

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU022221\
 Data File : VU042398.D
 Acq On : 22 Feb 2021 16:27
 Operator : SY/MD
 Sample : M1457-06
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 GWBR1-85-100-021921

Quant Time: Feb 22 16:50:02 2021
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U021721W.M
 Quant Title : SW846 8260
 QLast Update : Fri Feb 19 02:55:52 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.382	168	181931	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	273751	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	242522	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	114102	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	132066	47.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.74%	
35) Dibromofluoromethane	5.298	113	91619	47.42	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.84%	
50) Toluene-d8	7.906	98	333443	48.18	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.36%	
62) 4-Bromofluorobenzene	10.639	95	119039	43.64	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.28%	
Target Compounds						Qvalue
16) Acetone	2.646	43	5848	3.42	ug/l	94
17) Carbon Disulfide	2.800	76	2197	0.44	ug/l	99
21) trans-1,2-Dichloroethene	3.369	96	1464	0.81	ug/l	91
27) cis-1,2-Dichloroethene	4.674	96	14234	6.83	ug/l	90
30) Chloroform	5.099	83	11154	3.00	ug/l	97
40) Benzene	5.787	78	2430	0.32	ug/l #	83
44) Trichloroethene	6.549	130	54598	26.94	ug/l	96
52) Toluene	7.977	92	21143	4.50	ug/l	100

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

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