

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071624\
 Data File : VN082946.D
 Acq On : 16 Jul 2024 19:39
 Operator : JC\MD
 Sample : P3246-17
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-CTW2-BL-04

Quant Time: Jul 17 01:00:54 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N070824W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 08 13:46:04 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	85027	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	205446	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	181000	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	69311	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	79080	66.348	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 132.700%#	
35) Dibromofluoromethane	8.165	113	53846	40.845	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 81.700%	
50) Toluene-d8	10.571	98	227934	46.812	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 93.620%	
62) 4-Bromofluorobenzene	12.853	95	60662	34.683	ug/l	0.00
Spiked Amount	50.000	Range	64 - 133	Recovery	= 69.360%	
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	26924	64.160	ug/l #	87
20) Methylene Chloride	5.283	84	5913	5.835	ug/l #	80
25) 2-Butanone	7.494	43	4081	6.441	ug/l	93
43) Isopropyl Acetate	8.694	43	119307	33.196	ug/l #	89

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

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