

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU121518\
 Data File : VU028694.D
 Acq On : 14 Dec 2018 19:40
 Operator : MD/SY
 Sample : J6393-08
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MW-10C_R1

Quant Time: Dec 17 10:03:04 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	77241	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	137530	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	126561	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	50148	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	64964	55.11	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.22%	
35) Dibromofluoromethane	4.88	113	48439	55.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	111.94%	
50) Toluene-d8	7.57	98	178666	52.18	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.36%	
62) 4-Bromofluorobenzene	10.30	95	59304	46.34	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.68%	
Target Compounds						
27) cis-1,2-Dichloroethene	4.22	96	47374	43.227	ug/l	89
44) Trichloroethene	6.18	130	44181	44.829	ug/l	98
64) Tetrachloroethene	8.23	164	251226	297.593	ug/l	99
83) tert-Butylbenzene	11.10	119	4272	1.391	ug/l	88
91) 1,2-Dichlorobenzene	11.88	146	2187	1.287	ug/l	94

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

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