

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060122\
 Data File : VU048786.D
 Acq On : 01 Jun 2022 17:43
 Operator : SY/MD
 Sample : N3075-05
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 WELL-1

Quant Time: Jun 02 06:02:00 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	274295	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	468409	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	438673	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	222421	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	213888	51.922	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 103.840%	
35) Dibromofluoromethane	5.288	113	171374	51.283	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 102.560%	
50) Toluene-d8	7.899	98	603633	50.391	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 100.780%	
62) 4-Bromofluorobenzene	10.632	95	233827	51.614	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 103.220%	
Target Compounds						
					Qvalue	
16) Acetone	2.626	43	2146	1.022	ug/l	90
18) Methyl Acetate	2.957	43	1834	0.348	ug/l #	78
39) Methylcyclohexane	6.761	83	2496	0.488	ug/l	92
47) Bromodichloromethane	7.111	83	2179	0.398	ug/l #	88
52) Toluene	7.976	92	3806	0.442	ug/l	92
60) Dibromochloromethane	8.809	129	6900	1.619	ug/l	88
67) Ethyl Benzene	9.571	91	3868	0.247	ug/l	90
68) m/p-Xylenes	9.693	106	5128	0.838	ug/l	97
69) o-Xylene	10.105	106	1297	0.216	ug/l	84
71) Bromoform	10.295	173	12365	3.797	ug/l #	98
84) 1,2,4-Trimethylbenzene	11.471	105	6011	1.623	ug/l	99
95) Naphthalene	14.092	128	16499	3.427	ug/l	97

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

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