

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100418\
 Data File : VU027389.D
 Acq On : 04 Oct 2018 04:27
 Operator : MD/SY
 Sample : J5194-04
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 HR-06-100218-B

Quant Time: Oct 04 08:39:43 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
 Quant Title : SW846 8260
 QLast Update : Tue Oct 02 02:04:34 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	97368	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	166571	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	163417	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	85639	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	97806	60.16	ug/l	0.00
Spiked Amount 50.000			Recovery	=	120.32%	
35) Dibromofluoromethane	4.89	113	68560	60.06	ug/l	0.00
Spiked Amount 50.000			Recovery	=	120.12%	
50) Toluene-d8	7.57	98	261266	61.15	ug/l	0.00
Spiked Amount 50.000			Recovery	=	122.30%	
62) 4-Bromofluorobenzene	10.31	95	91731	57.77	ug/l	0.00
Spiked Amount 50.000			Recovery	=	115.54%	
Target Compounds						
16) Acetone	2.32	43	14003	14.579	ug/l	97
18) Methyl Acetate	2.62	43	3524	1.608	ug/l #	88
20) Methylene Chloride	2.70	84	27652	20.668	ug/l	98
30) Chloroform	4.68	83	2592	1.143	ug/l	89
43) Isopropyl Acetate	5.55	43	94783	26.109	ug/l	98
84) 1,2,4-Trimethylbenzene	11.15	105	6383	1.234	ug/l	94
95) Naphthalene	13.75	128	75863	12.988	ug/l	99

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

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