

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092618\
 Data File : VU027205.D
 Acq On : 27 Sep 2018 02:44
 Operator : MD/SY
 Sample : J5048-08
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 DUP-1-SEP18

Quant Time: Sep 27 04:32:29 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U092218W.M
 Quant Title : SW846 8260
 QLast Update : Sat Sep 22 01:31:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	115503	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	192144	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	181264	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	84705	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	110670	47.61	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.22%	
35) Dibromofluoromethane	4.88	113	78970	47.41	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.82%	
50) Toluene-d8	7.57	98	302792	47.70	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.40%	
62) 4-Bromofluorobenzene	10.31	95	99659	45.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.52%	
Target Compounds						
16) Acetone	2.32	43	161958	122.709	ug/l	99
27) cis-1,2-Dichloroethene	4.23	96	12035	6.257	ug/l	89
44) Trichloroethene	6.19	130	38279	23.065	ug/l	100
64) Tetrachloroethene	8.23	164	43601	31.216	ug/l	99

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

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