

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU071018\
 Data File : VU025291.D
 Acq On : 10 Jul 2018 16:04
 Operator : MD/SY
 Sample : J3905-06
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 SUMP-070518

Quant Time: Jul 11 03:27:06 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U070918W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jul 10 09:34:16 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	204910	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	307650	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	284859	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	145182	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	152796	48.00	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.00%	
35) Dibromofluoromethane	4.89	113	121244	46.72	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.44%	
50) Toluene-d8	7.57	98	439796	49.49	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.98%	
62) 4-Bromofluorobenzene	10.31	95	163072	45.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.16%	
Target Compounds						
16) Acetone	2.32	43	6750	8.94	ug/l	Qvalue 100
27) cis-1,2-Dichloroethene	4.23	96	16857	6.14	ug/l	94
30) Chloroform	4.68	83	2972	0.65	ug/l	89
44) Trichloroethene	6.20	130	2638	0.93	ug/l	87
64) Tetrachloroethene	8.23	164	120632	46.46	ug/l	98

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

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