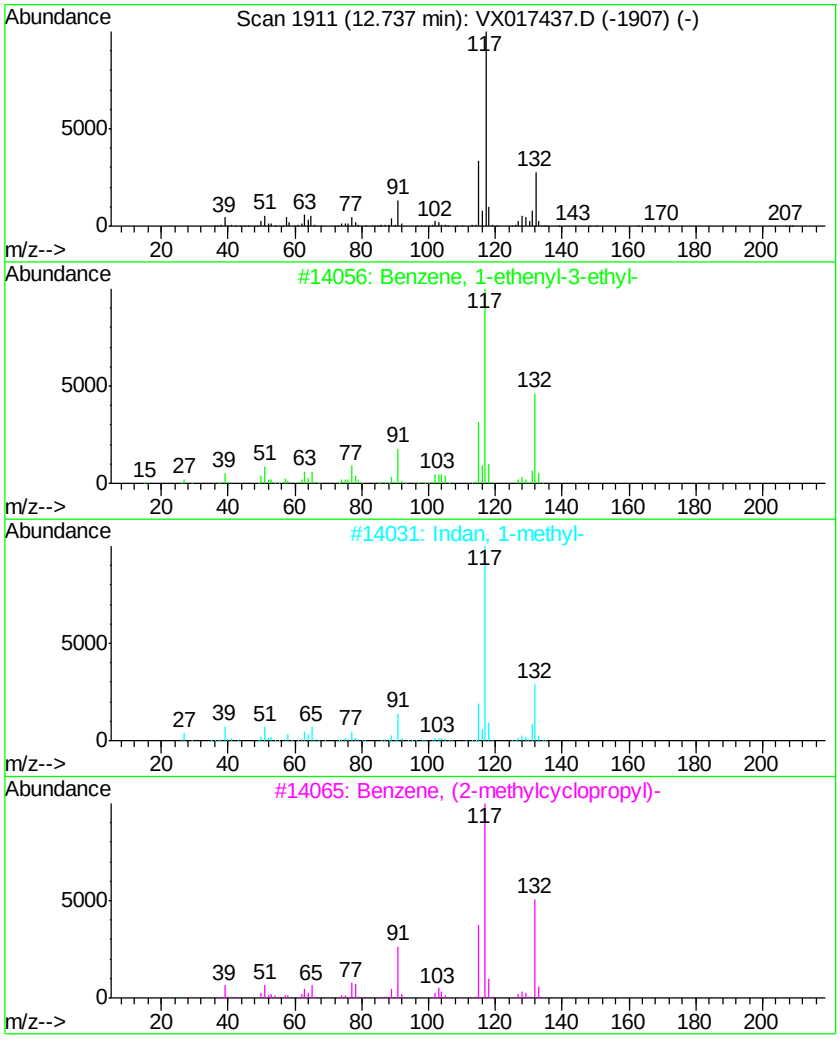
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

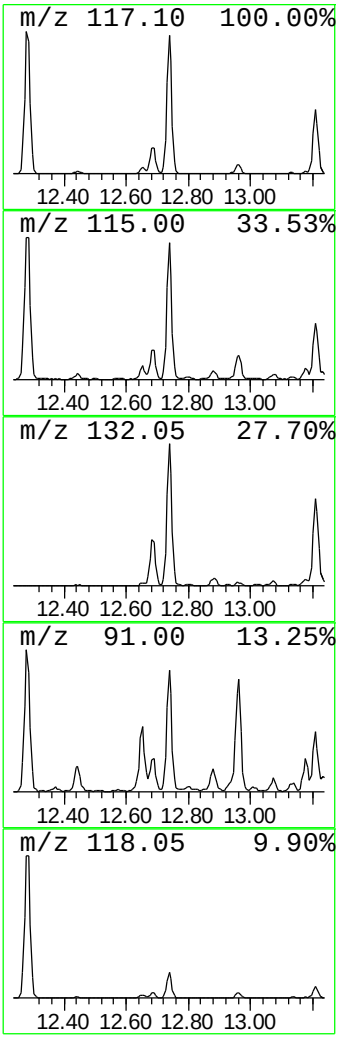
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.74	16.86 ug/l	523379	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
2	Indan, 1-methyl-	132	C10H12	000767-58-8	90
3	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	90
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	83





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX072020\
Data File : VX017437.D
Acq On : 20 Jul 2020 16:28
Operator : JC/SP
Sample : L3329-13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : LT-4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071320W.M
Quant Title : SW846 8260

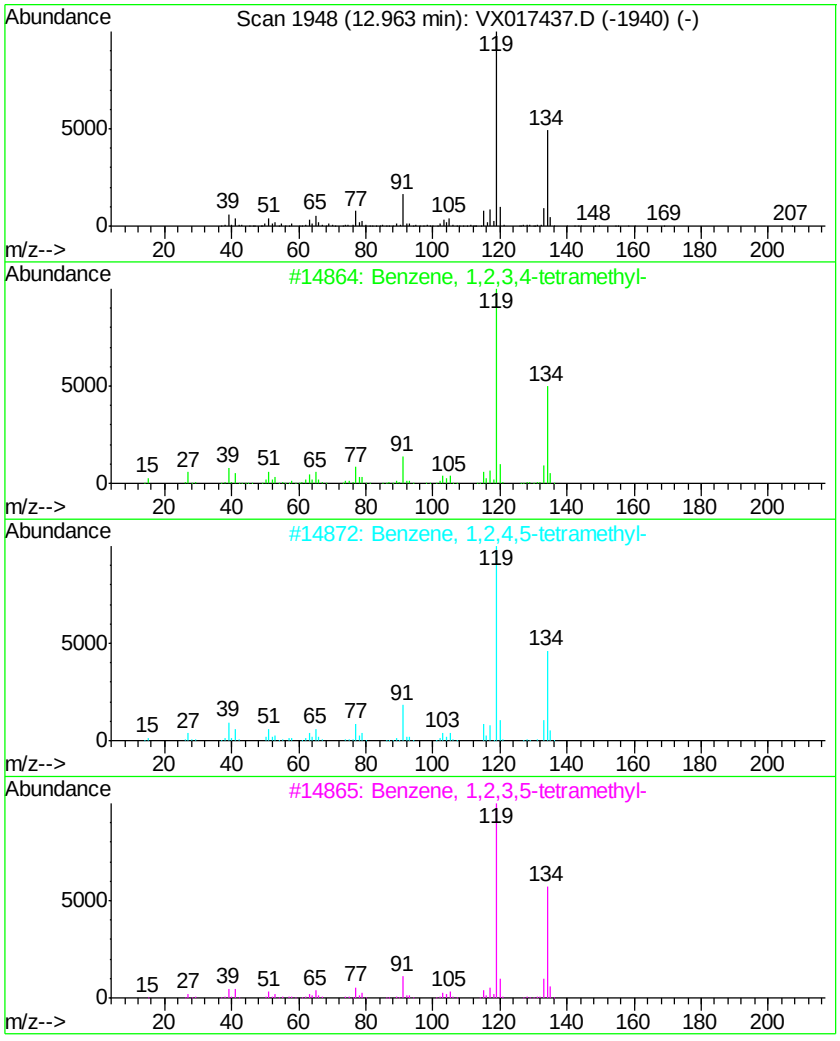
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

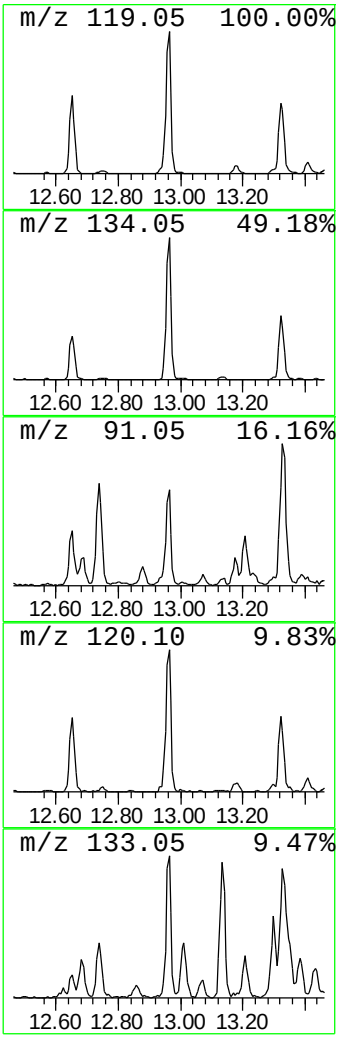
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.96	14.47 ug/l	449013	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	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	o-Cymene	134	C10H14	000527-84-4	94
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX072020\
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ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : LT-4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071320W.M
Quant Title : SW846 8260

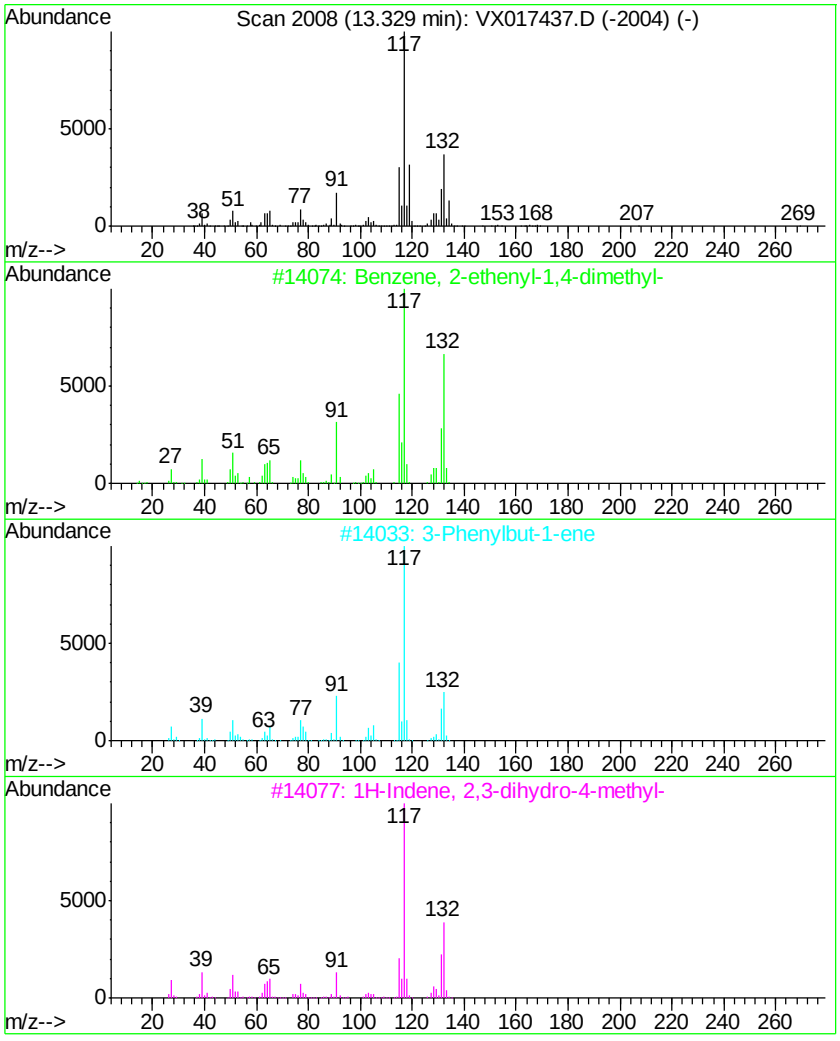
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

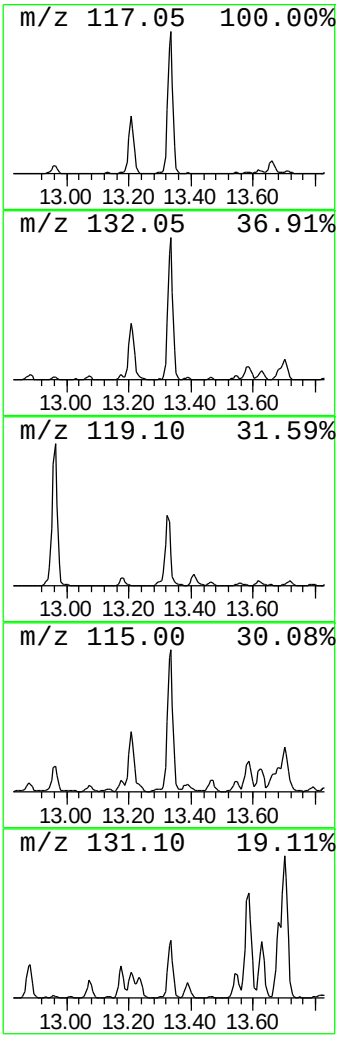
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.33	29.71 ug/l	922268	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	81
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	76
5	Indan, 1-methyl-	132	C10H12	000767-58-8	76





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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071320W.M
Quant Title : SW846 8260

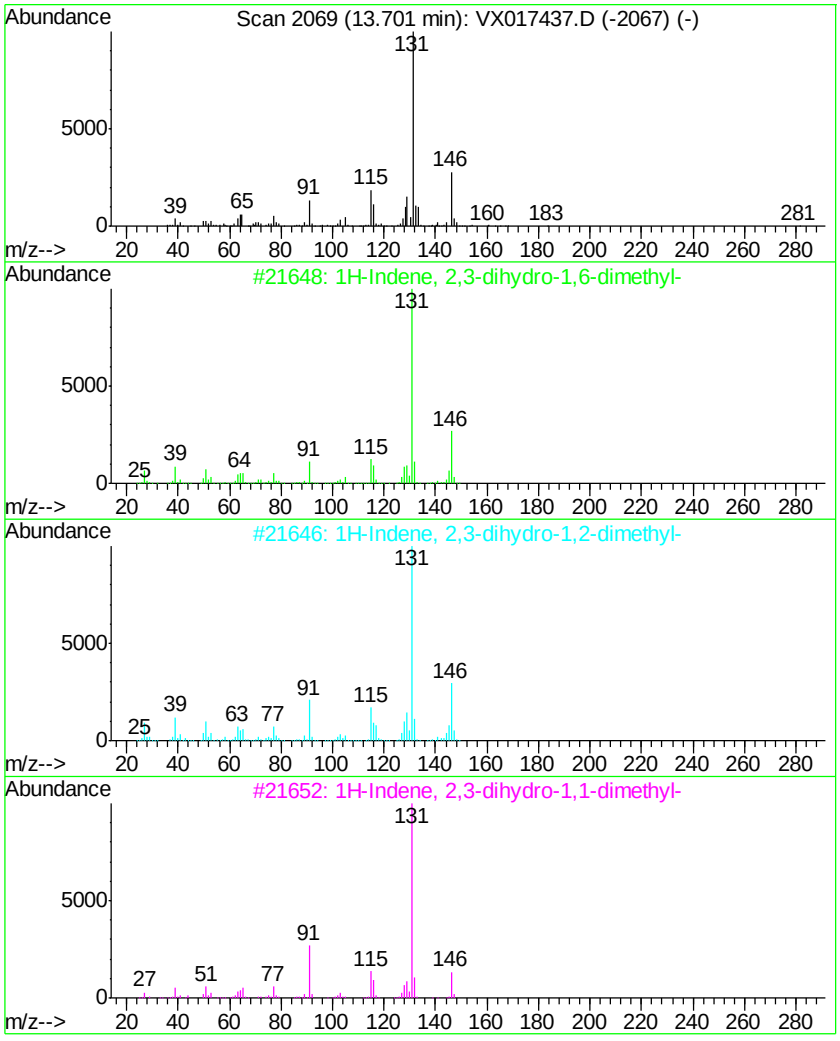
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

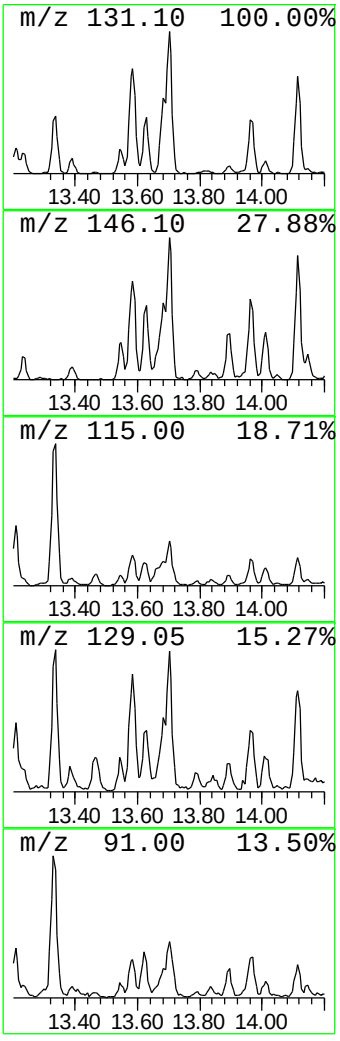
****

Peak Number 5 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.70	13.15 ug/l	408275	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	93
2	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	90
3	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	90
4	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	90
5	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	90





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ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : LT-4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071320W.M
Quant Title : SW846 8260

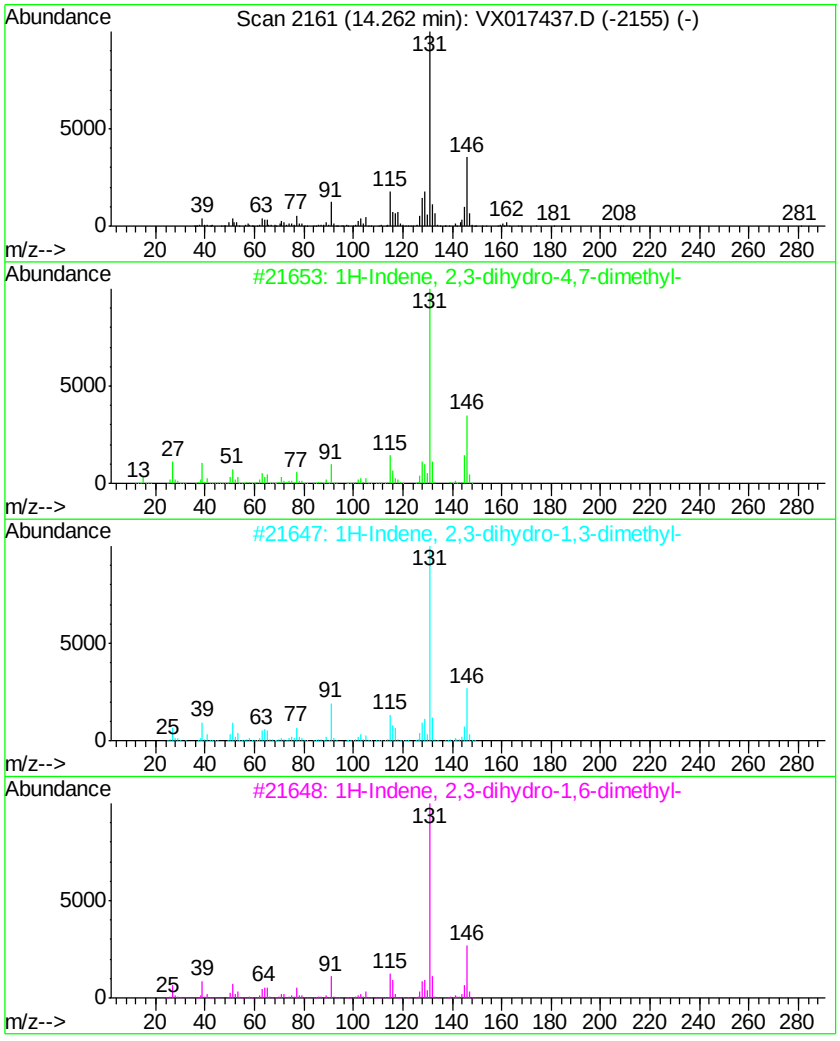
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

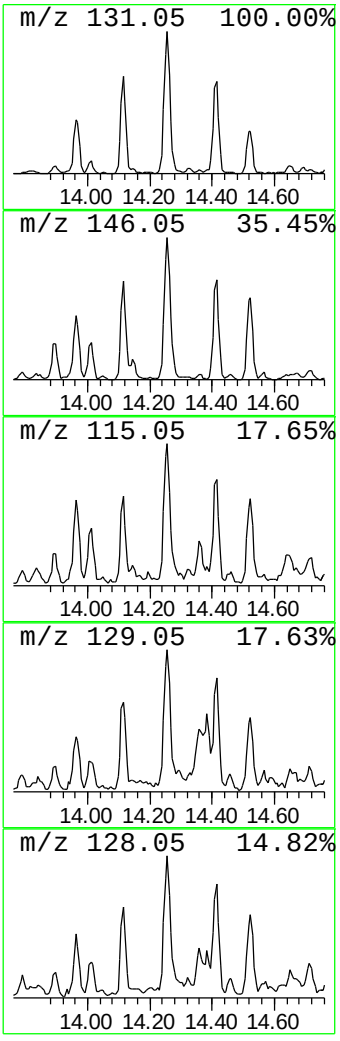
****

Peak Number 6 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.26	13.79 ug/l	428023	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	97
2	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
3	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	91
4	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	90
5	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	90





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ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : LT-4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071320W.M
Quant Title : SW846 8260

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

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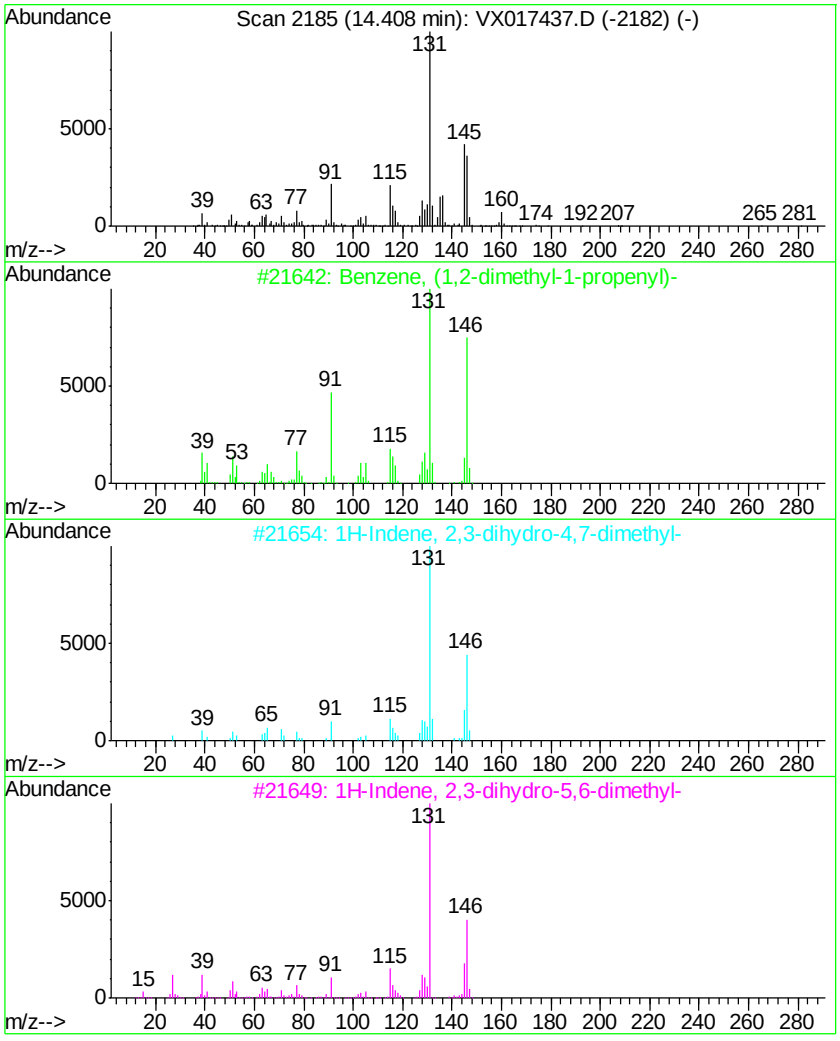
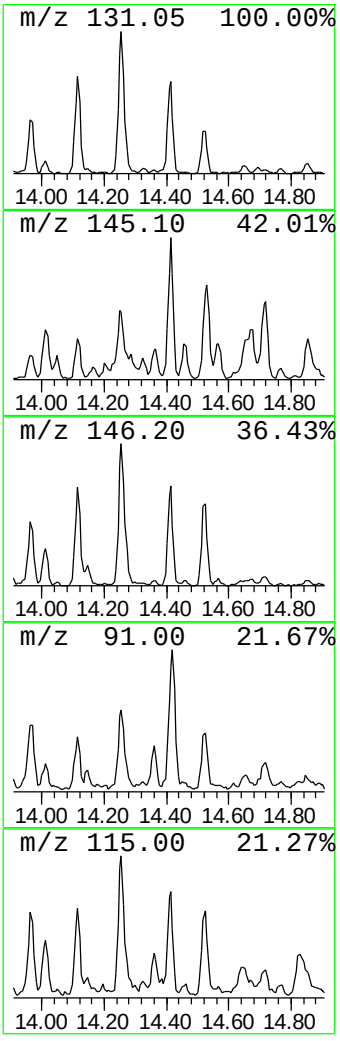
Peak Number 7 Benzene, (1,2-dimethyl-1-pr... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.41	14.62 ug/l	453892	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	89
2	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	55
3	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	50
4	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	50
5	1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	50





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX072020\
Data File : VX017437.D
Acq On : 20 Jul 2020 16:28
Operator : JC/SP
Sample : L3329-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : LT-4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071320W.M
Quant Title : SW846 8260

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

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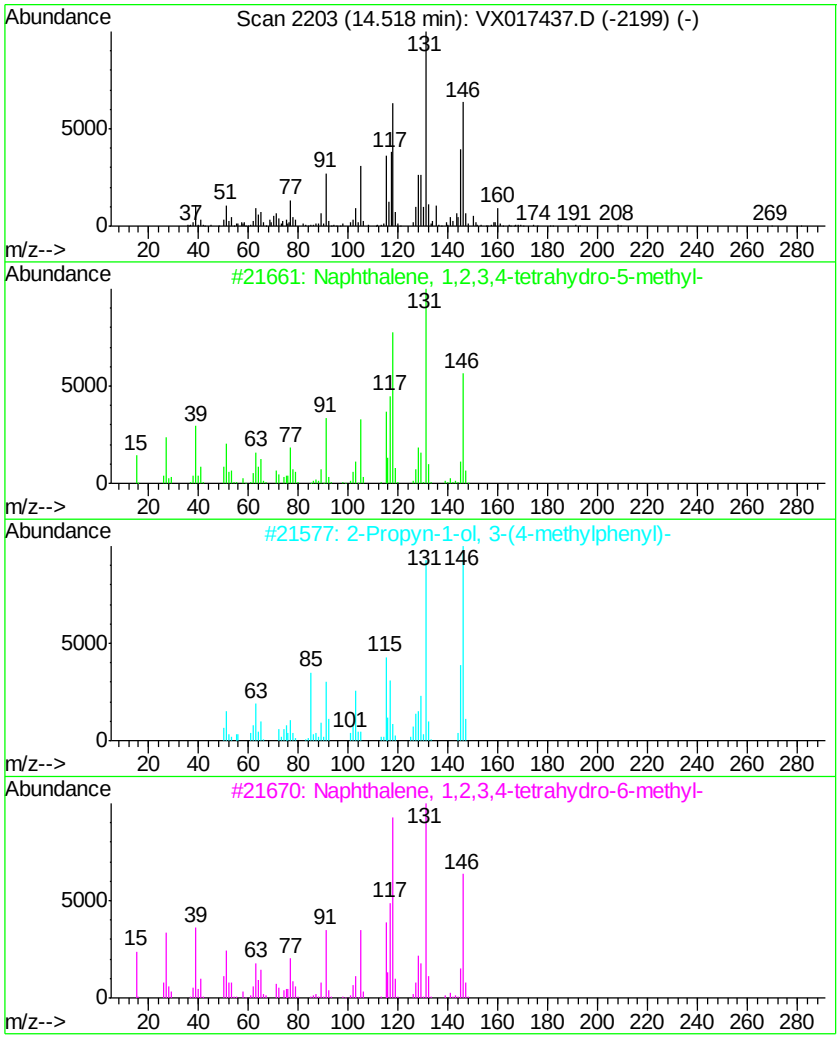
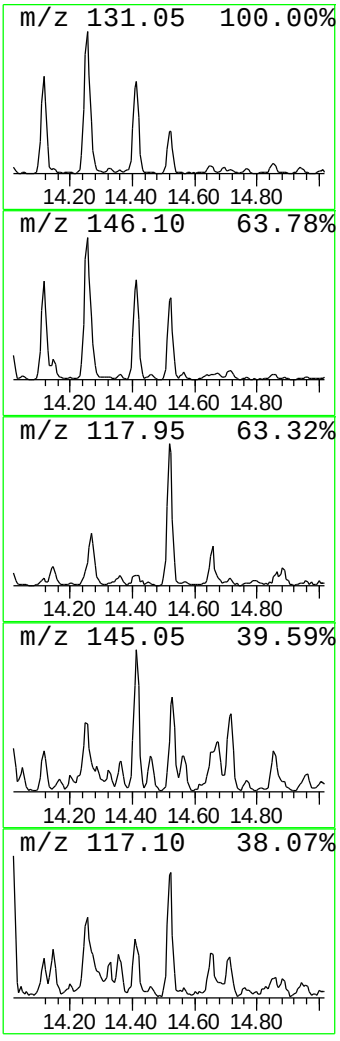
Peak Number 8 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.52	9.08 ug/l	281764	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	76
2	2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	70
3	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	64
4	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	60
5	1H-Benzimidazole, 5,6-dimethyl-	146	C9H10N2	000582-60-5	50





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX072020\
Data File : VX017437.D
Acq On : 20 Jul 2020 16:28
Operator : JC/SP
Sample : L3329-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : LT-4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071320W.M
Quant Title : SW846 8260

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

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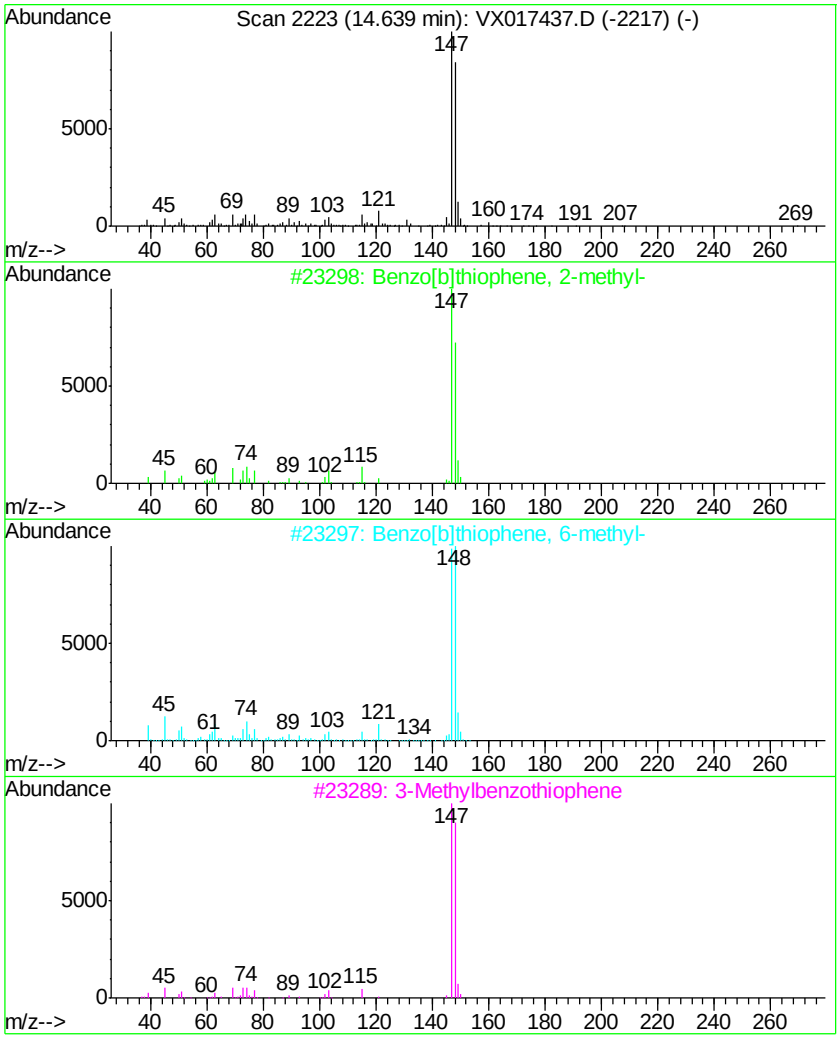
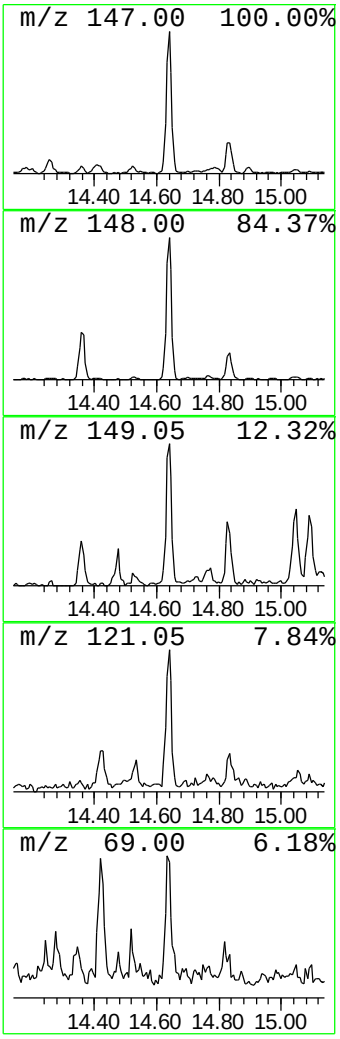
Peak Number 9 Benzo[b]thiophene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.64	8.79 ug/l	272852	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	91
2	Benzo[b]thiophene, 6-methyl-	148	C9H8S	016587-47-6	91
3	3-Methylbenzothiophene	148	C9H8S	001455-18-1	90
4	Benzo[b]thiophene, 5-methyl-	148	C9H8S	014315-14-1	87
5	Benzo[b]thiophene, 4-methyl-	148	C9H8S	014315-11-8	87





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX072020\
Data File : VX017437.D
Acq On : 20 Jul 2020 16:28
Operator : JC/SP
Sample : L3329-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : LT-4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071320W.M
Quant Title : SW846 8260

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

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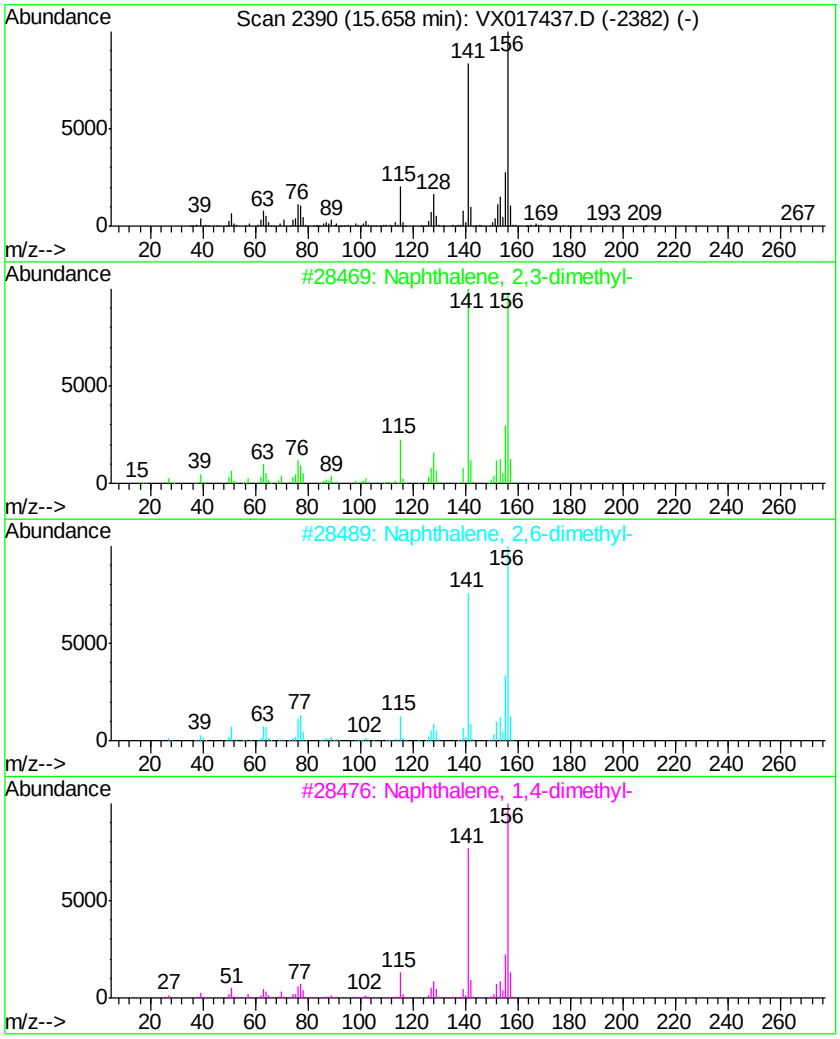
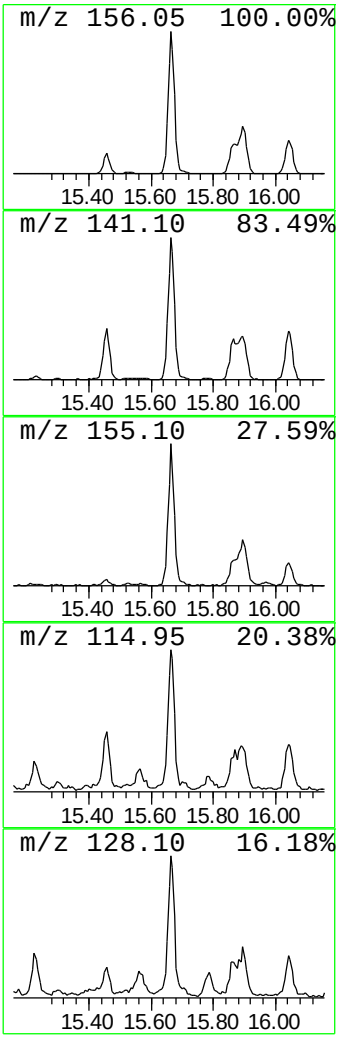
Peak Number 10 Naphthalene, 2,3-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.66	19.19 ug/l	595626	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
4	Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	96
5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX072020\
Data File : VX017437.D
Acq On : 20 Jul 2020 16:28
Operator : JC/SP
Sample : L3329-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : LT-4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071320W.M
Quant Title : SW846 8260

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Indane	12.28	20.1	ug/l	623282	4	12.06	1552050	50.0
Benzene, 1-etheny...	12.74	16.9	ug/l	523379	4	12.06	1552050	50.0
Benzene, 1,2,3,4-...	12.96	14.5	ug/l	449013	4	12.06	1552050	50.0
Benzene, 2-etheny...	13.33	29.7	ug/l	922268	4	12.06	1552050	50.0
1H-Indene, 2,3-di...	13.70	13.2	ug/l	408275	4	12.06	1552050	50.0
1H-Indene, 2,3-di...	14.26	13.8	ug/l	428023	4	12.06	1552050	50.0
Benzene, (1,2-dim...	14.41	14.6	ug/l	453892	4	12.06	1552050	50.0
Naphthalene, 1,2,...	14.52	9.1	ug/l	281764	4	12.06	1552050	50.0
Benzo[b]thiophene...	14.64	8.8	ug/l	272852	4	12.06	1552050	50.0
Naphthalene, 2,3-...	15.66	19.2	ug/l	595626	4	12.06	1552050	50.0