

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091421\
 Data File : VX024211.D
 Acq On : 14 Sep 2021 16:41
 Operator : JC/MD
 Sample : M3729-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-TT-VPB-RW7-GW-760-762

Manual Integrations
 APPROVED

MMDadoda
 9/15/2021 5:54:27 PM

Quant Time: Sep 15 01:57:24 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091421W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 14 12:18:04 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	137415	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	235422	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.061	117	218907	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	87873	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.970	65	101440	51.768	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 103.540%			
35) Dibromofluoromethane	5.397	113	78787	49.902	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 99.800%			
50) Toluene-d8	8.653	98	288280	50.121	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 100.240%			
62) 4-Bromofluorobenzene	11.085	95	103940	46.436	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 92.880%			
Target Compounds						Qvalue
3) Chloromethane	1.300	50	1017	0.737	ug/l	97
9) 1,1,2-Trichlorotrifluo...	2.331	101	1084	0.801	ug/l	94
12) 1,1-Dichloroethene	2.318	96	1107	0.874	ug/l #	93
16) Acetone	2.386	43	11432	13.683	ug/l	88
17) Carbon Disulfide	2.514	76	1541	0.485	ug/l #	84
27) cis-1,2-Dichloroethene	4.507	96	699m	0.436	ug/l	
30) Chloroform	5.098	83	1410	0.500	ug/l #	68
44) Trichloroethene	7.135	130	47931	27.479	ug/l	94

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

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