

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092218\
 Data File : VU027018.D
 Acq On : 22 Sep 2018 16:11
 Operator : MD/SY
 Sample : J4958-05
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BPSI-HN-MW24IR-20180914

Quant Time: Sep 24 04:59:22 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U092218W.M
 Quant Title : SW846 8260
 QLast Update : Sat Sep 22 01:31:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	96824	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	160378	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	154578	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	70954	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	96598	49.57	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.14%	
35) Dibromofluoromethane	4.88	113	67246	48.37	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.74%	
50) Toluene-d8	7.57	98	248212	46.85	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.70%	
62) 4-Bromofluorobenzene	10.31	95	83386	45.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.74%	
Target Compounds						
7) Trichlorofluoromethane	1.89	101	6106	2.642	ug/l	98
12) 1,1-Dichloroethene	2.29	96	1920	1.436	ug/l	94
19) Methyl tert-butyl Ether	3.00	73	7320	1.490	ug/l	94
44) Trichloroethene	6.19	130	10541	7.609	ug/l	95
64) Tetrachloroethene	8.23	164	7652	6.424	ug/l	93

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

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