

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN110724\
 Data File : VN084725.D
 Acq On : 07 Nov 2024 16:45
 Operator : JC\MD
 Sample : VIBLK
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VIBLK

Quant Time: Nov 08 00:29:10 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N103024W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 31 18:45:38 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.218	168	183285	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	323271	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	295746	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	139802	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	126157	47.656	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	95.320%	
35) Dibromofluoromethane	8.165	113	103897	47.482	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	94.960%	
50) Toluene-d8	10.565	98	378616	46.972	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	93.940%	
62) 4-Bromofluorobenzene	12.847	95	153696	51.020	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	=	102.040%	
Target Compounds						
					Qvalue	
67) Ethyl Benzene	11.959	91	11468	1.038	ug/l	96
78) n-propylbenzene	13.035	91	15311	1.334	ug/l	98
95) Naphthalene	15.641	128	12444	1.538	ug/l #	92

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN110724\
Data File : VN084725.D
Acq On : 07 Nov 2024 16:45
Operator : JC\MD
Sample : VIBLK
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VIBLK

Quant Time: Nov 08 00:29:10 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N103024W.M
Quant Title : SW846 8260
QLast Update : Thu Oct 31 18:45:38 2024
Response via : Initial Calibration

