

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU093021\
 Data File : VU045055.D
 Acq On : 30 Sep 2021 18:24
 Operator : SY/MD
 Sample : M3966-07
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 22 Sample Multiplier: 1

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

Quant Time: Oct 01 05:41:18 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	199185	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	358976	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	328962	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	144814	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.707	65	163521	49.317	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	98.640%
35) Dibromofluoromethane	5.295	113	112573	49.711	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.420%
50) Toluene-d8	7.903	98	449423	50.228	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	100.460%
62) 4-Bromofluorobenzene	10.636	95	155968	42.222	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	84.440%
Target Compounds						
					Qvalue	
16) Acetone	2.636	43	8525	4.332	ug/l	97
27) cis-1,2-Dichloroethene	4.675	96	1236	0.405	ug/l	91
44) Trichloroethene	6.549	130	1219	0.462	ug/l	98
65) Chlorobenzene	9.449	112	2409	0.349	ug/l	99
73) Isopropylbenzene	10.488	105	4588	0.425	ug/l	94
78) n-propylbenzene	10.909	91	14642	1.114	ug/l	97
84) 1,2,4-Trimethylbenzene	11.472	105	17754	1.900	ug/l	99
85) sec-Butylbenzene	11.649	105	3492	0.295	ug/l	90
95) Naphthalene	14.092	128	6242	0.584	ug/l	99

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

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