

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU110921\
 Data File : VU045665.D
 Acq On : 09 Nov 2021 17:51
 Operator : SY/MD
 Sample : M4594-04
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 22BR-20211108

Quant Time: Nov 10 02:01:07 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U110521W.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 05 13:35:16 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.375	168	139367	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	257888	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	266452	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	134372	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	113493	54.376	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 108.760%	
35) Dibromofluoromethane	5.292	113	87560	50.854	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 101.700%	
50) Toluene-d8	7.899	98	342186	53.446	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 106.900%	
62) 4-Bromofluorobenzene	10.636	95	132185	51.487	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 102.980%	
Target Compounds						
						Qvalue
16) Acetone	2.629	43	2313	2.177	ug/l	95
27) cis-1,2-Dichloroethene	4.671	96	47360	25.278	ug/l	91
44) Trichloroethene	6.546	130	30481	17.104	ug/l	95

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

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