

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU052722\
 Data File : VU048699.D
 Acq On : 27 May 2022 17:25
 Operator : SY/MD
 Sample : VIBLK
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 VIBLK

Quant Time: May 28 00:21:54 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U052622W.M
 Quant Title : SW846 8260
 QLast Update : Fri May 27 06:42:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.375	168	217957	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	360728	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	374887	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	212790	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	167819	51.269	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 102.540%		
35) Dibromofluoromethane	5.292	113	139411	54.171	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 108.340%		
50) Toluene-d8	7.899	98	432541	46.887	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 93.780%		
62) 4-Bromofluorobenzene	10.632	95	199965	57.315	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 114.640%		
Target Compounds						
						Qvalue
40) Benzene	5.777	78	4148	0.374	ug/l	# 89
67) Ethyl Benzene	9.571	91	44860	3.355	ug/l	97
68) m/p-Xylenes	9.697	106	12914	2.469	ug/l	96
73) Isopropylbenzene	10.484	105	3229	0.235	ug/l	99
78) n-propylbenzene	10.906	91	8919	0.545	ug/l	99
80) 1,3,5-Trimethylbenzene	11.092	105	5842	0.489	ug/l	96
84) 1,2,4-Trimethylbenzene	11.468	105	44721	4.578	ug/l	97

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

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