

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU111318\
 Data File : VU028101.D
 Acq On : 13 Nov 2018 18:55
 Operator : MD/SY
 Sample : J5909-12
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 SB-17-S

Quant Time: Nov 14 08:02:23 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	4.99	168	258509	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	479072	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	425783	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	176553	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	201000	52.30	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.60%	
35) Dibromofluoromethane	4.88	113	151244	50.07	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.14%	
50) Toluene-d8	7.57	98	610262	52.30	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.60%	
62) 4-Bromofluorobenzene	10.31	95	202301	47.00	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.00%	
Target Compounds						
16) Acetone	2.32	43	32083	13.454	ug/l	94
18) Methyl Acetate	2.62	43	10812	1.648	ug/l	96
20) Methylene Chloride	2.70	84	140702	40.413	ug/l	95
25) 2-Butanone	4.30	43	4710	1.242	ug/l	96
43) Isopropyl Acetate	5.55	43	254134	22.064	ug/l	98

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

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