

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX041819\
 Data File : VX009047.D
 Acq On : 18 Apr 2019 16:31
 Operator : JC/SP
 Sample : K2449-13
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SS037MW302-190416

Quant Time: Apr 19 09:25:00 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X041519W.M
 Quant Title : SW846 8260
 QLast Update : Wed Apr 17 06:47:58 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	228892	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	371668	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	304529	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	115094	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	155735	48.28	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.56%	
35) Dibromofluoromethane	5.49	113	113468	47.57	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.14%	
50) Toluene-d8	8.71	98	457173	48.06	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.12%	
62) 4-Bromofluorobenzene	11.13	95	138602	41.28	ug/l	0.00
Spiked Amount 50.000			Recovery	=	82.56%	
Target Compounds						
21) trans-1,2-Dichloroethene	3.17	96	2045	0.781	ug/l	84
27) cis-1,2-Dichloroethene	4.60	96	139633	46.974	ug/l	91
30) Chloroform	5.21	83	3028	0.648	ug/l	96
44) Trichloroethene	7.21	130	58149	20.929	ug/l	99
64) Tetrachloroethene	9.34	164	624000	269.727	ug/l	98

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

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