

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072423\
 Data File : VX036639.D
 Acq On : 24 Jul 2023 18:56
 Operator : JC/MD
 Sample : 03744-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 GWBR-04-440-455-072423

Quant Time: Jul 25 06:39:21 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 21 12:14:12 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	158771	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	283943	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	258256	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	114759	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	108179	46.048	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	92.100%		
35) Dibromofluoromethane	5.391	113	89976	46.102	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	92.200%		
50) Toluene-d8	8.647	98	341796	49.385	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	98.780%		
62) 4-Bromofluorobenzene	11.079	95	128232	49.109	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	98.220%		
Target Compounds					Qvalue	
12) 1,1-Dichloroethene	2.319	96	24337	13.689	ug/l	89
16) Acetone	2.386	43	2616	2.683	ug/l	92
24) 1,1-Dichloroethane	3.617	63	5152	1.382	ug/l #	90
27) cis-1,2-Dichloroethene	4.495	96	27736	11.752	ug/l	84
42) 1,2-Dichloroethane	6.092	62	2807	0.907	ug/l	97
44) Trichloroethene	7.129	130	21646	9.785	ug/l	98
51) 4-Methyl-2-Pentanone	8.574	43	2079	0.663	ug/l	93
52) Toluene	8.720	92	285360	53.810	ug/l	96

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

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