

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU111218\
 Data File : VU028076.D
 Acq On : 12 Nov 2018 19:46
 Operator : MD/SY
 Sample : J5887-12
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 W-2

Quant Time: Nov 13 03:58:23 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 02 02:28:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.98	168	169627	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	319933	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	283116	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	113239	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	148052	58.71	ug/l	0.00
Spiked Amount 50.000			Recovery	=	117.42%	
35) Dibromofluoromethane	4.89	113	106662	52.88	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.76%	
50) Toluene-d8	7.57	98	415673	53.34	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.68%	
62) 4-Bromofluorobenzene	10.31	95	134595	46.83	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.66%	
Target Compounds						
16) Acetone	2.32	43	8841	5.650	ug/l	100
18) Methyl Acetate	2.62	43	7975	1.853	ug/l	98
20) Methylene Chloride	2.70	84	25958	11.362	ug/l	98
43) Isopropyl Acetate	5.55	43	206531	26.850	ug/l	99

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

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