

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092018\
 Data File : VU026909.D
 Acq On : 19 Sep 2018 20:02
 Operator : MD/SY
 Sample : J4934-03
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 FT-PZ460I-20180912

Quant Time: Sep 20 07:15:11 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	79017	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	133962	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	132780	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	72419	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	66986	55.88	ug/l	0.00
Spiked Amount 50.000			Recovery	=	111.76%	
35) Dibromofluoromethane	4.89	113	59941	54.85	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.70%	
50) Toluene-d8	7.57	98	199854	49.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.90%	
62) 4-Bromofluorobenzene	10.31	95	66767	45.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.16%	
Target Compounds						
12) 1,1-Dichloroethene	2.28	96	98910	109.033	ug/l	99
16) Acetone	2.32	43	1886	2.779	ug/l	96
24) 1,1-Dichloroethane	3.45	63	81256	42.242	ug/l	98
27) cis-1,2-Dichloroethene	4.23	96	5669	5.147	ug/l	86
44) Trichloroethene	6.19	130	135751	117.428	ug/l	94
64) Tetrachloroethene	8.23	164	3624	3.429	ug/l	95

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

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