

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU121718\
 Data File : VU028727.D
 Acq On : 17 Dec 2018 17:55
 Operator : MD/SY
 Sample : J6393-08DL 4X
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MW-10C_R1DL

Quant Time: Dec 18 06:50:06 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U120618W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 07 00:39:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.98	168	83422	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	146795	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	129276	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	50839	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	66614	52.33	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.66%	
35) Dibromofluoromethane	4.88	113	49435	53.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.02%	
50) Toluene-d8	7.56	98	188923	51.70	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.40%	
62) 4-Bromofluorobenzene	10.31	95	61292	44.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.74%	
Target Compounds						
27) cis-1,2-Dichloroethene	4.23	96	11700	9.885	ug/l	89
44) Trichloroethene	6.18	130	10457	9.941	ug/l	100
64) Tetrachloroethene	8.22	164	57432	66.603	ug/l	97

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU121718\
Data File : VU028727.D
Acq On : 17 Dec 2018 17:55
Operator : MD/SY
Sample : J6393-08DL 4X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-10C_R1DL

Quant Time: Dec 18 06:50:06 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U120618W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 07 00:39:42 2018
Response via : Initial Calibration

