

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU062818\
 Data File : VU025030.D
 Acq On : 28 Jun 2018 19:07
 Operator : MD/SY
 Sample : J3755-01
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 PS-BASELINEWATER

Quant Time: Jun 29 11:09:37 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 13 13:55:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	221450	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	332890	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	302399	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	167605	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	5.31	65	168326	46.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.12%	
35) Dibromofluoromethane	4.89	113	129214	46.77	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.54%	
50) Toluene-d8	7.57	98	478792	47.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.02%	
62) 4-Bromofluorobenzene	10.31	95	178055	44.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.00%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
11) Tert butyl alcohol	2.82	59	28279	32.16	ug/l	# 92
16) Acetone	2.32	43	3393928	1747.24	ug/l	99
25) 2-Butanone	4.26	43	196315	75.24	ug/l	97
51) 4-Methyl-2-Pentanone	7.46	43	25848	5.58	ug/l	97

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

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