

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090922\
 Data File : VN074388.D
 Acq On : 09 Sep 2022 21:23
 Operator : JC\MD
 Sample : N4471-12
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BR-03D-479.5-499.5-083122-H

Quant Time: Sep 10 07:27:06 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N090922W.M
 Quant Title : SW846 8260
 QLast Update : Sat Sep 10 06:10:29 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	319509	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	571970	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	555694	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.615	152	242660	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	231903	41.951	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	83.900%
35) Dibromofluoromethane	7.951	113	192449	49.840	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.680%
50) Toluene-d8	10.380	98	646191	42.949	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	85.900%#
62) 4-Bromofluorobenzene	12.674	95	266062	48.707	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	97.420%
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	4.110	96	151850	38.125	ug/l	99
16) Acetone	4.228	43	14016	4.696	ug/l #	87
17) Carbon Disulfide	4.457	76	66939	6.184	ug/l	97
24) 1,1-Dichloroethane	6.304	63	30825	3.410	ug/l #	95
27) cis-1,2-Dichloroethene	7.251	96	170037	30.924	ug/l	86
30) Chloroform	7.745	83	3313	0.367	ug/l #	42
42) 1,2-Dichloroethane	8.457	62	14808	1.989	ug/l	99
44) Trichloroethene	9.151	130	104217	23.867	ug/l	95
49) 1,4-Dioxane	9.516	88	4538	32.378	ug/l #	84
52) Toluene	10.445	92	92928	7.715	ug/l	99

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

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