

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU102621\
 Data File : VU045407.D
 Acq On : 26 Oct 2021 17:46
 Operator : SY/MD
 Sample : M4360-04
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BP-VPB-RW9-GW-148-150

Quant Time: Oct 27 04:41:02 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	210780	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	399781	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	370361	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.815	152	165509	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	171496	48.877	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	97.760%
35) Dibromofluoromethane	5.292	113	124648	49.425	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.860%
50) Toluene-d8	7.899	98	505859	50.765	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	101.540%
62) 4-Bromofluorobenzene	10.635	95	180771	43.941	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	87.880%
Target Compounds						
					Qvalue	
3) Chloromethane	1.520	50	11261	3.331	ug/l	98
16) Acetone	2.639	43	6399	3.072	ug/l	95
18) Methyl Acetate	2.957	43	10392	1.707	ug/l	99

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

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