

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090222\
 Data File : VN074294.D
 Acq On : 02 Sep 2022 18:57
 Operator : JC\MD
 Sample : N4329-01DL 5X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-3RDL

Quant Time: Sep 03 04:40:56 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N090222W.M
 Quant Title : SW846 8260
 QLast Update : Sat Sep 03 04:35:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	270906	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	472656	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	452727	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	192194	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	187037	44.290	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	88.580%
35) Dibromofluoromethane	7.951	113	161552	51.580	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	103.160%
50) Toluene-d8	10.380	98	559828	45.329	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	90.660%
62) 4-Bromofluorobenzene	12.669	95	195525	43.648	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	87.300%
Target Compounds						
					Qvalue	
19) Methyl tert-butyl Ether	5.540	73	230450	21.803	ug/l	98
22) Diisopropyl ether	6.410	45	8084	0.803	ug/l #	88
27) cis-1,2-Dichloroethene	7.257	96	7685	1.977	ug/l #	16
40) Benzene	8.392	78	251135	17.908	ug/l	96
67) Ethyl Benzene	11.786	91	5052	0.277	ug/l #	86

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

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