

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062718\
 Data File : VX002969.D
 Acq On : 28 Jun 2018 06:38
 Operator : JC/MD
 Sample : J3718-10
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 W-33

Quant Time: Jul 07 02:07:15 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X062518W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jun 25 14:38:05 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	274858	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.87	114	386228	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	361159	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	211591	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	184060	48.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.90%	
35) Dibromofluoromethane	5.51	113	150117	49.08	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.16%	
50) Toluene-d8	8.72	98	543690	48.68	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.36%	
62) 4-Bromofluorobenzene	11.14	95	197804	45.92	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.84%	
Target Compounds						
10) Methyl Iodide	2.52	142	2843	7.034	ug/l	98
11) Tert butyl alcohol	3.07	59	7854	10.378	ug/l	97
16) Acetone	2.45	43	7657	4.867	ug/l	97
19) Methyl tert-butyl Ether	3.21	73	40850	4.192	ug/l	95
22) Diisopropyl ether	3.88	45	4698	0.460	ug/l #	77
65) Chlorobenzene	10.14	112	72776	7.933	ug/l	96

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

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