

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX081821\
 Data File : VX023813.D
 Acq On : 18 Aug 2021 13:52
 Operator : JC/MD
 Sample : M3433-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-TT-VPB-RW7-GW-150-152

Quant Time: Aug 18 14:15:16 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 17 11:57:51 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	132635	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	231457	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	218237	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	94433	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.971	65	95133	51.391	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 102.780%			
35) Dibromofluoromethane	5.404	113	74520	48.644	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 97.280%			
50) Toluene-d8	8.653	98	283045	48.544	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 97.080%			
62) 4-Bromofluorobenzene	11.085	95	100909	47.012	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 94.020%			
Target Compounds					Qvalue	
16) Acetone	2.386	43	10120	12.227	ug/l #	84
25) 2-Butanone	4.581	43	2620	2.090	ug/l	91
30) Chloroform	5.105	83	2253	0.818	ug/l	97
40) Benzene	6.050	78	1631	0.257	ug/l	93
52) Toluene	8.726	92	1055	0.257	ug/l	91
84) 1,2,4-Trimethylbenzene	11.756	105	1328	0.225	ug/l	99
95) Naphthalene	13.780	128	6571	2.209	ug/l	96

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

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