

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060524\
 Data File : VX041752.D
 Acq On : 05 Jun 2024 16:02
 Operator : JC/MD
 Sample : P2718-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-1

Quant Time: Jun 06 02:21:33 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X052924W.M
 Quant Title : SW846 8260
 QLast Update : Thu May 30 07:36:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	140547	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	217582	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	229947	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	133643	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	82760	48.011	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	96.020%
35) Dibromofluoromethane	5.385	113	69861	46.439	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	92.880%
50) Toluene-d8	8.646	98	258936	49.261	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	98.520%
62) 4-Bromofluorobenzene	11.079	95	110578	56.998	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	114.000%
Target Compounds						
					Qvalue	
16) Acetone	2.385	43	6181	5.478	ug/l	97
67) Ethyl Benzene	10.201	91	4169	0.530	ug/l	100
68) m/p-Xylenes	10.305	106	3719	1.189	ug/l	97
69) o-Xylene	10.646	106	781	0.251	ug/l	89
78) n-propylbenzene	11.311	91	2993	0.329	ug/l	93
84) 1,2,4-Trimethylbenzene	11.756	105	8614	1.228	ug/l	95

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

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