

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU102221\
 Data File : VU045364.D
 Acq On : 22 Oct 2021 15:00
 Operator : SY/MD
 Sample : M4315-01
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MH-D-S

Manual Integrations
 APPROVED

MMDadoda
 10/25/2021 10:26:51 AM

Quant Time: Oct 23 01:12:23 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 28 12:32:13 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.375	168	231090	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	437234	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.427	117	398296	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	169840	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	181786	47.257	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	94.520%		
35) Dibromofluoromethane	5.292	113	131609	47.715	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	95.440%		
50) Toluene-d8	7.902	98	545719	50.074	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	100.140%		
62) 4-Bromofluorobenzene	10.645	95	192696	42.827	ug/l	0.01
Spiked Amount 50.000	Range 83 - 123		Recovery =	85.660%		
Target Compounds						Qvalue
16) Acetone	2.652	43	14480	6.341	ug/l	98
18) Methyl Acetate	2.964	43	4393	0.658	ug/l	94
20) Methylene Chloride	3.038	84	17203	5.002	ug/l	95
29) Tetrahydrofuran	5.079	42	2297m	1.189	ug/l	
43) Isopropyl Acetate	5.925	43	33737	3.707	ug/l	100
68) m/p-Xylenes	9.703	106	2403	0.403	ug/l	94

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

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