

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091719\
 Data File : VX012451.D
 Acq On : 17 Sep 2019 23:01
 Operator : JC/SP
 Sample : K4858-04
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RE103D3-20190909

Quant Time: Sep 18 04:46:33 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091719W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 17 15:27:10 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.65	168	158076	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	254467	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	218339	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	86922	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.05	65	110357	48.71	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.42%	
35) Dibromofluoromethane	5.49	113	78147	50.74	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.48%	
50) Toluene-d8	8.71	98	297817	52.24	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.48%	
62) 4-Bromofluorobenzene	11.14	95	106011	44.79	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.58%	
Target Compounds						
					Qvalue	
9) 1,1,2-Trichlorotrifluoroet	2.37	101	6287	5.046	ug/l	96
27) cis-1,2-Dichloroethene	4.58	96	2050	1.330	ug/l	83
30) Chloroform	5.20	83	1859	0.624	ug/l	97
44) Trichloroethene	7.21	130	962041	743.654	ug/l	96
64) Tetrachloroethene	9.33	164	1332	1.331	ug/l #	81

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

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