

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042924\
 Data File : VX041280.D
 Acq On : 29 Apr 2024 17:04
 Operator : JC/MD
 Sample : P2270-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VNJ-240

Quant Time: Apr 30 02:13:08 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042224W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 23 05:30:28 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	173562	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	251411	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	234329	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	116339	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	89237	54.012	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 108.020%	
35) Dibromofluoromethane	5.385	113	74558	48.753	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 97.500%	
50) Toluene-d8	8.647	98	274094	53.122	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 106.240%	
62) 4-Bromofluorobenzene	11.079	95	99934	52.474	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 104.940%	
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	39843	66.252	ug/l	99
18) Methyl Acetate	2.709	43	1928	1.084	ug/l	94
20) Methylene Chloride	2.782	84	4516	3.075	ug/l	89
43) Isopropyl Acetate	6.349	43	83640	24.185	ug/l	99

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

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