

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091719\
 Data File : VN058144.D
 Acq On : 17 Sep 2019 13:12
 Operator : JC/SP
 Sample : K4666-18
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PL-02-091619

Quant Time: Sep 18 08:13:57 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	375865	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	621131	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	568240	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	310139	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	356770	46.86	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.72%	
35) Dibromofluoromethane	7.57	113	250400	43.68	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.36%	
50) Toluene-d8	10.09	98	950590	42.71	ug/l	0.00
Spiked Amount 50.000			Recovery	=	85.42%	
62) 4-Bromofluorobenzene	12.41	95	342864	40.42	ug/l	0.00
Spiked Amount 50.000			Recovery	=	80.84%	
Target Compounds						
16) Acetone	3.80	43	15868	6.358	ug/l	90
18) Methyl Acetate	4.31	43	7234	1.087	ug/l #	77
20) Methylene Chloride	4.53	84	58997	11.469	ug/l	96
43) Isopropyl Acetate	8.15	43	152545	12.188	ug/l #	85

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

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