

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062822\
 Data File : VN073211.D
 Acq On : 28 Jun 2022 21:02
 Operator : JC\MD
 Sample : N3494-30DL 10X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SP-EXC-1-2-GRDL

Quant Time: Jun 29 02:03:55 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062822W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 29 01:54:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	304853	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	504278	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.680	117	473903	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	223509	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	191521	43.519	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	87.040%
35) Dibromofluoromethane	7.951	113	159917	47.665	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	95.320%
50) Toluene-d8	10.374	98	535174	44.912	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	89.820%
62) 4-Bromofluorobenzene	12.668	95	235483	51.424	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	102.840%
Target Compounds						
					Qvalue	
16) Acetone	4.234	43	12407	5.022	ug/l #	86
20) Methylene Chloride	5.010	84	58648	13.509	ug/l	92
43) Isopropyl Acetate	8.486	43	14098	1.258	ug/l #	90
65) Chlorobenzene	11.710	112	228018	21.274	ug/l	98

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

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