

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU070318\
 Data File : VU025192.D
 Acq On : 03 Jul 2018 15:41
 Operator : MD/SY
 Sample : J3790-02
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 ZER-WC-L-01

Quant Time: Jul 05 09:03:20 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 13 13:55:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	206321	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	306868	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	277995	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	145448	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	155639	46.21	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.42%	
35) Dibromofluoromethane	4.89	113	120553	47.34	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.68%	
50) Toluene-d8	7.57	98	435143	46.84	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.68%	
62) 4-Bromofluorobenzene	10.31	95	161735	43.85	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.70%	
Target Compounds						
16) Acetone	2.32	43	18212	10.063	ug/l	97
17) Carbon Disulfide	2.48	76	4794	0.637	ug/l	95
30) Chloroform	4.68	83	79217	15.734	ug/l	97
47) Bromodichloromethane	6.76	83	4330	1.129	ug/l #	87
84) 1,2,4-Trimethylbenzene	11.16	105	3093	0.314	ug/l	88

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

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