

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU052722\
 Data File : VU048726.D
 Acq On : 28 May 2022 05:08
 Operator : SY/MD
 Sample : N2934-08
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 RB-220513

Quant Time: May 30 03:02:09 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	208073	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	348438	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	352489	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	193385	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	153138	49.006	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	98.020%
35) Dibromofluoromethane	5.292	113	130300	52.417	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	104.840%
50) Toluene-d8	7.899	98	396520	44.499	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	89.000%#
62) 4-Bromofluorobenzene	10.632	95	179726	53.331	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	106.660%
Target Compounds						
					Qvalue	
16) Acetone	2.633	43	2473	1.553	ug/l #	86
67) Ethyl Benzene	9.571	91	4236	0.337	ug/l	96
68) m/p-Xylenes	9.697	106	3194	0.649	ug/l	91
69) o-Xylene	10.102	106	977	0.202	ug/l	87
84) 1,2,4-Trimethylbenzene	11.468	105	5565	1.651	ug/l	97

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

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