

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060921\
 Data File : VX022614.D
 Acq On : 09 Jun 2021 15:21
 Operator : JC/MD
 Sample : M2652-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PDB-BLANK-20210608

Quant Time: Jun 10 05:42:52 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060721W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jun 07 14:27:50 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.568	168	59326	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.775	114	120549	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.061	117	105491	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	41508	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	50815	48.181	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	96.360%
35) Dibromofluoromethane	5.397	113	35206	46.678	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	93.360%
50) Toluene-d8	8.653	98	145952	48.498	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	97.000%
62) 4-Bromofluorobenzene	11.085	95	51148	45.725	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	91.460%
Target Compounds						
						Qvalue
16) Acetone	2.386	43	4919	10.380	ug/l	94
25) 2-Butanone	4.586	43	2121	2.708	ug/l	# 77
27) cis-1,2-Dichloroethene	4.501	96	8087	8.130	ug/l	94
37) Ethyl Acetate	4.733	43	3764	2.516	ug/l	# 91
44) Trichloroethene	7.135	130	5284	6.223	ug/l	76

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

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