

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU111021\
 Data File : VU045700.D
 Acq On : 10 Nov 2021 19:59
 Operator : SY/MD
 Sample : M4638-01
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 PILOT-PLANT-PROCESS-BLEED

Quant Time: Nov 11 02:32:45 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	124037	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	232471	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	237791	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	116808	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.706	65	101582	54.685	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 109.360%			
35) Dibromofluoromethane	5.295	113	79708	51.355	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 102.720%			
50) Toluene-d8	7.899	98	306709	53.143	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 106.280%			
62) 4-Bromofluorobenzene	10.635	95	116976	50.545	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 101.080%			
Target Compounds						
16) Acetone	2.626	43	7685	8.127	ug/l	99
30) Chloroform	5.092	83	49852	17.680	ug/l	97
42) 1,2-Dichloroethane	5.796	62	130020	53.636	ug/l	99
47) Bromodichloromethane	7.108	83	18812	8.058	ug/l	95
60) Dibromochloromethane	8.816	129	4677	2.638	ug/l	89

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

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