

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN041720\
 Data File : VN061080.D
 Acq On : 18 Apr 2020 3:07
 Operator : JC/MD
 Sample : L2338-12
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BP-MW202I-20200415

Quant Time: Apr 20 07:05:09 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N041620W.M
 Quant Title : SW846 8260
 QLast Update : Fri Apr 17 03:49:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.64	168	138894	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.56	114	240099	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	222086	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	96540	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	97118	54.65	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.30%	
35) Dibromofluoromethane	7.56	113	72779	51.91	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.82%	
50) Toluene-d8	10.08	98	297014	56.06	ug/l	0.00
Spiked Amount 50.000			Recovery	=	112.12%	
62) 4-Bromofluorobenzene	12.40	95	106563	51.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.64%	
Target Compounds						
7) Trichlorofluoromethane	3.01	101	11741	4.607	ug/l	99
12) 1,1-Dichloroethene	3.72	96	4913	3.232	ug/l	93
19) Methyl tert-butyl Ether	5.00	73	9858	1.860	ug/l	94
30) Chloroform	7.34	83	3097	1.029	ug/l	94
32) 1,1,1-Trichloroethane	7.54	97	4576	1.758	ug/l	100
44) Trichloroethene	8.81	130	8863	4.799	ug/l	97
64) Tetrachloroethene	10.62	164	9539	5.613	ug/l	92

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

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