

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100918\
 Data File : VX005175.D
 Acq On : 09 Oct 2018 22:30
 Operator : JC/MD
 Sample : J5284-14
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RE-104D1-20181003

Quant Time: Oct 10 07:57:57 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 03 04:00:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	174486	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	274863	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	275819	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	154163	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	142298	57.09	ug/l	0.00
Spiked Amount 50.000			Recovery	=	114.18%	
35) Dibromofluoromethane	5.50	113	114848	49.30	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.60%	
50) Toluene-d8	8.72	98	404020	52.17	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.34%	
62) 4-Bromofluorobenzene	11.14	95	142396	51.04	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.08%	
Target Compounds						
9) 1,1,2-Trichlorotrifluoroet	2.38	101	2541	1.331	ug/l	100
16) Acetone	2.45	43	6502	4.468	ug/l	98
27) cis-1,2-Dichloroethene	4.60	96	1447	0.626	ug/l	91
44) Trichloroethene	7.21	130	125012	51.107	ug/l	96
64) Tetrachloroethene	9.34	164	3157	1.163	ug/l	91

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

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