

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100419\
 Data File : VN058548.D
 Acq On : 4 Oct 2019 17:59
 Operator : JC/SP
 Sample : K5169-10
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-22D_100219

Quant Time: Oct 05 04:20:30 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N092719W.M
 Quant Title : SW846 8260
 QLast Update : Mon Sep 30 08:07:38 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	471082	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	754063	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	608245	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	248737	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	332017	48.23	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.46%	
35) Dibromofluoromethane	7.57	113	234623	50.77	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.54%	
50) Toluene-d8	10.09	98	881885	48.71	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.42%	
62) 4-Bromofluorobenzene	12.41	95	288440	44.06	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.12%	
Target Compounds						
3) Chloromethane	2.04	50	3909	0.688	ug/l	96
16) Acetone	3.80	43	4986	2.125	ug/l	96
19) Methyl tert-butyl Ether	5.02	73	8661	0.552	ug/l	94
27) cis-1,2-Dichloroethene	6.81	96	3633	0.629	ug/l	81
44) Trichloroethene	8.82	130	26688	5.215	ug/l	99
64) Tetrachloroethene	10.63	164	501845	129.864	ug/l	98

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

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