

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU082819\
 Data File : VU033991.D
 Acq On : 27 Aug 2019 22:41
 Operator : JC/SP
 Sample : K4516-28
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 T1-S-E-(0)

Quant Time: Aug 28 08:43:23 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U082819W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 28 02:42:03 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.96	168	435448	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.86	114	698233	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.07	117	679084	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.47	152	297212	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.29	65	282685	58.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	116.86%	
35) Dibromofluoromethane	4.86	113	231596	49.34	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.68%	
50) Toluene-d8	7.55	98	921810	53.25	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.50%	
62) 4-Bromofluorobenzene	10.29	95	278811	41.98	ug/l	0.00
Spiked Amount 50.000			Recovery	=	83.96%	
Target Compounds						
16) Acetone	2.31	43	8797	3.799	ug/l	99
18) Methyl Acetate	2.61	43	8709	1.527	ug/l	91
20) Methylene Chloride	2.69	84	17781	3.619	ug/l #	86
43) Isopropyl Acetate	5.53	43	85745	8.718	ug/l	96

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

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