

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX040422\
 Data File : VX027953.D
 Acq On : 04 Apr 2022 12:59
 Operator : JC/MD
 Sample : N2249-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-6S

Quant Time: Apr 05 05:48:23 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X032922W.M
 Quant Title : SW846 8260
 QLast Update : Wed Mar 30 04:39:27 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	93710	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	148913	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	119348	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	50585	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	65336	54.036	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 108.080%			
35) Dibromofluoromethane	5.391	113	48960	50.936	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 101.880%			
50) Toluene-d8	8.653	98	164963	48.778	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 97.560%			
62) 4-Bromofluorobenzene	11.085	95	67655	47.074	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 94.140%			
Target Compounds						
40) Benzene	6.050	78	2389	0.651	ug/l	93
67) Ethyl Benzene	10.201	91	2053	0.481	ug/l #	86
68) m/p-Xylenes	10.305	106	1814	1.156	ug/l	74
69) o-Xylene	10.640	106	633	0.413	ug/l	80
84) 1,2,4-Trimethylbenzene	11.756	105	1090	0.334	ug/l	94

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

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