

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092121\
 Data File : VX024391.D
 Acq On : 21 Sep 2021 17:12
 Operator : JC/MD
 Sample : M3835-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SW-CB-04-(1.0)

Manual Integrations
 APPROVED

MMDadoda
 9/23/2021 11:06:47 AM

Quant Time: Sep 22 04:51:37 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091421W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 21 10:52:48 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.568	168	151309	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	225875	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.061	117	219318	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	101656	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	103906	48.158	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	96.320%		
35) Dibromofluoromethane	5.397	113	89971	59.394	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	118.780%		
50) Toluene-d8	8.653	98	304916	55.254	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	110.500%		
62) 4-Bromofluorobenzene	11.085	95	106316	49.505	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	99.020%		
Target Compounds					Qvalue	
16) Acetone	2.386	43	69738	75.803	ug/l	91
25) 2-Butanone	4.574	43	10094	7.449	ug/l	90
40) Benzene	6.050	78	81003	13.600	ug/l	100
52) Toluene	8.726	92	64485	16.519	ug/l	97
65) Chlorobenzene	10.085	112	48784	10.812	ug/l	98
67) Ethyl Benzene	10.201	91	2271	0.288	ug/l	91
68) m/p-Xylenes	10.305	106	1892	0.626	ug/l	94
86) p-Isopropyltoluene	12.012	119	5626	0.869	ug/l	92
87) 1,3-Dichlorobenzene	11.975	146	1156	0.339	ug/l	96
88) 1,4-Dichlorobenzene	12.042	146	3921m	1.127	ug/l	
91) 1,2-Dichlorobenzene	12.335	146	4298	1.286	ug/l	98
95) Naphthalene	13.780	128	185196	26.122	ug/l	100

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

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