

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062722\
 Data File : VN073189.D
 Acq On : 28 Jun 2022 03:25
 Operator : JC\MD
 Sample : N3494-30
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 45 Sample Multiplier: 1

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

Quant Time: Jun 28 06:45:06 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	350608	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	566351	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.680	117	530933	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	273661	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	206310	43.524	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	87.040%
35) Dibromofluoromethane	7.951	113	174548	49.863	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.720%
50) Toluene-d8	10.375	98	598409	45.941	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	91.880%
62) 4-Bromofluorobenzene	12.668	95	273108	58.308	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	116.620%
Target Compounds						
					Qvalue	
16) Acetone	4.228	43	117195	44.978	ug/l	95
20) Methylene Chloride	5.016	84	632077	156.911	ug/l	90
25) 2-Butanone	7.263	43	22839	5.524	ug/l	90
37) Ethyl Acetate	7.339	43	46391	6.123	ug/l	95
40) Benzene	8.386	78	68936	4.631	ug/l	98
43) Isopropyl Acetate	8.492	43	170918	15.782	ug/l #	88
65) Chlorobenzene	11.710	112	2708401	261.206	ug/l	99
91) 1,2-Dichlorobenzene	13.921	146	18919	2.003	ug/l	96

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

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