

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX041619\
 Data File : VX008986.D
 Acq On : 16 Apr 2019 14:41
 Operator : JC/SP
 Sample : K2387-07
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 KY063MW003-190410

Quant Time: Apr 17 05:34:07 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X041519W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 16 09:35:21 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.66	168	261093	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	415914	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	359771	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	156728	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	168290	45.74	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.48%	
35) Dibromofluoromethane	5.50	113	123933	46.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.86%	
50) Toluene-d8	8.71	98	516497	48.52	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.04%	
62) 4-Bromofluorobenzene	11.13	95	175123	46.60	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.20%	
Target Compounds						
16) Acetone	2.44	43	7312	3.832	ug/l	99
21) trans-1,2-Dichloroethene	3.17	96	2639	0.884	ug/l	87
27) cis-1,2-Dichloroethene	4.60	96	557159	164.319	ug/l	90
30) Chloroform	5.21	83	6270	1.176	ug/l	96
44) Trichloroethene	7.21	130	518959	166.911	ug/l	99
64) Tetrachloroethene	9.34	164	1650281	603.811	ug/l	98

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

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