

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100419\
 Data File : VN058546.D
 Acq On : 4 Oct 2019 17:15
 Operator : JC/SP
 Sample : K5169-08
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-2R_100219

Quant Time: Oct 05 04:12:11 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	423657	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	681409	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	639210	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	226801	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	301887	48.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.52%	
35) Dibromofluoromethane	7.57	113	213196	51.05	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.10%	
50) Toluene-d8	10.09	98	856576	52.36	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.72%	
62) 4-Bromofluorobenzene	12.41	95	274764	46.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.90%	
Target Compounds						
3) Chloromethane	2.04	50	3005	0.588	ug/l	99
16) Acetone	3.80	43	15804	7.488	ug/l	94
27) cis-1,2-Dichloroethene	6.82	96	1949	0.375	ug/l	60
52) Toluene	10.15	92	8540	0.724	ug/l	91

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

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