

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100819\
 Data File : VX012920.D
 Acq On : 09 Oct 2019 04:18
 Operator : JC/SP
 Sample : K5154-24
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-101S

Quant Time: Oct 09 07:49:32 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X100819W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 09 01:51:23 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.65	168	179034	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	304627	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	260699	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	103779	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	118744	53.07	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.14%	
35) Dibromofluoromethane	5.49	113	90217	51.60	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.20%	
50) Toluene-d8	8.71	98	356399	52.26	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.52%	
62) 4-Bromofluorobenzene	11.14	95	122185	46.11	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.22%	
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	2.37	96	6861	3.792	ug/l	84
16) Acetone	2.44	43	4322	3.664	ug/l	89
24) 1,1-Dichloroethane	3.69	63	11346	3.267	ug/l	95
27) cis-1,2-Dichloroethene	4.59	96	9356	4.186	ug/l	89
30) Chloroform	5.21	83	4923	1.381	ug/l	89
32) 1,1,1-Trichloroethane	5.49	97	2638	0.874	ug/l #	46
44) Trichloroethene	7.21	130	38022	17.333	ug/l	88
64) Tetrachloroethene	9.33	164	21345	11.805	ug/l	98

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

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