

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU090619\
 Data File : VU034293.D
 Acq On : 05 Sep 2019 19:01
 Operator : JC/SP
 Sample : K4545-11
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 T16-17-E1-(2.25)

Quant Time: Sep 06 08:23:49 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	451053	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.87	114	707238	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.07	117	689592	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.46	152	339053	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.29	65	277058	40.60	ug/l	0.00
Spiked Amount 50.000			Recovery	=	81.20%	
35) Dibromofluoromethane	4.86	113	242675	36.88	ug/l	0.00
Spiked Amount 50.000			Recovery	=	73.76%	
50) Toluene-d8	7.55	98	962744	38.94	ug/l	0.00
Spiked Amount 50.000			Recovery	=	77.88%	
62) 4-Bromofluorobenzene	10.29	95	340485	35.67	ug/l	0.00
Spiked Amount 50.000			Recovery	=	71.34%	
Target Compounds						
16) Acetone	2.32	43	12374	3.504	ug/l	92
43) Isopropyl Acetate	5.53	43	76324	5.395	ug/l	99
95) Naphthalene	13.75	128	12604	4.479	ug/l #	88
96) 1,2,3-Trichlorobenzene	13.98	180	4269	3.160	ug/l	86

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

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