

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX052022\
 Data File : VX028863.D
 Acq On : 20 May 2022 20:03
 Operator : JC/MD
 Sample : N2958-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RW8-MW01D3-20220517

Quant Time: May 20 23:34:00 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
 Quant Title : SW846 8260
 QLast Update : Thu May 12 14:58:53 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	252508	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	446984	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	403756	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	176217	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	175071	50.580	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	101.160%
35) Dibromofluoromethane	5.391	113	142896	48.794	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	97.580%
50) Toluene-d8	8.653	98	528483	47.864	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	95.720%
62) 4-Bromofluorobenzene	11.079	95	186176	45.145	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	90.280%
Target Compounds						
						Qvalue
16) Acetone	2.380	43	4882	3.286	ug/l	# 81
29) Tetrahydrofuran	5.007	42	57183	37.728	ug/l	92
68) m/p-Xylenes	10.305	106	2872	0.480	ug/l	95
84) 1,2,4-Trimethylbenzene	11.756	105	3205	0.287	ug/l	99

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

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