

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN022422\
 Data File : VN071055.D
 Acq On : 24 Feb 2022 15:08
 Operator : JC\MD
 Sample : N1642-05 10X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 31457

Quant Time: Feb 25 02:43:55 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021522W.M
 Quant Title : SW846 8260
 QLast Update : Mon Feb 21 07:57:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.081	168	845282	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1344092	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1201747	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.674	152	421886	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	509932	52.811	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	105.620%
35) Dibromofluoromethane	8.016	113	394542	49.351	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	98.700%
50) Toluene-d8	10.439	98	1615525	50.319	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	100.640%
62) 4-Bromofluorobenzene	12.727	95	555543	48.409	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	96.820%
Target Compounds						
						Qvalue
16) Acetone	4.293	43	12847	2.416	ug/l	# 73
18) Methyl Acetate	4.863	43	21541	1.263	ug/l	91
40) Benzene	8.457	78	17348	0.448	ug/l	93
52) Toluene	10.498	92	9339	0.384	ug/l	92

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN022422\
Data File : VN071055.D
Acq On : 24 Feb 2022 15:08
Operator : JC\MD
Sample : N1642-05 10X
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
31457

Quant Time: Feb 25 02:43:55 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021522W.M
Quant Title : SW846 8260
QLast Update : Mon Feb 21 07:57:15 2022
Response via : Initial Calibration

