

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092618\
 Data File : VU027168.D
 Acq On : 26 Sep 2018 11:33
 Operator : MD/SY
 Sample : J4969-01DL 5X
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 SA-PZ171I-20180916DL

Quant Time: Sep 27 01:57:56 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	97197	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	166143	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	158408	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	71870	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	102188	52.24	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.48%	
35) Dibromofluoromethane	4.89	113	70881	49.22	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.44%	
50) Toluene-d8	7.57	98	265544	48.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.76%	
62) 4-Bromofluorobenzene	10.31	95	86478	45.92	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.84%	
Target Compounds						
6) Chloroethane	1.70	64	5266	4.035	ug/l	96
12) 1,1-Dichloroethene	2.28	96	4945	3.684	ug/l	94
24) 1,1-Dichloroethane	3.45	63	158403	49.385	ug/l	99
32) 1,1,1-Trichloroethane	4.92	97	13723	5.773	ug/l	96

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

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