

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062822\
 Data File : VN073216.D
 Acq On : 28 Jun 2022 23:04
 Operator : JC\MD
 Sample : N3494-32
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 27 Sample Multiplier: 1

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

Quant Time: Jun 29 02:04:41 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	360620	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	610383	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.680	117	564284	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	269272	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	229230	44.032	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	88.060%
35) Dibromofluoromethane	7.957	113	191545	47.167	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.340%
50) Toluene-d8	10.375	98	641776	44.496	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	89.000%
62) 4-Bromofluorobenzene	12.669	95	283237	51.100	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	102.200%
Target Compounds						
					Qvalue	
16) Acetone	4.222	43	69368	23.738	ug/l	90
20) Methylene Chloride	5.016	84	286655	55.819	ug/l	87
37) Ethyl Acetate	7.340	43	42106	4.460	ug/l #	94
43) Isopropyl Acetate	8.492	43	155272	11.443	ug/l #	88
65) Chlorobenzene	11.710	112	75948	5.951	ug/l	100

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

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