

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX040822\
 Data File : VX028025.D
 Acq On : 08 Apr 2022 14:24
 Operator : JC/MD
 Sample : N2333-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-9D

Quant Time: Apr 11 01:16:51 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X040722W.M
 Quant Title : SW846 8260
 QLast Update : Thu Apr 07 18:12:28 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	80463	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	142670	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	130744	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	61064	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	70957	58.093	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	116.180%
35) Dibromofluoromethane	5.391	113	50132	52.348	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	104.700%
50) Toluene-d8	8.653	98	169313	49.550	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	99.100%
62) 4-Bromofluorobenzene	11.079	95	71791	49.423	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	98.840%
Target Compounds						
					Qvalue	
3) Chloromethane	1.295	50	663	2.183	ug/l	99
11) Tert butyl alcohol	2.953	59	9497	44.129	ug/l	99
19) Methyl tert-butyl Ether	3.117	73	7483	2.506	ug/l	98
22) Diisopropyl ether	3.764	45	1319	0.444	ug/l #	80
40) Benzene	6.044	78	3622	0.921	ug/l #	82

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

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