

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX081119\
 Data File : VX011461.D
 Acq On : 12 Aug 2019 03:54
 Operator : JC/SP
 Sample : K4266-21
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 R1-R2(A1-A4)-(0-1)

Quant Time: Aug 12 09:13:44 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X080119W.M
 Quant Title : SW846 8260
 QLast Update : Fri Aug 02 01:55:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.66	168	196190	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.85	114	291676	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	270936	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	123218	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	105108	53.20	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.40%	
35) Dibromofluoromethane	5.49	113	90426	47.49	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.98%	
50) Toluene-d8	8.71	98	351667	48.92	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.84%	
62) 4-Bromofluorobenzene	11.13	95	116693	45.47	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.94%	
Target Compounds						
16) Acetone	2.43	43	6265	6.662	ug/l	96
20) Methylene Chloride	2.85	84	23788	11.550	ug/l	97
30) Chloroform	5.20	83	3821	1.140	ug/l	88
43) Isopropyl Acetate	6.45	43	7734	1.894	ug/l	97

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

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