

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101221\
 Data File : VN068864.D
 Acq On : 12 Oct 2021 14:20
 Operator : JC/MD
 Sample : IBLK
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 IBLK

Quant Time: Oct 13 05:52:53 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 13 05:44:40 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	486407	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	807471	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	712954	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	272888	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	302097	47.106	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	94.220%
35) Dibromofluoromethane	8.024	113	243267	51.782	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	103.560%
50) Toluene-d8	10.443	98	984358	52.367	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	104.740%
62) 4-Bromofluorobenzene	12.734	95	337085	47.531	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	95.060%

Target Compounds					Qvalue	
74) N-amyl acetate	12.396	43	4343	0.505	ug/l	# 86
82) 4-Chlorotoluene	13.106	91	4560	0.300	ug/l	95
87) 1,3-Dichlorobenzene	13.621	146	3002	0.346	ug/l	95
89) n-Butylbenzene	13.946	91	6750	0.443	ug/l	98
91) 1,2-Dichlorobenzene	13.991	146	3060	0.363	ug/l	92
92) 1,2-Dibromo-3-Chloropr...	14.611	75	662	0.523	ug/l	78
93) 1,2,4-Trichlorobenzene	15.268	180	5186	3.674	ug/l	99
95) Naphthalene	15.509	128	24868	5.626	ug/l	99
96) 1,2,3-Trichlorobenzene	15.702	180	5861	3.790	ug/l	96

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

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