

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060524\
 Data File : VN082476.D
 Acq On : 05 Jun 2024 17:27
 Operator : JC\MD
 Sample : P2704-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WCS-TPU

Manual Integrations
 APPROVED

Reviewed By : John Carlone 06/06/2024
 Supervised By : Mahesh Dadoda 06/06/2024

Quant Time: Jun 06 00:59:58 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	163401	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	304795	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	309579	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	128772	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	132409	51.318	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 102.640%			
35) Dibromofluoromethane	8.165	113	96694	51.571	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 103.140%			
50) Toluene-d8	10.571	98	378751	52.689	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 105.380%			
62) 4-Bromofluorobenzene	12.853	95	126924	49.469	ug/l	0.00
Spiked Amount 50.000	Range 64 - 133		Recovery = 98.940%			
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	99154	93.681	ug/l	89
20) Methylene Chloride	5.283	84	4078	1.910	ug/l	92
25) 2-Butanone	7.494	43	25272	15.748	ug/l #	87
43) Isopropyl Acetate	8.694	43	301398	48.334	ug/l #	70
51) 4-Methyl-2-Pentanone	10.447	43	24279m	6.853	ug/l	

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

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