

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061822\
 Data File : VN072927.D
 Acq On : 18 Jun 2022 10:04
 Operator : JC\MD
 Sample : N3334-20
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-14A-1A-5

Quant Time: Jun 20 08:02:04 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
 Quant Title : SW846 8260
 QLast Update : Sat Jun 18 06:22:45 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	301325	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.898	114	525632	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.680	117	459960	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	202435	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	173958	42.701	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	85.400%
35) Dibromofluoromethane	7.951	113	137295	42.259	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	84.520%
50) Toluene-d8	10.375	98	464662	38.436	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	76.880%#
62) 4-Bromofluorobenzene	12.669	95	162707	37.429	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	74.860%#

Target Compounds						Qvalue
16) Acetone	4.234	43	63480	28.347	ug/l	90
17) Carbon Disulfide	4.463	76	9948	1.346	ug/l #	90
25) 2-Butanone	7.269	43	28327	7.972	ug/l	90
43) Isopropyl Acetate	8.492	43	39222	3.902	ug/l #	88
69) o-Xylene	12.216	106	14956	2.413	ug/l	99
73) Isopropylbenzene	12.516	105	22025	1.406	ug/l	98
78) n-propylbenzene	12.857	91	20078	1.166	ug/l	100
80) 1,3,5-Trimethylbenzene	12.992	105	30522	2.373	ug/l	97
84) 1,2,4-Trimethylbenzene	13.298	105	89164	6.959	ug/l	99
95) Naphthalene	15.427	128	82638	5.284	ug/l	99

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

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