

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX051120\
 Data File : VX016145.D
 Acq On : 11 May 2020 13:48
 Operator : JC/SP
 Sample : L2552-08
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SS052OW377-200507

Quant Time: May 11 18:51:19 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X050420W.M
 Quant Title : SW846 8260
 QLast Update : Wed May 06 06:43:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.64	168	284328	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.84	114	463311	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	454517	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	236157	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.04	65	169310	45.25	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.50%	
35) Dibromofluoromethane	5.47	113	136317	46.44	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.88%	
50) Toluene-d8	8.70	98	570763	50.47	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.94%	
62) 4-Bromofluorobenzene	11.12	95	227527	52.89	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.78%	
Target Compounds						
16) Acetone	2.42	43	7443	4.458	ug/l	96
27) cis-1,2-Dichloroethene	4.57	96	6491	1.803	ug/l	77
30) Chloroform	5.18	83	2925	0.514	ug/l #	82
44) Trichloroethene	7.20	130	19150	4.039	ug/l	93
64) Tetrachloroethene	9.32	164	5486	0.983	ug/l	93

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

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