

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX122619\
 Data File : VX014281.D
 Acq On : 26 Dec 2019 19:50
 Operator : JC/SP
 Sample : K6405-01DL 20X
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 991-MW-01-(23)DL

Quant Time: Dec 27 07:26:58 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	525064	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.85	114	807481	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	734095	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	319659	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.05	65	294060	49.41	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.82%	
35) Dibromofluoromethane	5.49	113	240899	49.08	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.16%	
50) Toluene-d8	8.71	98	970563	50.79	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.58%	
62) 4-Bromofluorobenzene	11.14	95	323745	46.34	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.68%	
Target Compounds						
27) cis-1,2-Dichloroethene	4.59	96	25576	4.064	ug/l	93
44) Trichloroethene	7.21	130	30352	4.820	ug/l	96
64) Tetrachloroethene	9.33	164	267655	41.204	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122619\
Data File : VX014281.D
Acq On : 26 Dec 2019 19:50
Operator : JC/SP
Sample : K6405-01DL 20X
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
991-MW-01-(23)DL

Quant Time: Dec 27 07:26:58 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121319W.M
Quant Title : SW846 8260
QLast Update : Tue Dec 17 03:01:07 2019
Response via : Initial Calibration

