

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060222\
 Data File : VX029153.D
 Acq On : 02 Jun 2022 22:32
 Operator : JC/MD
 Sample : N3164-01 50X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 06BR-20220531

Quant Time: Jun 03 04:35:06 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X053122W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 01 04:45:42 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	317088	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	580175	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	564453	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	268536	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	200700	56.130	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	112.260%
35) Dibromofluoromethane	5.391	113	191030	49.571	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	99.140%
50) Toluene-d8	8.653	98	681089	48.323	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	96.640%
62) 4-Bromofluorobenzene	11.085	95	272475	50.012	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	100.020%
Target Compounds						
						Qvalue
17) Carbon Disulfide	2.514	76	2814	0.359	ug/l	# 89
20) Methylene Chloride	2.794	84	975	0.297	ug/l	# 76
27) cis-1,2-Dichloroethene	4.495	96	3475	0.904	ug/l	84
37) Ethyl Acetate	4.721	43	17394	3.620	ug/l	99
95) Naphthalene	13.780	128	4372	0.245	ug/l	96
96) 1,2,3-Trichlorobenzene	13.963	180	1423	0.259	ug/l	84

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

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