

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120922\
 Data File : VX033344.D
 Acq On : 09 Dec 2022 13:54
 Operator : JC/MD
 Sample : N5944-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 GW-BR-04-100-114-120722

Quant Time: Dec 09 14:42:47 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120722W.M
 Quant Title : SW846 8260
 QLast Update : Wed Dec 07 12:43:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	156126	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	248446	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	217900	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	97686	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	48797	48.719	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	97.440%
35) Dibromofluoromethane	5.385	113	53491	47.868	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	95.740%
50) Toluene-d8	8.647	98	112342	48.057	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	96.120%
62) 4-Bromofluorobenzene	11.079	95	84733	47.405	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	94.820%
Target Compounds						Qvalue
12) 1,1-Dichloroethene	2.325	96	5609	3.963	ug/l	97
16) Acetone	2.386	43	1933	2.608	ug/l	92
27) cis-1,2-Dichloroethene	4.495	96	24837	13.422	ug/l	90
40) Benzene	6.044	78	1442	0.227	ug/l #	87
44) Trichloroethene	7.129	130	39277	22.426	ug/l	99
52) Toluene	8.720	92	159072	38.731	ug/l	98
65) Chlorobenzene	10.073	112	1087	0.246	ug/l	97

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

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