

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX101718\
 Data File : VX005409.D
 Acq On : 18 Oct 2018 00:14
 Operator : JC/MD
 Sample : J5569-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RLF-GW-9S

Quant Time: Oct 19 02:27:33 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.M
 Quant Title : SW846 8260
 QLast Update : Fri Oct 12 11:28:43 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	275056	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	394563	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	354589	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	174392	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	154354	50.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.12%	
35) Dibromofluoromethane	5.50	113	125657	48.08	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.16%	
50) Toluene-d8	8.71	98	471554	52.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.12%	
62) 4-Bromofluorobenzene	11.14	95	155705	46.20	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.40%	
Target Compounds						
2) Dichlorodifluoromethane	1.20	85	1071	0.50	ug/l	98
8) Diethyl Ether	2.19	74	11091	6.17	ug/l	93
16) Acetone	2.44	43	3450	1.95	ug/l	97
65) Chlorobenzene	10.14	112	6147	0.81	ug/l	92

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

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