

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061022\
 Data File : VN072771.D
 Acq On : 10 Jun 2022 17:48
 Operator : JC\MD
 Sample : N3293-10
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 52322-C

Quant Time: Jun 11 01:55:44 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jun 09 05:17:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.081	168	162913	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	8.963	114	313060	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	290669	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.674	152	114595	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	243804	55.759	ug/l	0.00
Spiked Amount 50.000	Range 61	- 141	Recovery	=	111.520%	
35) Dibromofluoromethane	8.016	113	177134	57.183	ug/l	0.00
Spiked Amount 50.000	Range 69	- 133	Recovery	=	114.360%	
50) Toluene-d8	10.439	98	578965	46.845	ug/l	0.00
Spiked Amount 50.000	Range 65	- 126	Recovery	=	93.700%	
62) 4-Bromofluorobenzene	12.727	95	232799	54.893	ug/l	0.00
Spiked Amount 50.000	Range 58	- 135	Recovery	=	109.780%	
Target Compounds						
						Qvalue
11) Tert butyl alcohol	5.363	59	134671	129.113	ug/l	99
16) Acetone	4.281	43	101067	48.238	ug/l	92
25) 2-Butanone	7.334	43	53904	15.994	ug/l	# 84
51) 4-Methyl-2-Pentanone	10.328	43	104557	15.432	ug/l	91
52) Toluene	10.504	92	814836	100.020	ug/l	100
67) Ethyl Benzene	11.845	91	42715	2.933	ug/l	92
68) m/p-Xylenes	11.945	106	54340	10.499	ug/l	91
69) o-Xylene	12.280	106	21210	4.077	ug/l	85

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

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