

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN103024\
 Data File : VN084599.D
 Acq On : 31 Oct 2024 04:33
 Operator : JC\MD
 Sample : P4611-12
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-4

Quant Time: Nov 01 05:20:36 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	176854	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	330083	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	302284	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	128849	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.571	65	132308	51.797	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	103.600%	
35) Dibromofluoromethane	8.165	113	106378	47.612	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	95.220%	
50) Toluene-d8	10.565	98	389895	47.373	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	94.740%	
62) 4-Bromofluorobenzene	12.847	95	143236	46.567	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	=	93.140%	
Target Compounds						
						Qvalue
16) Acetone	4.436	43	12391	14.102	ug/l	97
18) Methyl Acetate	5.030	43	16162	2.178	ug/l	98
43) Isopropyl Acetate	8.688	43	185709	35.475	ug/l #	78

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

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