

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100419\
 Data File : VN058545.D
 Acq On : 4 Oct 2019 16:53
 Operator : JC/SP
 Sample : K5169-05
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-25S_100119

Quant Time: Oct 05 04:04:11 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	502924	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	765638	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	649601	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	267286	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	347891	47.34	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.68%	
35) Dibromofluoromethane	7.57	113	241725	51.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.02%	
50) Toluene-d8	10.08	98	916268	49.84	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.68%	
62) 4-Bromofluorobenzene	12.41	95	304172	45.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.52%	
Target Compounds						
3) Chloromethane	2.04	50	2077	0.343	ug/l	93
16) Acetone	3.80	43	9197	3.671	ug/l	92
30) Chloroform	7.36	83	21277	2.008	ug/l	93
64) Tetrachloroethene	10.63	164	18540	4.492	ug/l	92

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

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