

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071024\
 Data File : VX042274.D
 Acq On : 10 Jul 2024 17:54
 Operator : JC/MD
 Sample : P3184-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MOO-23-00199

Quant Time: Jul 11 01:01:29 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	5.550	168	183816	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	293135	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	270457	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	129446	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	121830	57.477	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 114.960%		
35) Dibromofluoromethane	5.385	113	117379	56.434	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 112.860%		
50) Toluene-d8	8.647	98	369392	50.553	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 101.100%		
62) 4-Bromofluorobenzene	11.079	95	139088	51.638	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 103.280%		
Target Compounds						
						Qvalue
11) Tert butyl alcohol	2.983	59	29673	73.764	ug/l	# 94
16) Acetone	2.386	43	48441	52.046	ug/l	# 88
18) Methyl Acetate	2.709	43	4078	1.606	ug/l	97
20) Methylene Chloride	2.788	84	13392	6.571	ug/l	# 84
25) 2-Butanone	4.580	43	11893	8.213	ug/l	98
43) Isopropyl Acetate	6.342	43	82585	18.700	ug/l	# 93

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

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