

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101422\
 Data File : VX032135.D
 Acq On : 14 Oct 2022 15:30
 Operator : JC/MD
 Sample : N5145-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-7-101322

Quant Time: Oct 17 02:25:34 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100622W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 06 14:28:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	88600	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	133452	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	137272	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	83832	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	98364	53.757	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 107.520%	
35) Dibromofluoromethane	5.385	113	72692	48.121	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 96.240%	
50) Toluene-d8	8.646	98	232774	42.823	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 85.640%#	
62) 4-Bromofluorobenzene	11.079	95	93636	48.727	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 97.460%	
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.989	59	29079	127.733	ug/l	100
16) Acetone	2.385	43	58357	93.841	ug/l	93
25) 2-Butanone	4.568	43	5620	5.781	ug/l	97
29) Tetrahydrofuran	5.019	42	4503	7.179	ug/l	98
30) Chloroform	5.098	83	6239	2.165	ug/l	99
51) 4-Methyl-2-Pentanone	8.579	43	3937	2.076	ug/l	96

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

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