

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071222\
 Data File : VN073440.D
 Acq On : 12 Jul 2022 16:35
 Operator : JC\MD
 Sample : N3636-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 GWOW-BR-03-150-170-070722

Quant Time: Jul 13 06:18:59 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	243867	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	416860	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	379694	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.616	152	180088	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	167952	47.707	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	95.420%
35) Dibromofluoromethane	7.951	113	138603	49.975	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.960%
50) Toluene-d8	10.380	98	439091	44.576	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	89.160%
62) 4-Bromofluorobenzene	12.669	95	186516	49.272	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	98.540%
Target Compounds						
					Qvalue	
16) Acetone	4.246	43	4842	2.450	ug/l #	76
40) Benzene	8.386	78	9710	0.737	ug/l #	86
51) 4-Methyl-2-Pentanone	10.275	43	5464	0.930	ug/l #	76
52) Toluene	10.439	92	6234	0.766	ug/l	93

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

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