

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN101918\
 Data File : VN051952.D
 Acq On : 19 Oct 2018 17:07
 Operator : MD\SY
 Sample : J5538-18
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 19 Sample Multiplier: 28

Instrument :
 MSVOA_N
 ClientSampleId :
 BL01

Quant Time: Oct 20 00:51:29 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N101718W.M
 Quant Title : SW846 8260
 QLast Update : Sat Oct 20 00:28:35 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	8.14	168	800846	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.91	114	1265882	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.48	117	1204227	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.36	152	597250	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.45	65	497969	67.07	ug/l	0.00
Spiked Amount 50.000			Recovery	=	134.14%	
35) Dibromofluoromethane	0.00	113	0	0.00	ug/l	
Spiked Amount 50.000			Recovery	=	0.00%	
50) Toluene-d8	10.25	98	1790382	61.13	ug/l	0.00
Spiked Amount 50.000			Recovery	=	122.26%	
62) 4-Bromofluorobenzene	12.44	95	654621	66.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	133.94%	
Target Compounds						
59) 2-Hexanone	10.85	43	31978	10.427	ug/l	83
93) 1,2,4-Trichlorobenzene	14.91	180	8875	4.904	ug/l	93
95) Naphthalene	15.14	128	159621	11.561	ug/l	99
96) 1,2,3-Trichlorobenzene	15.32	180	8028	3.725	ug/l #	88

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN101918\
Data File : VN051952.D
Acq On : 19 Oct 2018 17:07
Operator : MD\SY
Sample : J5538-18
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 19 Sample Multiplier: 28

Instrument :
MSVOA_N
ClientSampleId :
BL01

Quant Time: Oct 20 00:51:29 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N101718W.M
Quant Title : SW846 8260
QLast Update : Sat Oct 20 00:28:35 2018
Response via : Initial Calibration

