

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU040921\
 Data File : VU043003.D
 Acq On : 09 Apr 2021 15:32
 Operator : SY/MD
 Sample : M1977-01
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EB-01-040721

Manual Integrations
 APPROVED

MMDadoda
 4/12/2021 9:57:04 AM

Quant Time: Apr 09 16:26:17 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U040921W.M
 Quant Title : SW846 8260
 QLast Update : Fri Apr 09 09:26:47 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.382	168	118708	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	194304	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	192035	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	94517	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	98948	50.99	ug/l	0.00
Spiked Amount	50.000		Recovery	=	101.98%	
35) Dibromofluoromethane	5.298	113	68684	48.41	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.82%	
50) Toluene-d8	7.906	98	247669	48.45	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.90%	
62) 4-Bromofluorobenzene	10.639	95	92590	44.34	ug/l	0.00
Spiked Amount	50.000		Recovery	=	88.68%	
Target Compounds						
						Qvalue
16) Acetone	2.629	43	50067	46.01	ug/l	97
20) Methylene Chloride	3.054	84	2007	1.44	ug/l	81
25) 2-Butanone	4.710	43	11969	6.88	ug/l	96
29) Tetrahydrofuran	5.060	42	1542m	1.39	ug/l	
51) 4-Methyl-2-Pentanone	7.800	43	7623	11.38	ug/l	100

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

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