

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060324\
 Data File : VN082445.D
 Acq On : 03 Jun 2024 19:40
 Operator : JC\MD
 Sample : P2683-03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-4

Quant Time: Jun 04 00:35:04 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N051524W.M
 Quant Title : SW846 8260
 QLast Update : Thu May 16 05:21:06 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	149112	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	275557	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	273971	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	115210	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	131972	56.051	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	112.100%
35) Dibromofluoromethane	8.171	113	94781	55.915	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	111.820%
50) Toluene-d8	10.571	98	364950	56.156	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	112.320%
62) 4-Bromofluorobenzene	12.853	95	120817	52.085	ug/l	0.00
Spiked Amount	50.000	Range	64 - 133	Recovery	=	104.180%
Target Compounds						
						Qvalue
16) Acetone	4.430	43	94832	98.184	ug/l	90
20) Methylene Chloride	5.289	84	12955	6.648	ug/l #	78
25) 2-Butanone	7.494	43	14032	9.582	ug/l	90
43) Isopropyl Acetate	8.694	43	188612	33.456	ug/l #	74
51) 4-Methyl-2-Pentanone	10.447	43	16393	5.118	ug/l #	79

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

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