

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120222\
 Data File : VN075540.D
 Acq On : 02 Dec 2022 18:29
 Operator : JC\MD
 Sample : N5844-02 10X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MH-17

Quant Time: Dec 03 00:36:59 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112122W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 23 00:18:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	155100	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	9.106	114	254216	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	219214	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	92179	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	43131	51.806	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 103.620%		
35) Dibromofluoromethane	8.177	113	55603	55.678	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 111.360%		
50) Toluene-d8	10.571	98	112630	50.075	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 100.140%		
62) 4-Bromofluorobenzene	12.853	95	79854	48.100	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 96.200%		
Target Compounds						
						Qvalue
43) Isopropyl Acetate	8.700	43	9078	2.884	ug/l	# 91
52) Toluene	10.629	92	11188	2.604	ug/l	98
68) m/p-Xylenes	12.065	106	4455	1.420	ug/l	95

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120222\
Data File : VN075540.D
Acq On : 02 Dec 2022 18:29
Operator : JC\MD
Sample : N5844-02 10X
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MH-17

Quant Time: Dec 03 00:36:59 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112122W.M
Quant Title : SW846 8260
QLast Update : Wed Nov 23 00:18:03 2022
Response via : Initial Calibration

