

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX033123\
 Data File : VX034880.D
 Acq On : 31 Mar 2023 15:58
 Operator : JC/MD
 Sample : 02048-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-8

Quant Time: Apr 03 01:19:53 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X032923W.M
 Quant Title : SW846 8260
 QLast Update : Sat Apr 01 01:50:18 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	218360	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	372293	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	331031	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	146898	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	158178	47.366	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	=	94.740%	
35) Dibromofluoromethane	5.385	113	117495	47.715	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	95.440%	
50) Toluene-d8	8.647	98	450100	49.522	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	=	99.040%	
62) 4-Bromofluorobenzene	11.079	95	160479	45.627	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	91.260%	
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	3940	2.904	ug/l	98
17) Carbon Disulfide	2.514	76	3219	0.573	ug/l	100
19) Methyl tert-butyl Ether	3.111	73	96281	11.601	ug/l	100
40) Benzene	6.050	78	2114	0.202	ug/l #	86
52) Toluene	8.720	92	2162	0.317	ug/l	99
67) Ethyl Benzene	10.195	91	5459	0.438	ug/l	94
68) m/p-Xylenes	10.305	106	1666	0.350	ug/l	96
69) o-Xylene	10.646	106	957	0.205	ug/l	95
70) Styrene	10.659	104	2648	0.355	ug/l #	83

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

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