

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060822\
 Data File : VX029320.D
 Acq On : 08 Jun 2022 14:50
 Operator : JC/MD
 Sample : N3224-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-13

Quant Time: Jun 09 06:59:52 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 08 03:00:38 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	281376	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	473554	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	397372	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	165082	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	235721	55.217	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	110.440%
35) Dibromofluoromethane	5.391	113	192298	47.765	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	95.540%
50) Toluene-d8	8.653	98	704526	49.481	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	98.960%
62) 4-Bromofluorobenzene	11.085	95	257286	46.175	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	92.340%
Target Compounds						
					Qvalue	
5) Bromomethane	1.611	94	882	0.333	ug/l	94
16) Acetone	2.386	43	27118	11.914	ug/l	90
25) 2-Butanone	4.593	43	4572	1.270	ug/l	94
67) Ethyl Benzene	10.195	91	4961	0.258	ug/l #	87
68) m/p-Xylenes	10.305	106	1825	0.245	ug/l	94
78) n-propylbenzene	11.305	91	4291	0.291	ug/l	95

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

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