

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102722\
 Data File : VN075038.D
 Acq On : 27 Oct 2022 14:07
 Operator : JC\MD
 Sample : N5268-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 OK-03-102422

Quant Time: Oct 28 02:14:15 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N102622W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 27 09:48:49 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	121858	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	239045	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	215142	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	88261	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	97601	59.444	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 118.880%			
35) Dibromofluoromethane	8.171	113	73697	52.516	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 105.040%			
50) Toluene-d8	10.571	98	295816	54.333	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 108.660%			
62) 4-Bromofluorobenzene	12.853	95	92018	48.113	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 96.220%			
Target Compounds					Qvalue	
16) Acetone	4.448	43	18912	27.203	ug/l	94
20) Methylene Chloride	5.289	84	78907	50.905	ug/l	95
30) Chloroform	7.971	83	4072	1.424	ug/l	92
37) Ethyl Acetate	7.577	43	8121	3.408	ug/l	98
43) Isopropyl Acetate	8.694	43	32487	8.790	ug/l #	87

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

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