

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070324\
 Data File : VX042170.D
 Acq On : 03 Jul 2024 17:32
 Operator : JC/MD
 Sample : P3163-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TOWER-2

Quant Time: Jul 04 07:29:37 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	140807	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	218382	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	212972	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	101852	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	68633	42.270	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	84.540%
35) Dibromofluoromethane	5.391	113	67847	43.785	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	87.580%
50) Toluene-d8	8.647	98	252183	46.326	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	92.660%
62) 4-Bromofluorobenzene	11.079	95	92849	46.271	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	92.540%
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	9908	13.897	ug/l	# 83
20) Methylene Chloride	2.782	84	1165	0.746	ug/l	# 79
30) Chloroform	5.093	83	22896	9.639	ug/l	97
47) Bromodichloromethane	7.824	83	6409	3.511	ug/l	# 86
60) Dibromochloromethane	9.525	129	3344	2.091	ug/l	99
71) Bromoform	10.799	173	844	1.454	ug/l	# 93

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

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