

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU090719\
 Data File : VU034338.D
 Acq On : 06 Sep 2019 18:07
 Operator : JC/SP
 Sample : K4668-03
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 SU-04-090319

Quant Time: Sep 06 18:57:18 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U090619W.M
 Quant Title : SW846 8260
 QLast Update : Fri Sep 06 05:10:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.96	168	306444	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.86	114	469083	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.07	117	468727	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.46	152	283834	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.29	65	240695	51.91	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.82%	
35) Dibromofluoromethane	4.86	113	212527	48.69	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.38%	
50) Toluene-d8	7.55	98	813720	49.62	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.24%	
62) 4-Bromofluorobenzene	10.29	95	295070	46.60	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.20%	
Target Compounds						
16) Acetone	2.31	43	17022	7.094	ug/l	100
20) Methylene Chloride	2.69	84	27428	6.632	ug/l	94
43) Isopropyl Acetate	5.53	43	93344	9.948	ug/l	99
95) Naphthalene	13.73	128	83933	8.051	ug/l	97

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

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