

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN100721\
 Data File : VN068842.D
 Acq On : 7 Oct 2021 15:32
 Operator : JC/MD
 Sample : M4087-10
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-20211005-ROCK

Quant Time: Oct 08 05:59:19 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N100621W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 06 14:41:09 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	471597	50.00	ug/l	# 0.00
34) 1,4-Difluorobenzene	8.971	114	837113	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	812748	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	282170	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	330212	47.65	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	95.30%
35) Dibromofluoromethane	8.024	113	257450	46.82	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	93.64%
50) Toluene-d8	10.443	98	1072554	50.02	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	100.04%
62) 4-Bromofluorobenzene	12.734	95	343309	43.09	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	86.18%
Target Compounds						
						Qvalue
18) Methyl Acetate	4.859	43	9315	1.36	ug/l	# 60
20) Methylene Chloride	5.095	84	28338	3.39	ug/l	85
43) Isopropyl Acetate	8.560	43	60613	5.28	ug/l	# 83

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN100721\
Data File : VN068842.D
Acq On : 7 Oct 2021 15:32
Operator : JC/MD
Sample : M4087-10
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WC-20211005-ROCK

Quant Time: Oct 08 05:59:19 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N100621W.M
Quant Title : SW846 8260
QLast Update : Wed Oct 06 14:41:09 2021
Response via : Initial Calibration

