

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX122619\
 Data File : VX014280.D
 Acq On : 26 Dec 2019 19:27
 Operator : JC/SP
 Sample : K6420-04
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PW

Quant Time: Dec 27 07:23:36 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X121319W.M
 Quant Title : SW846 8260
 QLast Update : Tue Dec 17 03:01:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.65	168	521456	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.85	114	819147	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	730541	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	317615	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.05	65	291809	49.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.76%	
35) Dibromofluoromethane	5.49	113	243716	48.94	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.88%	
50) Toluene-d8	8.71	98	962611	49.65	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.30%	
62) 4-Bromofluorobenzene	11.13	95	326557	46.08	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.16%	
Target Compounds						
					Qvalue	
19) Methyl tert-butyl Ether	3.18	73	11072	0.672	ug/l #	85
30) Chloroform	5.20	83	14566	1.547	ug/l	98
44) Trichloroethene	7.21	130	4359	0.682	ug/l	74
47) Bromodichloromethane	7.89	83	7135	0.967	ug/l #	98
64) Tetrachloroethene	9.34	164	5754	0.890	ug/l #	94

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

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