

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092418\
 Data File : VU027100.D
 Acq On : 24 Sep 2018 11:53
 Operator : MD/SY
 Sample : J4969-02
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 SA-PZ172I-20180916

Quant Time: Sep 25 02:04:23 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	97340	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	165713	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	157997	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	73966	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	99904	50.99	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.98%	
35) Dibromofluoromethane	4.88	113	68899	47.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.94%	
50) Toluene-d8	7.57	98	266215	48.63	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.26%	
62) 4-Bromofluorobenzene	10.31	95	85773	45.66	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.32%	
Target Compounds						
6) Chloroethane	1.70	64	8398	6.426	ug/l	99
12) 1,1-Dichloroethene	2.29	96	4993	3.714	ug/l	88
24) 1,1-Dichloroethane	3.45	63	181499	56.502	ug/l	99
32) 1,1,1-Trichloroethane	4.92	97	6002	2.521	ug/l	97

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

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