LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX011719\
Data File : VX007088.D
Acq On : 17 Jan 2019 14:56
Operator : JC/SP
Sample : K1168-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038MW050-190115

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\82X010819W.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	1.617	67	76	80	rVB4	34593	94306	1.60%	0.129%
2	1.788	99	104	112	rVB	115848	159301	2.71%	0.217%
3	2.397	198	204	211	rBV2	42176	81329	1.38%	0.111%
4	2.842	268	277	291	rBV	462370	1036200	17.63%	1.415%
5	3.172	323	331	339	rBV2	55664	125804	2.14%	0.172%
6	4.141	477	490	500	rBV2	37045	112030	1.91%	0.153%
7	4.306	503	517	531	rBV	344399	1053199	17.92%	1.438%
8	4.556	545	558	570	rBV3	51912	195646	3.33%	0.267%
9	5.244	658	671	682	rVB4	49274	170798	2.91%	0.233%
10	5.507	700	714	725	rBV2	339388	984313	16.75%	1.344%
11	5.665	729	740	747	rBV2	579488	1694006	28.83%	2.312%
12	5.946	775	786	797	rBV3	194441	603566	10.27%	0.824%
13	6.067	797	806	817	rBV	348449	888955	15.13%	1.214%
14	6.348	840	852	866	rVB	1547519	4622013	78.66%	6.310%
15	6.860	925	936	947	rBV	877999	2045701	34.81%	2.793%
16	7.445	1023	1032	1042	rBV	258859	568961	9.68%	0.777%
17	7.573	1045	1053	1059	rBV	212691	441881	7.52%	0.603%
18	7.677	1059	1070	1078	rVB	245395	581842	9.90%	0.794%
19	7.848	1090	1098	1108	rVB	277783	547020	9.31%	0.747%
20	8.055	1121	1132	1141	rBV	1425948	2729488	46.45%	3.726%
21	8.177	1143	1152	1161	rVV	2548800	4955011	84.33%	6.764%
22	8.390	1178	1187	1192	rBV2	99361	191105	3.25%	0.261%
23	8.579	1210	1218	1227	rVB2	40595	78953	1.34%	0.108%
24	8.713	1227	1240	1254	rBV	1898250	3592075	61.13%	4.904%
25	8.835	1254	1260	1266	rVV	228074	416402	7.09%	0.568%
26	9.043	1288	1294	1301	rVB3	55807	112047	1.91%	0.153%
27	9.164	1307	1314	1319	rVB	59265	103972	1.77%	0.142%
28	9.567	1375	1380	1385	rBV3	49357	78556	1.34%	0.107%
29	9.677	1393	1398	1402	rBV	143318	219885	3.74%	0.300%
30	9.890	1426	1433	1439	rBV	81510	121899	2.07%	0.166%
31	10.109	1464	1469	1473	rBV	1766172	2923061	49.75%	3.990%
32	10.183	1478	1481	1486	rVB	149232	214676	3.65%	0.293%
33	10.335	1501	1506	1514	rVB2	143727	250342	4.26%	0.342%
34	10.408	1514	1518	1521	rBV	65104	86269	1.47%	0.118%

LSC Area Percent Report

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Operator : JC/SP
Sample : K1168-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038MW050-190115

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\82X010819W.M
Title : SW846 8260

35	10.512	1530	1535	1542	rVB	102343	145346	2.47%	0.198%
36	10.658	1545	1559	1564	rBV5	175345	446289	7.60%	0.609%
37	10.713	1564	1568	1574	rVB2	58539	83992	1.43%	0.115%
38	10.810	1574	1584	1585	rBV2	133388	213661	3.64%	0.292%
39	10.835	1585	1588	1593	rVB	366708	479488	8.16%	0.655%
40	10.914	1596	1601	1608	rBV4	130842	231429	3.94%	0.316%
41	11.018	1614	1618	1621	rBV	229754	300371	5.11%	0.410%
42	11.067	1621	1626	1631	rVB5	72038	150754	2.57%	0.206%
43	11.134	1632	1637	1642	rBV	1524528	2017648	34.34%	2.754%
44	11.182	1642	1645	1649	rVV	199542	255816	4.35%	0.349%
45	11.280	1657	1661	1666	rVB	320539	423147	7.20%	0.578%
46	11.390	1675	1679	1684	rBV5	43708	79709	1.36%	0.109%
47	11.445	1684	1688	1693	rVB3	113804	171717	2.92%	0.234%
48	11.579	1706	1710	1715	rVB4	47329	76483	1.30%	0.104%
49	11.670	1718	1725	1732	rBV	308469	455444	7.75%	0.622%
50	11.768	1733	1741	1745	rBV	389771	508831	8.66%	0.695%
51	11.920	1758	1766	1768	rBV	854245	1192848	20.30%	1.628%
52	11.944	1768	1770	1776	rVB	632890	709530	12.08%	0.969%
53	12.079	1787	1792	1798	rBV	2047047	2632840	44.81%	3.594%
54	12.176	1803	1808	1812	rVV3	51353	93881	1.60%	0.128%
55	12.231	1812	1817	1823	rVB	1273807	1530823	26.05%	2.090%
56	12.298	1823	1828	1834	rBV	1178422	1449555	24.67%	1.979%
57	12.365	1834	1839	1850	rVB	416386	604124	10.28%	0.825%
58	12.463	1850	1855	1861	rBV3	79936	140307	2.39%	0.192%
59	12.670	1878	1889	1892	rBV	3445635	4507280	76.71%	6.153%
60	12.700	1892	1894	1899	rVV	1726226	2057924	35.02%	2.809%
61	12.755	1899	1903	1910	rVV2	3722649	5875994	100.00%	8.021%
62	12.816	1910	1913	1918	rVV	266402	320874	5.46%	0.438%
63	12.896	1918	1926	1931	rVV	1082817	1542733	26.25%	2.106%
64	12.975	1931	1939	1944	rVV	1846234	2659675	45.26%	3.631%
65	13.030	1944	1948	1952	rVV	83311	118241	2.01%	0.161%
66	13.084	1952	1957	1963	rVB2	291237	455874	7.76%	0.622%
67	13.152	1963	1968	1971	rBV	312922	387788	6.60%	0.529%
68	13.194	1971	1975	1978	rVV	347460	451207	7.68%	0.616%
69	13.225	1978	1980	1982	rVV	160848	184755	3.14%	0.252%
70	13.249	1982	1984	1989	rVB	237615	263105	4.48%	0.359%
71	13.310	1989	1994	1995	rBV2	140869	178792	3.04%	0.244%

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Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038MW050-190115

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

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Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X010819W.M
Title : SW846 8260

72	13.347	1995	2000	2006	rVV2	2751377	3909471	66.53%	5.337%
73	13.402	2006	2009	2011	rVV2	223435	287562	4.89%	0.393%
74	13.426	2011	2013	2016	rVV	150554	180617	3.07%	0.247%
75	13.481	2019	2022	2029	rVB	124600	162433	2.76%	0.222%
76	13.560	2032	2035	2038	rBV	62255	81564	1.39%	0.111%
77	13.597	2038	2041	2044	rVV2	132903	185345	3.15%	0.253%
78	13.639	2044	2048	2052	rVV2	459089	661930	11.26%	0.904%
79	13.700	2052	2058	2059	rVV3	270193	517968	8.81%	0.707%
80	13.719	2059	2061	2070	rVB2	365056	496702	8.45%	0.678%
81	13.810	2070	2076	2079	rBV	109404	151971	2.59%	0.207%
82	13.853	2079	2083	2088	rVB2	87876	118045	2.01%	0.161%
83	13.914	2088	2093	2097	rBV	98894	128069	2.18%	0.175%
84	13.981	2097	2104	2109	rBV2	118237	195624	3.33%	0.267%
85	14.029	2109	2112	2115	rBV2	59600	69086	1.18%	0.094%
86	14.273	2147	2152	2158	rBV2	79918	155437	2.65%	0.212%
87	14.535	2190	2195	2201	rBV2	114025	175929	2.99%	0.240%
88	14.840	2238	2245	2250	rBV2	214023	297793	5.07%	0.407%

Sum of corrected areas: 73254433

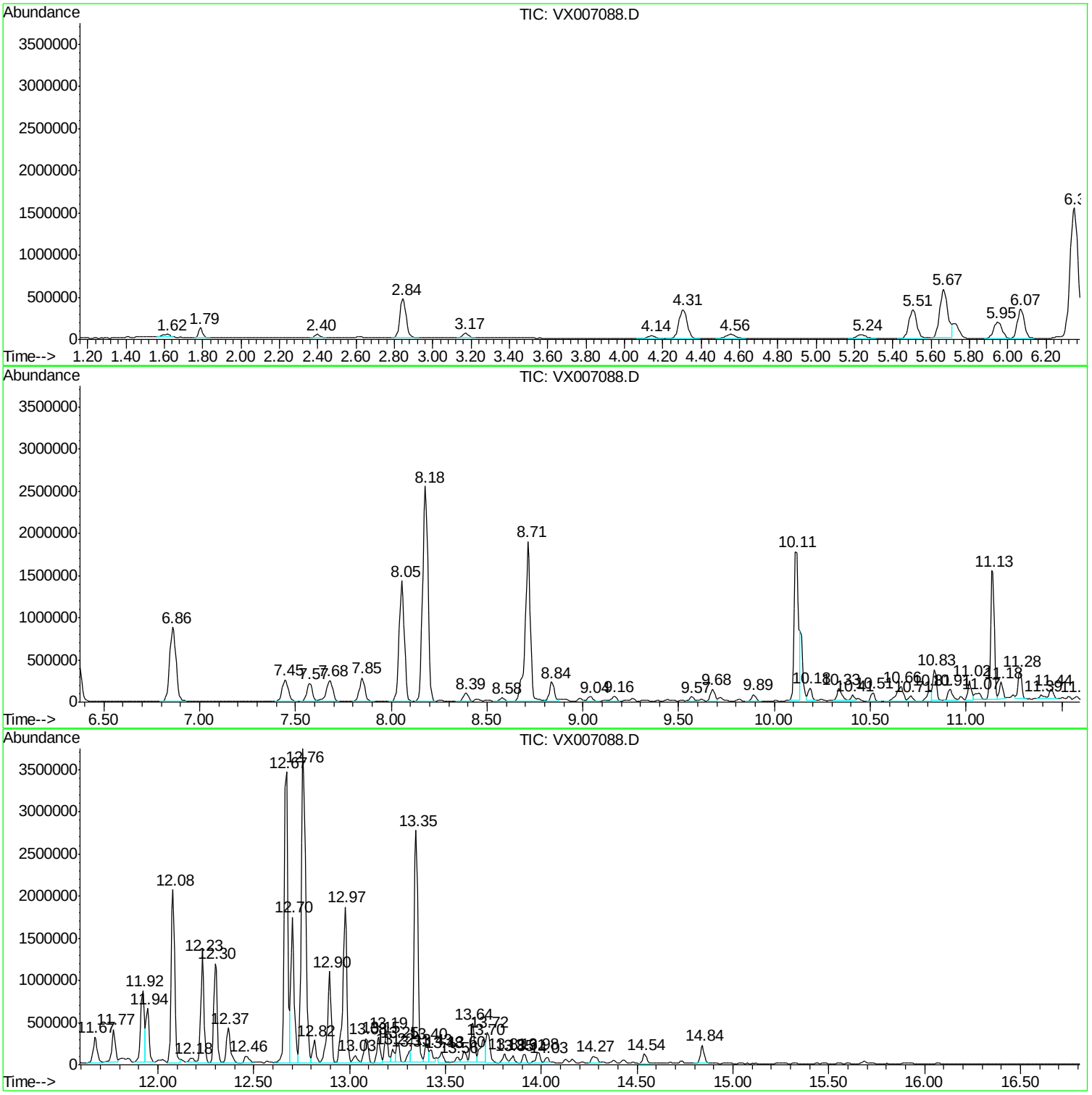
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX011719\
Data File : VX007088.D
Acq On : 17 Jan 2019 14:56
Operator : JC/SP
Sample : K1168-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038MW050-190115

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X010819W.M
Quant Title : SW846 8260

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX011719\
Data File : VX007088.D
Acq On : 17 Jan 2019 14:56
Operator : JC/SP
Sample : K1168-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038MW050-190115

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X010819W.M
Quant Title : SW846 8260

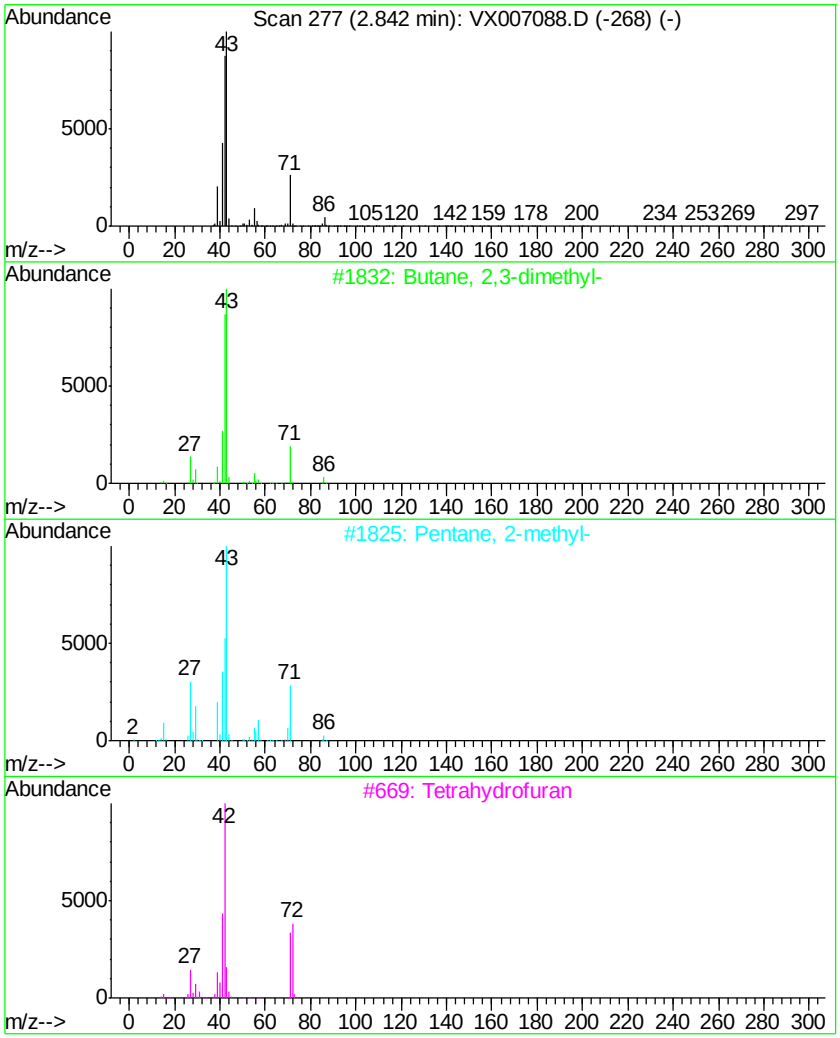
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

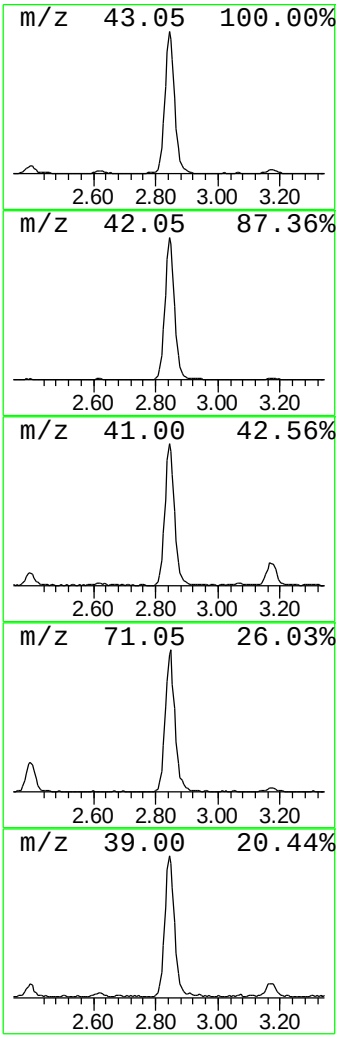
****

Peak Number 1 Butane, 2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.84	30.58 ug/l	1036200	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	86
2	Pentane, 2-methyl-	86	C6H14	000107-83-5	50
3	Tetrahydrofuran	72	C4H8O	000109-99-9	38
4	Pentane	72	C5H12	000109-66-0	38
5	Butane, 2-methyl-	72	C5H12	000078-78-4	36





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX011719\
Data File : VX007088.D
Acq On : 17 Jan 2019 14:56
Operator : JC/SP
Sample : K1168-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038MW050-190115

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X010819W.M
Quant Title : SW846 8260

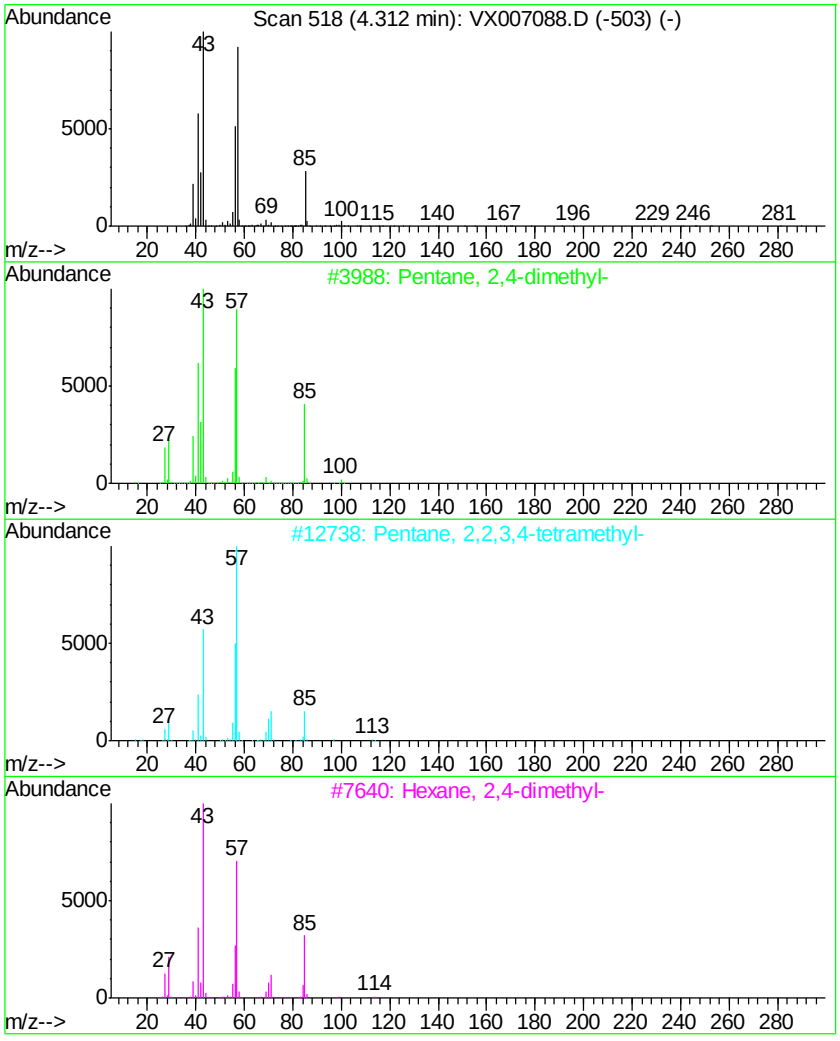
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

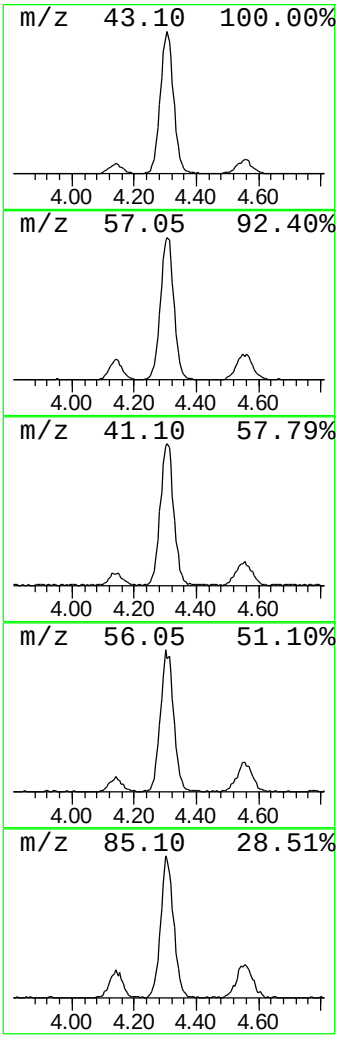
****

Peak Number 2 Pentane, 2,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.31	31.09 ug/l	1053200	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	94
2	Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	72
3	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	72
4	Oxalic acid, butyl isohexyl ester	230	C12H22O4	1000309-32-7	64
5	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	59





Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX011719\
Data File : VX007088.D
Acq On : 17 Jan 2019 14:56
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Sample : K1168-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038MW050-190115

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X010819W.M
Quant Title : SW846 8260

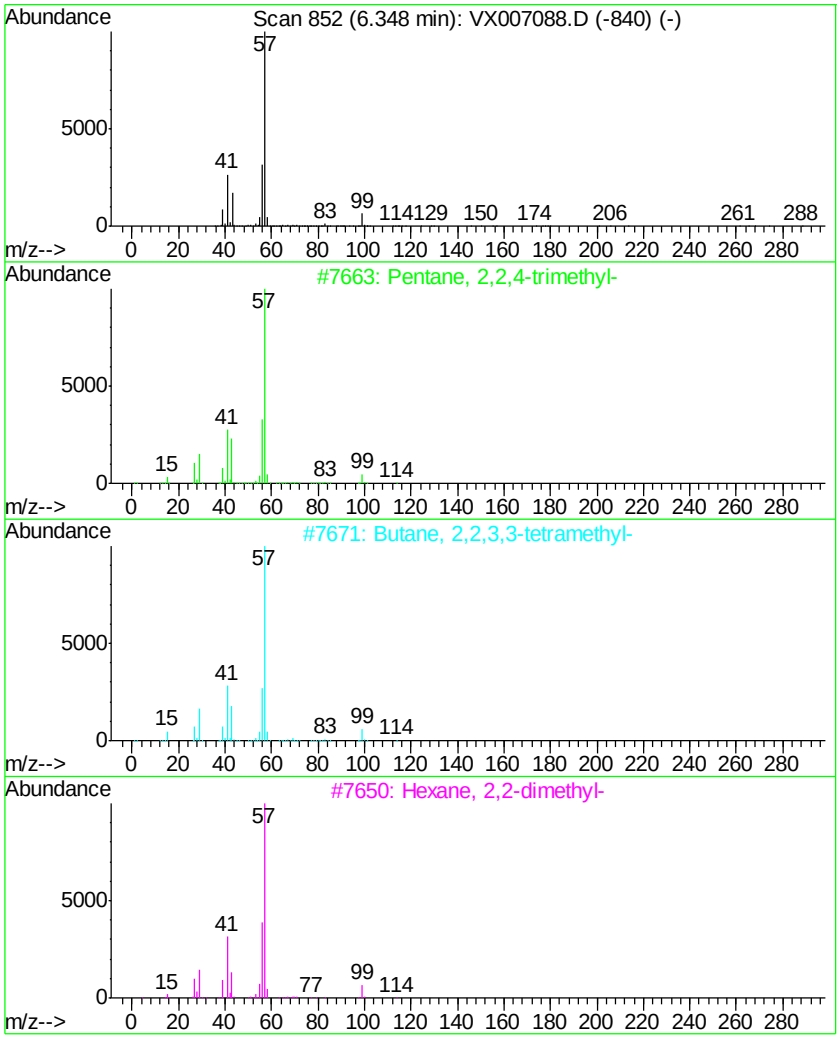
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

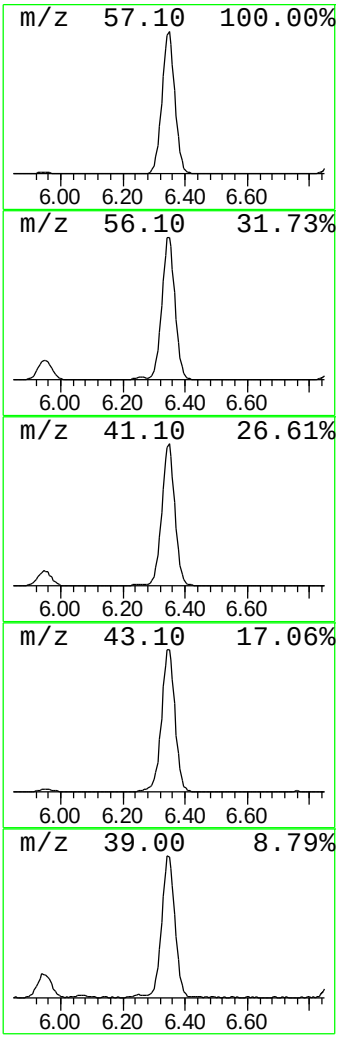
****

Peak Number 3 Pentane, 2,2,4-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.35	112.97 ug/l	4622010	1,4-Difluorobenzene	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	83
2	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	83
3	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	83
4	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	72
5	Hexane, 2,2,5-trimethyl-	128	C9H20	003522-94-9	50





Library Search Compound Report

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ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038MW050-190115

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X010819W.M
Quant Title : SW846 8260

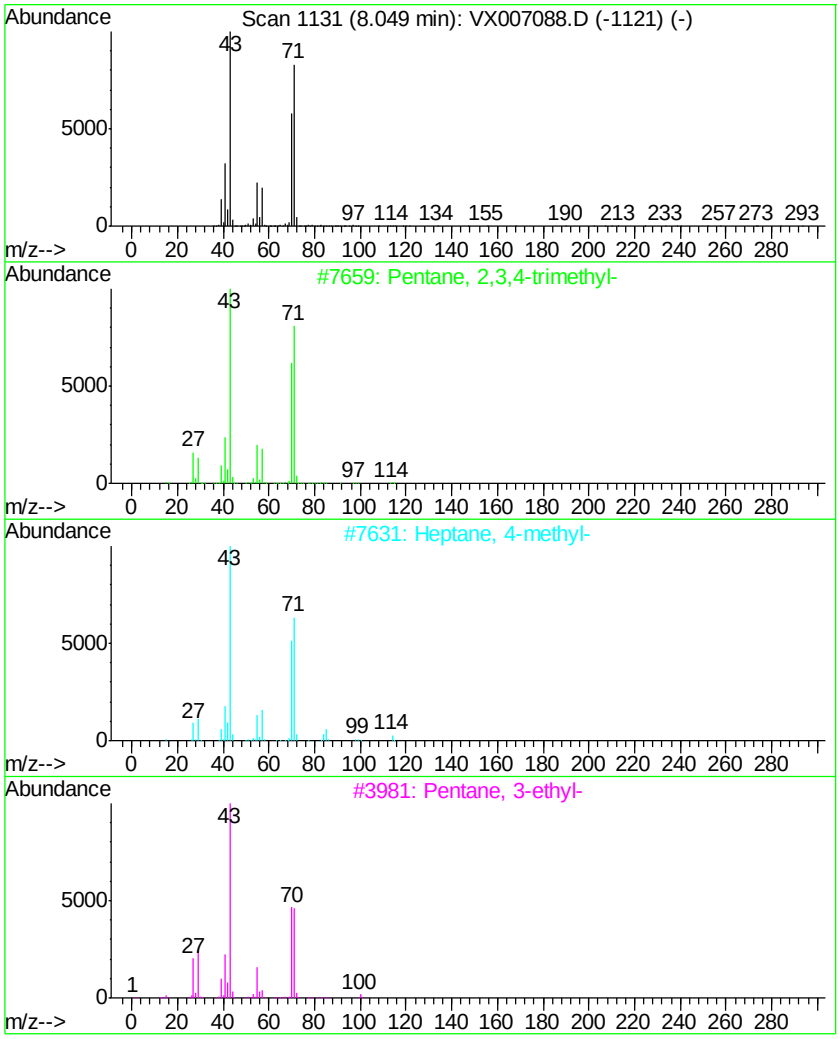
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

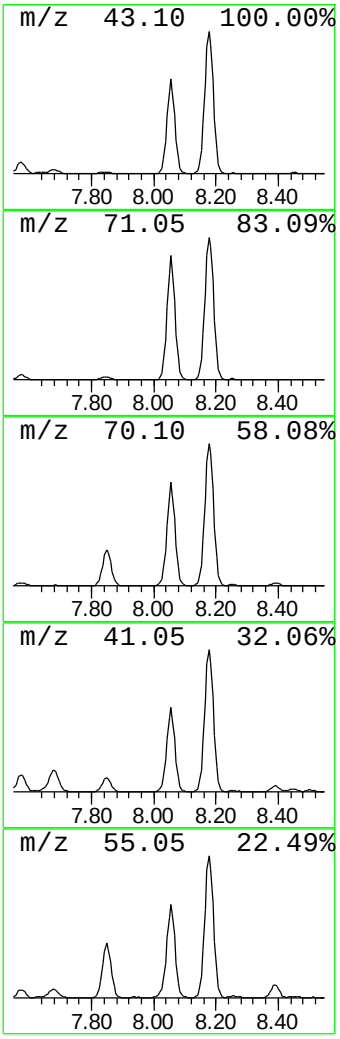
****

Peak Number 4 Pentane, 2,3,4-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.05	66.71 ug/l	2729490	1,4-Difluorobenzene	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2	Heptane, 4-methyl-	114	C8H18	000589-53-7	83
3	Pentane, 3-ethyl-	100	C7H16	000617-78-7	83
4	2,4,4-Trimethyl-1-hexene	126	C9H18	051174-12-0	78
5	Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	74





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Sample : K1168-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038MW050-190115

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X010819W.M
Quant Title : SW846 8260

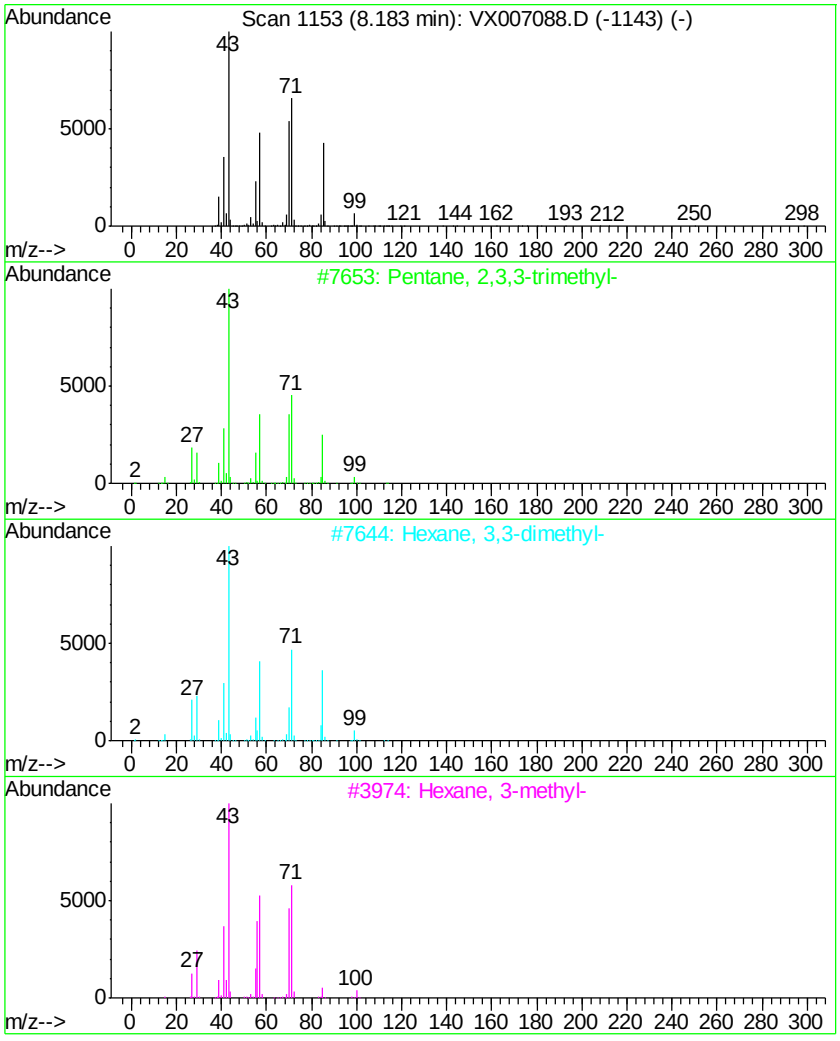
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

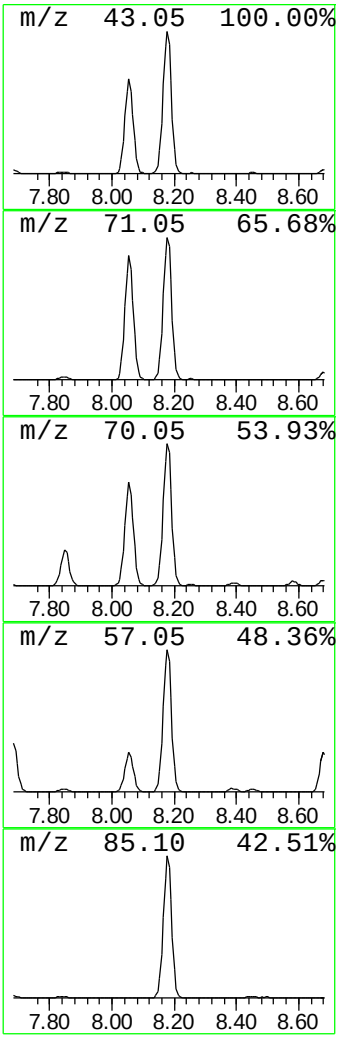
****

Peak Number 5 Pentane, 2,3,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.18	121.11 ug/l	4955010	1,4-Difluorobenzene	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
3	Hexane, 3-methyl-	100	C7H16	000589-34-4	53
4	Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	50
5	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	50





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Instrument : MSVOA_X
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X010819W.M
Quant Title : SW846 8260

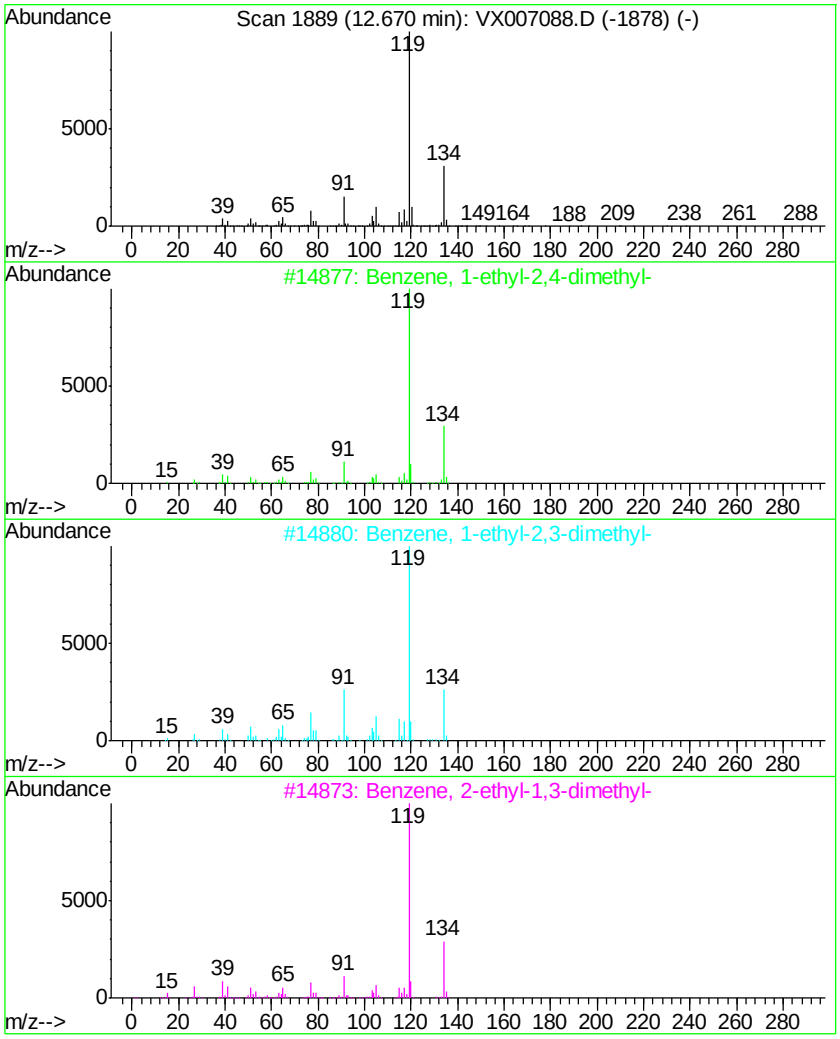
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

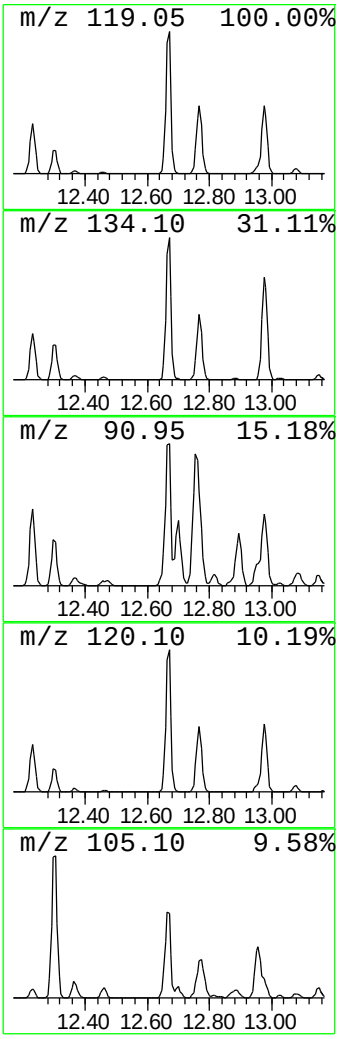
****

Peak Number 6 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	85.60 ug/l	4507280	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	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, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX011719\
Data File : VX007088.D
Acq On : 17 Jan 2019 14:56
Operator : JC/SP
Sample : K1168-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038MW050-190115

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X010819W.M
Quant Title : SW846 8260

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

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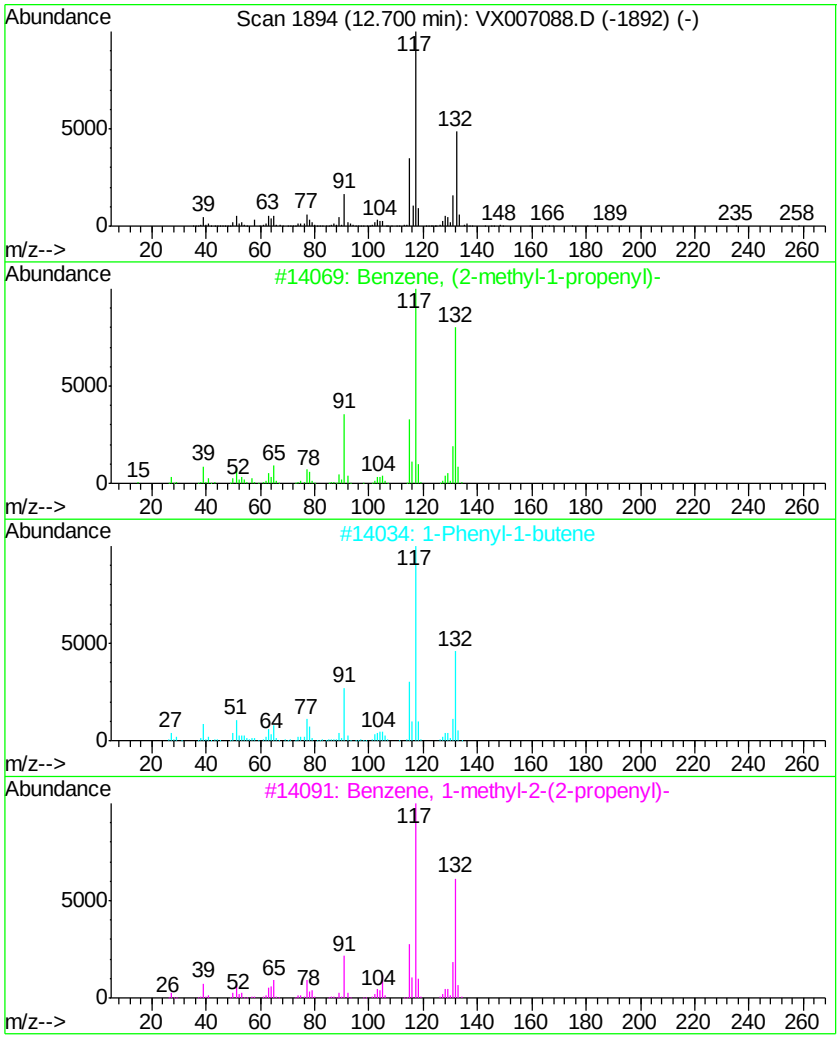
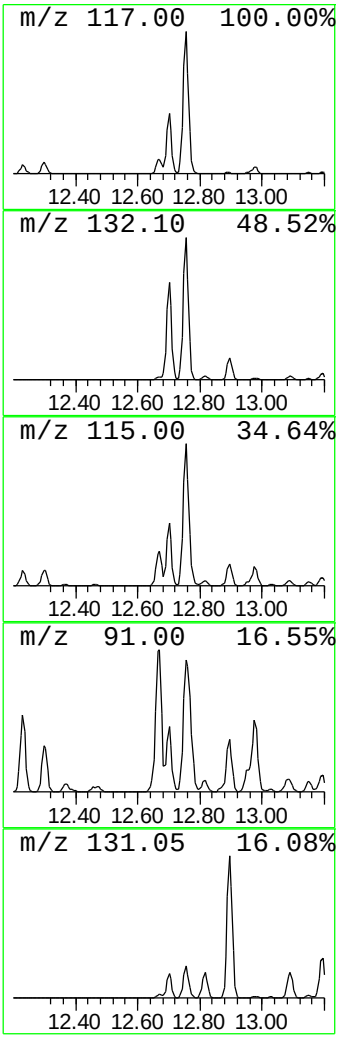
Peak Number 7 Benzene, (2-methyl-1-propen... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	39.08 ug/l	2057920	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	91
2	1-Phenyl-1-butene	132	C10H12	000824-90-8	91
3	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	91
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	91
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	91





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX011719\
Data File : VX007088.D
Acq On : 17 Jan 2019 14:56
Operator : JC/SP
Sample : K1168-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038MW050-190115

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X010819W.M
Quant Title : SW846 8260

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

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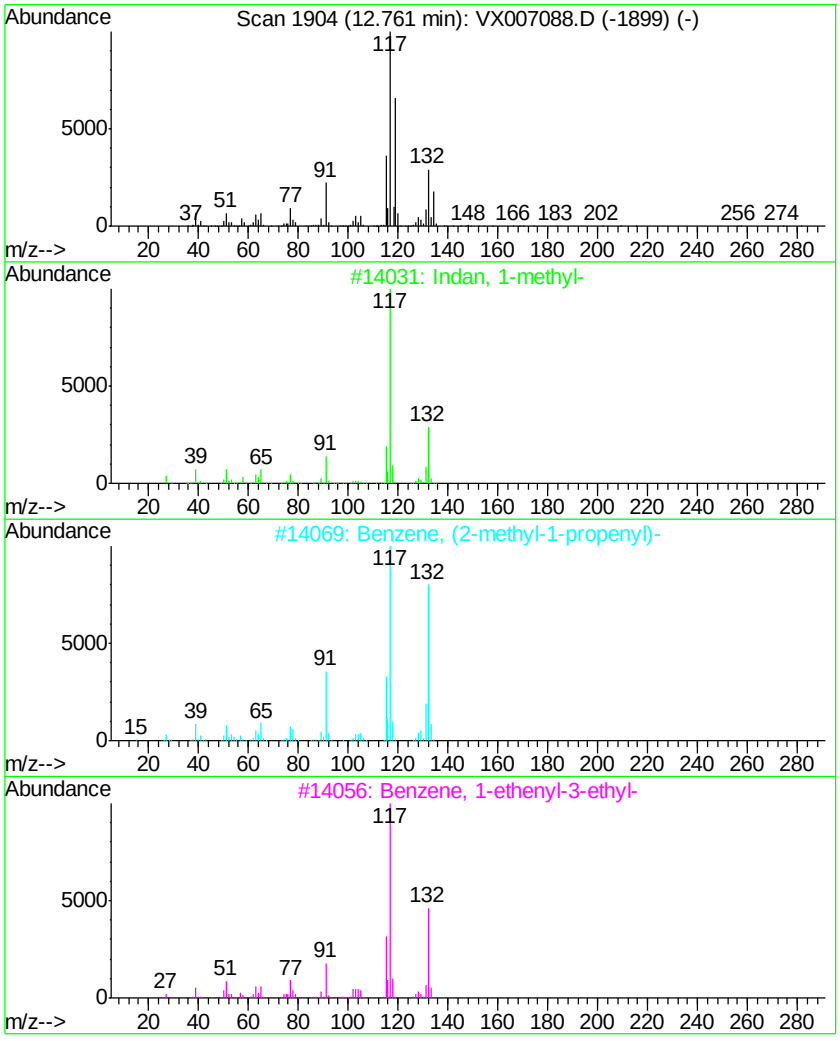
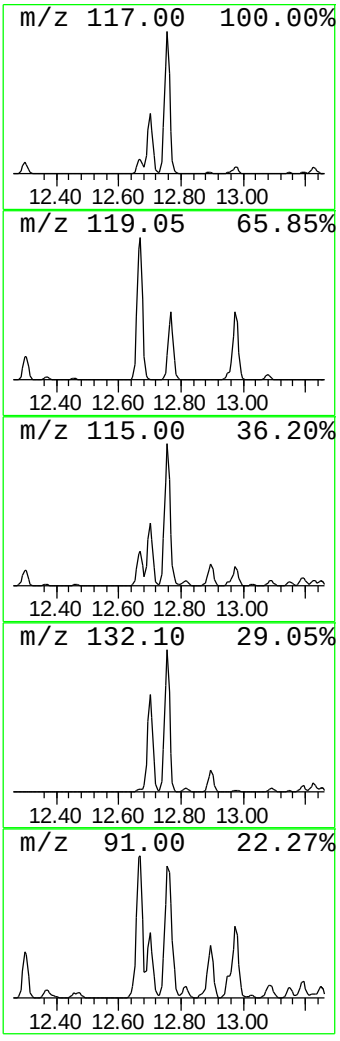
Peak Number 8 Indan, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	111.59 ug/l	5875990	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	94
2	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
4	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	80
5	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	76





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX011719\
Data File : VX007088.D
Acq On : 17 Jan 2019 14:56
Operator : JC/SP
Sample : K1168-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038MW050-190115

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X010819W.M
Quant Title : SW846 8260

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

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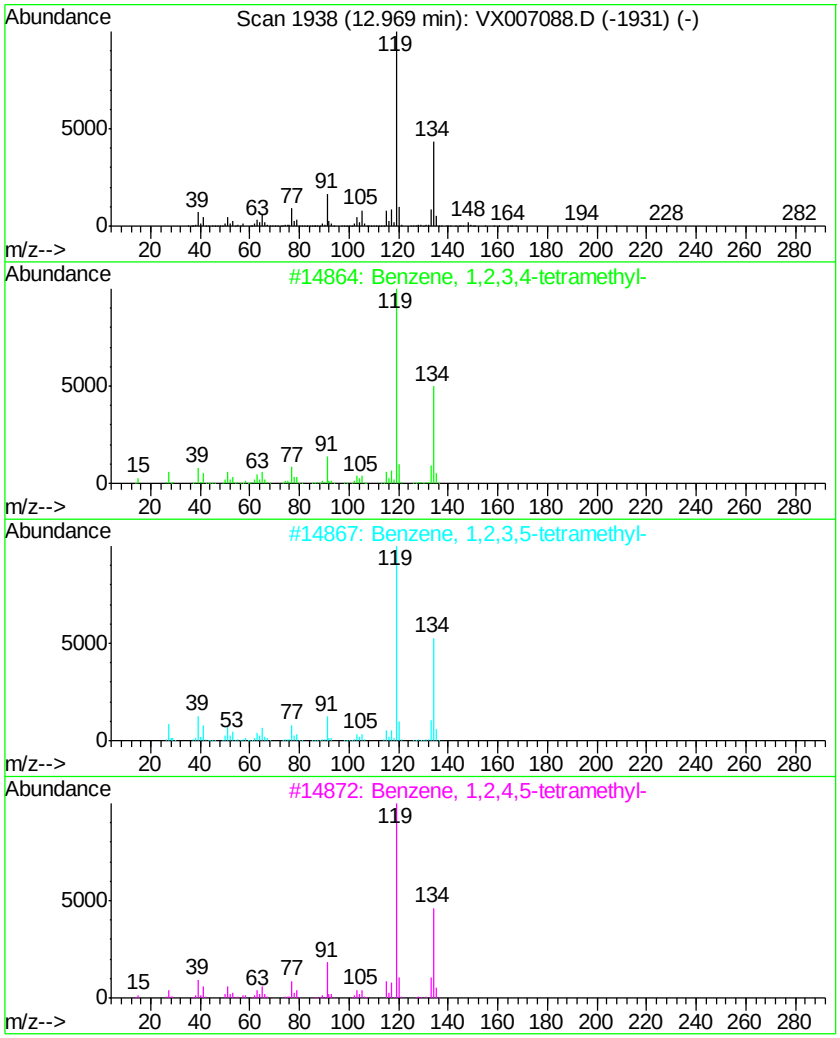
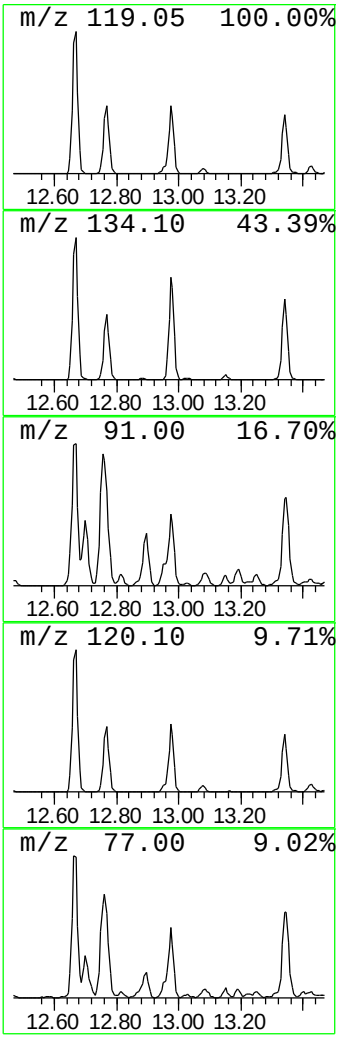
Peak Number 9 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	50.51 ug/l	2659680	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94





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Data File : VX007088.D
Acq On : 17 Jan 2019 14:56
Operator : JC/SP
Sample : K1168-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038MW050-190115

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X010819W.M
Quant Title : SW846 8260

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

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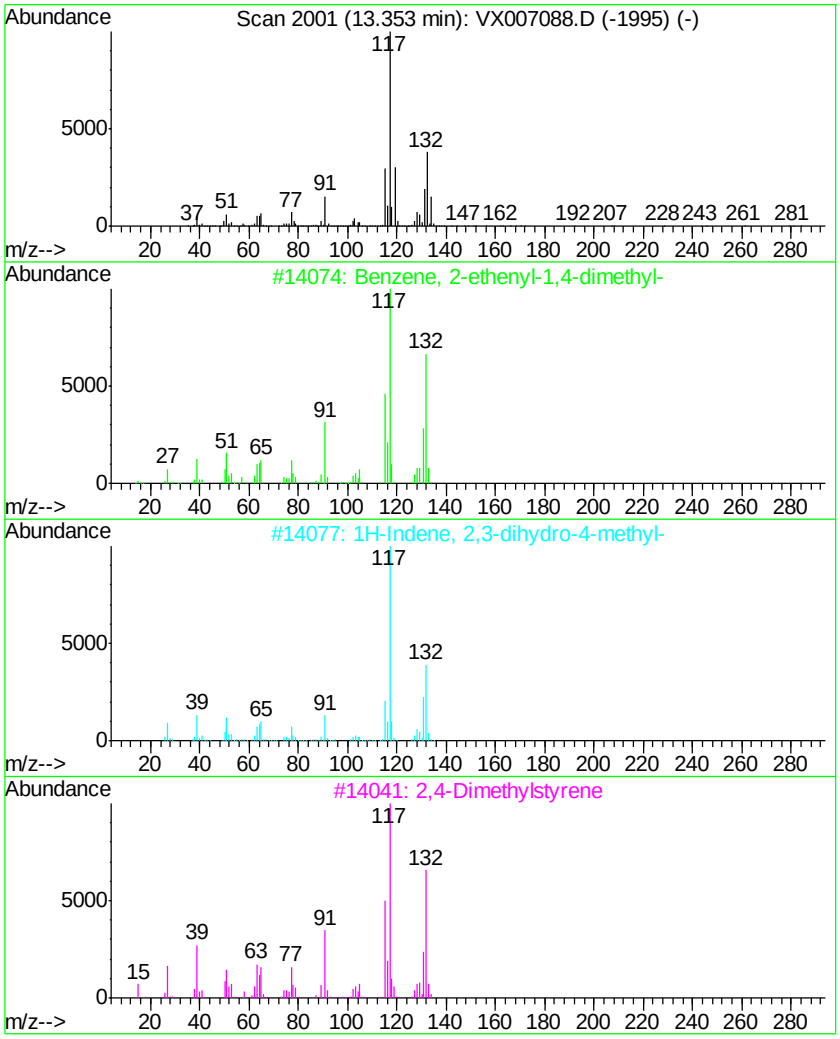
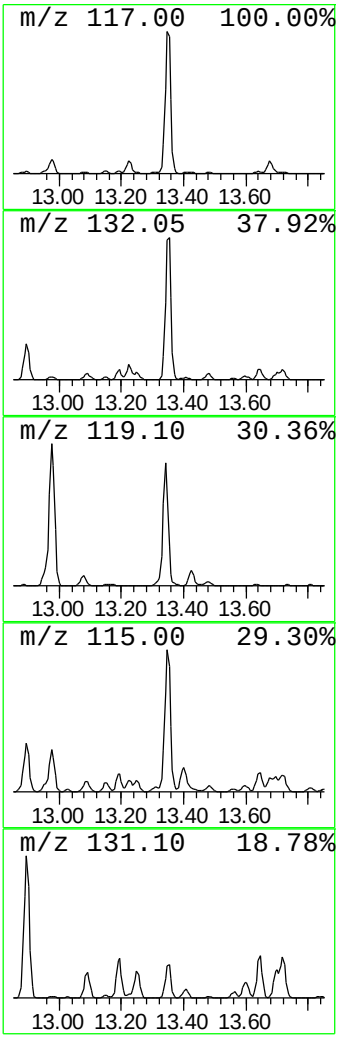
Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	74.24 ug/l	3909470	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	94
2	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
3	2,4-Dimethylstyrene	132	C10H12	002234-20-0	86
4	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	60
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	60





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX011719\
Data File : VX007088.D
Acq On : 17 Jan 2019 14:56
Operator : JC/SP
Sample : K1168-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS038MW050-190115

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X010819W.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
Butane, 2,3-dimet...	2.84	30.6	ug/l	1036200	1	5.67	1694010	50.0
Pentane, 2,4-dime...	4.31	31.1	ug/l	1053200	1	5.67	1694010	50.0
Pentane, 2,2,4-tr...	6.35	113.0	ug/l	4622010	2	6.86	2045700	50.0
Pentane, 2,3,4-tr...	8.05	66.7	ug/l	2729490	2	6.86	2045700	50.0
Pentane, 2,3,3-tr...	8.18	121.1	ug/l	4955010	2	6.86	2045700	50.0
Benzene, 1-ethyl-...	12.67	85.6	ug/l	4507280	4	12.08	2632840	50.0
Benzene, (2-methy...	12.70	39.1	ug/l	2057920	4	12.08	2632840	50.0
Indan, 1-methyl-	12.76	111.6	ug/l	5875990	4	12.08	2632840	50.0
Benzene, 1,2,3,4-...	12.97	50.5	ug/l	2659680	4	12.08	2632840	50.0
Benzene, 2-etheny...	13.35	74.2	ug/l	3909470	4	12.08	2632840	50.0