

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU062818\
 Data File : VU025045.D
 Acq On : 29 Jun 2018 01:59
 Operator : MD/SY
 Sample : J3669-31ME 10X
 Misc : 6.81g/5mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 WP021SB474-33-35-180620ME

Quant Time: Jun 29 08:15:36 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	4.99	168	220891	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	330348	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	307576	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	168238	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	166429	46.15	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.30%	
35) Dibromofluoromethane	4.89	113	127328	46.44	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.88%	
50) Toluene-d8	7.57	98	475035	47.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.00%	
62) 4-Bromofluorobenzene	10.31	95	179715	45.26	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.52%	
Target Compounds						
12) 1,1-Dichloroethene	2.28	96	6002	2.39	ug/l	95
27) cis-1,2-Dichloroethene	4.23	96	185729	58.32	ug/l	88
44) Trichloroethene	6.19	130	280275	89.03	ug/l	97
91) 1,2-Dichlorobenzene	11.88	146	3196	0.53	ug/l	96

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

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