

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU090419\
 Data File : VU034219.D
 Acq On : 03 Sep 2019 11:55
 Operator : JC/SP
 Sample : K4646-19
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EB-7-083019

Manual Integrations
 APPROVED

MMDadoda
 9/4/2019 4:57:34 PM

Quant Time: Sep 04 05:07:30 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U082919W.M
 Quant Title : SW846 8260
 QLast Update : Thu Aug 29 04:11:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.96	168	432589	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.87	114	698275	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.07	117	672388	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.46	152	336046	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.29	65	273003	49.25	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.50%	
35) Dibromofluoromethane	4.86	113	229961	47.48	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.96%	
50) Toluene-d8	7.55	98	888922	49.02	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.04%	
62) 4-Bromofluorobenzene	10.29	95	308214	43.69	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.38%	
Target Compounds						
16) Acetone	2.31	43	14207	5.023	ug/l #	84
25) 2-Butanone	4.27	43	15126m	3.844	ug/l	
51) 4-Methyl-2-Pentanone	7.45	43	4576	0.589	ug/l #	61

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

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