

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100821\
 Data File : VX024680.D
 Acq On : 08 Oct 2021 19:07
 Operator : JC/MD
 Sample : M4154-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SW-CB-32

Manual Integrations
 APPROVED

SAM
 10/22/2021 1:35:47 PM

Quant Time: Oct 11 05:50:02 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092321W.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 23 18:24:34 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	139710	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	238703	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	231661	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	102300	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	101373	51.48	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 102.96%			
35) Dibromofluoromethane	5.391	113	80054	50.48	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 100.96%			
50) Toluene-d8	8.653	98	293166	50.23	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 100.46%			
62) 4-Bromofluorobenzene	11.085	95	116055	50.57	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 101.14%			
Target Compounds						
						Qvalue
16) Acetone	2.386	43	16870	10.59	ug/l	94
17) Carbon Disulfide	2.581	76	84363m	25.67	ug/l	
25) 2-Butanone	4.574	43	6390	4.71	ug/l	97
52) Toluene	8.720	92	10373	2.45	ug/l	100
86) p-Isopropyltoluene	12.012	119	5881	0.91	ug/l	87

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

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