

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031221\
 Data File : VN066181.D
 Acq On : 12 Mar 2021 11:49
 Operator : JC/MD
 Sample : IBLK
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 IBLK

Quant Time: Mar 12 13:11:16 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N031221W.M
 Quant Title : SW846 8260
 QLast Update : Fri Mar 12 13:00:23 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.587	168	417960	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.518	114	638190	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.350	117	538115	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.286	152	196991	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.949	65	293429	43.51	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	87.02%
35) Dibromofluoromethane	7.509	113	206634	47.99	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	95.98%
50) Toluene-d8	10.031	98	774660	44.42	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	88.84%
62) 4-Bromofluorobenzene	12.345	95	240615	37.50	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	75.00%
Target Compounds						
					Qvalue	
13) Acrolein	3.535	56	10199	Below Cal		93
16) Acetone	3.747	43	5719	Below Cal		97
59) 2-Hexanone	10.698	43	15158	Below Cal		99
64) Tetrachloroethene	10.578	164	2027	0.43 ug/l #		77
65) Chlorobenzene	11.371	112	4527	0.41 ug/l		95

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

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