

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN092618\
 Data File : VN051499.D
 Acq On : 25 Sep 2018 22:01
 Operator : MD\SY
 Sample : J5065-02
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MH-169

Manual Integrations
 APPROVED

mmdadoda
 9/30/2018 9:53:56 AM

Quant Time: Sep 26 04:56:02 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N092318W.M
 Quant Title : SW846 8260
 QLast Update : Mon Sep 24 06:41:32 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	648220	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	960699	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	813302	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	333501	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	430623	53.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.06%	
35) Dibromofluoromethane	7.59	113	384308	50.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.06%	
50) Toluene-d8	10.09	98	1375772	47.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.12%	
62) 4-Bromofluorobenzene	12.40	95	383925	41.02	ug/l	0.00
Spiked Amount 50.000			Recovery	=	82.04%	
Target Compounds						
16) Acetone	3.82	43	107578	39.345	ug/l	98
25) 2-Butanone	6.85	43	9544	3.434	ug/l #	89
27) cis-1,2-Dichloroethene	6.83	96	3478m	0.430	ug/l	
30) Chloroform	7.37	83	203298	14.701	ug/l	98
37) Ethyl Acetate	6.94	43	18780	2.936	ug/l #	92
38) Carbon Tetrachloride	7.78	117	4108	0.353	ug/l #	83
60) Dibromochloromethane	10.90	129	2805	0.331	ug/l	97

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

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