

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072921\
 Data File : VN067972.D
 Acq On : 29 Jul 2021 18:41
 Operator : JC/MD
 Sample : M3147-19
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 152140-DDC-1-PDA

Quant Time: Jul 30 04:45:30 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072921W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jul 29 15:02:35 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	299246	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	570265	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	537743	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	174240	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	235881	55.680	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	111.360%
35) Dibromofluoromethane	8.024	113	174925	52.634	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	105.260%
50) Toluene-d8	10.446	98	722065	47.303	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	94.600%
62) 4-Bromofluorobenzene	12.733	95	225283	41.111	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	82.220%
Target Compounds						
						Qvalue
16) Acetone	4.304	43	4888	3.162	ug/l	99
44) Trichloroethene	9.223	130	1816	0.460	ug/l	70
64) Tetrachloroethene	10.979	164	59833	15.536	ug/l	92

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072921\
Data File : VN067972.D
Acq On : 29 Jul 2021 18:41
Operator : JC/MD
Sample : M3147-19
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
152140-DDC-1-PDA

Quant Time: Jul 30 04:45:30 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072921W.M
Quant Title : SW846 8260
QLast Update : Thu Jul 29 15:02:35 2021
Response via : Initial Calibration

