

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100121\
 Data File : VU045077.D
 Acq On : 01 Oct 2021 16:49
 Operator : SY/MD
 Sample : M3993-02
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 STEEL-WELL-GATE-1-GROUT

Quant Time: Oct 02 03:12:39 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	178949	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	329429	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	311052	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	141701	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	150630	50.567	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 101.140%		
35) Dibromofluoromethane	5.292	113	53781	25.879	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 51.760%#		
50) Toluene-d8	7.899	98	416880	50.770	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 101.540%		
62) 4-Bromofluorobenzene	10.636	95	160639	47.386	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 94.780%		
Target Compounds						
						Qvalue
16) Acetone	2.633	43	52538	29.713	ug/l	98
18) Methyl Acetate	2.951	43	35022	6.777	ug/l	98
20) Methylene Chloride	3.044	84	12931	4.855	ug/l	93
25) 2-Butanone	4.716	43	24571	9.781	ug/l	99
68) m/p-Xylenes	9.697	106	7557	1.621	ug/l	100
69) o-Xylene	10.105	106	5110	1.104	ug/l	93

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

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