

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122721\
 Data File : VN070317.D
 Acq On : 27 Dec 2021 18:35
 Operator : JC/MD
 Sample : IBLK
 Misc : 7.75g/10mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 IBLK

Quant Time: Dec 28 01:42:24 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
 Quant Title : SW846 8260
 QLast Update : Mon Dec 27 18:47:12 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.080	168	1108596	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.965	114	2128921	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	2342635	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	1084867	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.431	65	843586	52.92	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	105.84%
35) Dibromofluoromethane	8.016	113	609056	47.28	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	94.56%
50) Toluene-d8	10.440	98	2772762	52.69	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	105.38%
62) 4-Bromofluorobenzene	12.731	95	1222350	64.32	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	128.64%
Target Compounds						
					Qvalue	
74) N-amyl acetate	12.387	43	20775	0.49	ug/l	97
87) 1,3-Dichlorobenzene	13.616	146	8729	0.23	ug/l	85
89) n-Butylbenzene	13.943	91	24054	0.38	ug/l	96
91) 1,2-Dichlorobenzene	13.986	146	10318	0.27	ug/l	94
93) 1,2,4-Trichlorobenzene	15.262	180	18072	1.01	ug/l	96

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

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