

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062122\
 Data File : VX029671.D
 Acq On : 21 Jun 2022 18:53
 Operator : JC/MD
 Sample : N3342-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-TT-MW149S1-230-232

Quant Time: Jun 22 01:12:27 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jun 20 01:31:01 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	253996	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	430644	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	374174	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	165177	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	144281	42.913	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	85.820%
35) Dibromofluoromethane	5.391	113	128292	45.713	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	91.420%
50) Toluene-d8	8.653	98	484195	47.227	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	94.460%
62) 4-Bromofluorobenzene	11.085	95	177334	45.866	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	91.740%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	22894	17.083	ug/l	91
19) Methyl tert-butyl Ether	3.117	73	11895	1.332	ug/l #	89
20) Methylene Chloride	2.794	84	1457	0.488	ug/l #	91
25) 2-Butanone	4.568	43	13004	6.139	ug/l	96
30) Chloroform	5.111	83	1924	0.368	ug/l #	78
40) Benzene	6.050	78	3026	0.269	ug/l	99
44) Trichloroethene	7.129	130	5368	0.566	ug/l	90
52) Toluene	8.720	92	1500	0.208	ug/l	96
64) Tetrachloroethene	9.287	164	1180	0.436	ug/l #	58

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

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