

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX120919\
 Data File : VX0137863.D
 Acq On : 09 Dec 2019 20:04
 Operator : JC/SP
 Sample : K6185-11DL 5X
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :

Quant Time: Dec 10 04:49:14 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	610786	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.85	114	928712	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	829693	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	358729	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.05	65	343815	48.49	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.98%	
35) Dibromofluoromethane	5.49	113	279352	48.66	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.32%	
50) Toluene-d8	8.71	98	1111583	49.69	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.38%	
62) 4-Bromofluorobenzene	11.14	95	366965	45.49	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.98%	
Target Compounds						
9) 1,1,2-Trichlorotrifluoroet	2.37	101	13702	2.353	ug/l	98
16) Acetone	2.43	43	11918	3.405	ug/l	98
27) cis-1,2-Dichloroethene	4.59	96	6516	0.896	ug/l	88
44) Trichloroethene	7.21	130	316686	44.941	ug/l	98
64) Tetrachloroethene	9.33	164	10541	1.704	ug/l	98

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

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