

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081324\
 Data File : VN083266.D
 Acq On : 13 Aug 2024 18:33
 Operator : JC\MD
 Sample : P3548-11
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 289

Quant Time: Aug 14 01:47:51 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N080724W.M
 Quant Title : SW846 8260
 QLast Update : Thu Aug 08 06:30:41 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	134350	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	262098	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	283434	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	128932	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	109808	57.421	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	114.840%	
35) Dibromofluoromethane	8.171	113	86701	52.997	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	106.000%	
50) Toluene-d8	10.571	98	337172	55.251	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	110.500%	
62) 4-Bromofluorobenzene	12.853	95	142518	59.904	ug/l	0.00
Spiked Amount	50.000	Range 64 - 133	Recovery	=	119.800%	
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	23765	31.759	ug/l	97
20) Methylene Chloride	5.277	84	6279	3.646	ug/l	89
43) Isopropyl Acetate	8.688	43	161225	32.590	ug/l #	89

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081324\
Data File : VN083266.D
Acq On : 13 Aug 2024 18:33
Operator : JC\MD
Sample : P3548-11
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
289

Quant Time: Aug 14 01:47:51 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N080724W.M
Quant Title : SW846 8260
QLast Update : Thu Aug 08 06:30:41 2024
Response via : Initial Calibration

