

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN011922\
 Data File : VN070707.D
 Acq On : 19 Jan 2022 18:19
 Operator : JC/MD
 Sample : IBLK
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 IBLK

Quant Time: Jan 20 03:42:38 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N011822W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jan 18 13:32:07 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	1100147	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	1786412	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	1593748	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	605763	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	719486	45.633	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	91.260%
35) Dibromofluoromethane	8.024	113	507437	45.793	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	91.580%
50) Toluene-d8	10.440	98	2199480	47.116	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	94.240%
62) 4-Bromofluorobenzene	12.731	95	763139	45.876	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	91.760%
Target Compounds						
					Qvalue	
16) Acetone	4.296	43	20407	1.865	ug/l	99
18) Methyl Acetate	4.870	43	11134	0.517	ug/l #	57
52) Toluene	10.502	92	9416	0.272	ug/l	85

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

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