

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU112918\
 Data File : VU028387.D
 Acq On : 29 Nov 2018 12:41
 Operator : MD/SY
 Sample : J6124-01
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BW-2072-112818

Manual Integrations
 APPROVED

MMDadoda
 11/30/2018 11:25:41 AM

Quant Time: Nov 30 07:26:48 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U112818W.M
 Quant Title : SW846 8260
 QLast Update : Thu Nov 29 01:04:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.98	168	243675	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	460594	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	426619	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	193411	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	213788	52.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.00%	
35) Dibromofluoromethane	4.88	113	155543	50.79	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.58%	
50) Toluene-d8	7.57	98	607393	49.59	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.18%	
62) 4-Bromofluorobenzene	10.31	95	218478	49.49	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.98%	
Target Compounds						
16) Acetone	2.32	43	82868	33.264	ug/l	97
20) Methylene Chloride	2.70	84	4589	1.234	ug/l	97
25) 2-Butanone	4.29	43	8848m	2.478	ug/l	
29) Tetrahydrofuran	4.65	42	12807	5.525	ug/l	94
39) Methylcyclohexane	6.42	83	6138	1.033	ug/l	96
67) Ethyl Benzene	9.25	91	26851	1.609	ug/l	97
80) 1,3,5-Trimethylbenzene	10.77	105	50054	3.949	ug/l	98

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

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