Full wwPDB X-ray Structure Validation Report (i)

Jun 13, 2024 – 06:45 AM EDT

PDB ID	:	4G1O
Title	:	Crystal structure of Newcastle disease virus matrix protein
Authors	:	Meng, G.; Rossmann, M.G.
Deposited on	:	2012-07-11
Resolution	:	2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

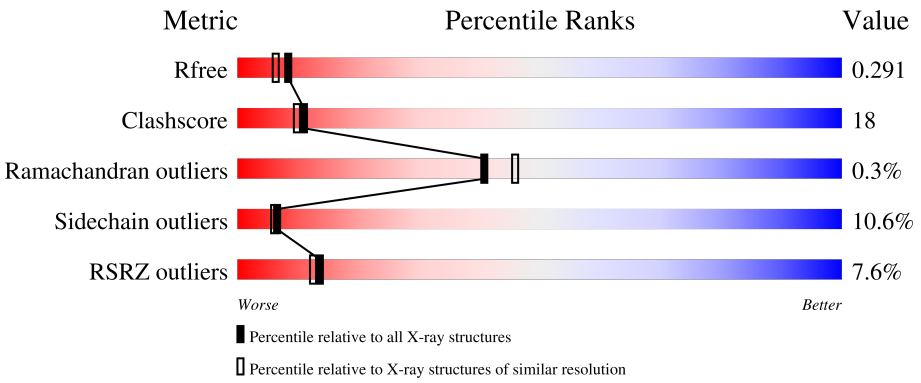
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	364	8% 61% 24% 5% 10%
1	B	364	5% 63% 20% 6% 10%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 5089 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Matrix protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	329	2517	1610	437	461	9	0	0	0
1	B	326	2501	1600	436	456	9	0	0	0

- Molecule 2 is water.

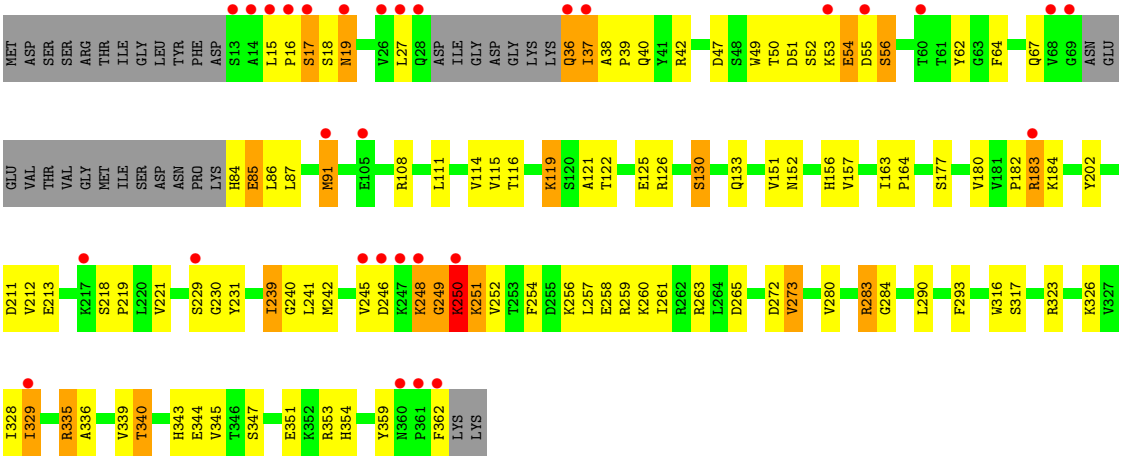
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	33	33	33	0	0
2	B	38	38	38	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

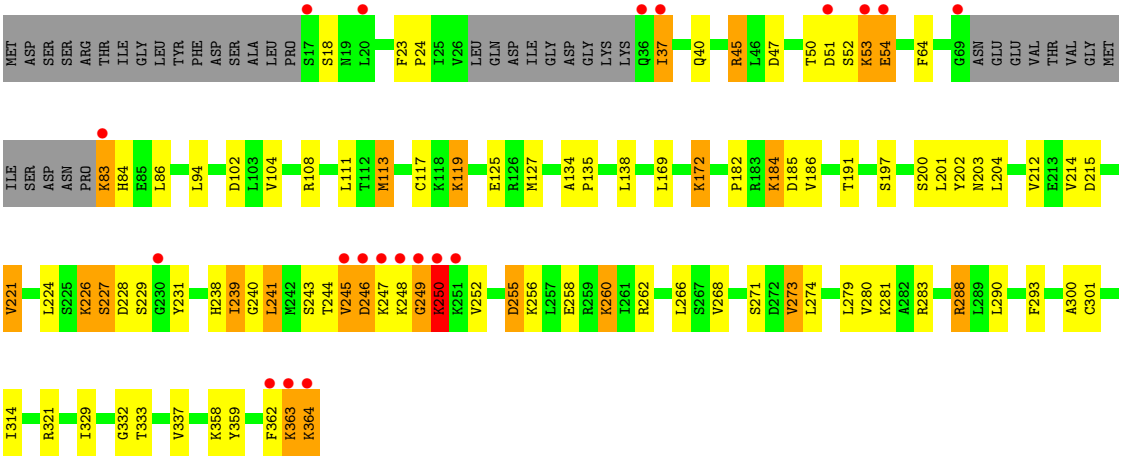
• Molecule 1: Matrix protein

Chain A:



• Molecule 1: Matrix protein

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	163.31Å 46.99Å 117.41Å 90.00° 132.05° 90.00°	Depositor
Resolution (Å)	30.00 – 2.20 27.73 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.4 (30.00-2.20) 99.5 (27.73-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	6.26 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.180 , 0.280 0.194 , 0.291	Depositor DCC
R_{free} test set	1680 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å^2)	47.2	Xtriage
Anisotropy	0.122	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 51.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.007 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	5089	wwPDB-VP
Average B, all atoms (Å^2)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.59% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#$\|Z\|$ >5	RMSZ	#$\|Z\|$ >5
1	A	0.68	2/2560 (0.1%)	0.89	8/3471 (0.2%)
1	B	0.68	0/2543	0.88	5/3443 (0.1%)
All	All	0.68	2/5103 (0.0%)	0.88	13/6914 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	316	TRP	CD2-CE2	5.39	1.47	1.41
1	A	49	TRP	CD2-CE2	5.14	1.47	1.41

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	56	SER	N-CA-CB	-8.11	98.34	110.50
1	B	250	LYS	N-CA-C	-7.56	90.60	111.00
1	A	250	LYS	N-CA-C	-7.18	91.61	111.00
1	B	273	VAL	CB-CA-C	-7.05	98.00	111.40
1	B	113	MET	CG-SD-CE	-6.75	89.41	100.20
1	B	249	GLY	N-CA-C	-6.72	96.30	113.10
1	A	17	SER	CB-CA-C	-6.68	97.41	110.10
1	A	335	ARG	NE-CZ-NH2	-6.26	117.17	120.30
1	A	230	GLY	C-N-CA	-5.49	107.98	121.70
1	A	230	GLY	O-C-N	5.28	131.15	122.70
1	B	246	ASP	CB-CG-OD2	5.20	122.98	118.30
1	A	335	ARG	NE-CZ-NH1	5.17	122.89	120.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	249	GLY	CA-C-N	-5.15	105.87	117.20

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	249	GLY	Mainchain
1	A	250	LYS	Peptide

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2517	0	2630	99	0
1	B	2501	0	2622	89	0
2	A	33	0	0	0	0
2	B	38	0	0	0	0
All	All	5089	0	5252	186	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 18.

All (186) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:PRO:HG2	1:A:19:ASN:OD1	1.47	1.13
1:B:288:ARG:HG2	1:B:288:ARG:HH11	1.02	1.11
1:B:182:PRO:HB3	1:B:184:LYS:HE2	1.34	1.04
1:A:54:GLU:OE1	1:A:180:VAL:HG12	1.56	1.04
1:A:15:LEU:HB3	1:A:16:PRO:HD2	1.40	1.00
1:A:245:VAL:HG12	1:A:251:LYS:H	1.25	1.00
1:A:183:ARG:HH11	1:A:183:ARG:HG3	1.32	0.91
1:A:259:ARG:HH11	1:A:263:ARG:HH12	1.02	0.91
1:B:50:THR:HG22	1:B:52:SER:H	1.37	0.90
1:B:247:LYS:O	1:B:248:LYS:HG2	1.71	0.90

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:288:ARG:HG2	1:B:288:ARG:NH1	1.80	0.89
1:A:259:ARG:HH11	1:A:263:ARG:NH1	1.72	0.87
1:A:16:PRO:O	1:A:17:SER:HB3	1.74	0.86
1:A:38:ALA:O	1:A:39:PRO:C	2.16	0.83
1:B:246:ASP:O	1:B:249:GLY:O	1.97	0.81
1:B:245:VAL:HG12	1:B:245:VAL:O	1.81	0.80
1:B:53:LYS:HD2	1:B:54:GLU:N	1.95	0.80
1:B:245:VAL:O	1:B:245:VAL:CG1	2.30	0.80
1:A:259:ARG:NH1	1:A:263:ARG:HH12	1.78	0.79
1:B:248:LYS:O	1:B:248:LYS:CG	2.30	0.79
1:A:16:PRO:O	1:A:17:SER:CB	2.30	0.78
1:A:50:THR:HG22	1:A:52:SER:H	1.48	0.78
1:B:111:LEU:HD11	1:B:362:PHE:CE2	2.19	0.77
1:B:248:LYS:O	1:B:248:LYS:HG3	1.84	0.77
1:B:184:LYS:HD3	1:B:184:LYS:H	1.50	0.75
1:A:64:PHE:HB3	1:A:86:LEU:HD11	1.69	0.75
1:A:16:PRO:HG2	1:A:19:ASN:CG	2.08	0.73
1:B:119:LYS:HE3	1:B:125:GLU:OE2	1.91	0.71
1:A:37:ILE:O	1:A:37:ILE:HG23	1.90	0.70
1:A:27:LEU:HB2	1:A:218:SER:HA	1.71	0.70
1:B:245:VAL:HG22	1:B:249:GLY:O	1.92	0.70
1:A:55:ASP:OD2	1:A:183:ARG:HB3	1.93	0.69
1:A:183:ARG:HG3	1:A:183:ARG:NH1	2.00	0.69
1:B:247:LYS:O	1:B:248:LYS:CG	2.40	0.69
1:A:336:ALA:O	1:A:340:THR:HG23	1.93	0.68
1:A:241:LEU:O	1:A:290:LEU:CD1	2.43	0.67
1:A:258:GLU:HG3	1:A:345:VAL:HG13	1.75	0.66
1:A:55:ASP:CG	1:A:183:ARG:H	1.98	0.65
1:A:242:MET:HG2	1:A:290:LEU:HD11	1.78	0.65
1:A:241:LEU:O	1:A:290:LEU:HD12	1.97	0.65
1:A:252:VAL:HG13	1:A:256:LYS:HB2	1.78	0.65
1:A:37:ILE:O	1:A:37:ILE:CG2	2.44	0.65
1:A:50:THR:HG22	1:A:51:ASP:N	2.11	0.64
1:A:39:PRO:HB3	1:A:111:LEU:HD11	1.80	0.64
1:A:259:ARG:NH1	1:A:263:ARG:NH1	2.43	0.64
1:A:15:LEU:HB3	1:A:16:PRO:CD	2.23	0.63
1:A:245:VAL:HG12	1:A:251:LYS:N	2.05	0.63
1:A:64:PHE:HB3	1:A:86:LEU:CD1	2.29	0.62
1:A:157:VAL:HG11	1:A:163:ILE:CD1	2.30	0.62
1:B:53:LYS:HD2	1:B:54:GLU:H	1.65	0.62
1:B:227:SER:O	1:B:228:ASP:C	2.34	0.62

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:246:ASP:OD1	1:B:247:LYS:N	2.34	0.60
1:A:157:VAL:HG11	1:A:163:ILE:HD12	1.83	0.60
1:B:255:ASP:N	1:B:255:ASP:OD1	2.34	0.60
1:B:363:LYS:O	1:B:364:LYS:C	2.39	0.60
1:B:50:THR:HG22	1:B:51:ASP:N	2.17	0.59
1:B:288:ARG:HH11	1:B:288:ARG:CG	1.94	0.59
1:A:67:GLN:HB3	1:A:85:GLU:HG2	1.85	0.59
1:A:265:ASP:O	1:A:283:ARG:HD3	2.02	0.59
1:B:108:ARG:HG2	1:B:362:PHE:CE1	2.38	0.58
1:A:39:PRO:CB	1:A:111:LEU:HD11	2.33	0.58
1:B:247:LYS:C	1:B:249:GLY:N	2.54	0.57
1:B:333:THR:O	1:B:337:VAL:HG23	2.04	0.57
1:A:55:ASP:C	1:A:56:SER:OG	2.39	0.57
1:A:256:LYS:O	1:A:260:LYS:HG3	2.05	0.57
1:A:38:ALA:O	1:A:40:GLN:N	2.38	0.56
1:A:42:ARG:HE	1:A:177:SER:HB2	1.70	0.56
1:A:50:THR:CG2	1:A:51:ASP:N	2.68	0.56
1:B:226:LYS:HD2	1:B:227:SER:N	2.19	0.56
1:B:245:VAL:HG22	1:B:250:LYS:HA	1.86	0.56
1:B:247:LYS:O	1:B:248:LYS:CB	2.51	0.56
1:B:247:LYS:O	1:B:247:LYS:CG	2.52	0.56
1:B:127:MET:HA	1:B:127:MET:HE2	1.89	0.55
1:A:239:ILE:HD11	1:A:280:VAL:HG21	1.87	0.55
1:B:247:LYS:O	1:B:247:LYS:HG2	2.06	0.55
1:B:50:THR:HG22	1:B:52:SER:N	2.17	0.54
1:B:37:ILE:O	1:B:37:ILE:HG23	2.07	0.54
1:A:16:PRO:CG	1:A:19:ASN:OD1	2.38	0.54
1:A:261:ILE:HD11	1:A:329:ILE:HD11	1.90	0.54
1:B:53:LYS:NZ	1:B:54:GLU:OE1	2.30	0.54
1:A:40:GLN:HB3	1:A:359:TYR:HB3	1.89	0.54
1:A:351:GLU:HG2	1:A:354:HIS:CE1	2.43	0.54
1:B:119:LYS:CE	1:B:125:GLU:OE2	2.55	0.54
1:B:23:PHE:CG	1:B:24:PRO:HA	2.43	0.53
1:B:247:LYS:C	1:B:249:GLY:H	2.12	0.52
1:A:202:TYR:CE2	1:A:257:LEU:HD11	2.45	0.52
1:A:15:LEU:HB3	1:A:19:ASN:HD21	1.73	0.52
1:B:256:LYS:O	1:B:260:LYS:HG3	2.09	0.52
1:B:64:PHE:CD1	1:B:86:LEU:HD11	2.45	0.52
1:A:18:SER:HG	1:A:359:TYR:HD2	1.56	0.52
1:B:184:LYS:H	1:B:184:LYS:CD	2.20	0.52
1:A:91:MET:O	1:A:91:MET:HG3	2.10	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:254:PHE:CE1	1:A:258:GLU:HB2	2.46	0.51
1:B:108:ARG:CG	1:B:362:PHE:CE1	2.93	0.51
1:B:119:LYS:HD3	1:B:127:MET:HE1	1.93	0.51
1:A:152:ASN:O	1:A:156:HIS:HD2	1.94	0.51
1:A:15:LEU:HB3	1:A:19:ASN:ND2	2.26	0.51
1:A:245:VAL:CG1	1:A:251:LYS:H	2.12	0.51
1:A:53:LYS:O	1:A:184:LYS:NZ	2.44	0.51
1:A:54:GLU:HA	1:A:182:PRO:HG3	1.94	0.50
1:B:203:ASN:OD1	1:B:238:HIS:ND1	2.37	0.50
1:B:226:LYS:HD2	1:B:226:LYS:C	2.31	0.50
1:A:241:LEU:O	1:A:290:LEU:HD11	2.12	0.50
1:B:40:GLN:HB3	1:B:359:TYR:HB3	1.94	0.50
1:A:55:ASP:OD1	1:A:56:SER:OG	2.26	0.49
1:B:221:VAL:HG22	1:B:231:TYR:CE1	2.47	0.49
1:B:50:THR:CG2	1:B:51:ASP:N	2.75	0.49
1:B:197:SER:HB2	1:B:332:GLY:O	2.13	0.49
1:A:339:VAL:HG23	1:A:340:THR:HG22	1.94	0.49
1:A:114:VAL:HG21	1:A:133:GLN:HE21	1.78	0.49
1:B:202:TYR:HB3	1:B:329:ILE:HG23	1.95	0.49
1:A:351:GLU:HG2	1:A:354:HIS:ND1	2.27	0.48
1:A:27:LEU:HB2	1:A:219:PRO:HD3	1.96	0.48
1:A:54:GLU:HG3	1:A:55:ASP:N	2.28	0.48
1:B:226:LYS:CD	1:B:227:SER:N	2.77	0.48
1:A:242:MET:CG	1:A:290:LEU:HD11	2.42	0.47
1:B:108:ARG:NH1	1:B:362:PHE:HD1	2.12	0.47
1:B:202:TYR:HB3	1:B:329:ILE:CG2	2.44	0.47
1:B:274:LEU:HD12	1:B:301:CYS:SG	2.55	0.47
1:A:36:GLN:O	1:A:38:ALA:HB2	2.15	0.47
1:A:116:THR:OG1	1:A:130:SER:HB2	2.16	0.46
1:A:353:ARG:HG3	1:A:354:HIS:CD2	2.50	0.46
1:B:119:LYS:HE3	1:B:125:GLU:CD	2.35	0.46
1:A:221:VAL:HG13	1:A:231:TYR:CE1	2.51	0.46
1:B:53:LYS:HD2	1:B:53:LYS:C	2.24	0.45
1:B:182:PRO:CB	1:B:184:LYS:HE2	2.25	0.45
1:B:250:LYS:HB3	1:B:250:LYS:HE3	1.48	0.45
1:A:328:ILE:HD13	1:A:344:GLU:HA	1.99	0.45
1:B:127:MET:HE3	1:B:127:MET:HB2	1.59	0.45
1:A:108:ARG:O	1:A:111:LEU:HB2	2.16	0.45
1:B:246:ASP:N	1:B:252:VAL:HG23	2.32	0.45
1:A:119:LYS:HE3	1:A:125:GLU:OE2	2.16	0.45
1:A:16:PRO:HG2	1:A:19:ASN:ND2	2.32	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:201:LEU:HD13	1:B:241:LEU:HG	1.99	0.44
1:B:221:VAL:HG22	1:B:231:TYR:CD1	2.52	0.44
1:A:115:VAL:HA	1:A:130:SER:O	2.17	0.44
1:A:183:ARG:HH11	1:A:183:ARG:CG	2.15	0.44
1:A:16:PRO:HD2	1:A:19:ASN:HD21	1.83	0.44
1:A:62:TYR:OH	1:A:272:ASP:OD1	2.36	0.44
1:A:18:SER:HB2	1:A:317:SER:O	2.18	0.44
1:B:172:LYS:HD3	1:B:172:LYS:HA	1.61	0.44
1:A:163:ILE:HA	1:A:164:PRO:HD3	1.92	0.43
1:B:268:VAL:HA	1:B:279:LEU:O	2.17	0.43
1:A:245:VAL:O	1:A:246:ASP:OD1	2.35	0.43
1:B:50:THR:HB	1:B:185:ASP:OD2	2.19	0.43
1:A:39:PRO:HB3	1:A:111:LEU:CD1	2.47	0.43
1:B:64:PHE:HB3	1:B:86:LEU:CD1	2.49	0.43
1:A:252:VAL:CG1	1:A:256:LYS:HB2	2.48	0.42
1:A:272:ASP:O	1:A:273:VAL:C	2.58	0.42
1:A:121:ALA:HA	1:A:126:ARG:HG3	2.00	0.42
1:B:51:ASP:N	1:B:51:ASP:OD2	2.48	0.42
1:B:280:VAL:HG22	1:B:300:ALA:HB3	2.02	0.42
1:B:83:LYS:HB3	1:B:84:HIS:H	1.72	0.42
1:B:239:ILE:HD13	1:B:239:ILE:HG21	1.57	0.42
1:A:55:ASP:OD1	1:A:56:SER:N	2.52	0.42
1:A:256:LYS:HA	1:A:256:LYS:HD3	1.87	0.42
1:A:261:ILE:CD1	1:A:329:ILE:HD11	2.50	0.42
1:B:111:LEU:HD11	1:B:362:PHE:HE2	1.78	0.42
1:B:117:CYS:O	1:B:169:LEU:HB2	2.19	0.42
1:B:200:SER:HB2	1:B:244:THR:OG1	2.19	0.42
1:B:224:LEU:HD21	1:B:314:ILE:HD13	2.02	0.42
1:B:246:ASP:O	1:B:249:GLY:C	2.59	0.42
1:B:247:LYS:O	1:B:248:LYS:C	2.55	0.42
1:B:266:LEU:HA	1:B:281:LYS:O	2.20	0.42
1:B:108:ARG:NH1	1:B:362:PHE:CD1	2.87	0.41
1:A:122:THR:O	1:B:238:HIS:CD2	2.73	0.41
1:A:248:LYS:HD3	1:A:248:LYS:HA	1.54	0.41
1:A:343:HIS:CD2	1:A:343:HIS:C	2.94	0.41
1:A:351:GLU:CG	1:A:354:HIS:ND1	2.83	0.41
1:B:271:SER:HB2	1:B:279:LEU:HD11	2.02	0.41
1:A:329:ILE:HD13	1:A:345:VAL:HG21	2.01	0.41
1:B:102:ASP:OD1	1:B:104:VAL:HB	2.19	0.41
1:B:240:GLY:HA2	1:B:293:PHE:CD1	2.56	0.41
1:A:16:PRO:C	1:A:18:SER:H	2.23	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:THR:O	1:B:238:HIS:HD2	2.03	0.41
1:B:214:VAL:HG22	1:B:215:ASP:O	2.20	0.41
1:A:67:GLN:HB2	1:A:87:LEU:HD11	2.03	0.41
1:A:151:VAL:HB	1:A:156:HIS:CD2	2.56	0.41
1:A:240:GLY:HA2	1:A:293:PHE:CD1	2.56	0.41
1:A:15:LEU:CB	1:A:16:PRO:HD2	2.24	0.40
1:B:113:MET:HG2	1:B:134:ALA:HA	2.03	0.40
1:B:45:ARG:HE	1:B:45:ARG:HB3	1.58	0.40
1:A:263:ARG:O	1:A:284:GLY:HA3	2.21	0.40
1:B:135:PRO:HD2	1:B:138:LEU:HD12	2.02	0.40
1:B:204:LEU:O	1:B:239:ILE:HG22	2.21	0.40
1:B:94:LEU:HD11	1:B:113:MET:HE1	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	323/364 (89%)	308 (95%)	14 (4%)	1 (0%)	41	46
1	B	320/364 (88%)	307 (96%)	12 (4%)	1 (0%)	41	46
All	All	643/728 (88%)	615 (96%)	26 (4%)	2 (0%)	41	46

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	19	ASN
1	B	37	ILE

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	283/317 (89%)	256 (90%)	27 (10%)	8 8
1	B	281/317 (89%)	248 (88%)	33 (12%)	5 4
All	All	564/634 (89%)	504 (89%)	60 (11%)	6 6

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	GLN
1	A	37	ILE
1	A	47	ASP
1	A	54	GLU
1	A	84	HIS
1	A	85	GLU
1	A	91	MET
1	A	119	LYS
1	A	130	SER
1	A	183	ARG
1	A	211	ASP
1	A	212	VAL
1	A	213	GLU
1	A	229	SER
1	A	239	ILE
1	A	248	LYS
1	A	250	LYS
1	A	251	LYS
1	A	273	VAL
1	A	283	ARG
1	A	323	ARG
1	A	326	LYS
1	A	329	ILE
1	A	335	ARG
1	A	340	THR
1	A	347	SER
1	A	362	PHE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	18	SER
1	B	45	ARG
1	B	47	ASP
1	B	53	LYS
1	B	54	GLU
1	B	83	LYS
1	B	119	LYS
1	B	172	LYS
1	B	184	LYS
1	B	186	VAL
1	B	191	THR
1	B	212	VAL
1	B	221	VAL
1	B	226	LYS
1	B	227	SER
1	B	229	SER
1	B	239	ILE
1	B	241	LEU
1	B	243	SER
1	B	245	VAL
1	B	250	LYS
1	B	255	ASP
1	B	258	GLU
1	B	260	LYS
1	B	262	ARG
1	B	273	VAL
1	B	283	ARG
1	B	288	ARG
1	B	290	LEU
1	B	321	ARG
1	B	358	LYS
1	B	363	LYS
1	B	364	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	133	GLN
1	A	156	HIS
1	A	174	ASN
1	A	310	GLN
1	A	343	HIS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	40	GLN
1	B	156	HIS
1	B	310	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	329/364 (90%)	0.40	30 (9%)	9	8	25, 59, 98, 134	0
1	B	326/364 (89%)	0.26	20 (6%)	21	20	25, 52, 89, 116	0
All	All	655/728 (89%)	0.33	50 (7%)	13	12	25, 56, 95, 134	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	15	LEU	8.4
1	A	69	GLY	7.0
1	B	248	LYS	6.8
1	B	247	LYS	6.5
1	A	248	LYS	6.5
1	B	246	ASP	6.2
1	B	364	LYS	6.0
1	B	245	VAL	5.6
1	A	246	ASP	5.4
1	A	14	ALA	5.2
1	A	28	GLN	5.1
1	B	36	GLN	4.9
1	B	249	GLY	4.8
1	A	27	LEU	4.8
1	B	362	PHE	4.7
1	A	26	VAL	4.6
1	A	13	SER	4.5
1	A	36	GLN	4.1
1	A	362	PHE	4.0
1	B	69	GLY	4.0
1	A	105	GLU	3.9
1	B	37	ILE	3.7
1	A	55	ASP	3.6
1	A	250	LYS	3.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	16	PRO	3.5
1	B	363	LYS	3.4
1	A	17	SER	3.3
1	B	20	LEU	3.2
1	B	53	LYS	3.2
1	A	361	PRO	3.1
1	A	360	ASN	3.0
1	A	37	ILE	3.0
1	A	53	LYS	3.0
1	A	245	VAL	3.0
1	B	250	LYS	2.9
1	B	230	GLY	2.8
1	B	54	GLU	2.8
1	A	229	SER	2.8
1	A	68	VAL	2.7
1	A	91	MET	2.7
1	A	247	LYS	2.6
1	A	19	ASN	2.6
1	A	60	THR	2.5
1	B	51	ASP	2.5
1	A	329	ILE	2.3
1	B	251	LYS	2.3
1	A	217	LYS	2.2
1	B	17	SER	2.2
1	A	183	ARG	2.1
1	B	83	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.