

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092922\
 Data File : VX031768.D
 Acq On : 29 Sep 2022 13:13
 Operator : JC/MD
 Sample : N4863-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-7

Quant Time: Sep 30 04:30:21 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092622W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 27 01:44:18 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	70855	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	108331	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	117001	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	74704	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	73927	49.159	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	98.320%
35) Dibromofluoromethane	5.391	113	62365	48.316	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	96.640%
50) Toluene-d8	8.653	98	217737	46.529	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	93.060%
62) 4-Bromofluorobenzene	11.079	95	76776	44.763	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	89.520%
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.983	59	257646	1475.969	ug/l	99
19) Methyl tert-butyl Ether	3.111	73	1016556	303.921	ug/l	96
22) Diisopropyl ether	3.757	45	17206	5.395	ug/l #	80
42) 1,2-Dichloroethane	6.098	62	1539	0.874	ug/l	83

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

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