

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092018\
 Data File : VU026910.D
 Acq On : 19 Sep 2018 20:25
 Operator : MD/SY
 Sample : J4934-04
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 FT-DUP03-20180912

Quant Time: Sep 20 07:16:50 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U091218W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 12 05:32:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	81585	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	137317	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	139935	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	75870	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	70039	56.59	ug/l	0.00
Spiked Amount 50.000			Recovery	=	113.18%	
35) Dibromofluoromethane	4.88	113	62131	55.47	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.94%	
50) Toluene-d8	7.57	98	208219	50.26	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.52%	
62) 4-Bromofluorobenzene	10.31	95	69984	46.61	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.22%	
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	2.28	96	98994	105.690	ug/l	100
16) Acetone	2.32	43	1600	2.283	ug/l	95
24) 1,1-Dichloroethane	3.45	63	80950	40.759	ug/l	97
27) cis-1,2-Dichloroethene	4.23	96	5715	5.025	ug/l	90
44) Trichloroethene	6.19	130	138814	117.144	ug/l	94
64) Tetrachloroethene	8.23	164	3772	3.387	ug/l	94

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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