

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU082819\
 Data File : VU034006.D
 Acq On : 28 Aug 2019 05:51
 Operator : JC/SP
 Sample : K4516-39
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 T2-S-D-(2.5)

Quant Time: Aug 28 08:45:29 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	437752	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.86	114	695766	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.07	117	677880	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.47	152	296264	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.29	65	299249	61.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	123.06%	
35) Dibromofluoromethane	4.86	113	235605	50.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.76%	
50) Toluene-d8	7.55	98	914852	53.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.06%	
62) 4-Bromofluorobenzene	10.29	95	299989	45.33	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.66%	
Target Compounds						
16) Acetone	2.31	43	12606	5.415	ug/l	100
18) Methyl Acetate	2.60	43	7594	1.324	ug/l	90
20) Methylene Chloride	2.68	84	11566	2.342	ug/l	89
43) Isopropyl Acetate	5.53	43	87573	8.936	ug/l	94

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

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