

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU052722\
 Data File : VU048725.D
 Acq On : 28 May 2022 04:45
 Operator : SY/MD
 Sample : N2934-07
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 RB-220512

Quant Time: May 30 03:01:53 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	213439	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	343295	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	349638	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	203401	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	162799	50.788	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 101.580%	
35) Dibromofluoromethane	5.292	113	138263	56.454	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 112.900%	
50) Toluene-d8	7.899	98	401297	45.709	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 91.420%#	
62) 4-Bromofluorobenzene	10.632	95	185305	55.811	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 111.620%	
Target Compounds						
					Qvalue	
16) Acetone	2.626	43	3764	2.304	ug/l	97
67) Ethyl Benzene	9.571	91	5703	0.457	ug/l	99
68) m/p-Xylenes	9.697	106	4635	0.950	ug/l	89
69) o-Xylene	10.105	106	1338	0.279	ug/l	97
84) 1,2,4-Trimethylbenzene	11.472	105	7705	1.798	ug/l	98

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

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