

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102924\
 Data File : VN084558.D
 Acq On : 29 Oct 2024 12:53
 Operator : JC\MD
 Sample : P4547-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BP-F-20

Quant Time: Oct 30 05:01:11 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N093024W.M
 Quant Title : SW846 8260
 QLast Update : Tue Oct 01 07:11:01 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.218	168	169528	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	295173	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	264179	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	103307	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	122489	48.731	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	97.460%
35) Dibromofluoromethane	8.165	113	96707	49.696	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.400%
50) Toluene-d8	10.565	98	365473	51.048	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	102.100%
62) 4-Bromofluorobenzene	12.847	95	123940	47.519	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	95.040%
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	31707	29.364	ug/l #	86
18) Methyl Acetate	5.030	43	3310	1.406	ug/l	93
20) Methylene Chloride	5.265	84	2627	1.197	ug/l #	91
43) Isopropyl Acetate	8.688	43	133121	23.437	ug/l #	86

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102924\
Data File : VN084558.D
Acq On : 29 Oct 2024 12:53
Operator : JC\MD
Sample : P4547-08
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
BP-F-20

Quant Time: Oct 30 05:01:11 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N093024W.M
Quant Title : SW846 8260
QLast Update : Tue Oct 01 07:11:01 2024
Response via : Initial Calibration

