

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU062718\
 Data File : VU024932.D
 Acq On : 26 Jun 2018 20:19
 Operator : MD/SY
 Sample : J3669-07ME
 Misc : 5.94g/5.0ml/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 WP021SB484-35-37-180620ME

Quant Time: Jun 27 04:36:20 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 13 13:55:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.98	168	222477	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	341959	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	329023	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	189071	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	168699	46.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.90%	
35) Dibromofluoromethane	4.89	113	129774	45.73	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.46%	
50) Toluene-d8	7.57	98	493452	47.67	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.34%	
62) 4-Bromofluorobenzene	10.32	95	204715	49.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.62%	
Target Compounds						
12) 1,1-Dichloroethene	2.25	96	3258	1.29	ug/l	93
27) cis-1,2-Dichloroethene	4.22	96	108219	33.74	ug/l	89
44) Trichloroethene	6.18	130	300003	92.06	ug/l	96
91) 1,2-Dichlorobenzene	11.89	146	5556	0.82	ug/l	96

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

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