

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU102621\
 Data File : VU045410.D
 Acq On : 26 Oct 2021 18:56
 Operator : SY/MD
 Sample : M4360-07
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BP-VPB-RW9-GW-238-240

Quant Time: Oct 27 04:41:26 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	5.376	168	211544	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	399389	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.421	117	359511	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.813	152	145736	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.704	65	170134	48.314	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	96.620%
35) Dibromofluoromethane	5.292	113	124598	49.454	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.900%
50) Toluene-d8	7.900	98	495543	49.779	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	99.560%
62) 4-Bromofluorobenzene	10.636	95	170055	41.377	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	82.760%#
Target Compounds						
					Qvalue	
3) Chloromethane	1.517	50	10350	3.050	ug/l	99
16) Acetone	2.636	43	5039	2.411	ug/l	97

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

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