

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN101819\
 Data File : VN058864.D
 Acq On : 18 Oct 2019 17:33
 Operator : JC/SP
 Sample : K5452-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 FA19-JPZ7807

Manual Integrations
 APPROVED

MMDadoda
 10/21/2019 10:42:58 AM

Quant Time: Oct 21 08:30:36 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N101819W.M
 Quant Title : SW846 8260
 QLast Update : Sat Oct 19 00:16:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	381203	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	606768	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	547451	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	205229	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	292374	52.29	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.58%	
35) Dibromofluoromethane	7.57	113	211274	52.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.92%	
50) Toluene-d8	10.08	98	780474	50.40	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.80%	
62) 4-Bromofluorobenzene	12.40	95	251630	44.25	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.50%	
Target Compounds						
3) Chloromethane	2.04	50	2837	0.482	ug/l	98
16) Acetone	3.80	43	9189	3.326	ug/l	99
44) Trichloroethene	8.82	130	2198m	0.495	ug/l	
64) Tetrachloroethene	10.63	164	102753	26.801	ug/l	99

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

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