

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU102621\
 Data File : VU045409.D
 Acq On : 26 Oct 2021 18:33
 Operator : SY/MD
 Sample : M4360-06
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BP-VPB-RW9-GW-218-220

Quant Time: Oct 27 04:41:18 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	207979	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	388276	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	348533	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.815	152	143904	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	166646	48.135	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	96.260%		
35) Dibromofluoromethane	5.292	113	121884	49.761	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	99.520%		
50) Toluene-d8	7.899	98	486473	50.266	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	100.540%		
62) 4-Bromofluorobenzene	10.635	95	164513	41.174	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	82.340%#		
Target Compounds						Qvalue
3) Chloromethane	1.520	50	8993	2.696	ug/l	98
16) Acetone	2.639	43	3424	1.666	ug/l	90
18) Methyl Acetate	2.957	43	6280	1.046	ug/l	94

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

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