

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071024\
 Data File : VX042278.D
 Acq On : 10 Jul 2024 19:27
 Operator : JC/MD
 Sample : P3183-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 N28660

Quant Time: Jul 11 01:03:31 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061824W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jun 21 23:51:58 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	137910	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	223069	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	196702	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	81755	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	85443	53.728	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 107.460%		
35) Dibromofluoromethane	5.385	113	81430	51.447	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 102.900%		
50) Toluene-d8	8.647	98	236118	42.464	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 84.920%#		
62) 4-Bromofluorobenzene	11.079	95	82409	40.205	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 80.420%#		
Target Compounds						
						Qvalue
11) Tert butyl alcohol	2.983	59	3285	10.884	ug/l	98
16) Acetone	2.386	43	74461	106.633	ug/l	91
18) Methyl Acetate	2.709	43	3445	1.808	ug/l	95
20) Methylene Chloride	2.788	84	6670	4.362	ug/l	86
25) 2-Butanone	4.581	43	11783	10.846	ug/l	91
40) Benzene	6.038	78	17928	3.163	ug/l	99
43) Isopropyl Acetate	6.355	43	9819	2.922	ug/l #	95

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

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