

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX102419\
 Data File : VX013197.D
 Acq On : 23 Oct 2019 19:18
 Operator : JC/SP
 Sample : K5451-10
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FA19-JPZ0341

Quant Time: Oct 24 04:24:57 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102419W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 24 03:56:38 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.65	168	82274	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.85	114	123923	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	120151	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	73440	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	90270	52.46	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.92%	
35) Dibromofluoromethane	5.49	113	69900	50.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.16%	
50) Toluene-d8	8.71	98	255059	49.80	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.60%	
62) 4-Bromofluorobenzene	11.14	95	93585	46.17	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.34%	
Target Compounds						
					Qvalue	
3) Chloromethane	1.32	50	796	1.548	ug/l #	86
16) Acetone	2.44	43	4718	5.820	ug/l #	88
27) cis-1,2-Dichloroethene	4.59	96	2321	1.352	ug/l	96
30) Chloroform	5.20	83	1612	0.571	ug/l	91
37) Ethyl Acetate	4.83	43	3337	1.616	ug/l #	88
44) Trichloroethene	7.21	130	9052	5.124	ug/l	99

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

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