

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092222\
 Data File : VX031583.D
 Acq On : 22 Sep 2022 20:14
 Operator : JC/MD
 Sample : N4751-09
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TB-9-21-22

Quant Time: Sep 23 01:51:10 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 20 09:45:29 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	83776	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	131666	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	116954	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	53112	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	82922	53.957	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 107.920%	
35) Dibromofluoromethane	5.391	113	69953	51.422	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 102.840%	
50) Toluene-d8	8.647	98	259656	49.499	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 99.000%	
62) 4-Bromofluorobenzene	11.079	95	79467	44.723	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 89.440%	
Target Compounds						
					Qvalue	
67) Ethyl Benzene	10.195	91	4900	0.769	ug/l	94
68) m/p-Xylenes	10.305	106	5248	2.051	ug/l	99
69) o-Xylene	10.640	106	1919	0.786	ug/l	79
78) n-propylbenzene	11.305	91	1562	0.342	ug/l	96
80) 1,3,5-Trimethylbenzene	11.451	105	2432	0.723	ug/l	95
84) 1,2,4-Trimethylbenzene	11.756	105	8143	2.456	ug/l	94

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

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