

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN110424\
 Data File : VN084660.D
 Acq On : 04 Nov 2024 16:17
 Operator : JC\MD
 Sample : P4660-07
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-WOOD-01-C

Quant Time: Nov 05 00:29:19 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N103024W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 31 18:45:38 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	176637	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	309644	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	292406	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	151986	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.576	65	121549	47.643	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	95.280%	
35) Dibromofluoromethane	8.159	113	98128	46.819	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	93.640%	
50) Toluene-d8	10.565	98	368332	47.707	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	95.420%	
62) 4-Bromofluorobenzene	12.847	95	157637	54.631	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	=	109.260%	
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	31850	40.832	ug/l	97
27) cis-1,2-Dichloroethene	7.477	96	6538	2.711	ug/l	84
43) Isopropyl Acetate	8.688	43	130590	26.346	ug/l #	87
44) Trichloroethene	9.347	130	17658	8.202	ug/l	97
67) Ethyl Benzene	11.965	91	11871	1.087	ug/l	94
68) m/p-Xylenes	12.065	106	11193	2.708	ug/l	94
69) o-Xylene	12.394	106	11044	2.856	ug/l	97
80) 1,3,5-Trimethylbenzene	13.170	105	24865	2.817	ug/l	96
84) 1,2,4-Trimethylbenzene	13.482	105	74948	8.226	ug/l	99

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

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