

Data Path : W:\HPCHEM1\MSVOA D\DATA\VD092115\
 Data File : VD046804.D
 Acq On : 21 Sep 2015 13:56
 Operator : FY/SY
 Sample : G3747-01
 Misc : 6.21gm/5mL/MSVOA D/SOIL
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 TRENCH-1-1

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 : W:\HPCHEM1\MSVOA_D\METHOD\82D091815S.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	1.236	69	72	73	rBV	14870	15675	2.05%	0.192%
2	1.275	73	76	85	rVV	372576	691325	90.32%	8.467%
3	1.384	85	87	94	rVB6	5386	14455	1.89%	0.177%
4	1.581	104	107	109	rBV3	3774	7706	1.01%	0.094%
5	1.650	111	114	118	rVV2	16400	26058	3.40%	0.319%
6	1.778	122	127	131	rVV	3270	9898	1.29%	0.121%
7	1.867	134	136	142	rBV6	5289	14699	1.92%	0.180%
8	2.941	240	245	249	rBV	5148	9274	1.21%	0.114%
9	3.493	297	301	308	rVB	15423	34493	4.51%	0.422%
10	5.386	489	493	500	rBV2	7213	22757	2.97%	0.279%
11	6.194	568	575	580	rBV	132224	332943	43.50%	4.078%
12	6.292	580	585	595	rVB	225084	526069	68.73%	6.443%
13	6.697	614	626	634	rBV2	103903	301356	39.37%	3.691%
14	7.416	693	699	707	rBV	287798	633525	82.77%	7.759%
15	8.678	826	827	829	rBV	14854	9834	1.28%	0.120%
16	9.605	915	921	927	rBV	356848	765448	100.00%	9.375%
17	11.566	1115	1120	1132	rVV	414564	754098	98.52%	9.236%
18	11.714	1132	1135	1140	rVB2	5340	10193	1.33%	0.125%
19	11.882	1147	1152	1156	rVB2	8972	18165	2.37%	0.222%
20	12.276	1189	1192	1199	rVB2	8931	16593	2.17%	0.203%
21	12.660	1224	1231	1235	rBV2	7207	13645	1.78%	0.167%
22	12.838	1245	1249	1254	rBV	345003	535992	70.02%	6.564%
23	12.936	1256	1259	1263	rVB2	8434	14208	1.86%	0.174%
24	13.064	1269	1272	1276	rVB	11199	19900	2.60%	0.244%
25	13.262	1288	1292	1297	rVB5	10525	20063	2.62%	0.246%
26	13.409	1304	1307	1312	rVB2	5051	9986	1.30%	0.122%
27	13.587	1321	1325	1330	rVV	141053	226244	29.56%	2.771%
28	13.725	1335	1339	1342	rBV2	9346	19653	2.57%	0.241%
29	13.902	1353	1357	1367	rBV	411779	727877	95.09%	8.915%
30	14.021	1367	1369	1374	rVB4	7183	12233	1.60%	0.150%
31	14.119	1374	1379	1383	rBV2	31479	61303	8.01%	0.751%
32	14.218	1383	1389	1393	rBV5	13956	51014	6.66%	0.625%
33	14.316	1397	1399	1402	rBV3	4184	7934	1.04%	0.097%
34	14.405	1405	1408	1410	rBV	24701	43315	5.66%	0.530%

Data Path : W:\HPCHEM1\MSVOA D\DATA\VD092115\
 Data File : VD046804.D
 Acq On : 21 Sep 2015 13:56
 Operator : FY/SY
 Sample : G3747-01
 Misc : 6.21qm/5mL/MSVOA D/SOIL
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 TRENCH-1-1

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 : W:\HPCHEM1\MSVOA_D\METHOD\82D091815S.M
 Title : SW846 8260

35	14.435	1410	1411	1415	rVV	33034	48267	6.31%	0.591%
36	14.494	1415	1417	1420	rVB	48007	71467	9.34%	0.875%
37	14.553	1420	1423	1425	rVB2	11753	15449	2.02%	0.189%
38	14.602	1425	1428	1432	rVB2	37914	66189	8.65%	0.811%
39	14.691	1433	1437	1439	rBV4	7037	11841	1.55%	0.145%
40	14.740	1439	1442	1443	rBV2	8744	12814	1.67%	0.157%
41	14.819	1448	1450	1452	rBV	27922	32758	4.28%	0.401%
42	14.858	1452	1454	1457	rVB	66578	97623	12.75%	1.196%
43	14.918	1457	1460	1465	rVB4	18794	48131	6.29%	0.589%
44	14.987	1465	1467	1469	rBV	7839	10460	1.37%	0.128%
45	15.046	1469	1473	1475	rVB4	8165	16212	2.12%	0.199%
46	15.095	1475	1478	1481	rBV	26191	54219	7.08%	0.664%
47	15.154	1481	1484	1486	rVB2	23240	38205	4.99%	0.468%
48	15.213	1486	1490	1494	rBV2	92974	195529	25.54%	2.395%
49	15.282	1494	1497	1501	rVB3	50594	90409	11.81%	1.107%
50	15.381	1501	1507	1510	rBV4	23495	58452	7.64%	0.716%
51	15.499	1513	1519	1521	rBV2	45868	111611	14.58%	1.367%
52	15.617	1527	1531	1534	rBV4	48208	100363	13.11%	1.229%
53	15.716	1539	1541	1544	rVB4	19314	30302	3.96%	0.371%
54	15.785	1544	1548	1553	rBV3	57716	124288	16.24%	1.522%
55	15.874	1553	1557	1560	rBV2	18408	42266	5.52%	0.518%
56	15.933	1560	1563	1565	rBV3	19917	41322	5.40%	0.506%
57	16.120	1578	1582	1587	rBV5	43285	120293	15.72%	1.473%
58	16.199	1587	1590	1593	rVV4	22822	47437	6.20%	0.581%
59	16.288	1594	1599	1602	rVV3	27964	83221	10.87%	1.019%
60	16.337	1602	1604	1608	rVV3	28745	53561	7.00%	0.656%
61	16.426	1609	1613	1616	rVV5	14139	30131	3.94%	0.369%
62	16.505	1616	1621	1626	rBV5	20854	66822	8.73%	0.818%
63	16.652	1631	1636	1640	rBV2	37879	110079	14.38%	1.348%
64	16.771	1644	1648	1652	rVB2	42288	75999	9.93%	0.931%
65	16.869	1653	1658	1662	rBV3	28876	67442	8.81%	0.826%
66	16.928	1662	1664	1666	rBV2	11251	21494	2.81%	0.263%
67	17.135	1679	1685	1694	rBV2	57809	192446	25.14%	2.357%

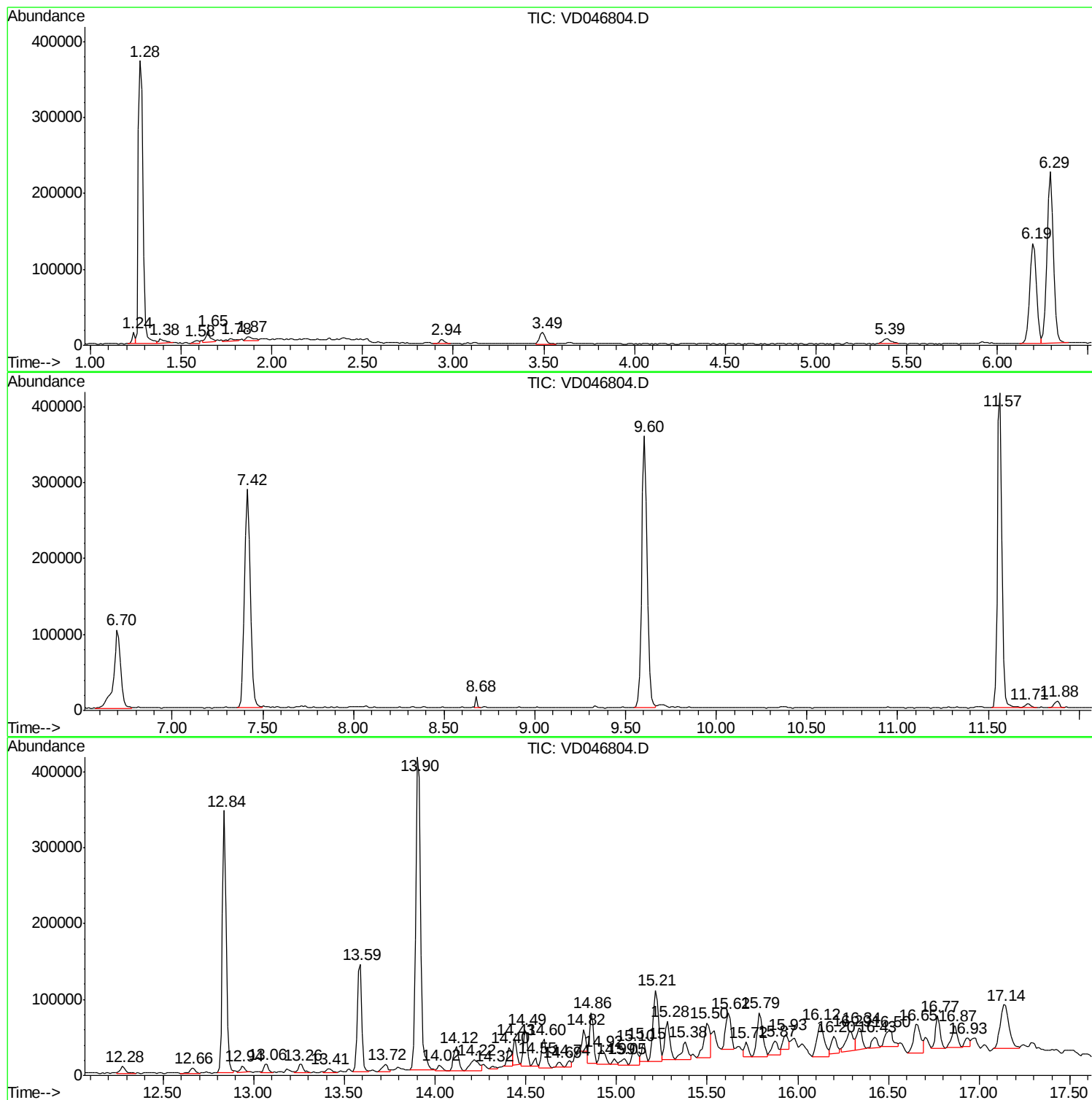
Sum of corrected areas: 8165033

Data Path : W:\HPCHEM1\MSVOA D\DATA\VD092115\
Data File : VD046804.D
Acq On : 21 Sep 2015 13:56
Operator : FY/SY
Sample : G3747-01
Misc : 6.21µm/5mL/MSVOA D/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
TRENCH-1-1

Quant Method : W:\HPCHEM1\MSVOA_D\METHOD\82D091815S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



Data Path : W:\HPCHEM1\MSVOA_D\DATA\VD092115\
Data File : VD046804.D
Acq On : 21 Sep 2015 13:56
Operator : FY/SY
Sample : G3747-01
Misc : 6.21gm/5mL/MSVOA_D/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
TRENCH-1-1

Quant Method : W:\HPCHEM1\MSVOA_D\METHOD\82D091815S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.86	6.71 ug/l	97623	1,4-Dichlorobenzene-d4	13.91
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9 90
2		Benzene, 1-methyl-3-(1-methyleth...	134 C10H14	000535-77-3 90
3		Benzene, 2-ethyl-1,4-dimethyl-	134 C10H14	001758-88-9 90
4		Benzene, 1-methyl-2-(1-methyleth...	134 C10H14	000527-84-4 90
5		Benzene, 4-ethyl-1,2-dimethyl-	134 C10H14	000934-80-5 90

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.21	13.43 ug/l	195529	1,4-Dichlorobenzene-d4	13.91
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-methyl-4-(2-propenyl)-	132 C10H12	003333-13-9 90
2		Benzene, 1-methyl-2-(2-propenyl)-	132 C10H12	001587-04-8 87
3		Benzene, (1-methyl-1-propenyl)-,...	132 C10H12	000767-99-7 87
4		1-Phenyl-1-butene	132 C10H12	000824-90-8 50
5		3-Phenylbut-1-ene	132 C10H12	000934-10-1 50

Peak Number 3 unknown15.28 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.28	6.21 ug/l	90409	1,4-Dichlorobenzene-d4	13.91
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (1,1-dimethylpropyl)-	148 C11H16	002049-95-8 46
2		Benzene, 1-methyl-4-(1-methylpro...	148 C11H16	001595-16-0 43
3		Bis[3,4-dimethylbenzyl]sulfone	302 C18H22O2S	034277-83-3 38
4		Benzene, 1-ethyl-2,3-dimethyl-	134 C10H14	000933-98-2 38
5		o-Toluylic acid, tridec-2-ynyl e...	314 C21H30O2	1000292-51-7 38

Peak Number 4 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
------	---------	------	------------------	------

Data Path : W:\HPCHEM1\MSVOA_D\DATA\VD092115\
 Data File : VD046804.D
 Acq On : 21 Sep 2015 13:56
 Operator : FY/SY
 Sample : G3747-01
 Misc : 6.21gm/5mL/MSVOA_D/SOIL
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 TRENCH-1-1

Quant Method : W:\HPCHEM1\MSVOA_D\METHOD\82D091815S.M
 Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

15.50	7.67 ug/l	111611	1,4-Dichlorobenzene-d4	13.91			
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	87
2			Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	55
3			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	55
4			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	55
5			.alpha.,.beta.,.beta.-Trimethyls...	146	C11H14	000769-57-3	55

 Peak Number 5 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.			
15.62	6.89 ug/l	100363	1,4-Dichlorobenzene-d4	13.91			
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70
2			1H-Indene,2,3-dihydro-2,2-dimethyl-	146	C11H14	020836-11-7	64
3			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	64
4			Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	62
5			1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	58

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.12	8.26 ug/l	120293	1,4-Dichlorobenzene-d4	13.91		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14		001685-82-1	93
2	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14		006682-71-9	93
3	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14		001075-22-5	91
4	.alpha.,.beta.,.beta.-Trimethyls...	146	C11H14		000769-57-3	91
5	Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14		052161-57-6	90

 Peak Number 7 unknown16.65 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.65	7.56 ug/l	110079	1,4-Dichlorobenzene-d4	13.91		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual

Data Path : W:\HPCHEM1\MSVOA_D\DATA\VD092115\
Data File : VD046804.D
Acq On : 21 Sep 2015 13:56
Operator : FY/SY
Sample : G3747-01
Misc : 6.21gm/5mL/MSVOA_D/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
TRENCH-1-1

Quant Method : W:\HPCHEM1\MSVOA_D\METHOD\82D091815S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

1	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	45
2	2-Propenal, 3-(4-methylphenyl)-	146	C10H10O	001504-75-2	43
3	Naphthalene, 1,2,3,4-tetrahydro...	160	C12H16	001076-61-5	43
4	Benzonitrile, 4-(dimethylamino)-	146	C9H10N2	001197-19-9	38
5	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	38

Peak Number 8 1H-Indene, 1-ethylidene- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.14	13.22 ug/l	192446	1,4-Dichlorobenzene-d4	13.91

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	95
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	93
3			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	89
4			Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	76
5			Benzocycloheptatriene	142	C11H10	000264-09-5	76

Data Path : W:\HPCHEM1\MSVOA_D\DATA\VD092115\
Data File : VD046804.D
Acq On : 21 Sep 2015 13:56
Operator : FY/SY
Sample : G3747-01
Misc : 6.21gm/5mL/MSVOA_D/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
TRENCH-1-1

Quant Method : W:\HPCHEM1\MSVOA_D\METHOD\82D091815S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-ethyl-...	14.86	6.7	ug/l	97623	4	13.91	727877	50.0
Benzene, 1-methyl...	15.21	13.4	ug/l	195529	4	13.91	727877	50.0
unknown15.28	15.28	6.2	ug/l	90409	4	13.91	727877	50.0
1H-Indene, 2,3-di...	15.50	7.7	ug/l	111611	4	13.91	727877	50.0
1H-Indene, 2,3-di...	15.62	6.9	ug/l	100363	4	13.91	727877	50.0
1H-Indene, 2,3-di...	16.12	8.3	ug/l	120293	4	13.91	727877	50.0
unknown16.65	16.65	7.6	ug/l	110079	4	13.91	727877	50.0
1H-Indene, 1-ethy...	17.14	13.2	ug/l	192446	4	13.91	727877	50.0