

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX011020\
 Data File : VX014543.D
 Acq On : 10 Jan 2020 17:31
 Operator : JC/SP
 Sample : L1065-02
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 6-HER-002-003-02

Quant Time: Jan 11 00:01:30 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X010720W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jan 07 15:54:41 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.65	168	487326	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.85	114	749944	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	670369	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	310041	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.05	65	269754	49.59	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.18%	
35) Dibromofluoromethane	5.49	113	226928	49.99	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.98%	
50) Toluene-d8	8.71	98	889920	50.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.02%	
62) 4-Bromofluorobenzene	11.13	95	299131	46.25	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.50%	
Target Compounds						
16) Acetone	2.43	43	17793	2.384	ug/l	93
30) Chloroform	5.20	83	68699	7.914	ug/l	95
80) 1,3,5-Trimethylbenzene	11.50	105	7087	0.392	ug/l	96
84) 1,2,4-Trimethylbenzene	11.81	105	7876	0.431	ug/l	87
86) p-Isopropyltoluene	12.06	119	36271	1.841	ug/l	92

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

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