

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU071218\
 Data File : VU025371.D
 Acq On : 13 Jul 2018 04:51
 Operator : MD/SY
 Sample : J3828-04
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 43 Sample Multiplier: 1

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

Quant Time: Jul 13 05:59:37 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	189923	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	292558	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	271405	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	143610	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	146804	49.75	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.50%	
35) Dibromofluoromethane	4.89	113	113798	46.11	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.22%	
50) Toluene-d8	7.57	98	422198	49.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.92%	
62) 4-Bromofluorobenzene	10.31	95	154188	45.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.64%	
Target Compounds						
16) Acetone	2.32	43	19221	17.384	ug/l	99
18) Methyl Acetate	2.62	43	18577	6.149	ug/l	98
20) Methylene Chloride	2.71	84	135642	59.290	ug/l	99
95) Naphthalene	13.75	128	18002	4.038	ug/l	97

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

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