

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100121\
 Data File : VU045066.D
 Acq On : 01 Oct 2021 12:32
 Operator : SY/MD
 Sample : M3966-07
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 S-830-J-GW-2.7-3.7-092721

Quant Time: Oct 02 03:10:25 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	5.379	168	197335	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	357405	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	319591	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	133576	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.707	65	160334	48.810	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	=	97.620%	
35) Dibromofluoromethane	5.295	113	111918	49.639	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	99.280%	
50) Toluene-d8	7.903	98	444503	49.897	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	=	99.800%	
62) 4-Bromofluorobenzene	10.636	95	151781	41.269	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	82.540%#	
Target Compounds						
					Qvalue	
16) Acetone	2.639	43	12460	6.390	ug/l	93
25) 2-Butanone	4.729	43	3380	1.220	ug/l	94
27) cis-1,2-Dichloroethene	4.674	96	2484	0.822	ug/l	96
44) Trichloroethene	6.546	130	2522	0.961	ug/l	96
65) Chlorobenzene	9.449	112	2340	0.349	ug/l	99
95) Naphthalene	14.092	128	10763	1.092	ug/l #	96

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

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