

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX103122\
 Data File : VX032489.D
 Acq On : 31 Oct 2022 15:43
 Operator : JC/MD
 Sample : N5340-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-8-102722

Quant Time: Nov 01 05:11:08 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 27 08:28:06 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	124415	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	202764	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	180282	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	84577	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	82199	47.549	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	95.100%
35) Dibromofluoromethane	5.391	113	66317	45.852	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	91.700%
50) Toluene-d8	8.647	98	242258	48.170	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	96.340%
62) 4-Bromofluorobenzene	11.079	95	94462	48.084	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	96.160%
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.971	59	48404	192.375	ug/l	98
16) Acetone	2.386	43	9407	14.972	ug/l	93
19) Methyl tert-butyl Ether	3.111	73	3004	0.685	ug/l	93
29) Tetrahydrofuran	5.013	42	14966	24.111	ug/l	100
30) Chloroform	5.099	83	5211	1.919	ug/l	98
40) Benzene	6.044	78	1300	0.228	ug/l	93

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

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