

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU071218\
 Data File : VU025380.D
 Acq On : 13 Jul 2018 08:28
 Operator : MD/SY
 Sample : J3825-04
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 CL-02-071118-B

Quant Time: Jul 13 09:12:29 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	187732	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	293420	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	276447	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	145339	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	146721	50.30	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.60%	
35) Dibromofluoromethane	4.89	113	114091	46.09	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.18%	
50) Toluene-d8	7.57	98	422225	49.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.62%	
62) 4-Bromofluorobenzene	10.31	95	154167	45.18	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.36%	
Target Compounds						
16) Acetone	2.32	43	20197	18.173	ug/l	97
18) Methyl Acetate	2.62	43	19512	6.534	ug/l	99
20) Methylene Chloride	2.71	84	11723	5.815	ug/l	96
95) Naphthalene	13.75	128	137448	14.672	ug/l	100

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

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