

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081823\
 Data File : VN078957.D
 Acq On : 18 Aug 2023 21:13
 Operator : JC\MD
 Sample : 04043-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 E-4

Quant Time: Aug 18 23:50:17 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081523W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 16 02:42:02 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	202640	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	398097	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	377236	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	120436	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.589	65	162394	51.501	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	103.000%	
35) Dibromofluoromethane	8.171	113	132258	49.354	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	98.700%	
50) Toluene-d8	10.571	98	488294	49.074	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	98.140%	
62) 4-Bromofluorobenzene	12.853	95	145882	44.854	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	89.700%	
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	37608	28.905	ug/l	88
20) Methylene Chloride	5.277	84	8354	3.013	ug/l #	78
25) 2-Butanone	7.489	43	19434	9.316	ug/l #	69
43) Isopropyl Acetate	8.694	43	224634	28.179	ug/l #	71
51) 4-Methyl-2-Pentanone	10.447	43	66555	14.077	ug/l #	81

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

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