

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070124\
 Data File : VX042110.D
 Acq On : 01 Jul 2024 13:35
 Operator : JC/MD
 Sample : P3073-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW14-20240625

Quant Time: Jul 02 02:04:30 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jun 21 23:51:58 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	5.550	168	140588	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.757	114	215496	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	215533	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	122242	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	68508	42.259	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	84.520%
35) Dibromofluoromethane	5.385	113	66814	43.696	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	87.400%
50) Toluene-d8	8.647	98	251420	46.805	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	93.600%
62) 4-Bromofluorobenzene	11.079	95	100539	50.774	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	101.540%
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	2258	3.172	ug/l	# 79
17) Carbon Disulfide	2.502	76	1153	0.466	ug/l	# 90
20) Methylene Chloride	2.782	84	2152	1.381	ug/l	90
40) Benzene	6.044	78	1471	0.269	ug/l	98
51) 4-Methyl-2-Pentanone	8.647	43	880	0.393	ug/l	# 1
67) Ethyl Benzene	10.195	91	7705	1.097	ug/l	98
69) o-Xylene	10.640	106	1580	0.563	ug/l	90
73) Isopropylbenzene	10.963	105	25866	3.480	ug/l	96
78) n-propylbenzene	11.305	91	20059	2.487	ug/l	97
80) 1,3,5-Trimethylbenzene	11.451	105	6722	1.086	ug/l	90
84) 1,2,4-Trimethylbenzene	11.750	105	242607	39.060	ug/l	91
86) p-Isopropyltoluene	12.012	119	3864	0.576	ug/l	# 72
95) Naphthalene	13.780	128	8359543	940.307	ug/l	# 92

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

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