

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU071218\
 Data File : VU025372.D
 Acq On : 13 Jul 2018 05:15
 Operator : MD/SY
 Sample : J3828-06
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EO-01-071018-C

Quant Time: Jul 13 06:03:30 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	186869	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	287033	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	262851	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	139179	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	144881	49.90	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.80%	
35) Dibromofluoromethane	4.89	113	112460	46.44	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.88%	
50) Toluene-d8	7.57	98	410364	49.49	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.98%	
62) 4-Bromofluorobenzene	10.31	95	151245	45.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.62%	
Target Compounds						
16) Acetone	2.32	43	22674	19.874	ug/l	99
18) Methyl Acetate	2.62	43	17831	5.999	ug/l	99
20) Methylene Chloride	2.71	84	10158	5.152	ug/l	93
95) Naphthalene	13.75	128	11808	3.513	ug/l	96

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

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