

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX052424\
 Data File : VX041649.D
 Acq On : 24 May 2024 12:57
 Operator : JC/MD
 Sample : P2602-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-16B-67-052224

Quant Time: May 25 04:22:40 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X043024W.M
 Quant Title : SW846 8260
 QLast Update : Wed May 08 06:03:16 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	134015	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	201198	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	197106	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	96595	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	76389	52.359	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	104.720%
35) Dibromofluoromethane	5.385	113	63711	46.727	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	93.460%
50) Toluene-d8	8.647	98	234977	54.450	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	108.900%
62) 4-Bromofluorobenzene	11.079	95	87999	52.677	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	105.360%
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	4823	3.653	ug/l	91
30) Chloroform	5.099	83	20567	9.463	ug/l	98
51) 4-Methyl-2-Pentanone	8.580	43	4906	2.673	ug/l	92
52) Toluene	8.720	92	2483	0.808	ug/l	100
68) m/p-Xylenes	10.299	106	880	0.333	ug/l	96

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

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