

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091419\
 Data File : VN058026.D
 Acq On : 13 Sep 2019 19:01
 Operator : JC/SP
 Sample : K4859-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 IDW-DECON-WATER-091119

Quant Time: Sep 14 05:07:50 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091019W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 10 03:02:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	318934	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	527488	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	472795	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	241835	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	348556	53.96	ug/l	0.00
Spiked Amount	50.000		Recovery	=	107.92%	
35) Dibromofluoromethane	7.58	113	240198	49.34	ug/l	0.00
Spiked Amount	50.000		Recovery	=	98.68%	
50) Toluene-d8	10.09	98	951392	50.33	ug/l	0.00
Spiked Amount	50.000		Recovery	=	100.66%	
62) 4-Bromofluorobenzene	12.40	95	310708	43.13	ug/l	0.00
Spiked Amount	50.000		Recovery	=	86.26%	
Target Compounds						
16) Acetone	3.80	43	70559	33.320	ug/l	94
30) Chloroform	7.36	83	50722	5.695	ug/l	92
44) Trichloroethene	8.83	130	5558	1.230	ug/l	87
47) Bromodichloromethane	9.39	83	7889	1.220	ug/l #	96
51) 4-Methyl-2-Pentanone	9.98	43	33128	5.267	ug/l	100
64) Tetrachloroethene	10.63	164	14881	4.263	ug/l	97
69) o-Xylene	11.95	106	8489	1.075	ug/l	93

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

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