

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN102518\
 Data File : VN052009.D
 Acq On : 25 Oct 2018 14:01
 Operator : MD\SY
 Sample : J5651-03
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 12 Sample Multiplier: 28

Instrument :
 MSVOA_N
 ClientSampleId :
 RBR-200030

Quant Time: Oct 26 02:10:19 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N102318W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 24 10:12:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	1148572	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	1681972	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	1576892	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	736775	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	580843	56.78	ug/l	0.00
Spiked Amount 50.000			Recovery	=	113.56%	
35) Dibromofluoromethane	7.59	113	613074	57.12	ug/l	0.00
Spiked Amount 50.000			Recovery	=	114.24%	
50) Toluene-d8	10.09	98	2297186	59.23	ug/l	0.00
Spiked Amount 50.000			Recovery	=	118.46%	
62) 4-Bromofluorobenzene	12.40	95	725254	54.40	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.80%	
Target Compounds						
7) Trichlorofluoromethane	3.02	101	16985	1.084	ug/l	96
16) Acetone	3.82	43	16125	9.568	ug/l	92
20) Methylene Chloride	4.55	84	11579	1.171	ug/l #	87
43) Isopropyl Acetate	8.17	43	289220	26.766	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN102518\
Data File : VN052009.D
Acq On : 25 Oct 2018 14:01
Operator : MD\SY
Sample : J5651-03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 12 Sample Multiplier: 28

Instrument :
MSVOA_N
ClientSampleId :
RBR-200030

Quant Time: Oct 26 02:10:19 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N102318W.M
Quant Title : SW846 8260
QLast Update : Wed Oct 24 10:12:55 2018
Response via : Initial Calibration

