

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX122619\
 Data File : VX014283.D
 Acq On : 26 Dec 2019 20:37
 Operator : JC/SP
 Sample : K6405-05DL 5X
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 995-MW-11-(15)DL

Quant Time: Dec 27 07:32:34 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	530079	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.85	114	816484	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	736656	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	320708	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.05	65	300079	49.95	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.90%	
35) Dibromofluoromethane	5.49	113	248434	50.05	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.10%	
50) Toluene-d8	8.71	98	982658	50.85	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.70%	
62) 4-Bromofluorobenzene	11.13	95	329402	46.63	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.26%	
Target Compounds						
27) cis-1,2-Dichloroethene	4.59	96	6125	0.964	ug/l	85
44) Trichloroethene	7.21	130	9225	1.449	ug/l	90
64) Tetrachloroethene	9.33	164	282687	43.367	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX122619\
Data File : VX014283.D
Acq On : 26 Dec 2019 20:37
Operator : JC/SP
Sample : K6405-05DL 5X
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
995-MW-11-(15)DL

Quant Time: Dec 27 07:32:34 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

