LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100818\
Data File : VX005133.D
Acq On : 08 Oct 2018 17:49
Operator : JC/MD
Sample : J5268-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-10

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\82X092618W.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	1.404	37	41	46	rBV	292189	285872	2.30%	0.206%
2	1.782	97	103	119	rVB	4498253	6221881	49.98%	4.485%
3	1.995	133	138	152	rVB	1240481	1846419	14.83%	1.331%
4	2.117	152	158	164	rBV	195616	292008	2.35%	0.210%
5	2.263	177	182	193	rVB	323181	511067	4.11%	0.368%
6	2.391	195	203	210	rBV	218139	426204	3.42%	0.307%
7	2.775	257	266	270	rBV3	101067	251862	2.02%	0.182%
8	2.843	270	277	280	rVV	624246	1404905	11.29%	1.013%
9	2.885	280	284	288	rVV	1251439	2542857	20.43%	1.833%
10	2.928	288	291	304	rVB2	676598	1601834	12.87%	1.155%
11	3.166	320	330	353	rVB	1056345	2405653	19.32%	1.734%
12	3.391	358	367	378	rBV	148457	365907	2.94%	0.264%
13	3.513	379	387	400	rVB	174553	399750	3.21%	0.288%
14	3.659	401	411	417	rBV2	93308	241976	1.94%	0.174%
15	3.745	417	425	429	rVV	94351	239780	1.93%	0.173%
16	3.806	429	435	445	rVB	359017	873415	7.02%	0.630%
17	3.915	445	453	460	rBV	244940	651570	5.23%	0.470%
18	3.989	460	465	475	rVV2	135973	463234	3.72%	0.334%
19	4.074	476	479	487	rVV	66036	140684	1.13%	0.101%
20	4.190	489	498	508	rVB	271809	699812	5.62%	0.504%
21	4.397	521	532	544	rVB	1581123	4595711	36.92%	3.313%
22	4.519	544	552	567	rVB	117109	365506	2.94%	0.263%
23	5.299	653	680	697	rBV2	211193	819850	6.59%	0.591%
24	5.501	697	713	717	rBV	162988	477857	3.84%	0.344%
25	5.574	717	725	734	rVB	746768	2134231	17.14%	1.538%
26	5.665	734	740	745	rBV	134965	335543	2.70%	0.242%
27	5.927	771	783	796	rBV4	154800	529341	4.25%	0.382%
28	6.068	796	806	819	rVB	150571	389100	3.13%	0.280%
29	6.257	819	837	847	rBV3	198377	926248	7.44%	0.668%
30	6.354	847	853	862	rVV2	255624	774734	6.22%	0.558%
31	6.452	862	869	881	rVB2	189344	549937	4.42%	0.396%
32	6.860	927	936	943	rBV2	327278	932236	7.49%	0.672%
33	6.939	946	949	963	rVB5	199594	479474	3.85%	0.346%
34	7.256	992	1001	1010	rBV	171831	372758	2.99%	0.269%

LSC Area Percent Report

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Operator : JC/MD
Sample : J5268-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-10

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\82X092618W.M
Title : SW846 8260

35	7.464	1023	1035	1047	rBV3	710025	1848926	14.85%	1.333%
36	7.732	1073	1079	1092	rVB	98397	201047	1.61%	0.145%
37	7.933	1104	1112	1113	rBV	111704	197037	1.58%	0.142%
38	7.957	1113	1116	1123	rVV2	138921	275748	2.21%	0.199%
39	8.055	1123	1132	1144	rVB2	84670	253888	2.04%	0.183%
40	8.177	1144	1152	1155	rBV	163524	341282	2.74%	0.246%
41	8.213	1155	1158	1163	rVV	226324	376385	3.02%	0.271%
42	8.573	1211	1217	1226	rVB3	286249	570718	4.58%	0.411%
43	8.671	1226	1233	1235	rBV3	79246	151473	1.22%	0.109%
44	8.713	1235	1240	1247	rVV	727210	1200714	9.64%	0.865%
45	8.787	1247	1252	1258	rVV	498908	747000	6.00%	0.538%
46	8.915	1269	1273	1280	rVB	108187	172724	1.39%	0.125%
47	9.037	1288	1293	1299	rVB4	60941	128679	1.03%	0.093%
48	9.226	1319	1324	1334	rVB	128814	229432	1.84%	0.165%
49	10.116	1460	1470	1475	rBV	668625	977529	7.85%	0.705%
50	10.256	1487	1493	1504	rBV	9406562	12449242	100.00%	8.974%
51	10.366	1505	1511	1526	rVB	9024850	12149484	97.59%	8.757%
52	10.701	1560	1566	1577	rVB	3459266	4418040	35.49%	3.185%
53	11.018	1612	1618	1625	rBV	597854	786249	6.32%	0.567%
54	11.140	1633	1638	1643	rBV	696451	892002	7.17%	0.643%
55	11.359	1666	1674	1681	rBV	1297932	1594176	12.81%	1.149%
56	11.439	1681	1687	1688	rBV	2184431	2815613	22.62%	2.030%
57	11.506	1694	1698	1704	rVB	1634387	1955484	15.71%	1.410%
58	11.670	1719	1725	1735	rBV	2544191	3139694	25.22%	2.263%
59	11.810	1743	1748	1754	rBV	6627588	7799981	62.65%	5.622%
60	11.945	1768	1770	1777	rVB	108011	129631	1.04%	0.093%
61	12.079	1787	1792	1798	rVB	942135	1180582	9.48%	0.851%
62	12.146	1798	1803	1808	rBV	1765240	2173895	17.46%	1.567%
63	12.298	1823	1828	1836	rBV	9343743	12317917	98.95%	8.879%
64	12.371	1836	1840	1850	rVB3	701713	1035129	8.31%	0.746%
65	12.469	1850	1856	1860	rBV2	300459	428945	3.45%	0.309%
66	12.518	1860	1864	1870	rVB	331696	388958	3.12%	0.280%
67	12.585	1870	1875	1878	rBV	680758	917820	7.37%	0.662%
68	12.670	1884	1889	1892	rBV	1431792	1716130	13.79%	1.237%
69	12.701	1892	1894	1898	rVV	634093	704168	5.66%	0.508%
70	12.755	1898	1903	1910	rVB	2432474	3085393	24.78%	2.224%
71	12.902	1922	1927	1932	rVB2	212759	285630	2.29%	0.206%

LSC Area Percent Report

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ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-10

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\82X092618W.M
Title : SW846 8260

72	12.975	1932	1939	1943	rBV	770968	975321	7.83%	0.703%
73	13.018	1943	1946	1952	rVV	668054	767089	6.16%	0.553%
74	13.085	1952	1957	1962	rVV4	136496	285572	2.29%	0.206%
75	13.194	1966	1975	1976	rVV3	155012	229451	1.84%	0.165%
76	13.225	1976	1980	1990	rVV	1902437	2411467	19.37%	1.738%
77	13.310	1990	1994	1996	rVV2	122077	177611	1.43%	0.128%
78	13.353	1996	2001	2006	rVV2	3688649	4875678	39.16%	3.514%
79	13.402	2006	2009	2018	rVV3	358695	588324	4.73%	0.424%
80	13.481	2018	2022	2028	rVB	614077	777399	6.24%	0.560%
81	13.560	2030	2035	2038	rBV2	178871	248257	1.99%	0.179%
82	13.603	2038	2042	2045	rVV	518218	704018	5.66%	0.507%
83	13.645	2045	2049	2053	rVV2	370064	676665	5.44%	0.488%
84	13.700	2054	2058	2059	rVV	262477	356987	2.87%	0.257%
85	13.719	2059	2061	2067	rVB2	511101	714998	5.74%	0.515%
86	13.834	2072	2080	2089	rBV	4851477	5852627	47.01%	4.219%
87	13.926	2089	2095	2100	rBV	358694	482118	3.87%	0.348%
88	13.981	2100	2104	2108	rBV2	168712	219495	1.76%	0.158%
89	14.030	2108	2112	2116	rVB	159283	178592	1.43%	0.129%
90	14.133	2124	2129	2135	rBV	289052	366559	2.94%	0.264%
91	14.273	2148	2152	2158	rBV2	196801	328657	2.64%	0.237%
92	14.432	2174	2178	2183	rVB	169895	210182	1.69%	0.152%
93	14.657	2209	2215	2218	rBV	158992	222315	1.79%	0.160%
94	14.694	2218	2221	2232	rVB	583637	780498	6.27%	0.563%
95	14.840	2240	2245	2255	rBV2	524966	716769	5.76%	0.517%

Sum of corrected areas: 138732610

LSC Report - Integrated Chromatogram

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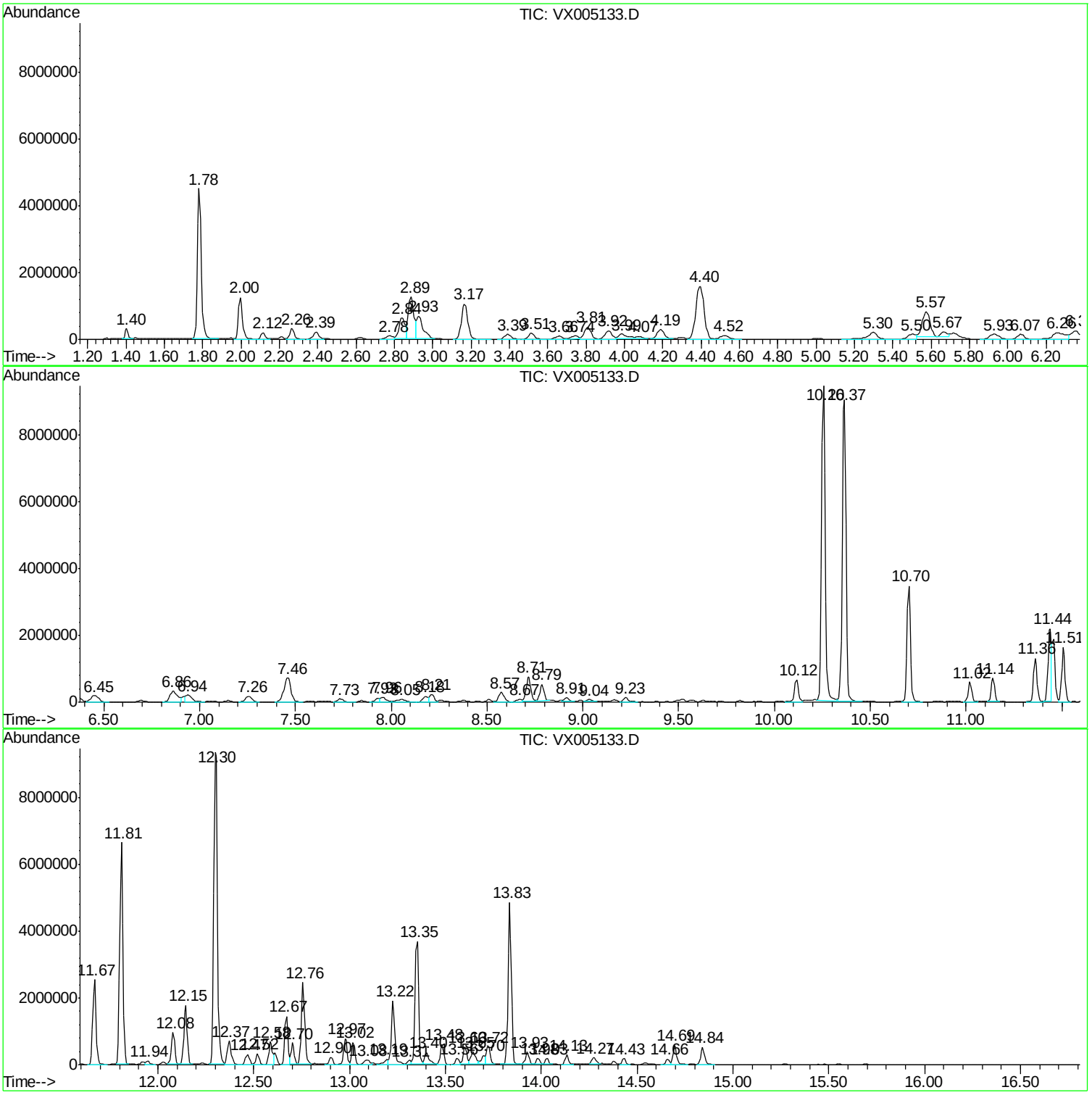
Data File : VX005133.D
Acq On : 08 Oct 2018 17:49
Operator : JC/MD
Sample : J5268-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.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\VX100818\
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Operator : JC/MD
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Instrument : MSVOA_X
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

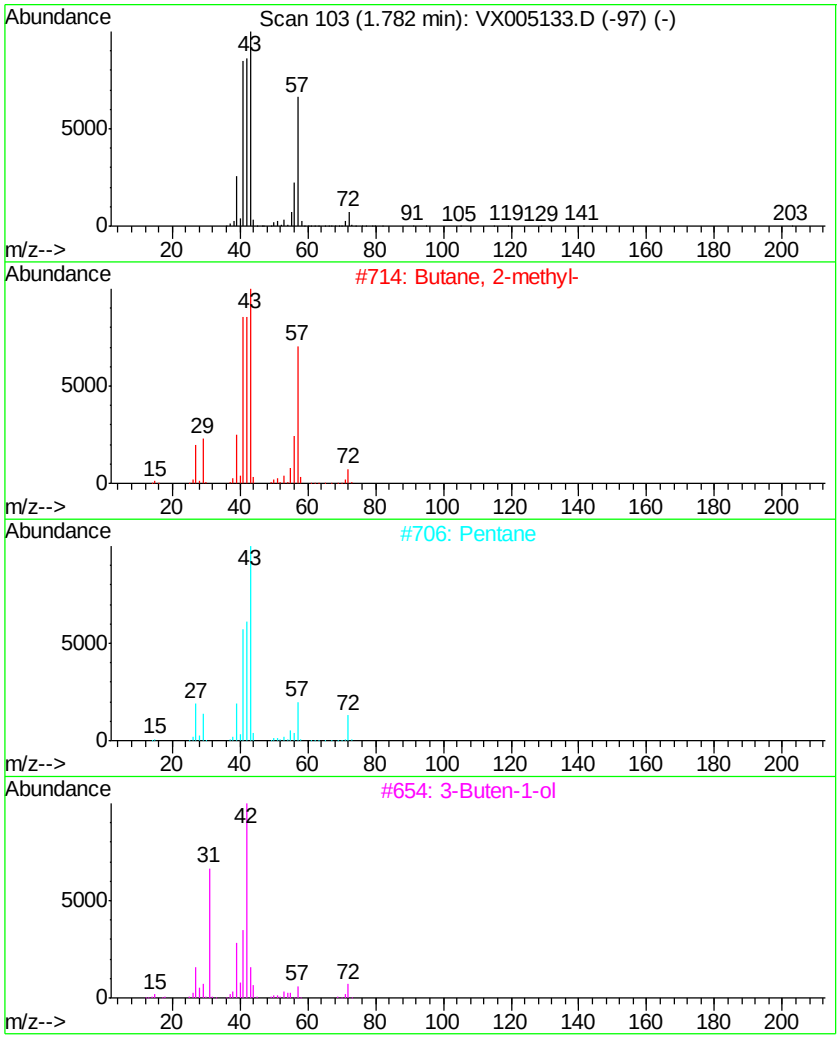
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

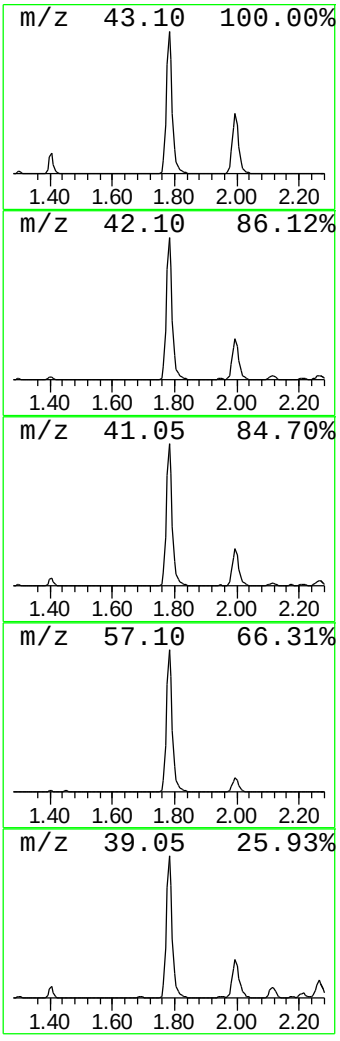
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.78	927.14 ug/l	6221880	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	94
2	Pentane	72	C5H12	000109-66-0	36
3	3-Buten-1-ol	72	C4H8O	000627-27-0	28
4	Methane, isocyanato-	57	C2H3NO	000624-83-9	9
5	1-Propene, 2-methyl-	56	C4H8	000115-11-7	9





Library Search Compound Report

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Data File : VX005133.D
Acq On : 08 Oct 2018 17:49
Operator : JC/MD
Sample : J5268-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

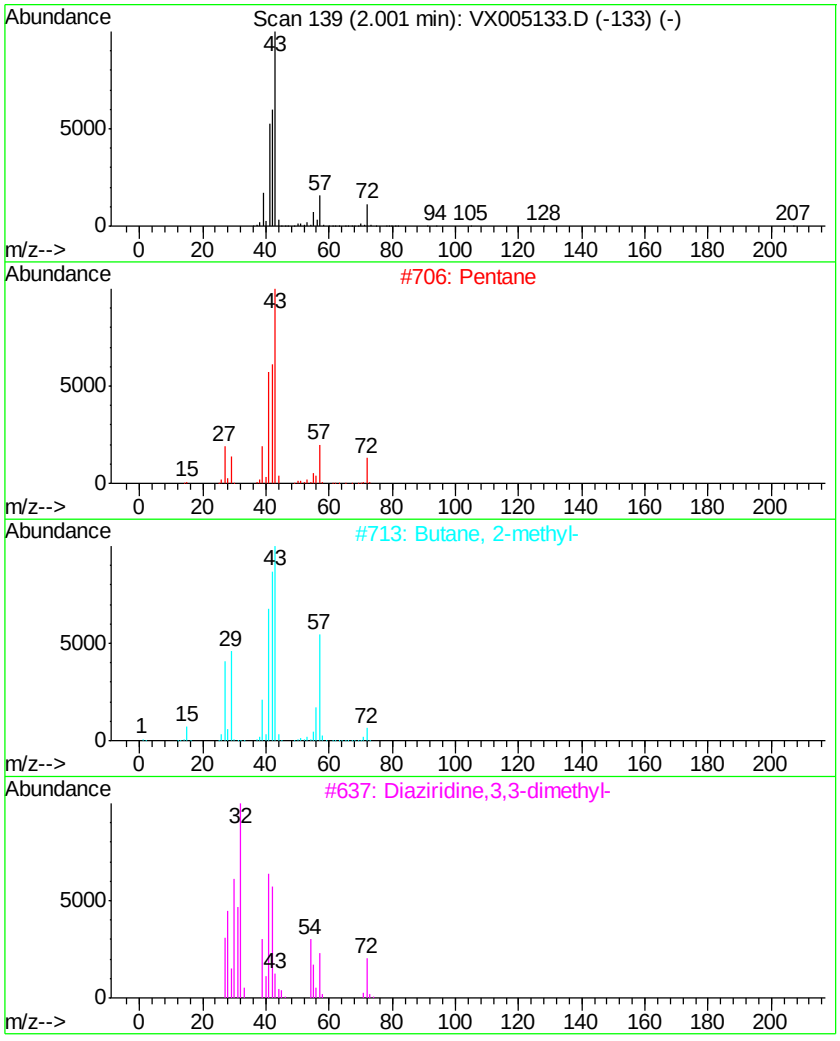
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

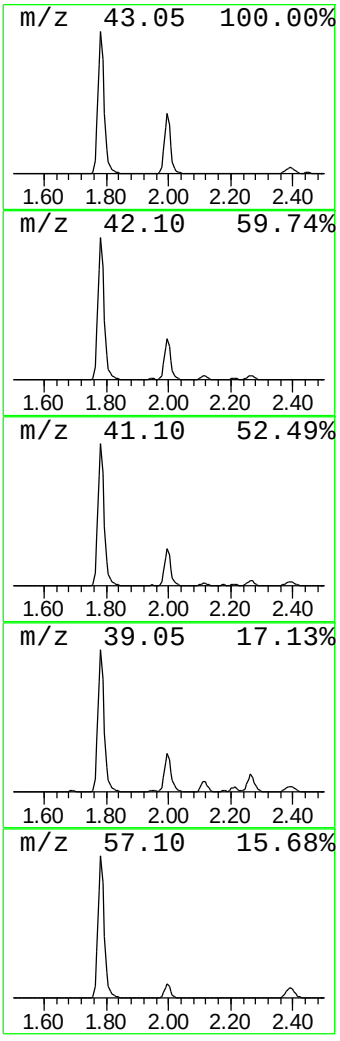
****

Peak Number 2 Pentane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.00	275.14 ug/l	1846420	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	91
2	Butane, 2-methyl-	72	C5H12	000078-78-4	38
3	Diaziridine,3,3-dimethyl-	72	C3H8N2	004901-76-2	9
4	Oxirane, ethyl-	72	C4H8O	000106-88-7	9
5	Butanal, 2,2-dimethyl-	100	C6H12O	002094-75-9	9





Library Search Compound Report

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ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

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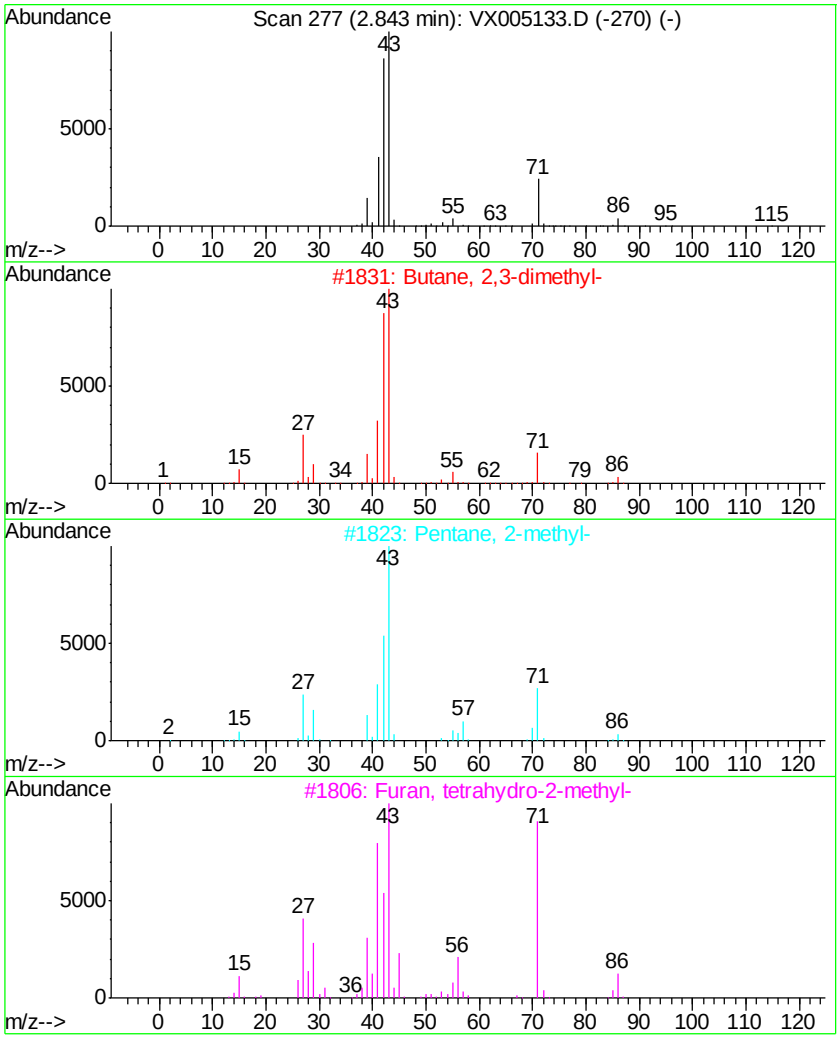
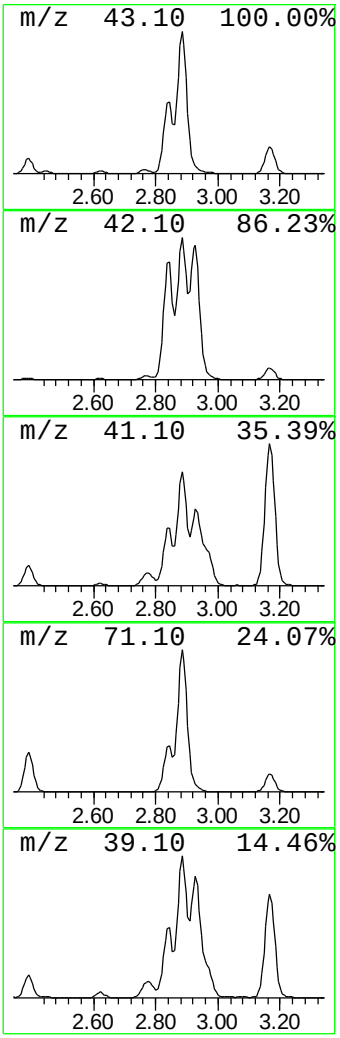
Peak Number 3 Butane, 2,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.84	209.35 ug/l	1404910	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	80
2	Pentane, 2-methyl-	86	C6H14	000107-83-5	40
3	Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	9
4	Propanoic acid, 2-methyl-, 2-pro...	128	C7H12O2	015727-77-2	9
5	Pentane	72	C5H12	000109-66-0	9





Library Search Compound Report

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Sample : J5268-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

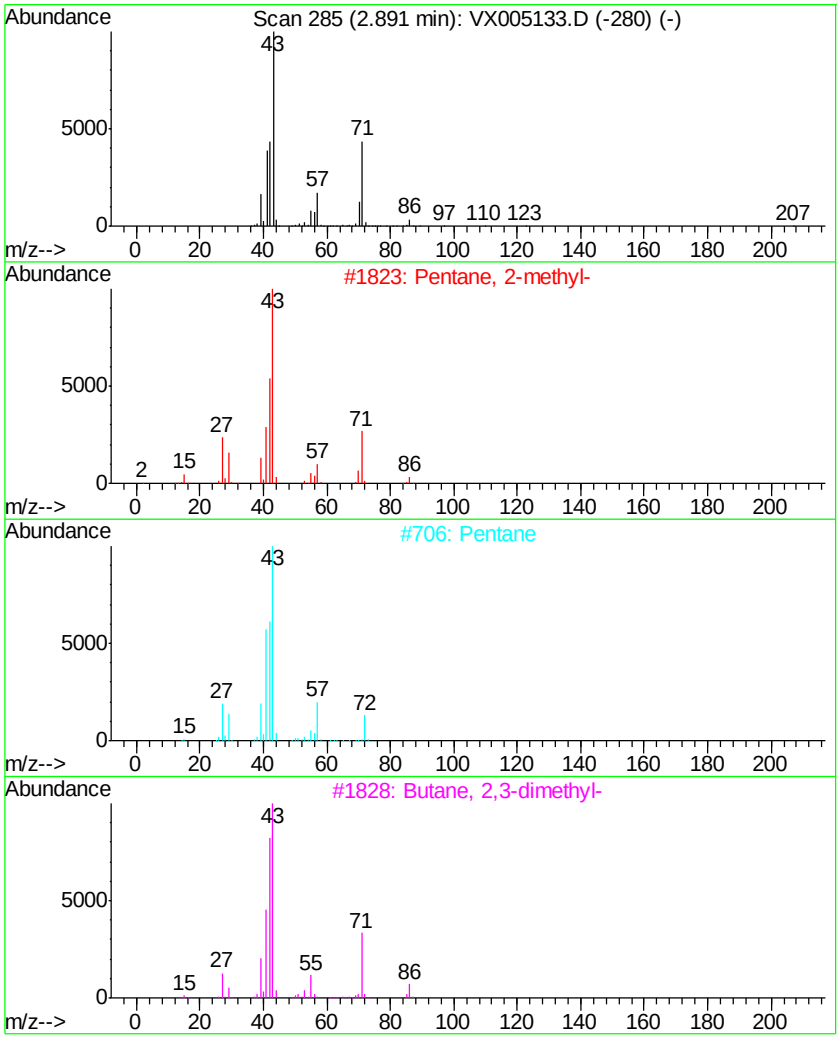
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

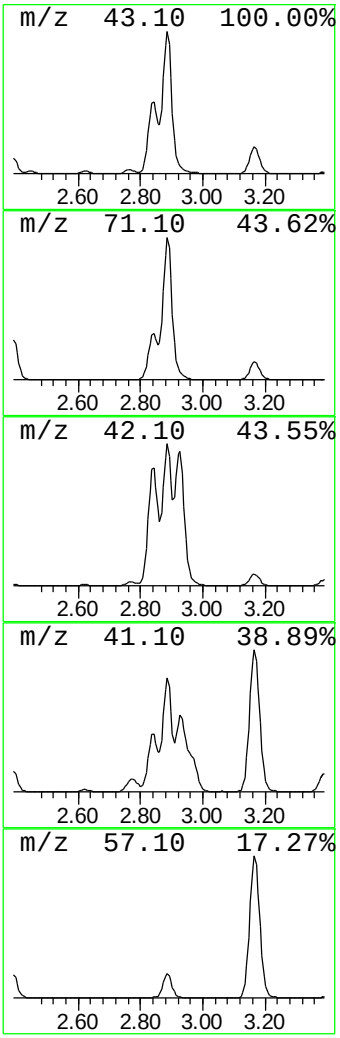
****

Peak Number 4 Pentane, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.89	378.92 ug/l	2542860	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2-methyl-	86	C6H14	000107-83-5	90
2	Pentane	72	C5H12	000109-66-0	43
3	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	43
4	Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	28
5	Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	23





Library Search Compound Report

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Sample : J5268-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

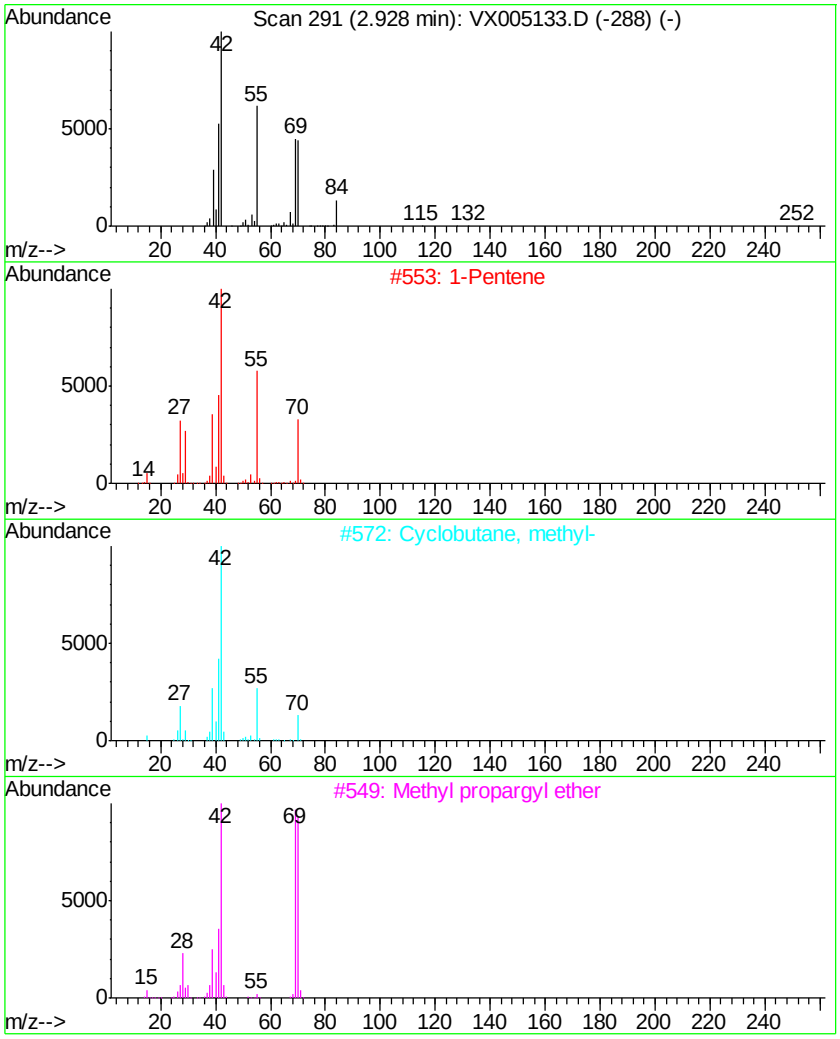
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

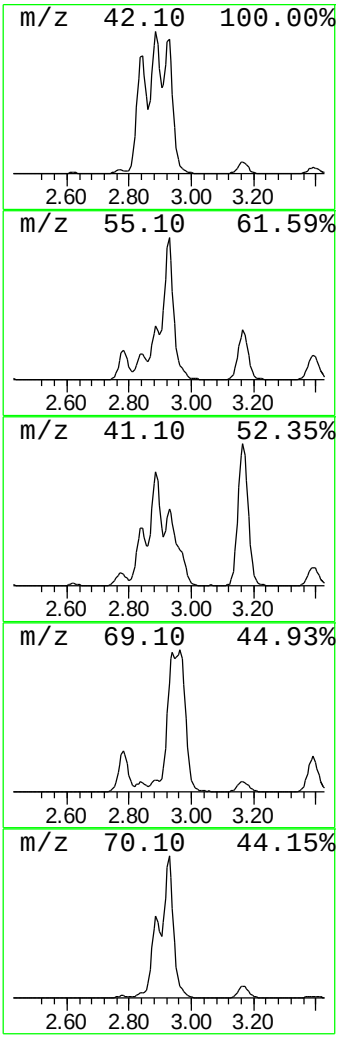
****

Peak Number 5 1-Pentene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.93	238.69 ug/l	1601830	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentene	70	C5H10	000109-67-1	80
2	Cyclobutane, methyl-	70	C5H10	000598-61-8	72
3	Methyl propargyl ether	70	C4H6O	000627-41-8	37
4	Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	32
5	(Z)-1,3-Butadien-1-ol	70	C4H6O	070415-58-6	27





Library Search Compound Report

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ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

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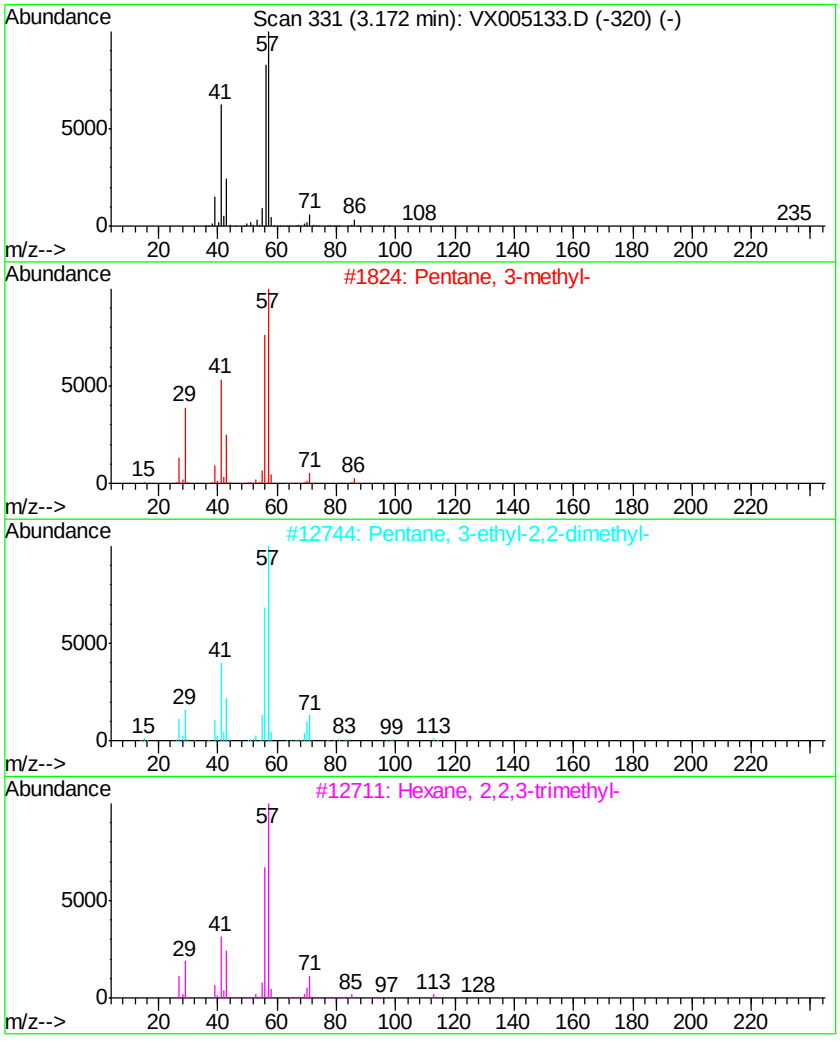
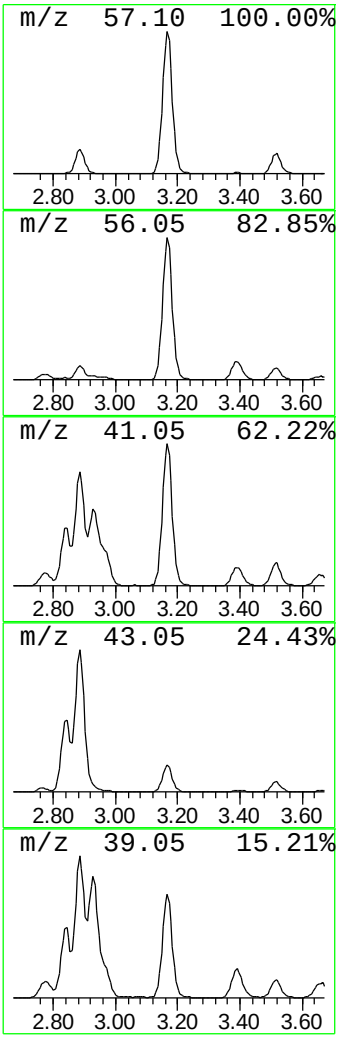
Peak Number 6 Pentane, 3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.17	358.47 ug/l	2405650	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64
3	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	64
4	Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	50
5	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	45





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100818\
Data File : VX005133.D
Acq On : 08 Oct 2018 17:49
Operator : JC/MD
Sample : J5268-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

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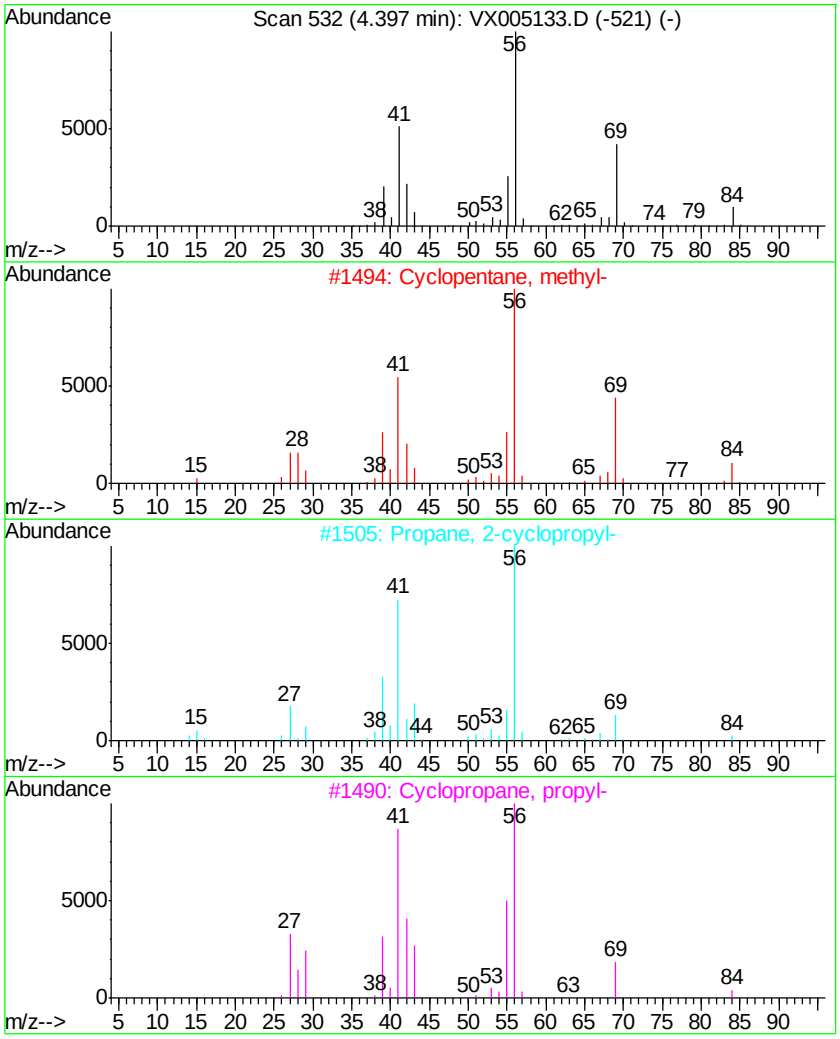
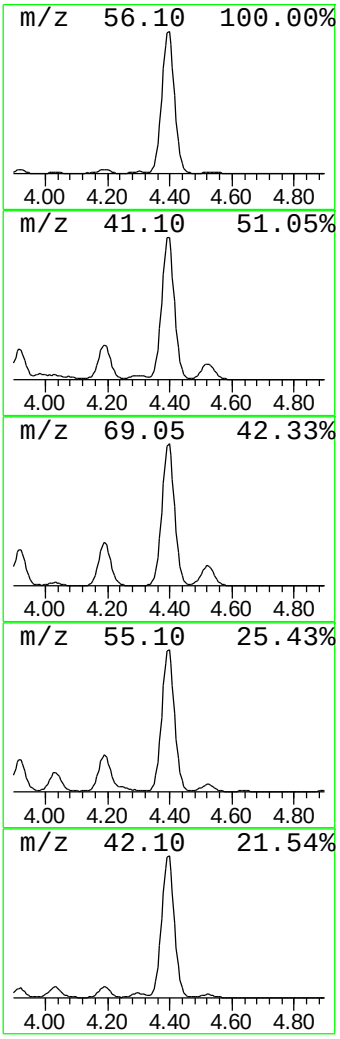
Peak Number 7 Cyclopentane, methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.40	684.82 ug/l	4595710	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2	Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	80
3	Cyclopropane, propyl-	84	C6H12	002415-72-7	78
4	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
5	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100818\
Data File : VX005133.D
Acq On : 08 Oct 2018 17:49
Operator : JC/MD
Sample : J5268-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

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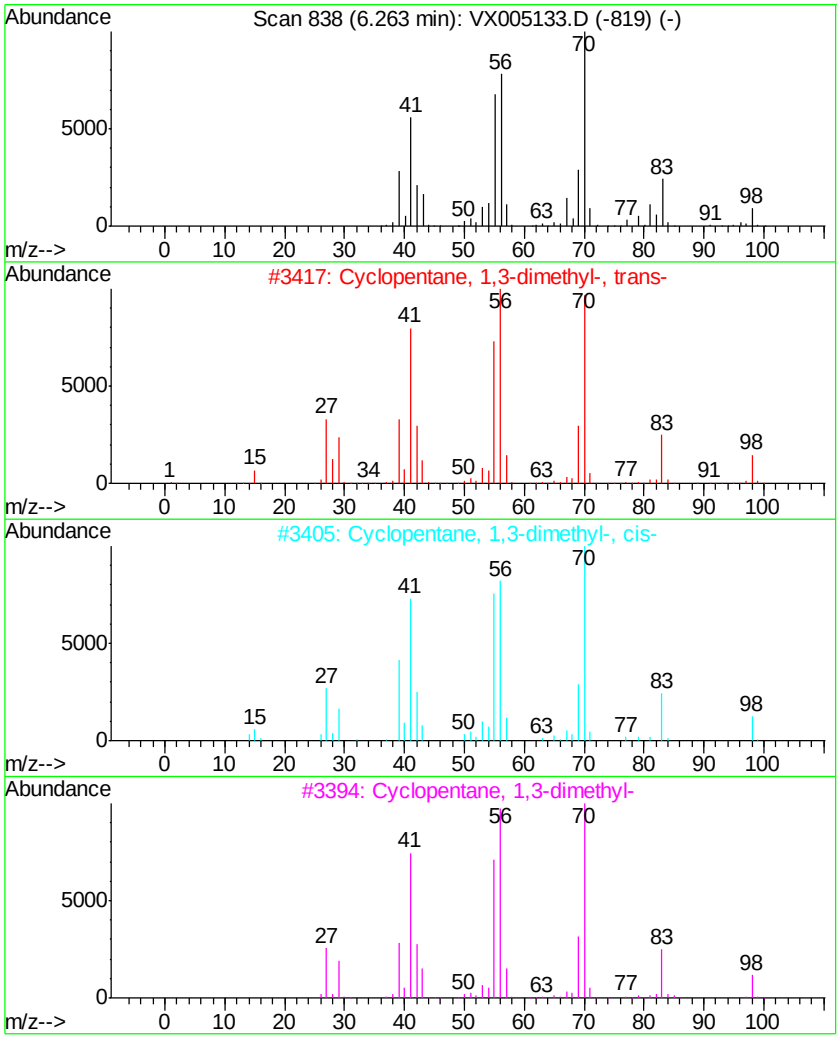
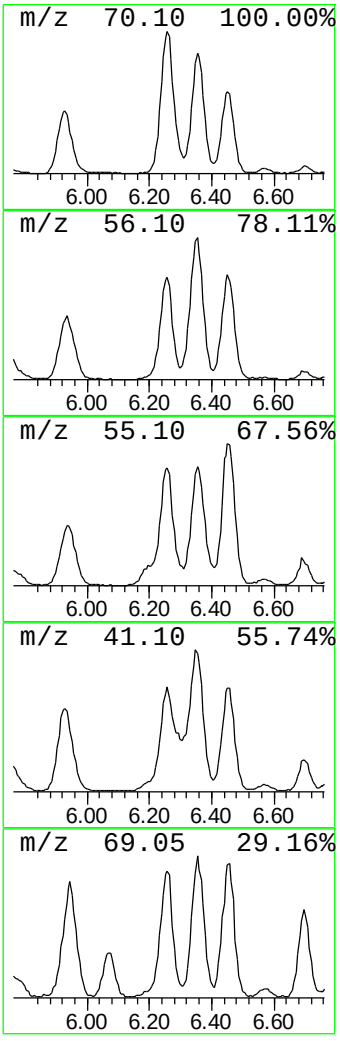
Peak Number 8 Cyclopentane, 1,3-dimethyl-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.26	138.02 ug/l	926248	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	94
2	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	94
3	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	93
4	Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	87
5	1-Pentene, 3,4-dimethyl-	98	C7H14	007385-78-6	64





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100818\
Data File : VX005133.D
Acq On : 08 Oct 2018 17:49
Operator : JC/MD
Sample : J5268-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

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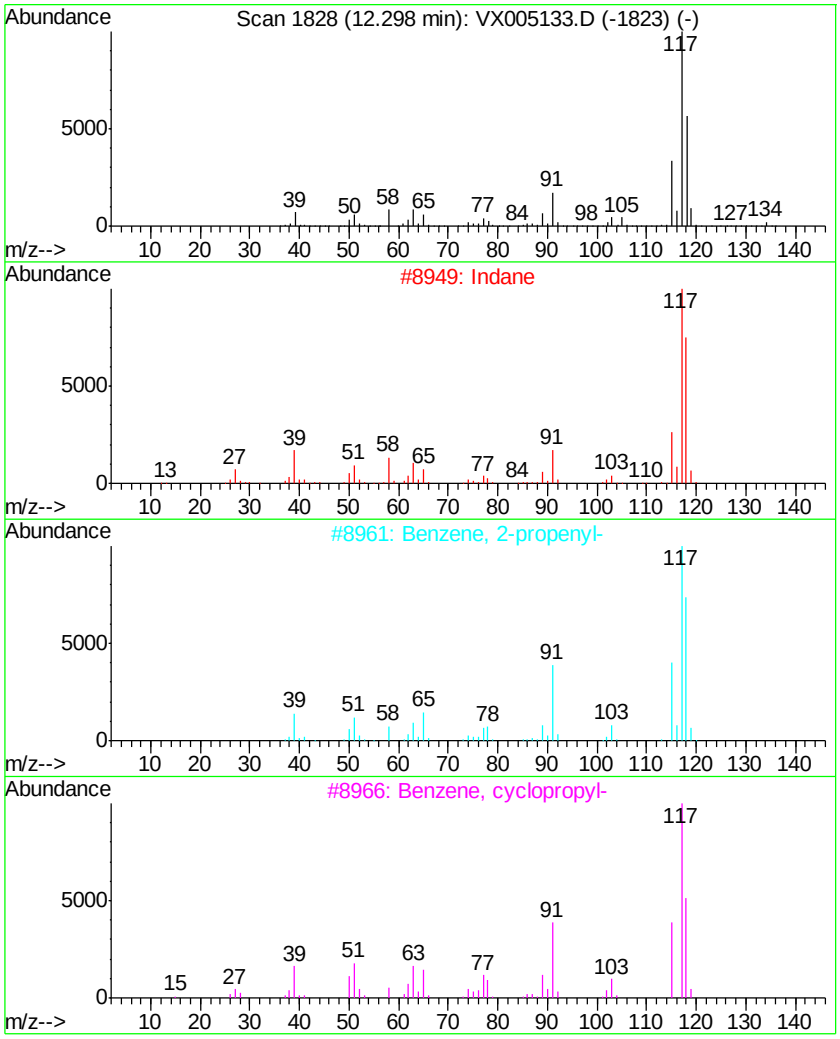
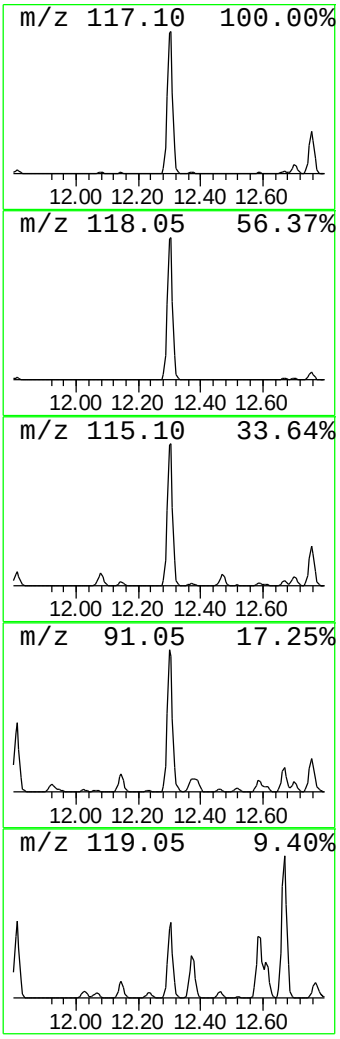
Peak Number 9 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	521.69 ug/l	12317900	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	91
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	87
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	81
4	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	74
5	Deltacyclene	118	C9H10	007785-10-6	64





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100818\
Data File : VX005133.D
Acq On : 08 Oct 2018 17:49
Operator : JC/MD
Sample : J5268-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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

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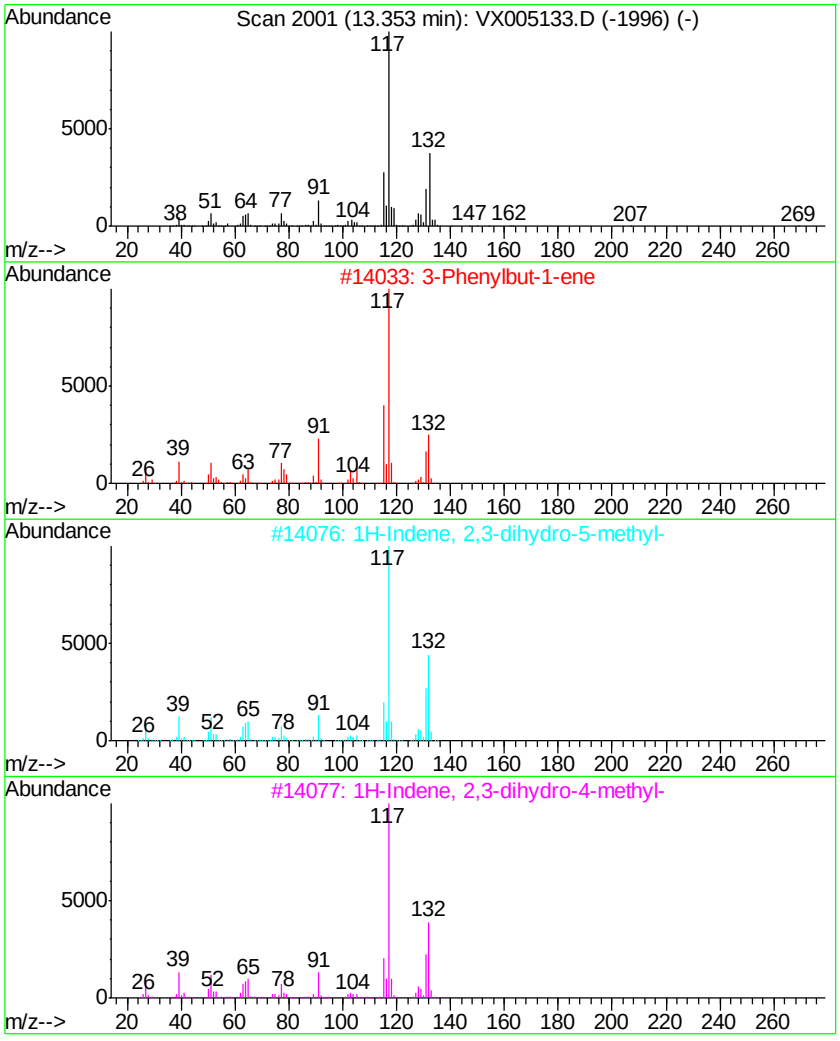
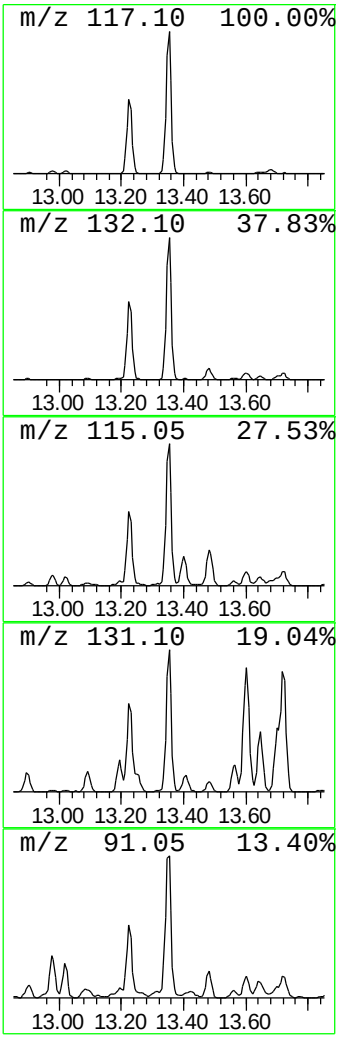
Peak Number 10 3-Phenylbut-1-ene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	206.50 ug/l	4875680	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	86
5	Indan, 1-methyl-	132	C10H12	000767-58-8	81





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100818\
Data File : VX005133.D
Acq On : 08 Oct 2018 17:49
Operator : JC/MD
Sample : J5268-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.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-	1.78	927.1	ug/l	6221880	1	5.67	335543	50.0
Pentane	2.00	275.1	ug/l	1846420	1	5.67	335543	50.0
Butane, 2,3-dimet...	2.84	209.3	ug/l	1404910	1	5.67	335543	50.0
Pentane, 2-methyl-	2.89	378.9	ug/l	2542860	1	5.67	335543	50.0
1-Pentene	2.93	238.7	ug/l	1601830	1	5.67	335543	50.0
Pentane, 3-methyl-	3.17	358.5	ug/l	2405650	1	5.67	335543	50.0
Cyclopentane, met...	4.40	684.8	ug/l	4595710	1	5.67	335543	50.0
Cyclopentane, 1,3...	6.26	138.0	ug/l	926248	1	5.67	335543	50.0
Indane	12.30	521.7	ug/l	12317900	4	12.08	1180580	50.0
3-Phenylbut-1-ene	13.35	206.5	ug/l	4875680	4	12.08	1180580	50.0