

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU052722\
 Data File : VU048723.D
 Acq On : 28 May 2022 03:58
 Operator : SY/MD
 Sample : N3033-12
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 TW-22

Quant Time: May 30 03:01:18 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	225214	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	374378	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	377300	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	227542	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	168359	49.777	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	99.560%
35) Dibromofluoromethane	5.292	113	143324	53.661	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	107.320%
50) Toluene-d8	7.899	98	438797	45.831	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	91.660%#
62) 4-Bromofluorobenzene	10.632	95	210396	58.106	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	116.220%
Target Compounds						
					Qvalue	
16) Acetone	2.630	43	3636	2.109	ug/l	98
30) Chloroform	5.093	83	2146	0.417	ug/l #	77
67) Ethyl Benzene	9.571	91	14386	1.069	ug/l	95
68) m/p-Xylenes	9.697	106	11273	2.142	ug/l	100
69) o-Xylene	10.102	106	3624	0.701	ug/l	92
84) 1,2,4-Trimethylbenzene	11.468	105	16695	2.371	ug/l	95

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

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