

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU110218\
 Data File : VU027926.D
 Acq On : 02 Nov 2018 15:14
 Operator : MD/SY
 Sample : J5741-10
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 TP-2

Quant Time: Nov 03 05:39:10 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 02 02:28:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	187584	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	330858	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	295490	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	121838	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	145321	52.11	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.22%	
35) Dibromofluoromethane	4.89	113	107734	51.64	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.28%	
50) Toluene-d8	7.57	98	423116	52.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.02%	
62) 4-Bromofluorobenzene	10.31	95	138146	46.48	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.96%	
Target Compounds						
16) Acetone	2.32	43	6660	3.849	ug/l	96
18) Methyl Acetate	2.62	43	5699	1.197	ug/l	92
20) Methylene Chloride	2.70	84	24985	9.890	ug/l	99
43) Isopropyl Acetate	5.55	43	208846	26.255	ug/l	100
95) Naphthalene	13.75	128	31618	2.993	ug/l	99

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

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