

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU102318\
 Data File : VU027758.D
 Acq On : 23 Oct 2018 11:21
 Operator : MD/SY
 Sample : J5026-06DL 4X
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 1721DL

Quant Time: Oct 24 05:42:45 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA U\METHOD\524U0102318DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Tue Oct 23 09:12:15 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
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1) Fluorobenzene	5.75	96	13315	1.00	ug/l	0.00
System Monitoring Compounds						
57) 4-Bromofluorobenzene	10.31	95	5123	1.01	ug/l	0.00
Spiked Amount 1.000			Recovery	=	101.00%	
68) 1,2-Dichlorobenzene-d4	11.86	152	4837	1.01	ug/l	0.00
Spiked Amount 1.000			Recovery	=	101.00%	
Target Compounds					Ovalue	
35) Benzene	5.39	78	12647	0.820	ug/l	100
49) Toluene	7.64	92	29320	3.202	ug/l	100
59) Chlorobenzene	9.12	112	30971	3.156	ug/l	99
63) Ethyl Benzene	9.25	91	77964	4.194	ug/l	99
64) m/p-Xylenes	9.38	106	16840	2.539	ug/l	99
65) o-Xylene	9.78	106	16528	2.513	ug/l	99
66) Styrene	9.79	104	47499	4.377	ug/l	96
69) Isopropylbenzene	10.17	105	28118	1.562	ug/l	99
73) n-propylbenzene	10.59	120	15751	3.392	ug/l	95
75) 1,3,5-Trimethylbenzene	10.77	105	68452	4.557	ug/l	99
77) tert-Butylbenzene	11.10	119	39903	2.701	ug/l	92
78) 1,2,4-Trimethylbenzene	11.15	105	18660	1.209	ug/l	100
79) sec-Butylbenzene	11.32	105	84289	4.308	ug/l	99
81) p-Isopropyltoluene	11.48	119	19613	1.226	ug/l	99
82) 1,3-Dichlorobenzene	11.42	146	34617	4.361	ug/l	99
83) 1,4-Dichlorobenzene	11.51	146	28466	3.579	ug/l	98
85) 1,2-Dichlorobenzene	11.88	146	28289	3.735	ug/l	99
87) 1,2,4-Trichlorobenzene	13.50	180	24353	4.425	ug/l	99
90) 1,2,3-Trichlorobenzene	13.99	180	21143	4.037	ug/l	99
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(#) = qualifier out of range (m) = manual integration (+) = signals summed

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