

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061421\
 Data File : VU044215.D
 Acq On : 15 Jun 2021 05:07
 Operator : SY/MD
 Sample : M2706-01
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 CIEW_WC

Manual Integrations
 APPROVED

MMDadoda
 6/15/2021 9:51:09 AM

Quant Time: Jun 15 05:39:26 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U060921W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jun 10 05:11:37 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.385	168	177180	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.259	114	312839	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	298557	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	156706	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.713	65	131838	46.791	ug/l	0.00
Spiked Amount	50.000		Recovery	=	93.580%	
35) Dibromofluoromethane	5.298	113	104527	48.032	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.060%	
50) Toluene-d8	7.906	98	390806	46.760	ug/l	0.00
Spiked Amount	50.000		Recovery	=	93.520%	
62) 4-Bromofluorobenzene	10.639	95	142779	42.629	ug/l	0.00
Spiked Amount	50.000		Recovery	=	85.260%	
Target Compounds					Qvalue	
16) Acetone	2.658	43	11468m	5.791	ug/l	
27) cis-1,2-Dichloroethene	4.677	96	10907	4.063	ug/l	87
30) Chloroform	5.099	83	8808	1.939	ug/l	97
44) Trichloroethene	6.552	130	13494	5.678	ug/l	99
80) 1,3,5-Trimethylbenzene	11.095	105	12488	1.338	ug/l	99
84) 1,2,4-Trimethylbenzene	11.475	105	19918	2.124	ug/l	99

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

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