

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX080719\
 Data File : VX011376.D
 Acq On : 07 Aug 2019 20:25
 Operator : JC/SP
 Sample : K4220-10
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 DUP-01_080719

Quant Time: Aug 08 07:58:09 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X080119W.M
 Quant Title : SW846 8260
 QLast Update : Fri Aug 02 01:55:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.65	168	154741	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.85	114	253536	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	245866	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	113096	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	89978	57.74	ug/l	0.00
Spiked Amount 50.000			Recovery	=	115.48%	
35) Dibromofluoromethane	5.49	113	84340	50.95	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.90%	
50) Toluene-d8	8.71	98	310551	49.70	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.40%	
62) 4-Bromofluorobenzene	11.13	95	109824	49.23	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.46%	
Target Compounds						
17) Carbon Disulfide	2.56	76	46175	13.976	ug/l	99
19) Methyl tert-butyl Ether	3.19	73	3906	0.941	ug/l	92
27) cis-1,2-Dichloroethene	4.59	96	1156	0.674	ug/l	93
44) Trichloroethene	7.21	130	18534	9.270	ug/l	97
64) Tetrachloroethene	9.34	164	528478	253.104	ug/l	96

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

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