

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU091019\
 Data File : VU034445.D
 Acq On : 10 Sep 2019 12:52
 Operator : JC/SP
 Sample : K4602-35
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 PAR-D2-(0)

Quant Time: Sep 11 06:38:27 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	240776	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.87	114	373885	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.07	117	391381	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.46	152	214429	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.29	65	190755	52.36	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.72%	
35) Dibromofluoromethane	4.86	113	167747	48.22	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.44%	
50) Toluene-d8	7.54	98	619200	47.37	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.74%	
62) 4-Bromofluorobenzene	10.29	95	210753	41.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	83.52%	
Target Compounds						
16) Acetone	2.31	43	12459	6.609	ug/l	97
20) Methylene Chloride	2.69	84	6078	1.581	ug/l	95
43) Isopropyl Acetate	5.53	43	60153	8.043	ug/l	99
95) Naphthalene	13.75	128	9400	4.572	ug/l #	89

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

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