

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU061818\
 Data File : VU024706.D
 Acq On : 18 Jun 2018 21:04
 Operator : MD/SY
 Sample : J3477-11
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 S36-EB-2018-1

Quant Time: Jun 19 07:29:36 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	193070	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	307218	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	286901	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	154286	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	140262	44.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.00%	
35) Dibromofluoromethane	4.89	113	120240	47.16	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.32%	
50) Toluene-d8	7.57	98	449096	48.29	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.58%	
62) 4-Bromofluorobenzene	10.31	95	161446	43.72	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.44%	
Target Compounds						
16) Acetone	2.32	43	13152	7.77	ug/l	# 88
20) Methylene Chloride	2.71	84	112225	42.23	ug/l	97
25) 2-Butanone	4.27	43	27344	12.02	ug/l	94
29) Tetrahydrofuran	4.64	42	1168404	829.42	ug/l	98

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

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