

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX121019\
 Data File : VX013901.D
 Acq On : 10 Dec 2019 16:23
 Operator : JC/SP
 Sample : K6185-03
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RE103D3-20191204

Quant Time: Dec 11 06:26:27 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 26 15:53:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.65	168	579971	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.85	114	887932	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	785994	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	340497	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.05	65	322700	47.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.86%	
35) Dibromofluoromethane	5.49	113	270572	49.29	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.58%	
50) Toluene-d8	8.71	98	1046835	48.95	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.90%	
62) 4-Bromofluorobenzene	11.14	95	342107	44.36	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.72%	
Target Compounds						
9) 1,1,2-Trichlorotrifluoroet	2.37	101	13233	2.393	ug/l	94
27) cis-1,2-Dichloroethene	4.58	96	4486	0.650	ug/l	85
44) Trichloroethene	7.21	130	2692772	399.681	ug/l	96
64) Tetrachloroethene	9.33	164	4375	0.746	ug/l #	80

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX121019\
Data File : VX013901.D
Acq On : 10 Dec 2019 16:23
Operator : JC/SP
Sample : K6185-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
RE103D3-20191204

Quant Time: Dec 11 06:26:27 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 26 15:53:00 2019
Response via : Initial Calibration

