

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091919\
 Data File : VN058209.D
 Acq On : 19 Sep 2019 17:25
 Operator : JC/SP
 Sample : K4896-17
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 EB01-20190912

Manual Integrations
 APPROVED

MMDadoda
 9/20/2019 3:37:40 PM

Quant Time: Sep 20 08:08:45 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091819W.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 19 09:27:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	506306	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	830864	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	701839	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	289735	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	365340	51.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.62%	
35) Dibromofluoromethane	7.58	113	252418	49.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.94%	
50) Toluene-d8	10.08	98	1037572	52.72	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.44%	
62) 4-Bromofluorobenzene	12.40	95	332217	45.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.90%	
Target Compounds						
16) Acetone	3.80	43	34562	14.439	ug/l	92
20) Methylene Chloride	4.54	84	5220m	1.028	ug/l	
52) Toluene	10.15	92	15154	1.321	ug/l	98

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

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