

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU110918\
 Data File : VU028042.D
 Acq On : 09 Nov 2018 14:14
 Operator : MD/SY
 Sample : J5885-01
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 RW-2

Quant Time: Nov 12 11:37:39 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 02 02:28:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	192833	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	374302	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	334233	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	143950	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	162985	56.85	ug/l	0.00
Spiked Amount	50.000		Recovery	=	113.70%	
35) Dibromofluoromethane	4.89	113	120960	51.25	ug/l	0.00
Spiked Amount	50.000		Recovery	=	102.50%	
50) Toluene-d8	7.57	98	485466	53.25	ug/l	0.00
Spiked Amount	50.000		Recovery	=	106.50%	
62) 4-Bromofluorobenzene	10.31	95	160194	47.64	ug/l	0.00
Spiked Amount	50.000		Recovery	=	95.28%	
Target Compounds						
27) cis-1,2-Dichloroethene	4.23	96	12917	4.939	ug/l	94
30) Chloroform	4.68	83	4893	1.051	ug/l	94
44) Trichloroethene	6.19	130	7982	3.098	ug/l	92
64) Tetrachloroethene	8.23	164	2748720	1268.294	ug/l	99
78) n-propylbenzene	10.59	91	5388	0.404	ug/l	93
80) 1,3,5-Trimethylbenzene	10.77	105	5809	0.629	ug/l	95
84) 1,2,4-Trimethylbenzene	11.15	105	14041	1.509	ug/l	98

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

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