

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062722\
 Data File : VN073190.D
 Acq On : 28 Jun 2022 03:49
 Operator : JC\MD
 Sample : N3494-32
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 46 Sample Multiplier: 1

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

Quant Time: Jun 28 06:45:18 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	369190	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	594942	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.680	117	555257	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	268700	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	218904	43.857	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	87.720%
35) Dibromofluoromethane	7.951	113	188406	51.235	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	102.480%
50) Toluene-d8	10.374	98	633595	46.304	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	92.600%
62) 4-Bromofluorobenzene	12.668	95	276210	56.137	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	112.280%
Target Compounds						
					Qvalue	
16) Acetone	4.234	43	72313	26.356	ug/l	100
20) Methylene Chloride	5.016	84	316253	74.430	ug/l	94
37) Ethyl Acetate	7.339	43	43630	5.482	ug/l	98
43) Isopropyl Acetate	8.498	43	169924	14.936	ug/l #	87
65) Chlorobenzene	11.710	112	92301	8.512	ug/l	98

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

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