

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU111221\
 Data File : VU045773.D
 Acq On : 12 Nov 2021 16:44
 Operator : SY/MD
 Sample : M4668-03
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BP-VPB-RW9-GW-638-640

Quant Time: Nov 15 05:21:22 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U110521W.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 05 13:35:16 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.378	168	117196	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	219805	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	231034	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.815	152	116177	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	96638	55.060	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 110.120%	
35) Dibromofluoromethane	5.292	113	77379	52.728	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 105.460%	
50) Toluene-d8	7.902	98	291544	53.426	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 106.860%	
62) 4-Bromofluorobenzene	10.635	95	110330	50.420	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 100.840%	
Target Compounds						
						Qvalue
16) Acetone	2.626	43	6598	7.385	ug/l	98
17) Carbon Disulfide	2.793	76	1561	0.432	ug/l #	93
25) 2-Butanone	4.706	43	1902	1.535	ug/l	99

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

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