LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX011719\
Data File : VX007087.D
Acq On : 17 Jan 2019 14:29
Operator : JC/SP
Sample : K1168-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS037MW266-190114

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	1.623	65	77	84	rBV2	34871	92601	3.18%	0.324%
2	1.788	100	104	110	rBV	80380	110202	3.79%	0.386%
3	2.001	135	139	147	rBV2	20578	36040	1.24%	0.126%
4	2.617	234	240	247	rBV	30254	53837	1.85%	0.189%
5	2.843	271	277	281	rBV	98739	206854	7.11%	0.725%
6	2.885	281	284	289	rVV2	65972	144168	4.95%	0.505%
7	2.934	289	292	298	rVB	40775	71253	2.45%	0.250%
8	3.172	321	331	342	rBV	141654	327207	11.24%	1.146%
9	4.306	507	517	523	rBV2	28677	82284	2.83%	0.288%
10	4.403	523	533	543	rVB	121665	360234	12.38%	1.262%
11	4.598	551	565	579	rVB	119290	351507	12.08%	1.231%
12	5.501	701	713	720	rBV2	337092	961255	33.03%	3.367%
13	5.574	721	725	732	rVV2	113481	314652	10.81%	1.102%
14	5.665	732	740	759	rVB	589712	1764314	60.63%	6.180%
15	5.946	774	786	796	rBV5	46797	164711	5.66%	0.577%
16	6.068	796	806	824	rVV	346876	922607	31.70%	3.232%
17	6.257	827	837	843	rVV2	75352	219553	7.54%	0.769%
18	6.348	843	852	862	rVV3	189409	592476	20.36%	2.075%
19	6.452	862	869	879	rVB2	60066	161868	5.56%	0.567%
20	6.860	924	936	949	rBV	874200	2055087	70.62%	7.198%
21	7.214	984	994	1004	rBV	208613	446664	15.35%	1.565%
22	7.458	1023	1034	1043	rVB2	130897	333997	11.48%	1.170%
23	7.574	1047	1053	1059	rVV4	14588	30587	1.05%	0.107%
24	7.641	1059	1064	1067	rVV6	14964	34272	1.18%	0.120%
25	7.683	1067	1071	1075	rVV3	18464	35071	1.21%	0.123%
26	7.738	1075	1080	1087	rVB2	16093	31410	1.08%	0.110%
27	7.848	1090	1098	1104	rBV3	23266	43568	1.50%	0.153%
28	8.055	1121	1132	1141	rBV2	97402	241433	8.30%	0.846%
29	8.177	1144	1152	1160	rBV	186532	359715	12.36%	1.260%
30	8.665	1225	1232	1234	rBV2	51168	86331	2.97%	0.302%
31	8.714	1234	1240	1255	rVV	1822207	2903677	99.78%	10.171%
32	8.835	1255	1260	1265	rVB3	20868	32013	1.10%	0.112%
33	9.043	1288	1294	1301	rVB2	41848	73607	2.53%	0.258%
34	9.165	1307	1314	1319	rBV2	21707	38519	1.32%	0.135%

LSC Area Percent Report

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Sample : K1168-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS037MW266-190114

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	9.335	1335	1342	1351	rBV	564166	829730	28.51%	2.906%
36	9.622	1384	1389	1394	rVB	46105	64922	2.23%	0.227%
37	9.677	1394	1398	1402	rBV2	19732	30123	1.04%	0.106%
38	10.110	1464	1469	1473	rBV	1799633	2910038	100.00%	10.193%
39	10.183	1479	1481	1486	rVB2	38140	47343	1.63%	0.166%
40	10.335	1500	1506	1509	rBV	35022	58118	2.00%	0.204%
41	10.372	1509	1512	1516	rVB3	25218	34939	1.20%	0.122%
42	10.512	1529	1535	1542	rBV	29699	43965	1.51%	0.154%
43	10.658	1544	1559	1564	rBV4	64378	174717	6.00%	0.612%
44	10.707	1564	1567	1574	rVB2	19997	30965	1.06%	0.108%
45	10.811	1576	1584	1585	rBV3	33550	64576	2.22%	0.226%
46	10.835	1585	1588	1593	rVB	77646	101969	3.50%	0.357%
47	10.914	1593	1601	1608	rBV5	65743	120488	4.14%	0.422%
48	11.018	1614	1618	1621	rBV	77948	99532	3.42%	0.349%
49	11.061	1621	1625	1628	rVV6	19164	33994	1.17%	0.119%
50	11.134	1632	1637	1642	rVV	1565929	1998512	68.68%	7.000%
51	11.183	1642	1645	1649	rVB2	58559	77185	2.65%	0.270%
52	11.280	1657	1661	1665	rVB	81980	102830	3.53%	0.360%
53	11.359	1670	1674	1678	rBV	63971	79221	2.72%	0.277%
54	11.445	1684	1688	1692	rBV4	33306	54363	1.87%	0.190%
55	11.573	1706	1709	1714	rVB4	18808	29269	1.01%	0.103%
56	11.670	1719	1725	1731	rVV2	115820	175260	6.02%	0.614%
57	11.768	1737	1741	1744	rBV	72138	83599	2.87%	0.293%
58	11.914	1758	1765	1768	rBV2	93702	165031	5.67%	0.578%
59	11.945	1768	1770	1775	rVB	101203	108532	3.73%	0.380%
60	12.079	1786	1792	1802	rVV	2022516	2613850	89.82%	9.156%
61	12.170	1804	1807	1812	rVV3	21749	37883	1.30%	0.133%
62	12.231	1812	1817	1823	rVV	257239	320770	11.02%	1.124%
63	12.298	1823	1828	1835	rVV	326826	407147	13.99%	1.426%
64	12.365	1835	1839	1850	rVB2	155208	268293	9.22%	0.940%
65	12.457	1850	1854	1860	rBV2	72332	106336	3.65%	0.372%
66	12.585	1869	1875	1877	rBV4	22011	35954	1.24%	0.126%
67	12.664	1884	1888	1891	rBV	168947	203938	7.01%	0.714%
68	12.701	1891	1894	1899	rVV2	212149	335349	11.52%	1.175%
69	12.762	1899	1904	1910	rVV3	467437	832675	28.61%	2.917%
70	12.817	1910	1913	1917	rVV2	50079	62397	2.14%	0.219%
71	12.871	1917	1922	1923	rVV	79988	100718	3.46%	0.353%

LSC Area Percent Report

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ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS037MW266-190114

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

72	12.896	1923	1926	1931	rVB	237518	286164	9.83%	1.002%
73	12.975	1936	1939	1943	rVB	184654	228100	7.84%	0.799%
74	13.018	1943	1946	1952	rVB2	67505	99290	3.41%	0.348%
75	13.085	1952	1957	1963	rVB3	87209	127073	4.37%	0.445%
76	13.152	1964	1968	1971	rBV	28908	36889	1.27%	0.129%
77	13.194	1971	1975	1977	rBV2	74764	88600	3.04%	0.310%
78	13.225	1977	1980	1982	rVV	36710	43464	1.49%	0.152%
79	13.249	1982	1984	1988	rVB	56047	62399	2.14%	0.219%
80	13.304	1988	1993	1996	rBV4	34183	56498	1.94%	0.198%
81	13.347	1996	2000	2005	rVB2	309290	416465	14.31%	1.459%
82	13.481	2015	2022	2030	rVB3	69488	130001	4.47%	0.455%
83	13.560	2030	2035	2038	rBV2	21381	32781	1.13%	0.115%
84	13.597	2038	2041	2044	rVV	34792	46179	1.59%	0.162%
85	13.640	2044	2048	2052	rVV2	82958	115223	3.96%	0.404%
86	13.694	2052	2057	2059	rVV2	57454	105136	3.61%	0.368%
87	13.719	2059	2061	2068	rVB2	57887	71264	2.45%	0.250%
88	13.810	2071	2076	2078	rBV	26521	38854	1.34%	0.136%
89	13.835	2078	2080	2088	rVB3	18218	36364	1.25%	0.127%

Sum of corrected areas: 28548849

LSC Report - Integrated Chromatogram

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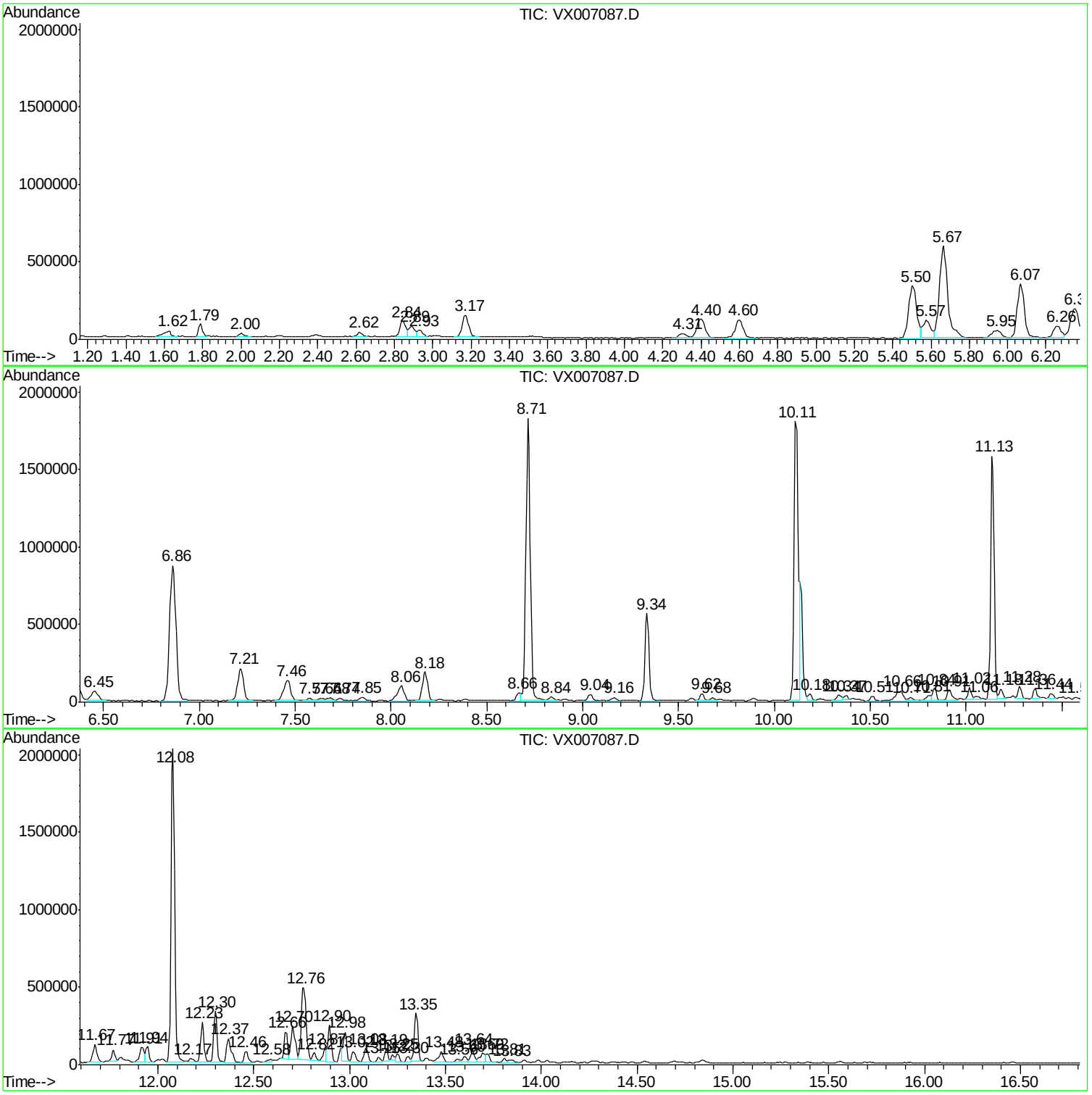
Operator : JC/SP
Sample : K1168-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS037MW266-190114

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 : VX007087.D
Acq On : 17 Jan 2019 14:29
Operator : JC/SP
Sample : K1168-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS037MW266-190114

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

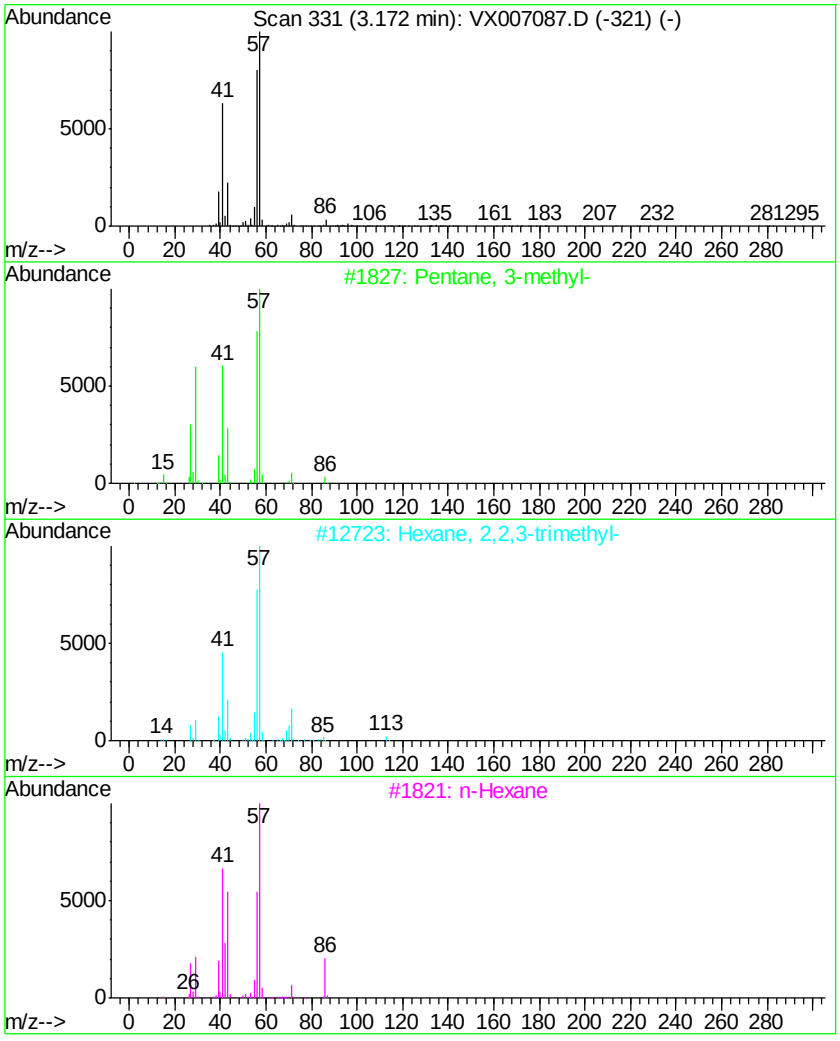
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

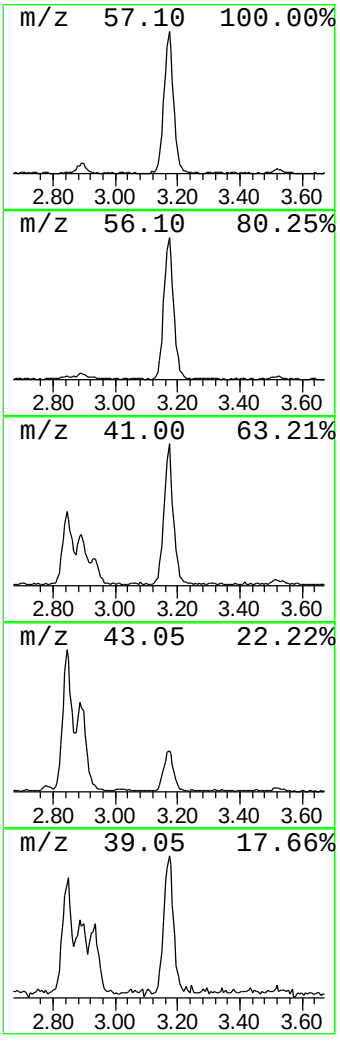
****

Peak Number 1 Pentane, 3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.17	9.27 ug/l	327207	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	86
2	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	72
3	n-Hexane	86	C6H14	000110-54-3	50
4	Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	45
5	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	40





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX011719\
Data File : VX007087.D
Acq On : 17 Jan 2019 14:29
Operator : JC/SP
Sample : K1168-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS037MW266-190114

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

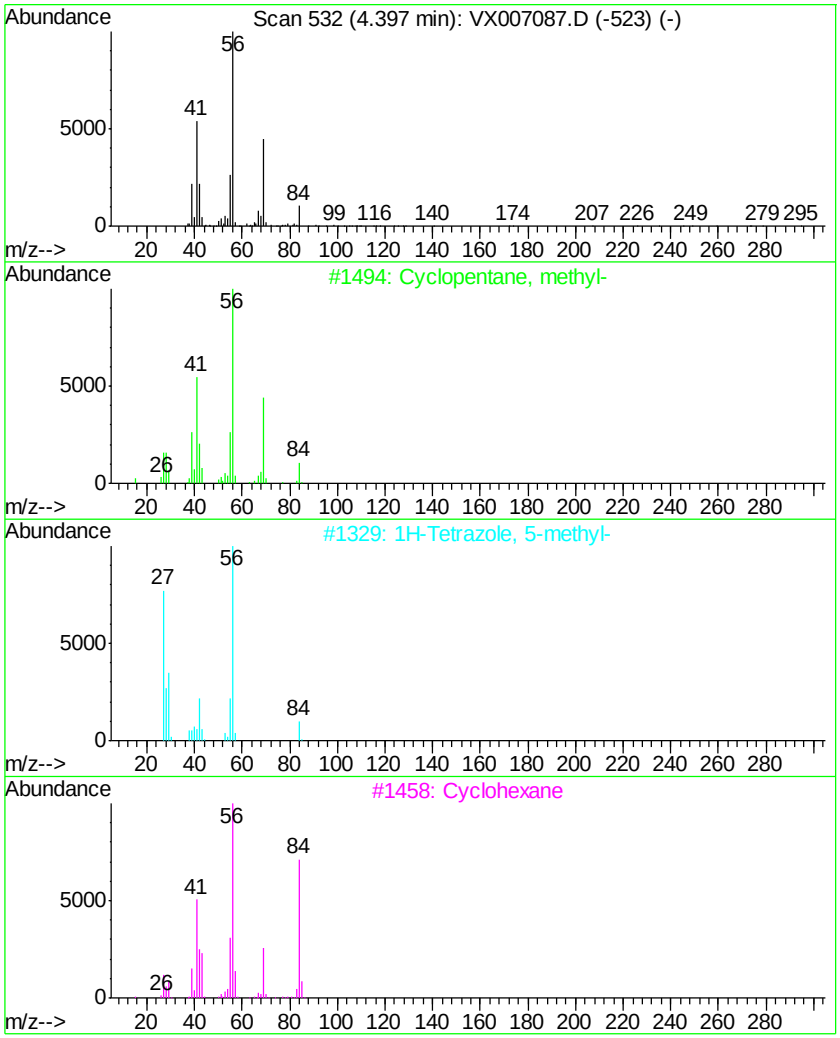
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

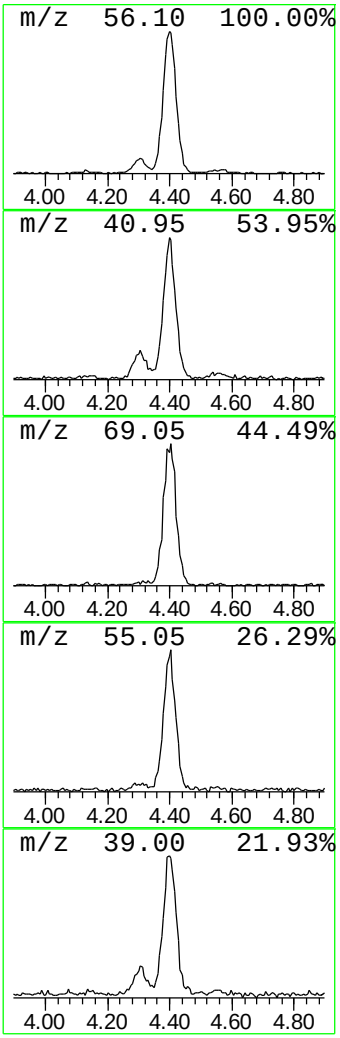
****

Peak Number 2 Cyclopentane, methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.40	10.21 ug/l	360234	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
3	Cyclohexane	84	C6H12	000110-82-7	64
4	1-Butanol	74	C4H10O	000071-36-3	53
5	Cyclopropane, propyl-	84	C6H12	002415-72-7	50





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX011719\
Data File : VX007087.D
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS037MW266-190114

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

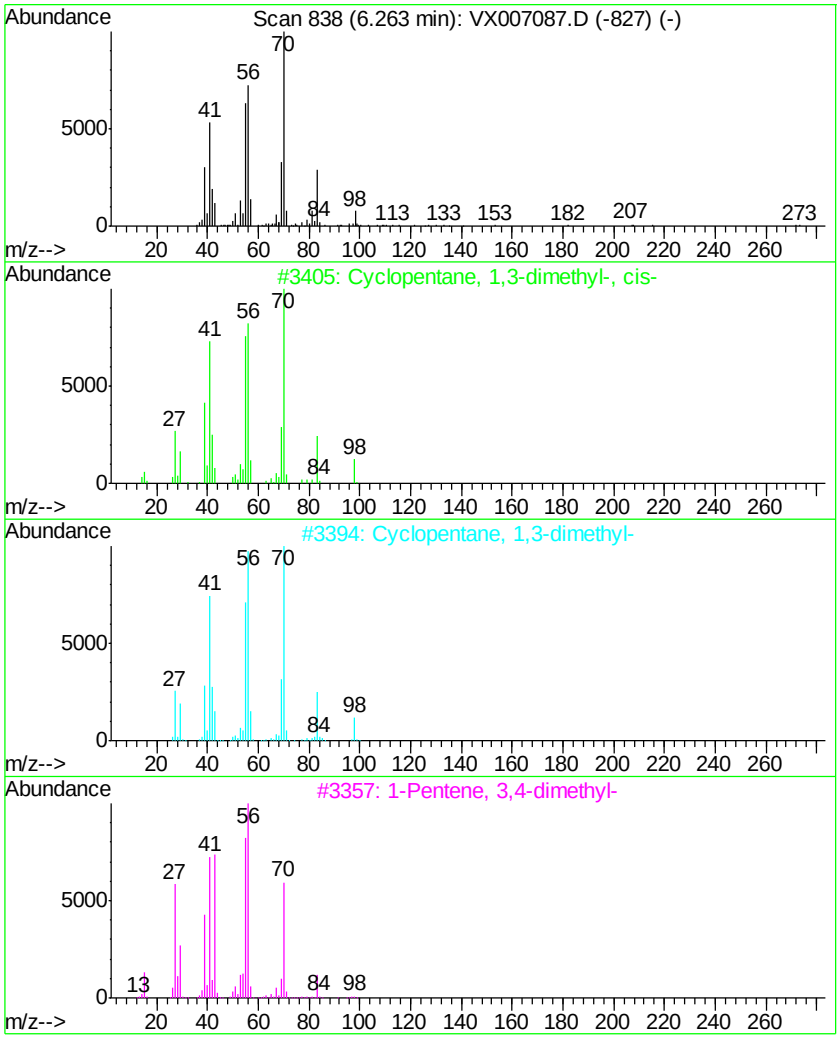
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

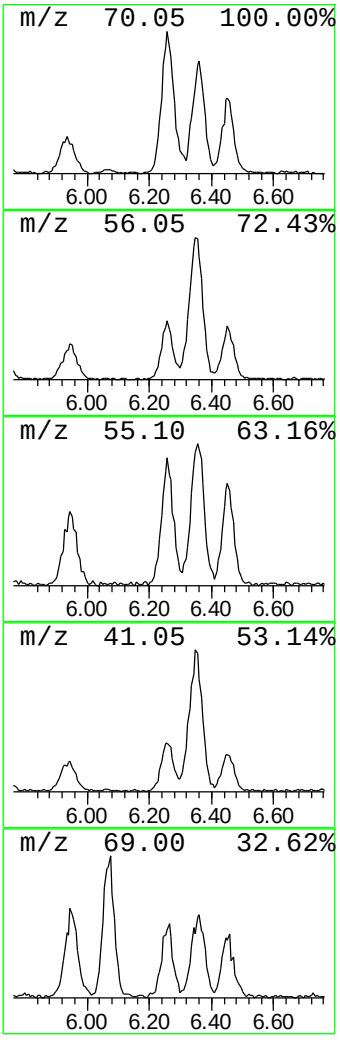
****

Peak Number 3 Cyclopentane, 1,3-dimethyl-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.26	6.22 ug/l	219553	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	91
2	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	90
3	1-Pentene, 3,4-dimethyl-	98	C7H14	007385-78-6	72
4	3-Methylcyclohexylamine,c&t	113	C7H15N	006850-35-7	64
5	Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	52





Library Search Compound Report

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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS037MW266-190114

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

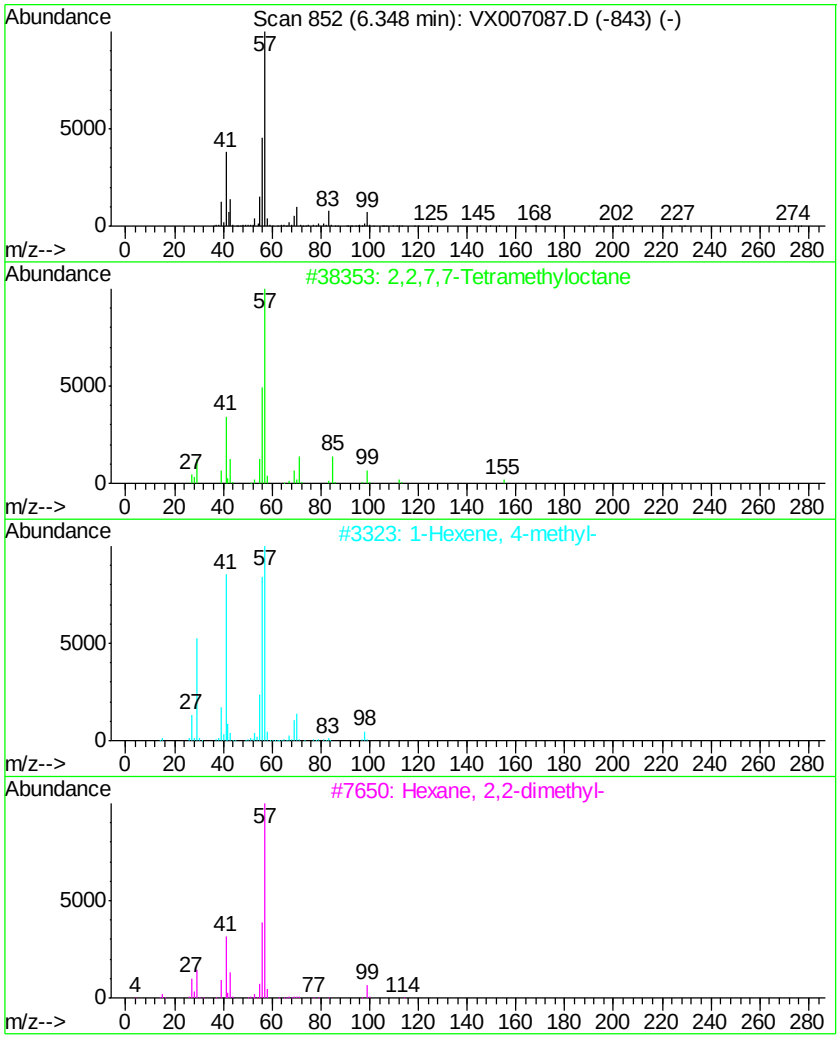
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

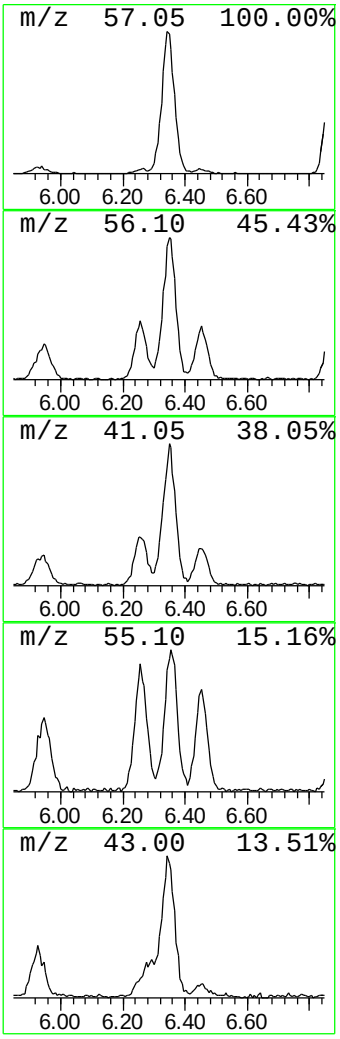
****

Peak Number 4 2,2,7,7-Tetramethyloctane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.35	14.41 ug/l	592476	1,4-Difluorobenzene	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2,7,7-Tetramethyloctane	170	C12H26	001071-31-4	72
2	1-Hexene, 4-methyl-	98	C7H14	003769-23-1	64
3	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	42
4	Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	39
5	Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	39





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Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS037MW266-190114

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

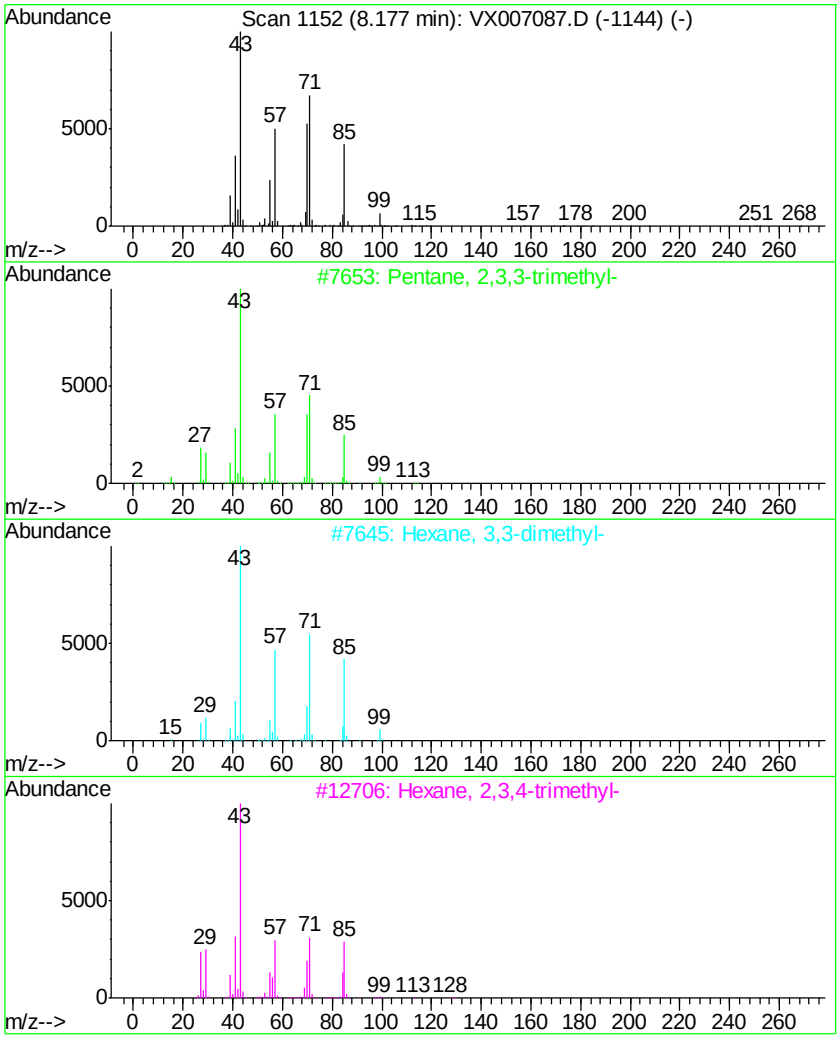
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

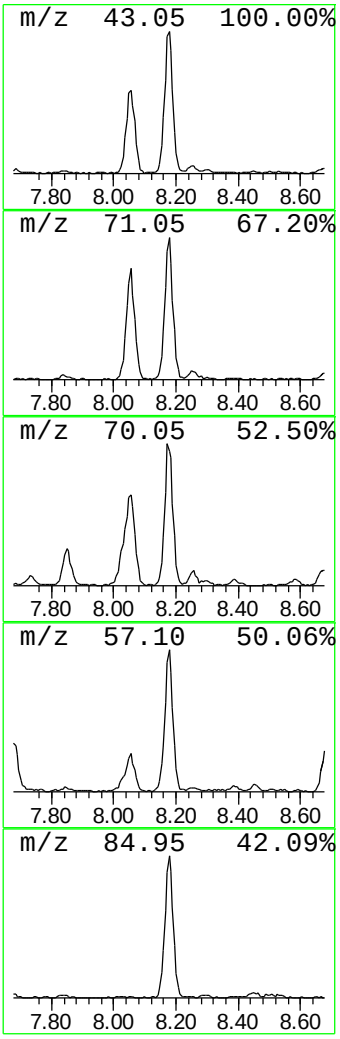
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.18	8.75 ug/l	359715	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, 2,3,4-trimethyl-	128	C9H20	000921-47-1	64
4	Undecane, 4,8-dimethyl-	184	C13H28	017301-33-6	59
5	Heptane, 4-methyl-	114	C8H18	000589-53-7	53





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ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS037MW266-190114

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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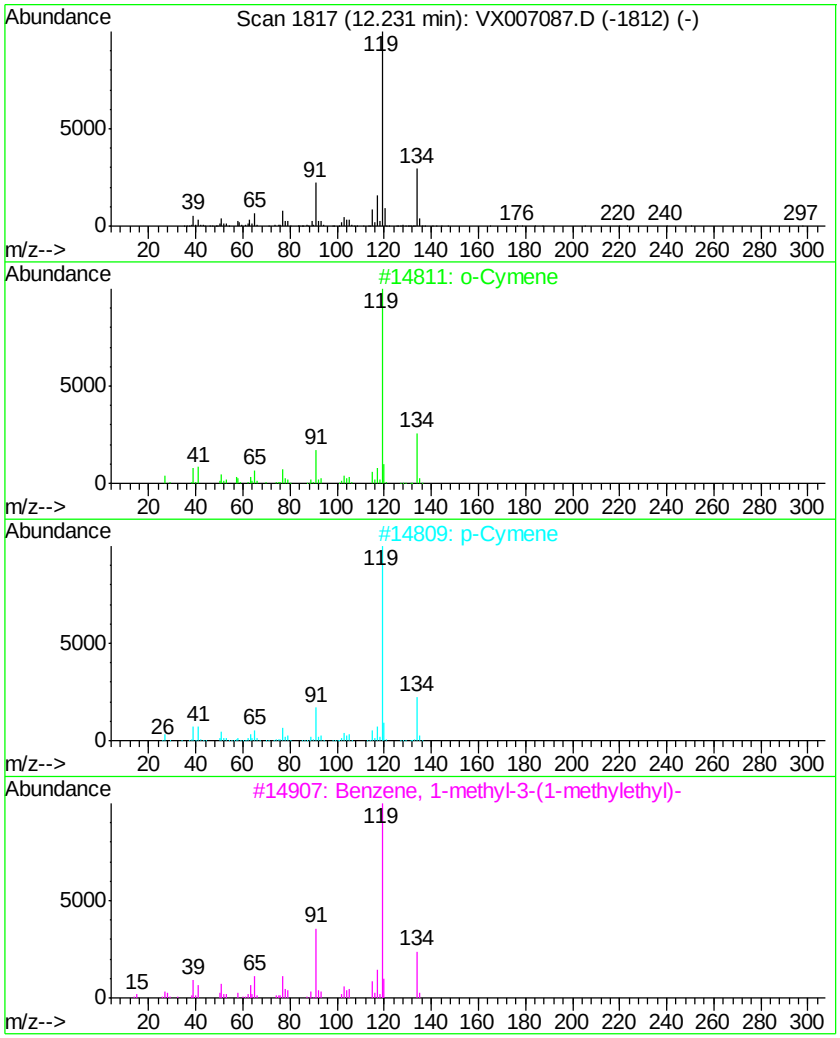
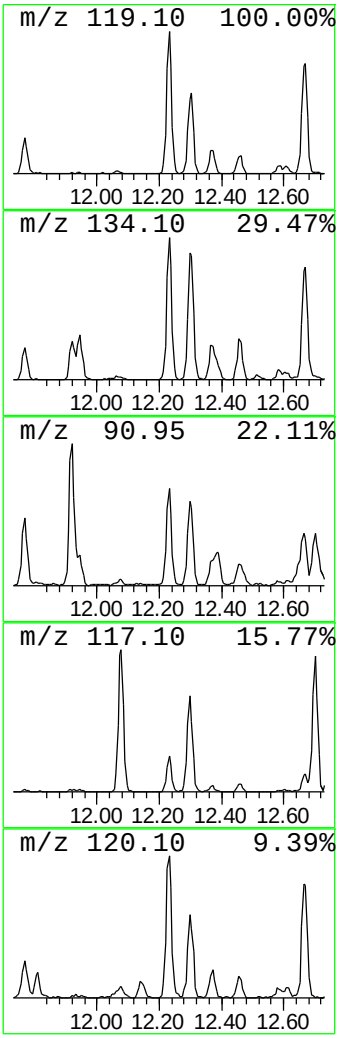
Peak Number 6 o-Cymene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.23	6.14 ug/l	320770	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	97
2	p-Cymene	134	C10H14	000099-87-6	97
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	91
5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX011719\
Data File : VX007087.D
Acq On : 17 Jan 2019 14:29
Operator : JC/SP
Sample : K1168-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS037MW266-190114

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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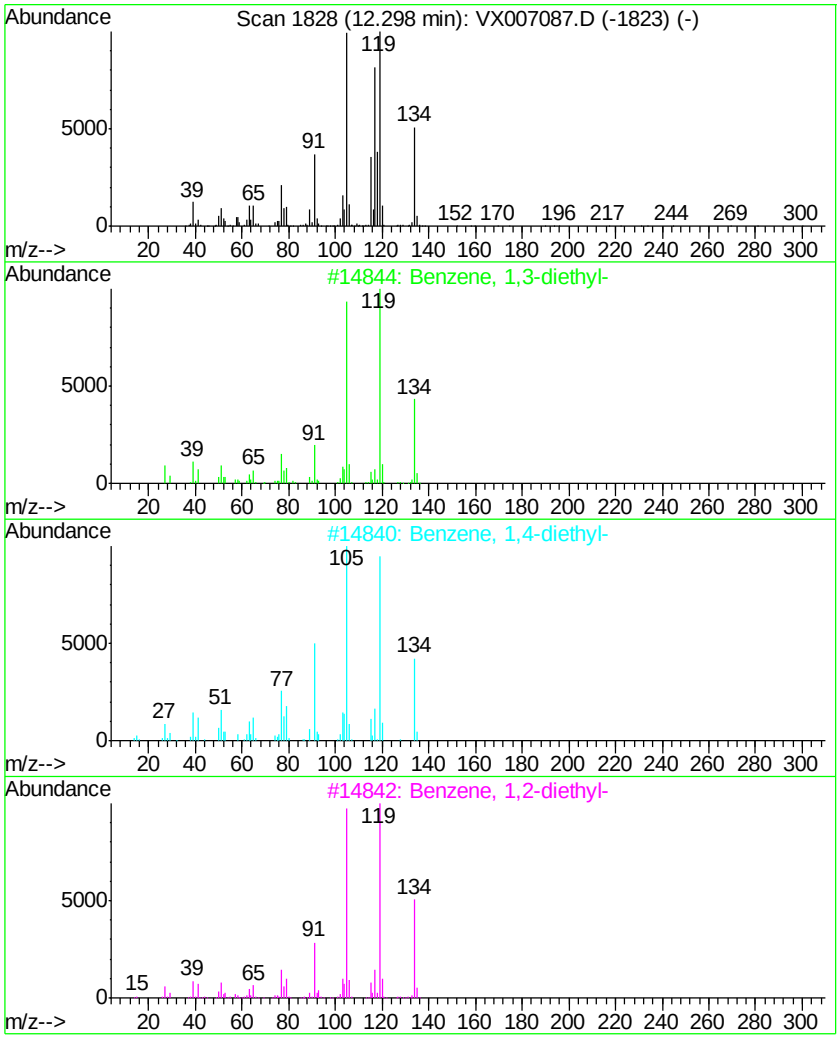
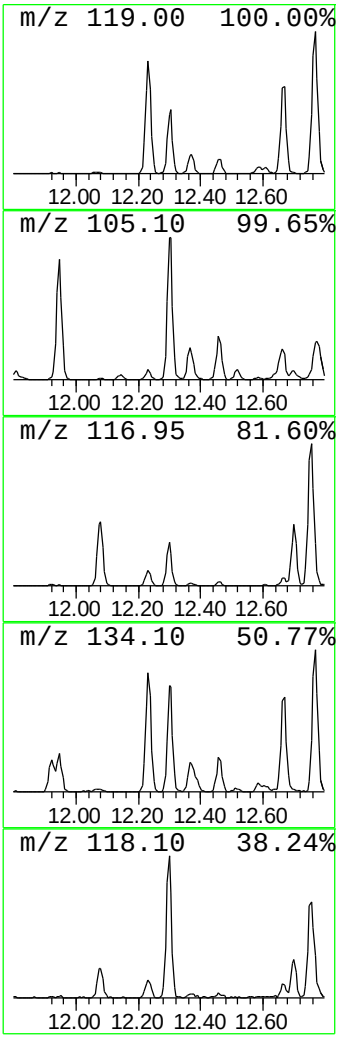
Peak Number 7 Benzene, 1,3-diethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	7.79 ug/l	407147	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	90
2	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	60
3	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	60
4	Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	51
5	1,3-Methanopentalene, 1,2,3,5-te...	118	C9H10	128600-88-4	42





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX011719\
Data File : VX007087.D
Acq On : 17 Jan 2019 14:29
Operator : JC/SP
Sample : K1168-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS037MW266-190114

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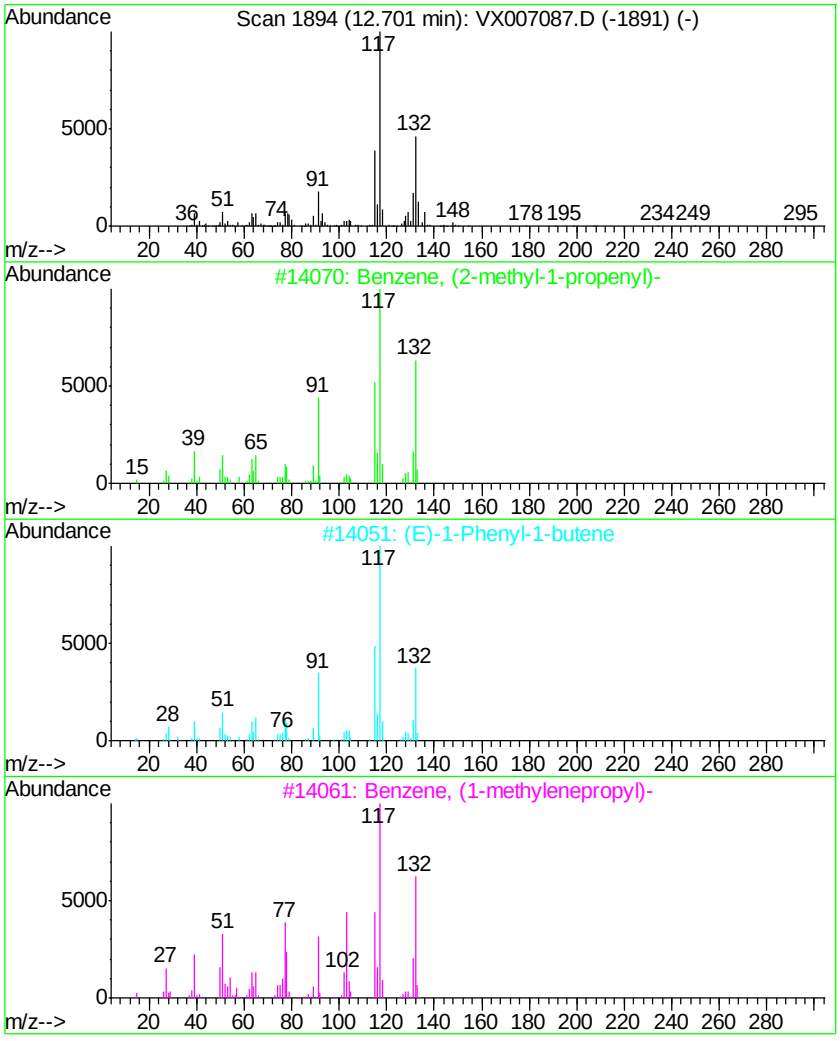
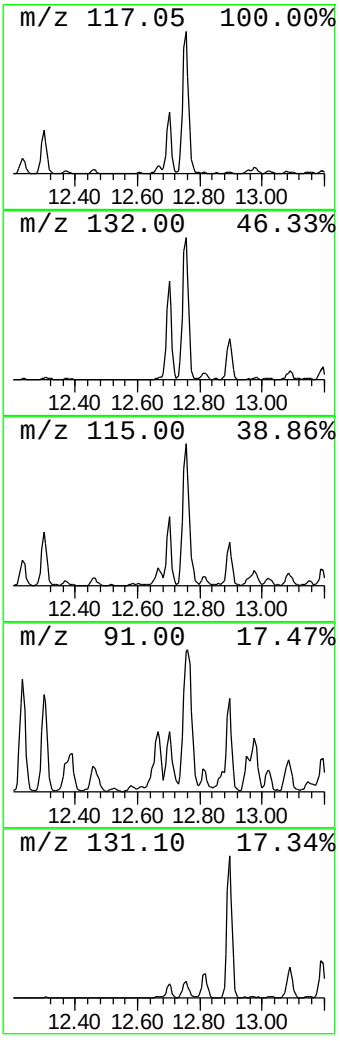
Peak Number 8 Benzene, (2-methyl-1-propen... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	6.41 ug/l	335349	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	90
2	(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	90
3	Benzene, (1-methylenepropyl)-	132	C10H12	002039-93-2	90
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	87





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX011719\
Data File : VX007087.D
Acq On : 17 Jan 2019 14:29
Operator : JC/SP
Sample : K1168-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS037MW266-190114

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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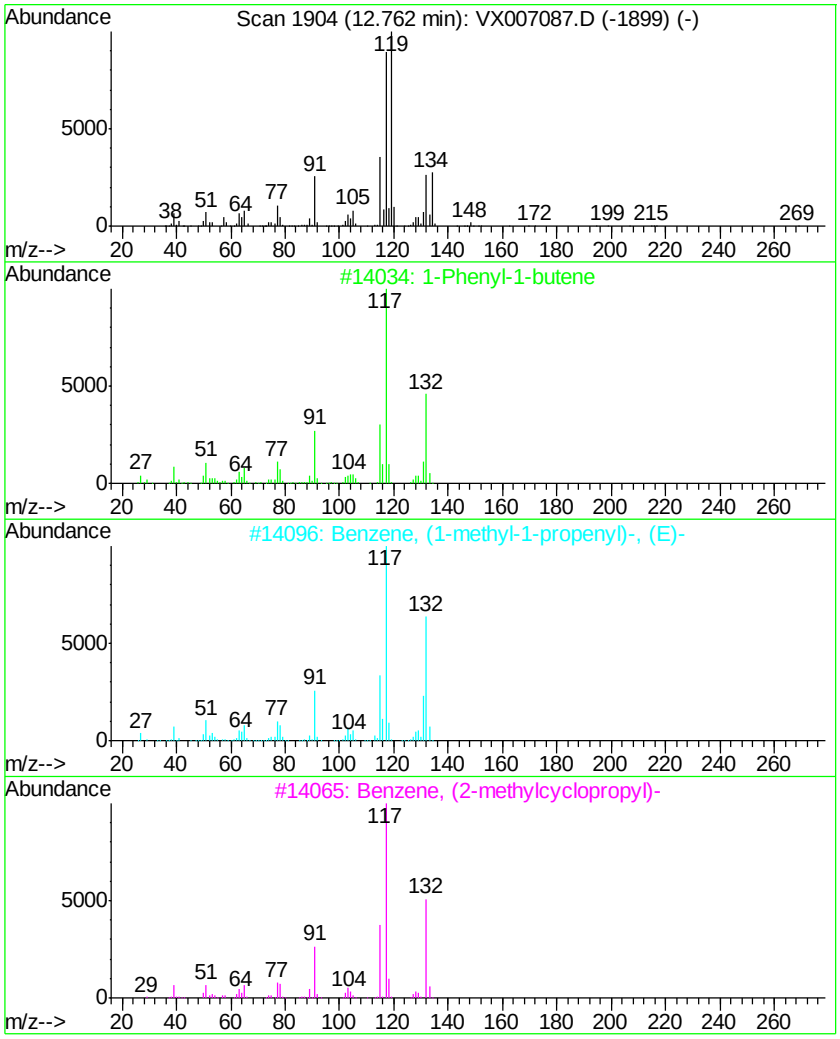
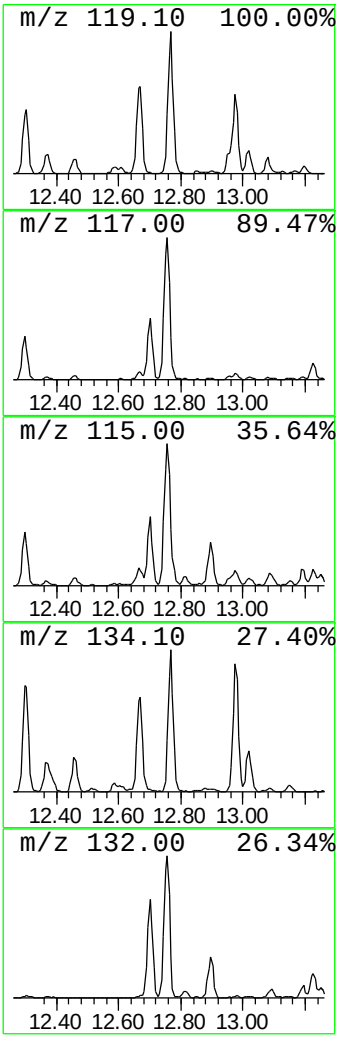
Peak Number 9 1-Phenyl-1-butene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	15.93 ug/l	832675	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	55
3	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	50
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	50
5	(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	49





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX011719\
Data File : VX007087.D
Acq On : 17 Jan 2019 14:29
Operator : JC/SP
Sample : K1168-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS037MW266-190114

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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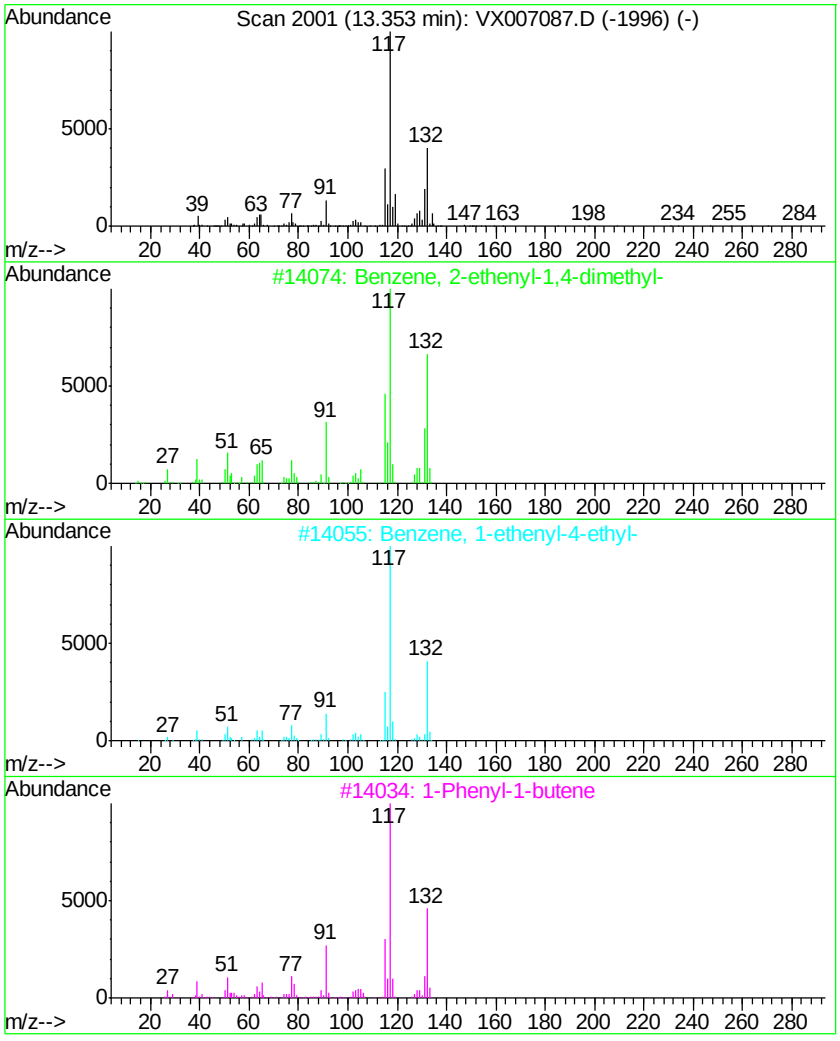
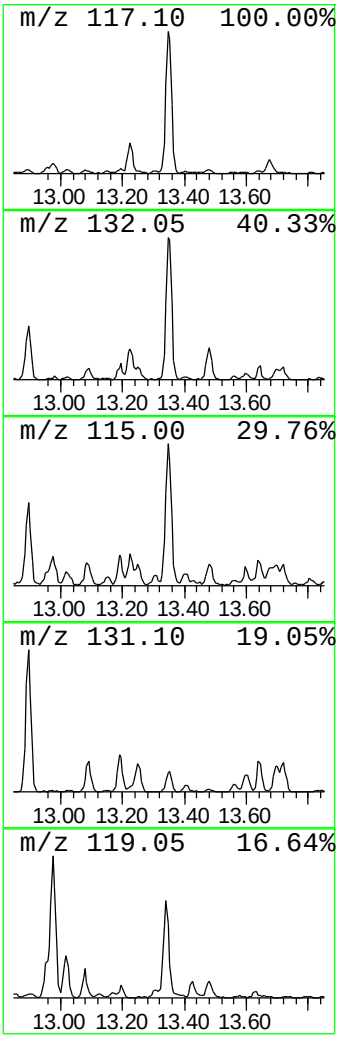
Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	7.97 ug/l	416465	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	95
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	89
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	76
5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	76





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX011719\
Data File : VX007087.D
Acq On : 17 Jan 2019 14:29
Operator : JC/SP
Sample : K1168-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : SS037MW266-190114

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
Pentane, 3-methyl-	3.17	9.3	ug/l	327207	1	5.67	1764310	50.0
Cyclopentane, met...	4.40	10.2	ug/l	360234	1	5.67	1764310	50.0
Cyclopentane, 1,3...	6.26	6.2	ug/l	219553	1	5.67	1764310	50.0
2,2,7,7-Tetrameth...	6.35	14.4	ug/l	592476	2	6.86	2055090	50.0
Pentane, 2,3,3-tr...	8.18	8.8	ug/l	359715	2	6.86	2055090	50.0
o-Cymene	12.23	6.1	ug/l	320770	4	12.08	2613850	50.0
Benzene, 1,3-diet...	12.30	7.8	ug/l	407147	4	12.08	2613850	50.0
Benzene, (2-methy...	12.70	6.4	ug/l	335349	4	12.08	2613850	50.0
1-Phenyl-1-butene	12.76	15.9	ug/l	832675	4	12.08	2613850	50.0
Benzene, 2-etheny...	13.35	8.0	ug/l	416465	4	12.08	2613850	50.0