

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101321\
 Data File : VN068896.D
 Acq On : 13 Oct 2021 14:32
 Operator : JC/MD
 Sample : M4188-01
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PP-PROCESS-BLEED

Quant Time: Oct 14 06:41:58 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 13 05:44:40 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	8.088	168	417222	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	677646	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	601895	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	242117	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	242519	44.09	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	88.18%
35) Dibromofluoromethane	8.024	113	203036	51.50	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	103.00%
50) Toluene-d8	10.443	98	813892	51.59	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	103.18%
62) 4-Bromofluorobenzene	12.733	95	292509	49.15	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	98.30%
Target Compounds						
					Qvalue	
16) Acetone	4.296	43	16436	8.01	ug/l	97
29) Tetrahydrofuran	7.689	42	4054	2.01	ug/l	93
30) Chloroform	7.820	83	154199	17.92	ug/l	98
42) 1,2-Dichloroethane	8.533	62	196511	31.49	ug/l	98
47) Bromodichloromethane	9.764	83	59410	9.18	ug/l	100
60) Dibromochloromethane	11.239	129	20431	4.05	ug/l	95
95) Naphthalene	15.509	128	3563	4.10	ug/l #	94

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

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