

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU102621\
 Data File : VU045406.D
 Acq On : 26 Oct 2021 17:22
 Operator : SY/MD
 Sample : M4360-03
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BP-VPB-RW9-GW-98-100

Quant Time: Oct 27 04:40:53 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	5.378	168	210457	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	399478	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	370577	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.815	152	167559	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.706	65	172233	49.163	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	98.320%		
35) Dibromofluoromethane	5.295	113	125180	49.674	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	99.340%		
50) Toluene-d8	7.899	98	500874	50.303	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	100.600%		
62) 4-Bromofluorobenzene	10.635	95	182102	44.298	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	88.600%		
Target Compounds						Qvalue
3) Chloromethane	1.520	50	6925	2.051	ug/l	93
16) Acetone	2.636	43	4353	2.093	ug/l	91
24) 1,1-Dichloroethane	3.874	63	8957	1.655	ug/l	93

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

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