LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX022719\
Data File : VX007746.D
Acq On : 27 Feb 2019 20:44
Operator : JC/SP
Sample : K1653-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-1R

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\82X022519W.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	1.794	99	105	117	rVB	290049	424315	12.51%	1.727%
2	2.398	197	204	209	rBV	40345	78559	2.32%	0.320%
3	2.843	269	277	288	rBV	214973	529208	15.61%	2.154%
4	2.934	288	292	306	rVB3	27635	65101	1.92%	0.265%
5	3.172	321	331	349	rVB2	258577	733973	21.65%	2.988%
6	4.135	476	489	496	rBV3	11875	41153	1.21%	0.168%
7	4.306	505	517	526	rBV	96547	291521	8.60%	1.187%
8	4.403	526	533	543	rVB	22825	65782	1.94%	0.268%
9	4.537	543	555	557	rBV4	15033	38248	1.13%	0.156%
10	5.238	655	670	690	rVB5	21649	97485	2.87%	0.397%
11	5.507	702	714	724	rBV	187065	543017	16.01%	2.210%
12	5.665	730	740	745	rBV	301650	876729	25.86%	3.569%
13	5.726	745	750	770	rVB	289804	844505	24.91%	3.438%
14	5.946	774	786	796	rBV2	24524	81222	2.40%	0.331%
15	6.068	796	806	814	rBV	198174	514362	15.17%	2.094%
16	6.147	814	819	827	rVV2	28893	74115	2.19%	0.302%
17	6.263	827	838	843	rVV3	111976	352151	10.39%	1.434%
18	6.348	843	852	863	rVV	495906	1556520	45.90%	6.336%
19	6.452	863	869	880	rVB3	35222	101843	3.00%	0.415%
20	6.860	925	936	946	rBV	483564	1141456	33.66%	4.647%
21	7.256	993	1001	1008	rVB2	18131	40456	1.19%	0.165%
22	7.439	1022	1031	1044	rVB4	62065	156936	4.63%	0.639%
23	7.573	1044	1053	1058	rBV2	47152	98724	2.91%	0.402%
24	7.634	1058	1063	1067	rVV	76800	157137	4.63%	0.640%
25	7.677	1067	1070	1077	rVB2	38590	73664	2.17%	0.300%
26	7.848	1090	1098	1105	rVB3	51231	107350	3.17%	0.437%
27	7.988	1105	1121	1123	rBV6	14971	46704	1.38%	0.190%
28	8.055	1123	1132	1143	rVB	311375	644951	19.02%	2.625%
29	8.177	1143	1152	1160	rBV	452323	906793	26.74%	3.691%
30	8.256	1160	1165	1169	rVB	51210	82162	2.42%	0.334%
31	8.378	1179	1185	1192	rBV4	24785	55568	1.64%	0.226%
32	8.451	1192	1197	1201	rVV	25571	43130	1.27%	0.176%
33	8.506	1201	1206	1214	rVV2	52056	127619	3.76%	0.520%
34	8.579	1214	1218	1225	rVB2	38089	70111	2.07%	0.285%

LSC Area Percent Report

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Operator : JC/SP
Sample : K1653-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-1R

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\82X022519W.M
Title : SW846 8260

35	8.671	1225	1233	1235	rBV2	94422	165582	4.88%	0.674%
36	8.713	1235	1240	1249	rVB	1021824	1590702	46.91%	6.475%
37	8.835	1256	1260	1264	rVB	28487	39049	1.15%	0.159%
38	9.037	1288	1293	1298	rVB4	32579	59565	1.76%	0.242%
39	9.165	1306	1314	1319	rVB2	61851	117360	3.46%	0.478%
40	9.219	1319	1323	1327	rBV	47710	74125	2.19%	0.302%
41	9.439	1354	1359	1363	rBV	29808	45678	1.35%	0.186%
42	9.518	1363	1372	1377	rVV	556921	792240	23.36%	3.225%
43	9.567	1377	1380	1386	rVB3	33451	58129	1.71%	0.237%
44	10.006	1447	1452	1458	rBV	30183	43180	1.27%	0.176%
45	10.110	1464	1469	1475	rBV	999258	1337422	39.44%	5.444%
46	11.067	1622	1626	1631	rBV	29177	42574	1.26%	0.173%
47	11.134	1632	1637	1643	rBV	855347	1115410	32.89%	4.541%
48	11.664	1721	1724	1733	rVB	39150	60569	1.79%	0.247%
49	11.768	1733	1741	1750	rBV5	24316	44836	1.32%	0.183%
50	12.079	1786	1792	1799	rBV	1065966	1324734	39.07%	5.393%
51	12.231	1813	1817	1822	rVB	32703	37975	1.12%	0.155%
52	12.298	1822	1828	1835	rVV	2844571	3390896	100.00%	13.803%
53	12.365	1835	1839	1848	rVB	145071	183254	5.40%	0.746%
54	12.457	1848	1854	1859	rBV	42606	60705	1.79%	0.247%
55	12.530	1859	1866	1870	rVB2	64232	95479	2.82%	0.389%
56	12.664	1882	1888	1891	rBV2	113930	192264	5.67%	0.783%
57	12.701	1891	1894	1898	rVV	189447	232978	6.87%	0.948%
58	12.755	1898	1903	1910	rVB	376790	522944	15.42%	2.129%
59	12.896	1919	1926	1931	rVB2	34223	54832	1.62%	0.223%
60	12.981	1931	1940	1946	rBV2	39962	86442	2.55%	0.352%
61	13.079	1951	1956	1962	rBV2	63241	97954	2.89%	0.399%
62	13.225	1977	1980	1989	rVB	151481	202697	5.98%	0.825%
63	13.347	1995	2000	2005	rBV	663313	807959	23.83%	3.289%
64	13.481	2018	2022	2027	rVB	38628	52725	1.55%	0.215%
65	13.603	2038	2042	2045	rBV	33201	42316	1.25%	0.172%
66	13.639	2045	2048	2052	rBV	51269	65895	1.94%	0.268%
67	13.719	2058	2061	2068	rVB	115455	151587	4.47%	0.617%
68	13.981	2098	2104	2108	rBV	47830	62103	1.83%	0.253%
69	14.030	2108	2112	2115	rBV	32781	42429	1.25%	0.173%
70	14.133	2123	2129	2133	rBV	33998	46710	1.38%	0.190%
71	14.267	2146	2151	2156	rBV2	25957	40368	1.19%	0.164%

LSC Area Percent Report

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Operator : JC/SP
Sample : K1653-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-1R

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\82X022519W.M
Title : SW846 8260

72	14.432	2174	2178	2183	rVB2	29723	41370	1.22%	0.168%
73	14.542	2191	2196	2200	rBV2	22989	35680	1.05%	0.145%
74	14.840	2240	2245	2250	rBV2	25851	44439	1.31%	0.181%

Sum of corrected areas: 24565726

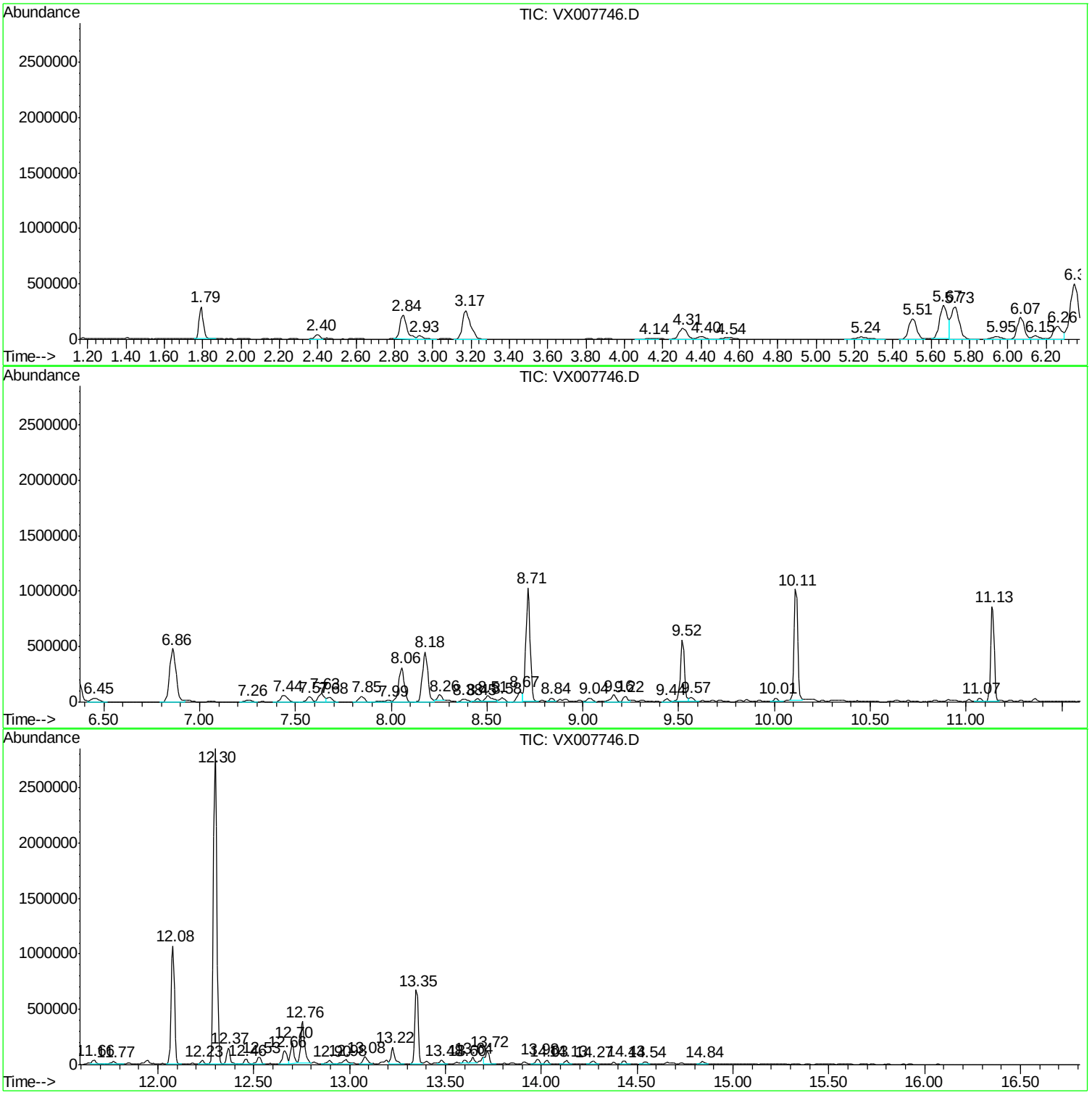
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX022719\
Data File : VX007746.D
Acq On : 27 Feb 2019 20:44
Operator : JC/SP
Sample : K1653-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-1R

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X022519W.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\VX022719\
Data File : VX007746.D
Acq On : 27 Feb 2019 20:44
Operator : JC/SP
Sample : K1653-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-1R

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X022519W.M
Quant Title : SW846 8260

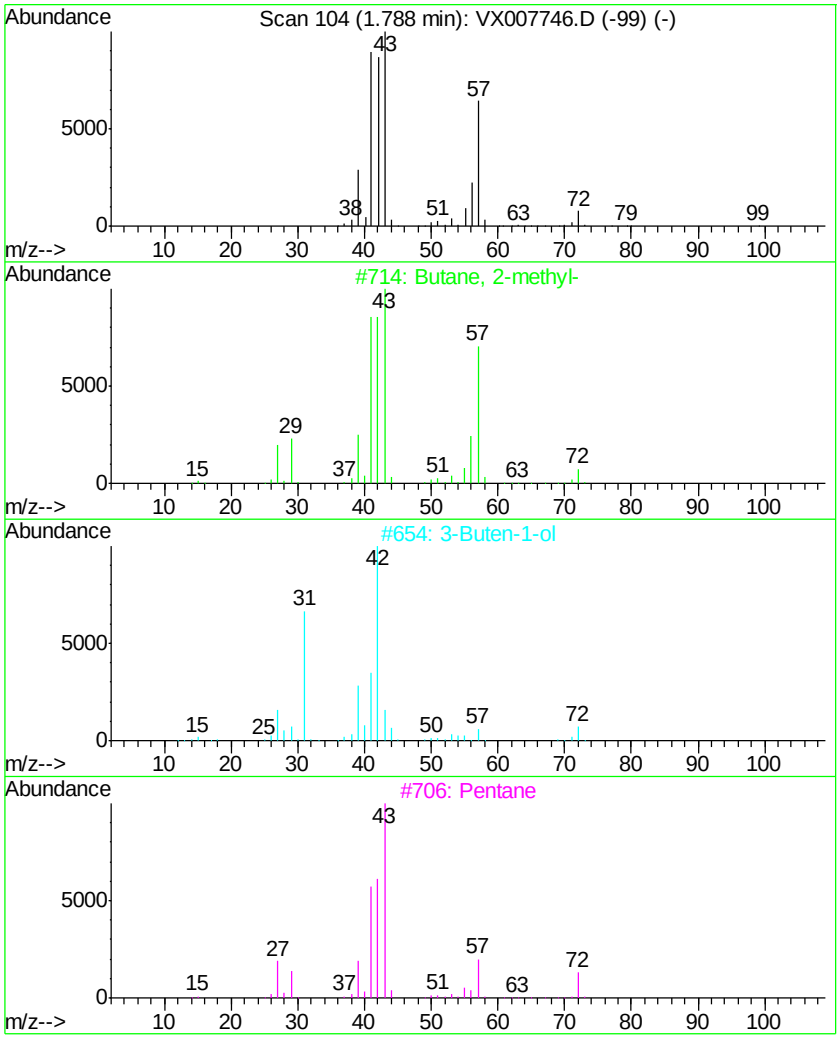
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

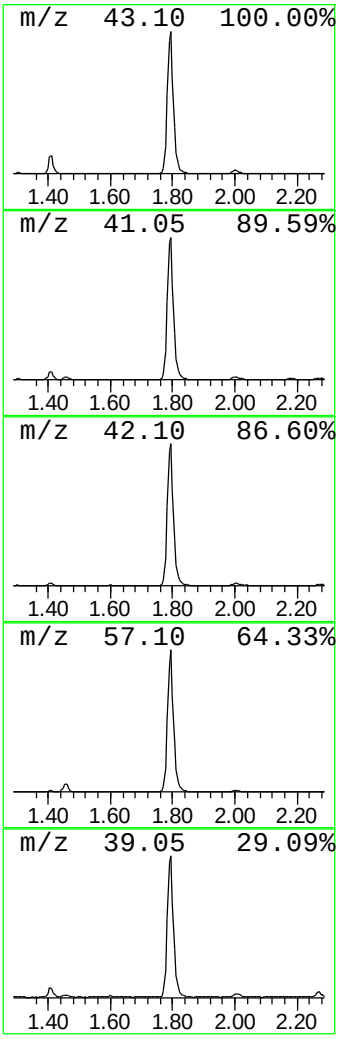
****

Peak Number 1 Butane, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.79	24.20 ug/l	424315	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	94
2	3-Buten-1-ol	72	C4H8O	000627-27-0	47
3	Pentane	72	C5H12	000109-66-0	36
4	1-Propene, 2-methyl-	56	C4H8	000115-11-7	10
5	1-Butene	56	C4H8	000106-98-9	9





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX022719\
Data File : VX007746.D
Acq On : 27 Feb 2019 20:44
Operator : JC/SP
Sample : K1653-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-1R

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X022519W.M
Quant Title : SW846 8260

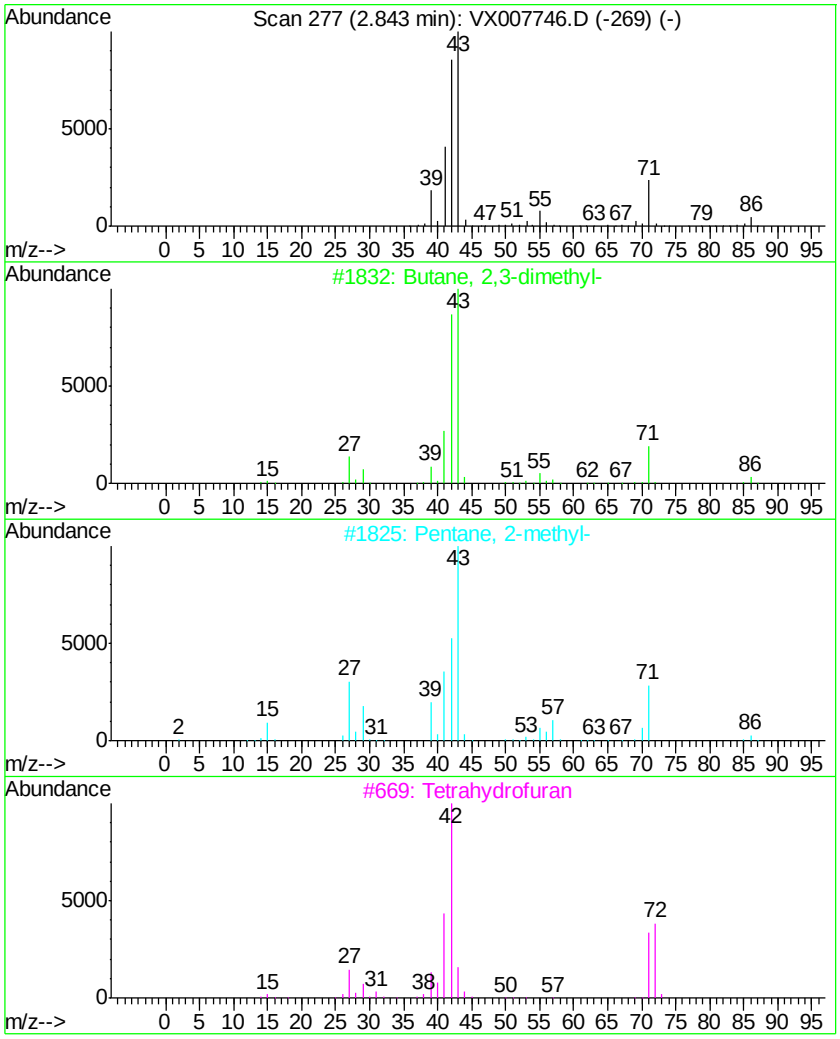
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

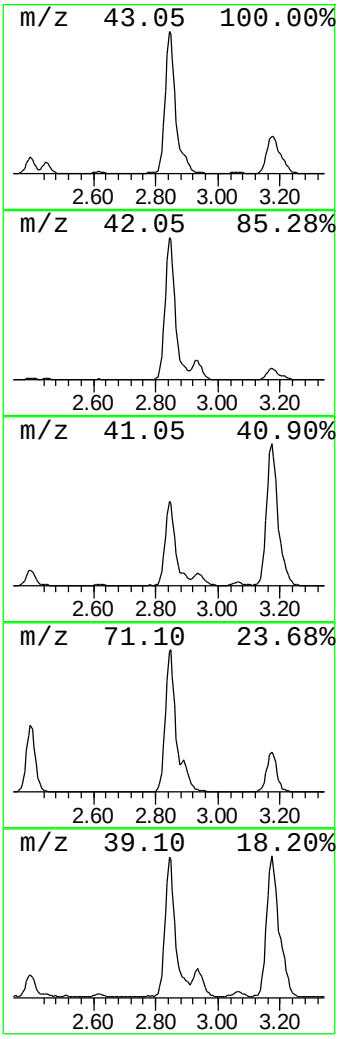
****

Peak Number 2 Butane, 2,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.84	30.18 ug/l	529208	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	Butane, 2-methyl-	72	C5H12	000078-78-4	36
5	Cyclobutane, methyl-	70	C5H10	000598-61-8	28





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX022719\
Data File : VX007746.D
Acq On : 27 Feb 2019 20:44
Operator : JC/SP
Sample : K1653-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-1R

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X022519W.M
Quant Title : SW846 8260

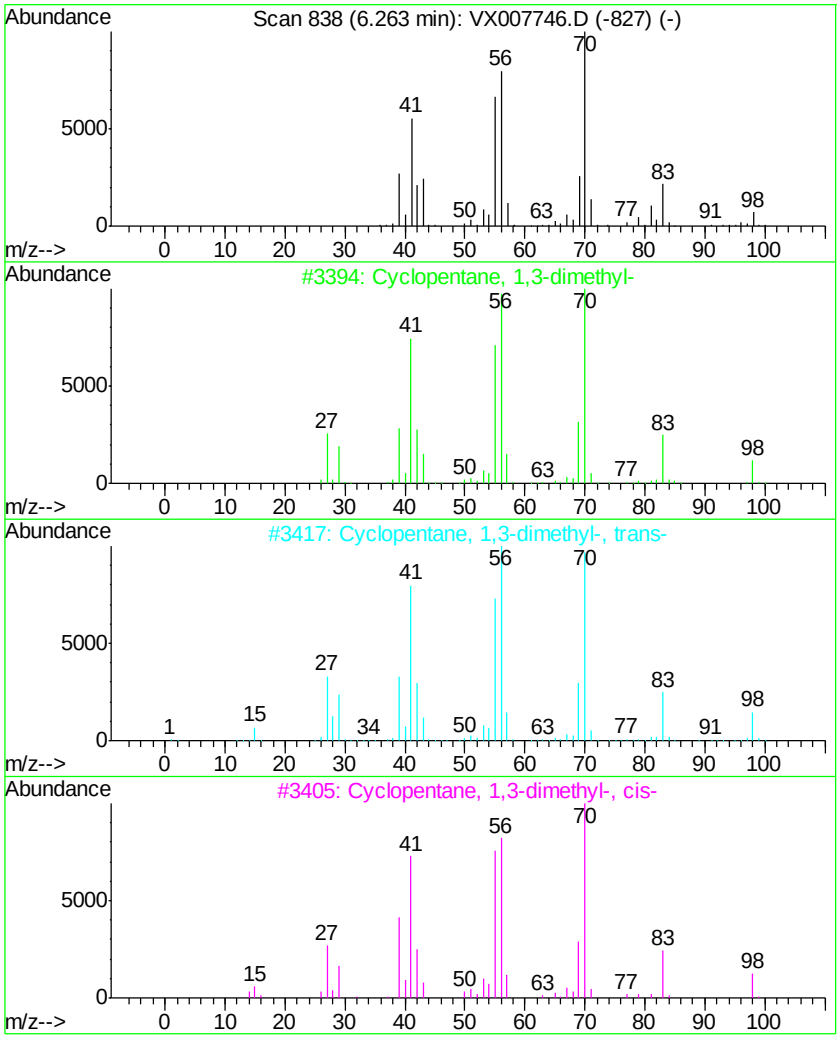
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

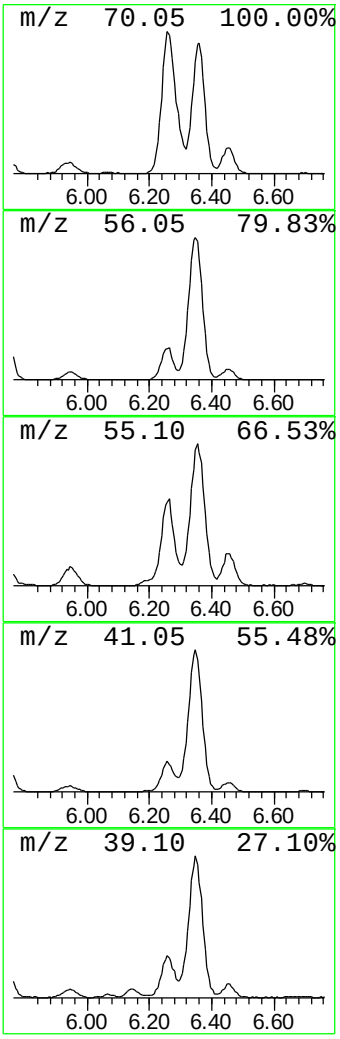
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.26	20.08 ug/l	352151	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	90
2	Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	90
3	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	90
4	Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	87
5	2H-Pyran-2,6(3H)-dione, dihydro-...	142	C7H10O3	004160-82-1	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX022719\
Data File : VX007746.D
Acq On : 27 Feb 2019 20:44
Operator : JC/SP
Sample : K1653-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-1R

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X022519W.M
Quant Title : SW846 8260

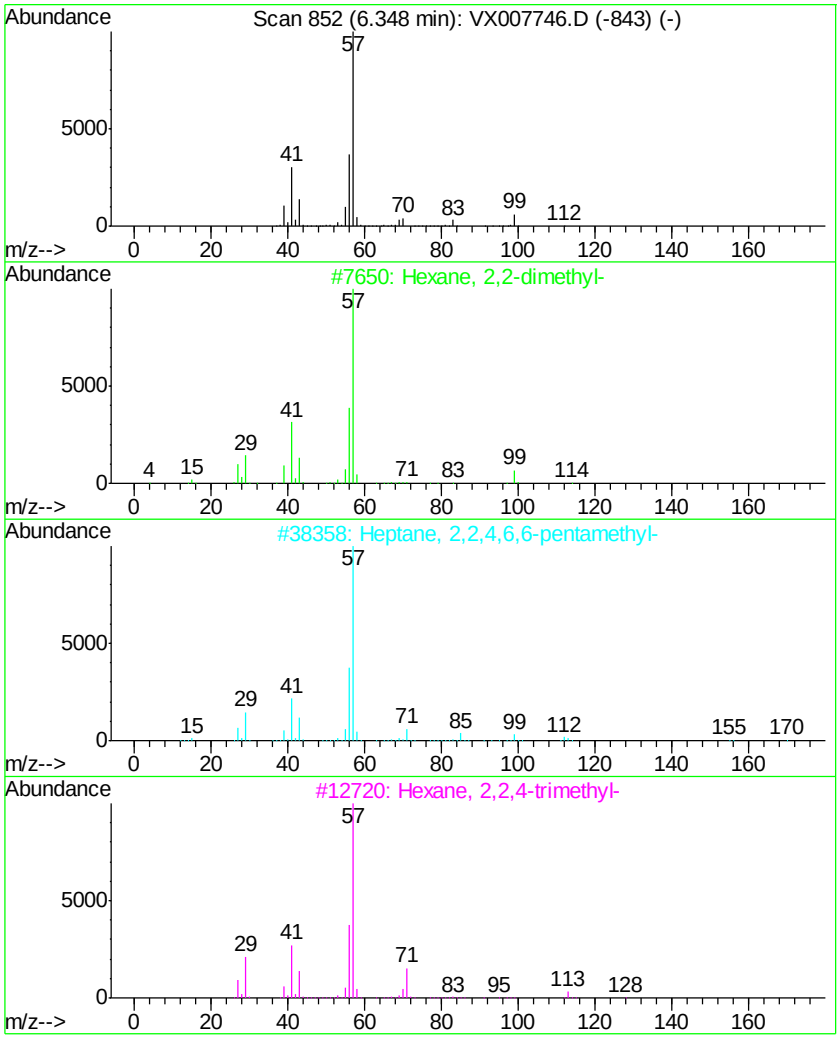
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

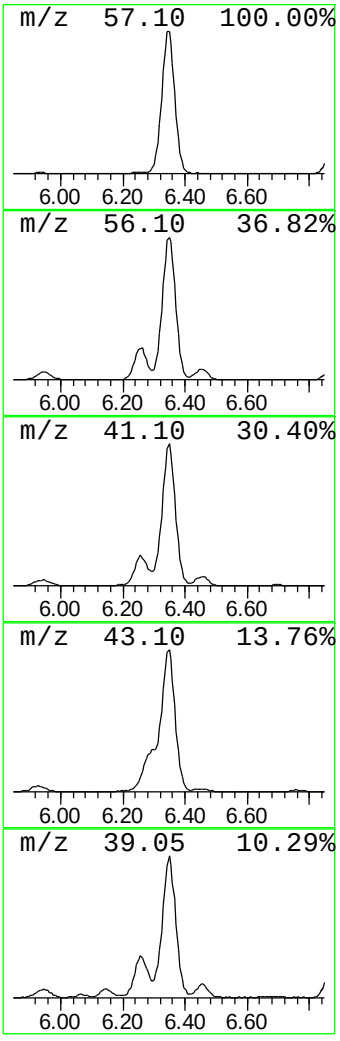
****

Peak Number 4 Hexane, 2,2-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.35	68.18 ug/l	1556520	1,4-Difluorobenzene	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	59
2	Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	59
3	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	59
4	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	53
5	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	50





Library Search Compound Report

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Data File : VX007746.D
Acq On : 27 Feb 2019 20:44
Operator : JC/SP
Sample : K1653-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-1R

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X022519W.M
Quant Title : SW846 8260

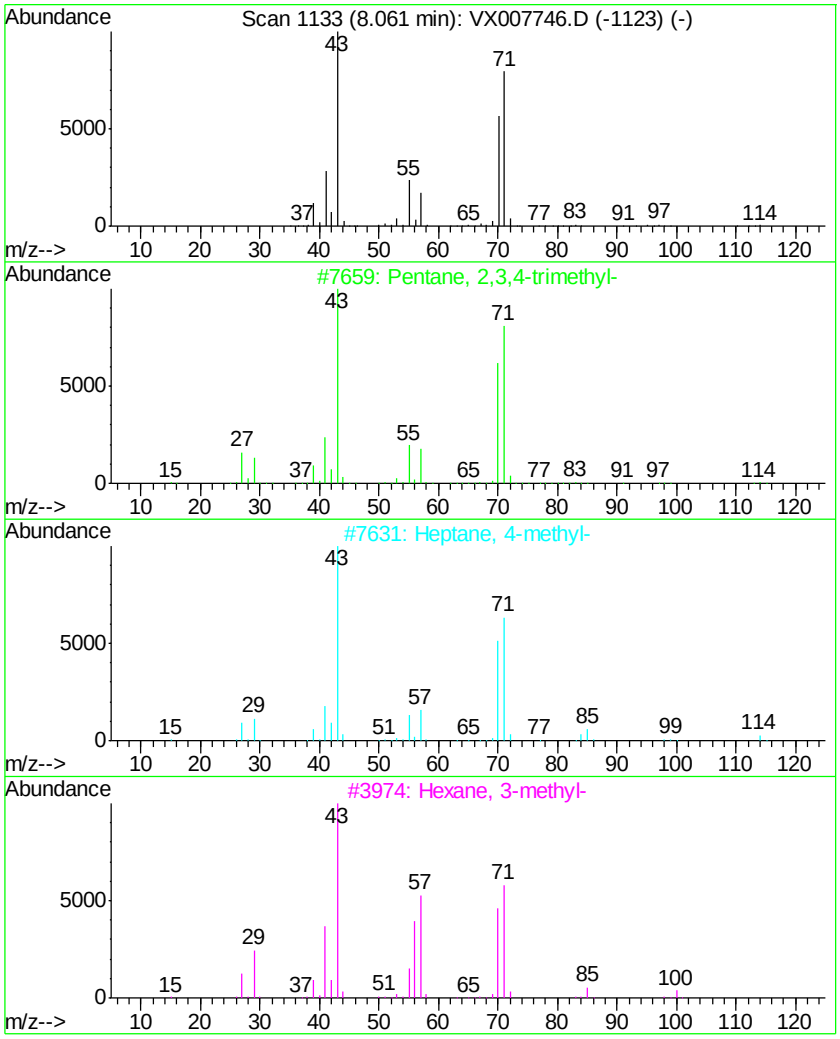
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

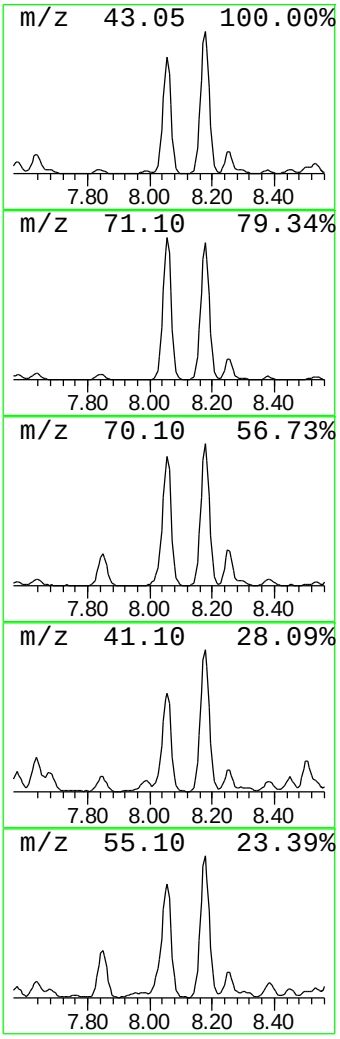
****

Peak Number 5 Pentane, 2,3,4-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.06	28.25 ug/l	644951	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	Hexane, 3-methyl-	100	C7H16	000589-34-4	72
4	Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	64
5	Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX022719\
Data File : VX007746.D
Acq On : 27 Feb 2019 20:44
Operator : JC/SP
Sample : K1653-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-1R

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X022519W.M
Quant Title : SW846 8260

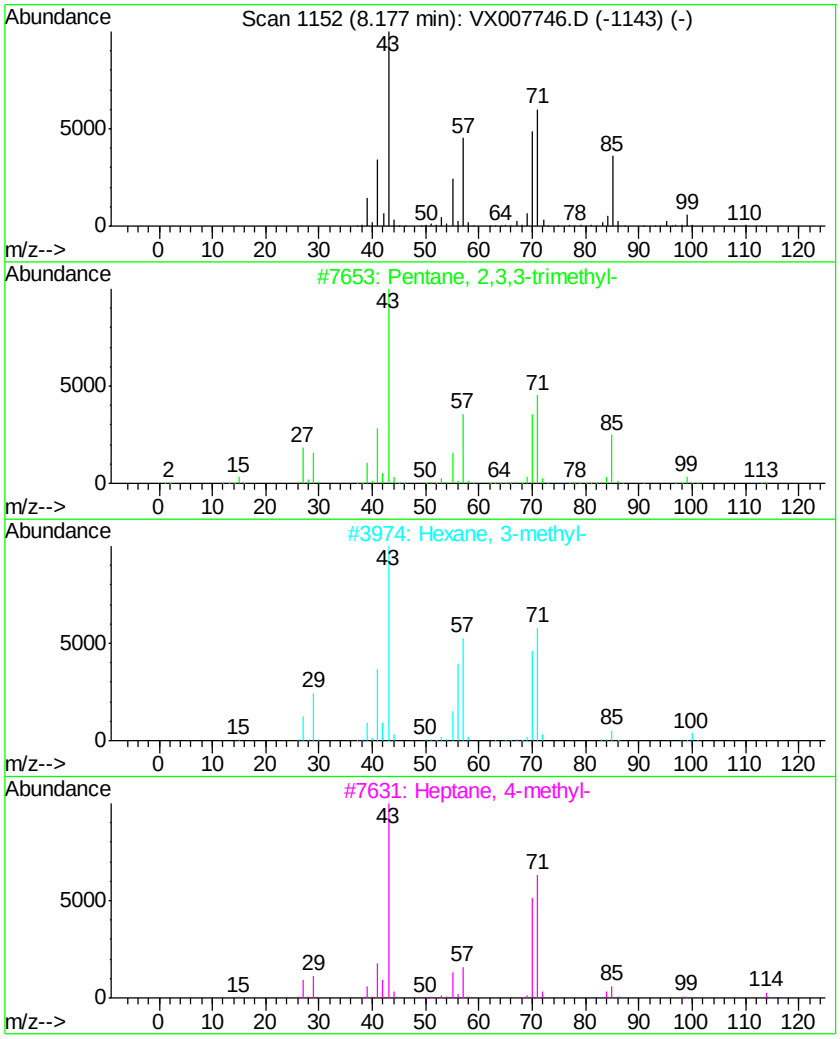
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

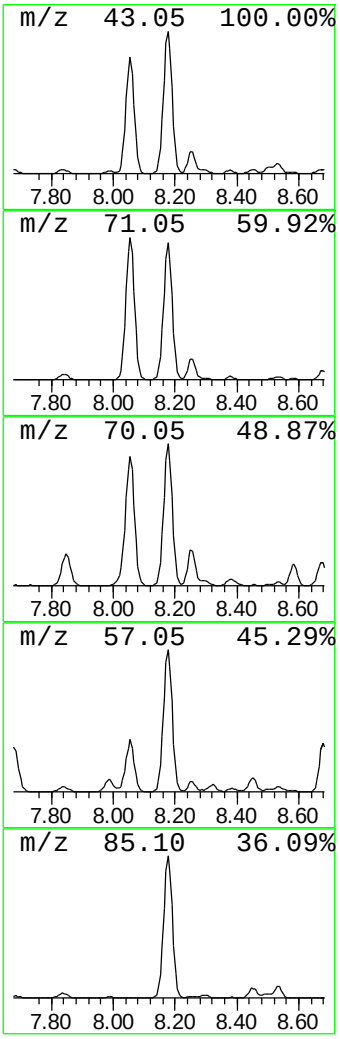
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.18	39.72 ug/l	906793	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-methyl-	100	C7H16	000589-34-4	64
3	Heptane, 4-methyl-	114	C8H18	000589-53-7	64
4	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	59
5	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	59





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX022719\
Data File : VX007746.D
Acq On : 27 Feb 2019 20:44
Operator : JC/SP
Sample : K1653-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-1R

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X022519W.M
Quant Title : SW846 8260

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

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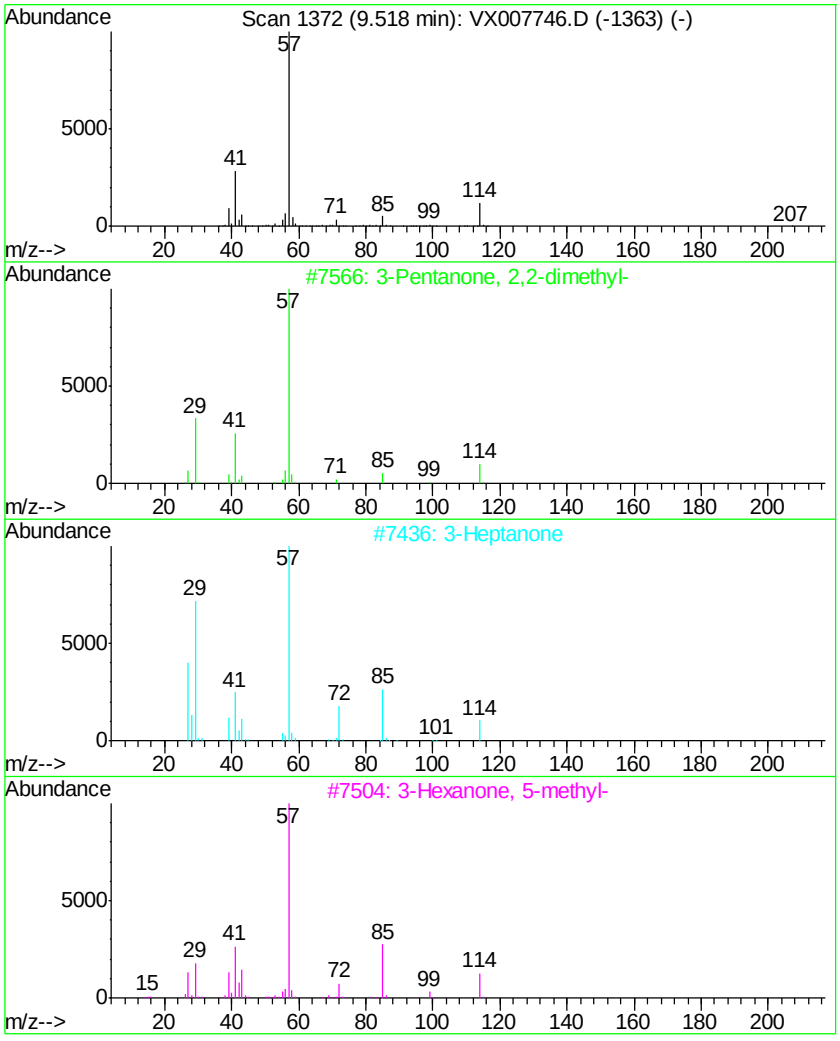
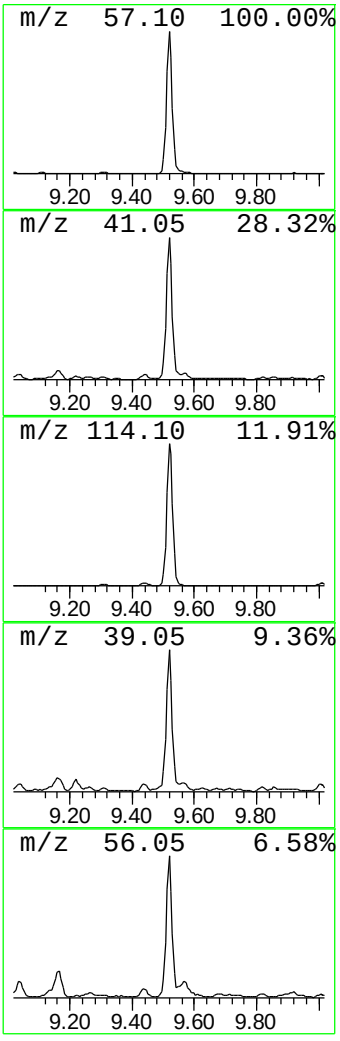
Peak Number 7 3-Pentanone, 2,2-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.52	29.62 ug/l	792240	Chlorobenzene-d5	10.11

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Pentanone, 2,2-dimethyl-	114	C7H14O	000564-04-5	90
2	3-Heptanone	114	C7H14O	000106-35-4	59
3	3-Hexanone, 5-methyl-	114	C7H14O	000623-56-3	59
4	3,4-Hexanedione	114	C6H10O2	004437-51-8	52
5	3-Hexanone, 4-methyl-	114	C7H14O	017042-16-9	38





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX022719\
Data File : VX007746.D
Acq On : 27 Feb 2019 20:44
Operator : JC/SP
Sample : K1653-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-1R

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X022519W.M
Quant Title : SW846 8260

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

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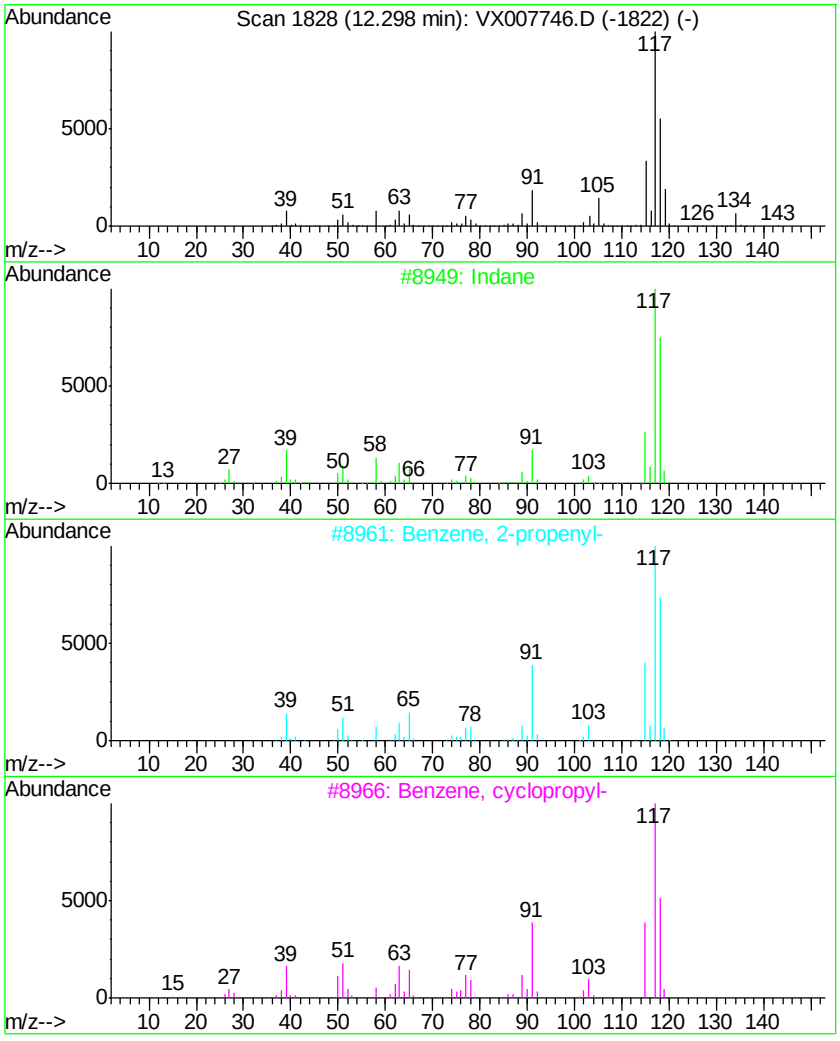
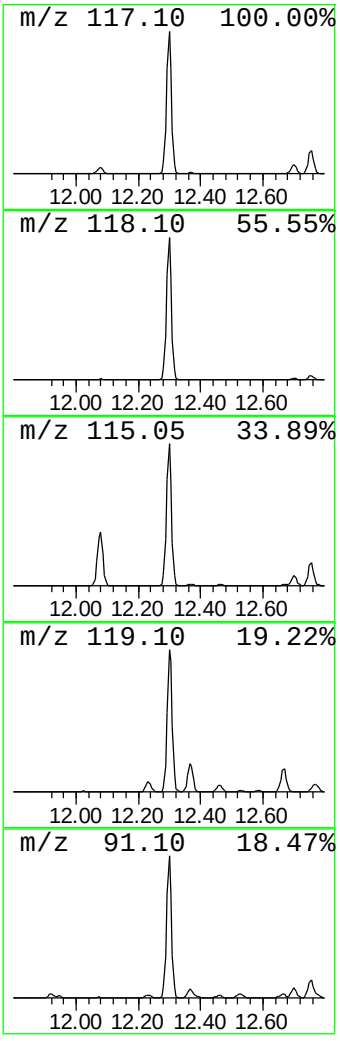
Peak Number 8 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	127.98 ug/l	3390900	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	76
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	76
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
4	Deltacyclene	118	C9H10	007785-10-6	58
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	55





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX022719\
Data File : VX007746.D
Acq On : 27 Feb 2019 20:44
Operator : JC/SP
Sample : K1653-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-1R

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X022519W.M
Quant Title : SW846 8260

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

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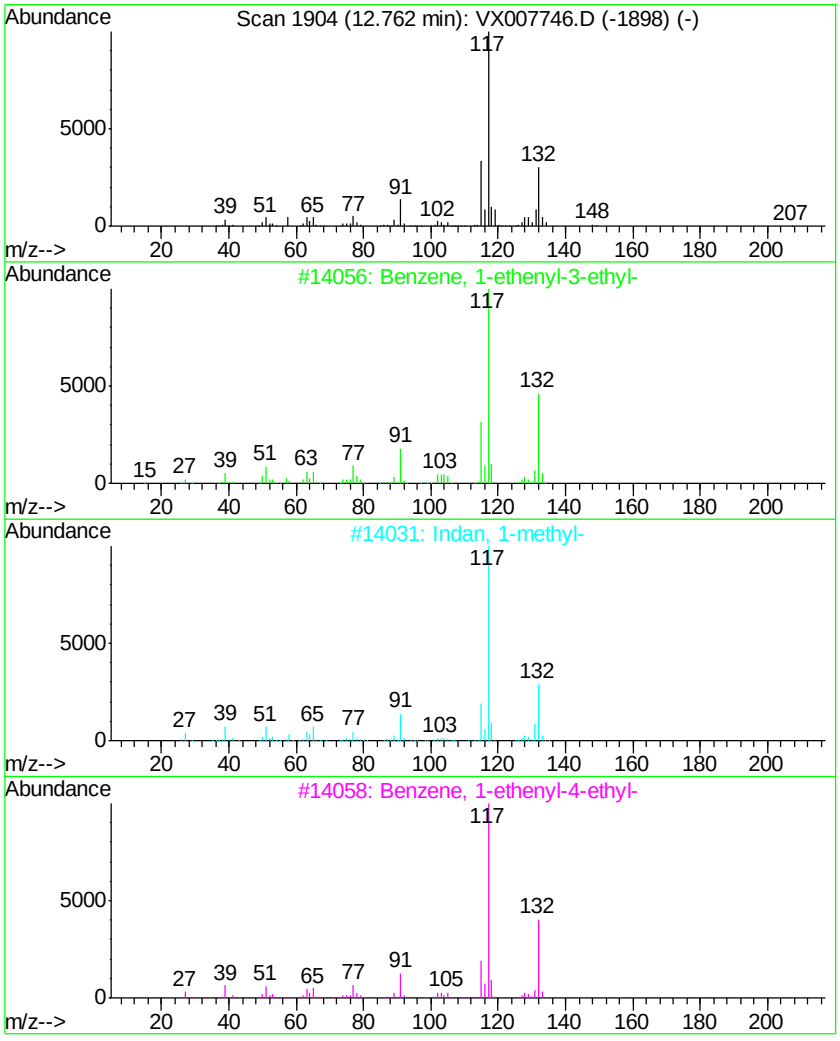
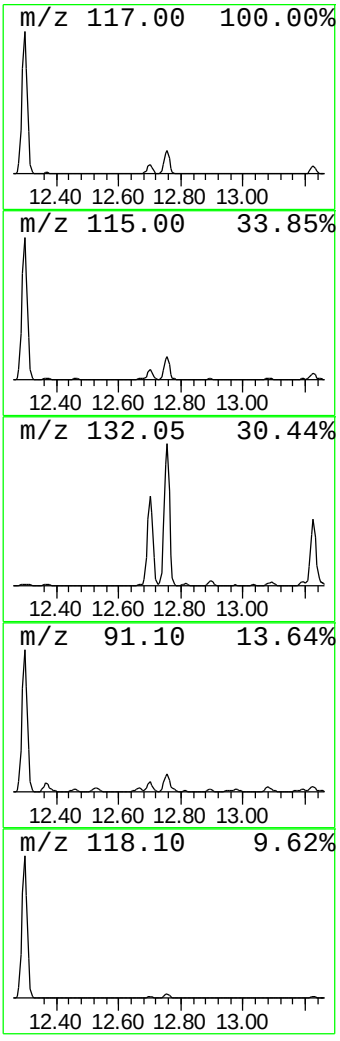
Peak Number 9 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	19.74 ug/l	522944	1,4-Dichlorobenzene-d4	12.08

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, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	86
5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	86





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX022719\
Data File : VX007746.D
Acq On : 27 Feb 2019 20:44
Operator : JC/SP
Sample : K1653-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-1R

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X022519W.M
Quant Title : SW846 8260

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

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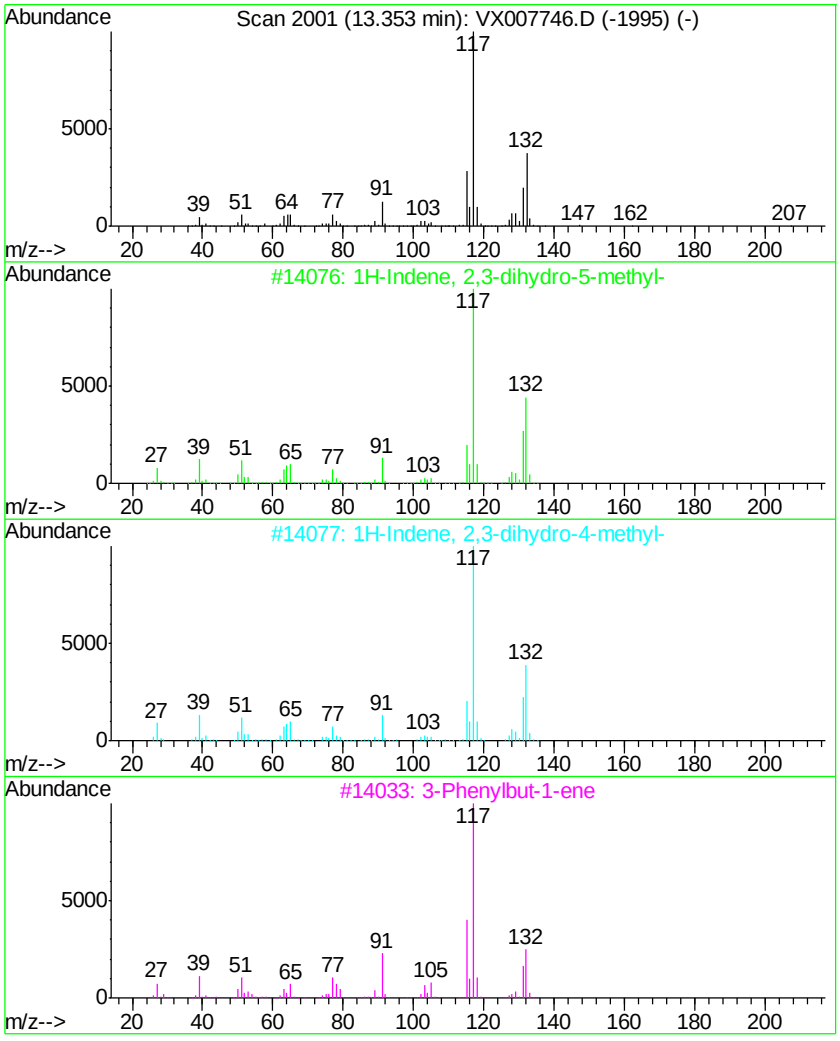
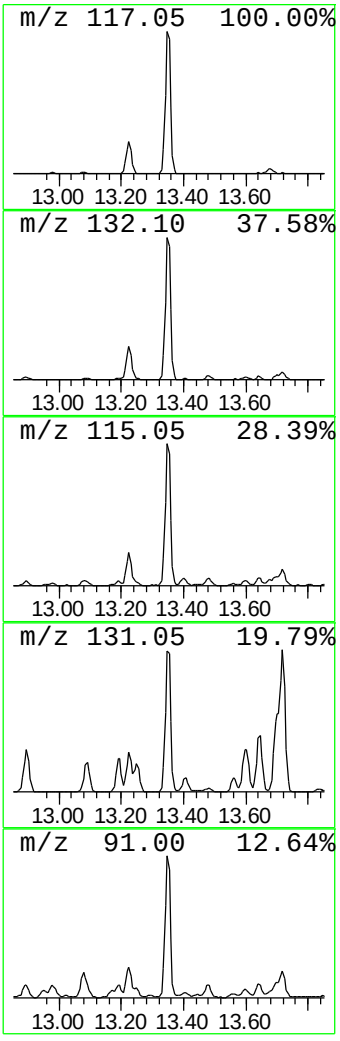
Peak Number 10 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	30.50 ug/l	807959	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
2	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4	Indan, 1-methyl-	132	C10H12	000767-58-8	87
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX022719\
Data File : VX007746.D
Acq On : 27 Feb 2019 20:44
Operator : JC/SP
Sample : K1653-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-1R

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X022519W.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-methyl-	1.79	24.2	ug/l	424315	1	5.67	876729	50.0
Butane, 2,3-dimet...	2.84	30.2	ug/l	529208	1	5.67	876729	50.0
Cyclopentane, 1,3...	6.26	20.1	ug/l	352151	1	5.67	876729	50.0
Hexane, 2,2-dimet...	6.35	68.2	ug/l	1556520	2	6.86	1141460	50.0
Pentane, 2,3,4-tr...	8.06	28.3	ug/l	644951	2	6.86	1141460	50.0
Pentane, 2,3,3-tr...	8.18	39.7	ug/l	906793	2	6.86	1141460	50.0
3-Pentanone, 2,2-...	9.52	29.6	ug/l	792240	3	10.11	1337420	50.0
Indane	12.30	128.0	ug/l	3390900	4	12.08	1324730	50.0
Benzene, 1-etheny...	12.76	19.7	ug/l	522944	4	12.08	1324730	50.0
1H-Indene, 2,3-di...	13.35	30.5	ug/l	807959	4	12.08	1324730	50.0