

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX051120\
 Data File : VX016149.D
 Acq On : 11 May 2020 15:19
 Operator : JC/SP
 Sample : L2552-12
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 Z4-DUP-0552-200507

Quant Time: May 11 18:57:09 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	5.63	168	286020	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.84	114	469838	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	457868	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	236598	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.04	65	174225	46.29	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.58%	
35) Dibromofluoromethane	5.47	113	137201	46.09	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.18%	
50) Toluene-d8	8.70	98	578418	50.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.86%	
62) 4-Bromofluorobenzene	11.12	95	235134	53.90	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.80%	
Target Compounds						
16) Acetone	2.42	43	4771	2.841	ug/l	93
27) cis-1,2-Dichloroethene	4.57	96	6057	1.672	ug/l	87
30) Chloroform	5.17	83	2885	0.504	ug/l	98
44) Trichloroethene	7.19	130	18750	3.900	ug/l	95
64) Tetrachloroethene	9.32	164	5198	0.925	ug/l	97

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

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