

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072123\
 Data File : VN078611.D
 Acq On : 21 Jul 2023 14:11
 Operator : JC\MD
 Sample : 03709-03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 GWBR-04-400-415-072023

Quant Time: Jul 21 23:22:57 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071023W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 10 16:13:53 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	419023	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	754724	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	706281	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	243958	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	282909	50.363	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 100.720%	
35) Dibromofluoromethane	8.177	113	282833	55.205	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 110.420%	
50) Toluene-d8	10.571	98	898532	46.425	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 92.860%	
62) 4-Bromofluorobenzene	12.853	95	298355	50.047	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 100.100%	
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	4.348	96	61323	13.869	ug/l	90
24) 1,1-Dichloroethane	6.565	63	13736	1.507	ug/l	97
27) cis-1,2-Dichloroethene	7.494	96	71127	12.340	ug/l	96
42) 1,2-Dichloroethane	8.677	62	7344	0.985	ug/l	96
44) Trichloroethene	9.365	130	61950	10.488	ug/l	88
51) 4-Methyl-2-Pentanone	10.453	43	3160	0.523	ug/l #	85
52) Toluene	10.635	92	743730	54.755	ug/l	95

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

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