

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071423\
 Data File : VN078525.D
 Acq On : 14 Jul 2023 16:24
 Operator : JC\MD
 Sample : 03632-09
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 RINSATE-BLANK

Quant Time: Jul 15 04:04:10 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	8.236	168	386005	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	711037	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	652236	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	222298	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	264666	51.145	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	102.300%
35) Dibromofluoromethane	8.177	113	271239	56.195	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	112.400%
50) Toluene-d8	10.577	98	830757	45.561	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	91.120%
62) 4-Bromofluorobenzene	12.859	95	285815	50.889	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	101.780%
Target Compounds						
						Qvalue
3) Chloromethane	2.377	50	2073	0.403	ug/l	94
16) Acetone	4.453	43	92078	57.122	ug/l	89
25) 2-Butanone	7.512	43	33448	13.182	ug/l	# 80
51) 4-Methyl-2-Pentanone	10.453	43	92183	16.203	ug/l	90

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

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