

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120424\
 Data File : VN085092.D
 Acq On : 04 Dec 2024 14:25
 Operator : JC\MD
 Sample : P5052-14
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 OR-640-COMP-62

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/05/2024
 Supervised By : Mahesh Dadoda 12/05/2024

Quant Time: Dec 04 22:06:51 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112224W.M
 Quant Title : SW846 8260
 QLast Update : Sat Nov 23 02:54:10 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	159336	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	287542	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	252903	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	95200	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	113352	49.943	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	99.880%		
35) Dibromofluoromethane	8.165	113	92932	47.926	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	95.860%		
50) Toluene-d8	10.565	98	337228	48.527	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	97.060%		
62) 4-Bromofluorobenzene	12.847	95	116360	44.775	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery =	89.540%		
Target Compounds						
16) Acetone	4.430	43	41164	62.275	ug/l	Qvalue 93
43) Isopropyl Acetate	8.688	43	163770m	33.020	ug/l	
51) 4-Methyl-2-Pentanone	10.441	43	9732m	4.296	ug/l	

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

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