

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU052722\
 Data File : VU048692.D
 Acq On : 27 May 2022 14:41
 Operator : SY/MD
 Sample : N3056-03
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MW-13D

Quant Time: May 28 00:19:06 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	205035	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	356600	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	367001	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	208491	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	167043	54.248	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 108.500%	
35) Dibromofluoromethane	5.292	113	143646	56.463	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 112.920%	
50) Toluene-d8	7.899	98	428471	46.984	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 93.960%	
62) 4-Bromofluorobenzene	10.632	95	196823	57.068	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 114.140%	
Target Compounds						
					Qvalue	
16) Acetone	2.633	43	2615	1.666	ug/l #	79
30) Chloroform	5.089	83	3619	0.772	ug/l	96
64) Tetrachloroethene	8.555	164	3708	0.354	ug/l	90
67) Ethyl Benzene	9.571	91	20612	1.574	ug/l	100
73) Isopropylbenzene	10.488	105	6299	0.468	ug/l	100
78) n-propylbenzene	10.906	91	8282	0.516	ug/l	99
84) 1,2,4-Trimethylbenzene	11.468	105	17047	2.506	ug/l	99

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

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