



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 11:28 pm GMT

PDB ID : 4G1G
Title : Crystal structure of Newcastle disease virus matrix protein
Authors : Meng, G.; Rossmann, M.G.
Deposited on : 2012-07-10
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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

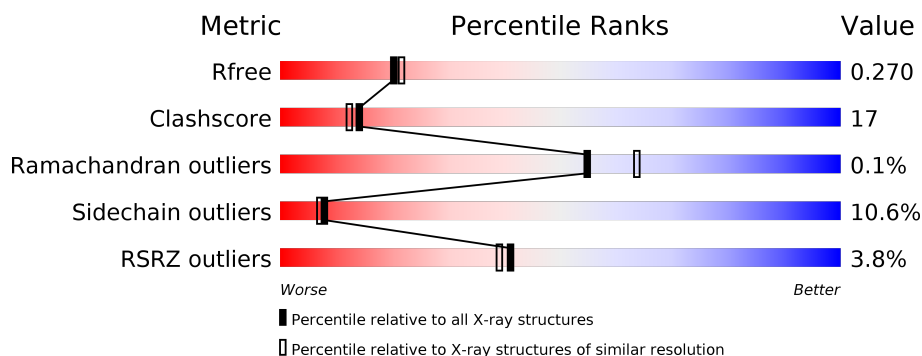
1 Overall quality at a glance

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}	100719	4002 (2.20-2.20)
Clashscore	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)
RSRZ outliers	101464	4033 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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 $\leq 5\%$. 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	
1	B	364	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5421 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
1	A	329	Total	C	N	O	S	0	0	0
			2504	1599	436	460	9			
1	B	354	Total	C	N	O	S	0	0	0
			2709	1730	468	501	10			

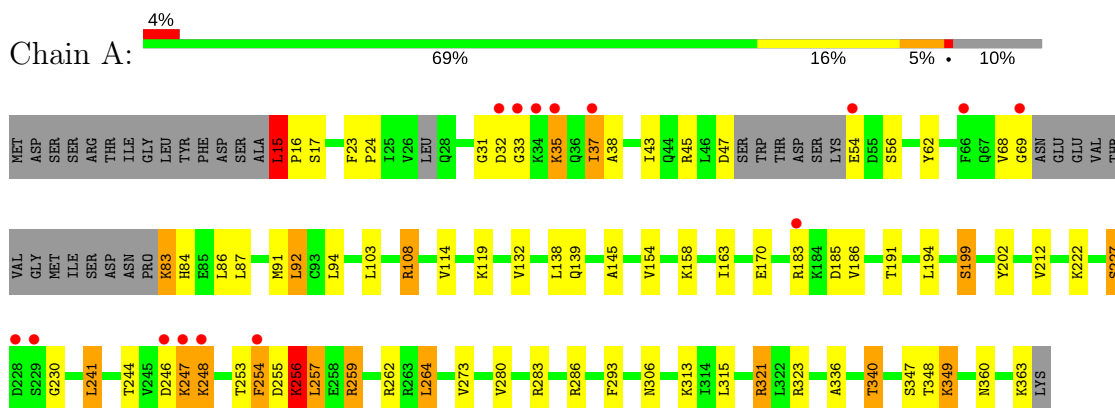
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	98	Total	O	0	0
			98	98		
2	B	110	Total	O	0	0
			110	110		

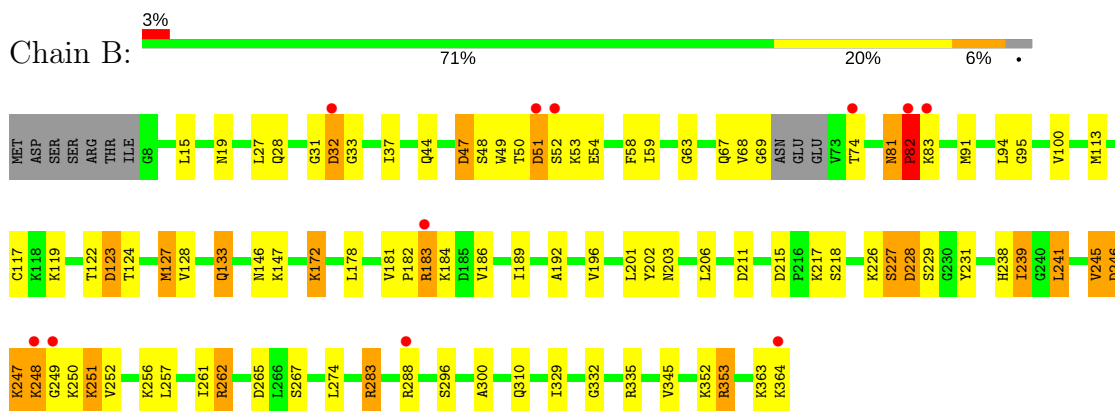
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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



• Molecule 1: Matrix protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2 ₁	Depositor
Cell constants a, b, c, α , β , γ	46.84Å 112.05Å 141.65Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.20 29.29 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.9 (30.00-2.20) 99.0 (29.29-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	9.37 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.220 , 0.270 0.221 , 0.270	Depositor DCC
R_{free} test set	1917 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	27.1	Xtriage
Anisotropy	0.056	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 32.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5421	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.53% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0.89	1/2544 (0.0%)	0.97	5/3445 (0.1%)
1	B	0.97	2/2756 (0.1%)	0.99	7/3733 (0.2%)
All	All	0.94	3/5300 (0.1%)	0.98	12/7178 (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	3
1	B	0	1
All	All	0	4

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	67	GLN	C-N	-5.59	1.21	1.34
1	B	123	ASP	CB-CG	-5.31	1.40	1.51
1	A	15	LEU	C-N	5.21	1.44	1.34

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	92	LEU	CA-CB-CG	6.69	130.70	115.30
1	A	264	LEU	CB-CG-CD1	-6.53	99.89	111.00
1	B	123	ASP	N-CA-CB	-6.39	99.10	110.60
1	B	74	THR	N-CA-C	6.23	127.83	111.00
1	A	108	ARG	NE-CZ-NH2	-6.10	117.25	120.30
1	A	185	ASP	CB-CG-OD1	5.94	123.65	118.30
1	A	185	ASP	CB-CG-OD2	-5.94	112.96	118.30
1	B	241	LEU	CA-CB-CG	5.86	128.78	115.30
1	B	32	ASP	CB-CG-OD2	5.50	123.25	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	211	ASP	CB-CG-OD1	5.29	123.06	118.30
1	B	32	ASP	CB-CA-C	5.16	120.72	110.40
1	B	82	PRO	CB-CA-C	-5.04	99.40	112.00

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	15	LEU	Mainchain
1	A	256	LYS	Peptide
1	A	35	LYS	Peptide
1	B	31	GLY	Peptide

5.2 Too-close contacts ⓘ

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	2504	0	2612	71	0
1	B	2709	0	2824	123	0
2	A	98	0	0	4	0
2	B	110	0	0	6	0
All	All	5421	0	5436	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 17.

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:B:69:GLY:N	1:B:83:LYS:HB3	1.27	1.41
1:B:69:GLY:C	1:B:83:LYS:HD2	1.41	1.40
1:B:69:GLY:CA	1:B:83:LYS:HB3	1.59	1.31
1:B:81:ASN:CG	1:B:82:PRO:HD3	1.49	1.31
1:B:69:GLY:HA2	1:B:83:LYS:CG	1.61	1.28
1:A:37:ILE:HD12	1:A:38:ALA:N	1.44	1.28

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:81:ASN:ND2	1:B:82:PRO:HD3	1.48	1.25
1:B:48:SER:OG	2:B:507:HOH:O	1.52	1.23
1:B:247:LYS:HD3	1:B:247:LYS:N	1.46	1.20
1:B:283:ARG:HG2	1:B:283:ARG:HH11	1.04	1.17
1:B:69:GLY:HA2	1:B:83:LYS:CB	1.73	1.17
1:B:47:ASP:HA	1:B:54:GLU:CD	1.63	1.16
1:A:37:ILE:CD1	1:A:38:ALA:H	1.61	1.11
1:B:247:LYS:H	1:B:247:LYS:CD	1.64	1.10
1:B:245:VAL:HG22	1:B:250:LYS:HA	1.35	1.07
1:A:306:ASN:O	1:B:124:THR:HG22	1.51	1.07
1:B:69:GLY:H	1:B:83:LYS:CB	1.67	1.06
1:B:27:LEU:HD22	1:B:217:LYS:HG3	1.38	1.05
1:B:247:LYS:H	1:B:247:LYS:HD3	0.96	1.03
1:B:262:ARG:HH11	1:B:262:ARG:HG3	0.90	1.02
1:B:47:ASP:HA	1:B:54:GLU:OE2	1.58	1.02
1:B:262:ARG:NH1	1:B:262:ARG:HG3	1.60	1.02
1:B:69:GLY:N	1:B:83:LYS:CB	2.23	1.01
1:B:81:ASN:CG	1:B:82:PRO:CD	2.28	1.01
1:B:69:GLY:CA	1:B:83:LYS:CB	2.32	1.00
1:B:69:GLY:HA2	1:B:83:LYS:HG3	1.38	0.99
1:A:248:LYS:HE3	1:A:248:LYS:HA	1.46	0.97
1:B:262:ARG:HH11	1:B:262:ARG:CG	1.77	0.95
1:B:69:GLY:C	1:B:83:LYS:CD	2.34	0.95
1:B:81:ASN:ND2	1:B:82:PRO:CD	2.29	0.94
1:B:50:THR:HG22	1:B:51:ASP:N	1.82	0.94
1:A:17:SER:CB	1:A:37:ILE:HG12	1.98	0.92
1:B:82:PRO:HG2	1:B:83:LYS:N	1.86	0.91
1:A:336:ALA:O	1:A:340:THR:HG23	1.70	0.91
1:A:108:ARG:HH22	1:A:363:LYS:HD3	1.34	0.91
1:B:283:ARG:HG2	1:B:283:ARG:NH1	1.79	0.90
1:B:69:GLY:CA	1:B:83:LYS:CG	2.48	0.90
1:A:17:SER:HB3	1:A:37:ILE:HG12	1.53	0.87
1:A:37:ILE:CD1	1:A:38:ALA:N	2.30	0.85
1:B:50:THR:HG22	1:B:51:ASP:H	1.42	0.84
1:B:251:LYS:HE3	2:B:436:HOH:O	1.79	0.82
1:B:119:LYS:HD3	1:B:127:MET:HE3	1.62	0.80
1:A:253:THR:C	1:A:255:ASP:H	1.83	0.79
1:B:172:LYS:NZ	1:B:172:LYS:HB3	1.98	0.79
1:B:69:GLY:CA	1:B:83:LYS:HD2	2.15	0.77
1:A:255:ASP:O	1:A:256:LYS:HD3	1.84	0.76
1:B:245:VAL:HG22	1:B:250:LYS:CA	2.16	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:69:GLY:O	1:A:83:LYS:HB2	1.86	0.75
1:B:119:LYS:CD	1:B:127:MET:HE3	2.17	0.75
1:B:82:PRO:CG	1:B:83:LYS:H	1.99	0.73
1:A:227:SER:HB3	1:A:230:GLY:H	1.54	0.73
1:A:37:ILE:HD12	1:A:38:ALA:H	0.68	0.72
1:A:256:LYS:HD2	1:A:259:ARG:H	1.55	0.71
1:B:82:PRO:HG2	1:B:83:LYS:H	1.54	0.71
1:B:69:GLY:H	1:B:83:LYS:HB3	0.88	0.71
1:B:47:ASP:CA	1:B:54:GLU:OE2	2.37	0.70
1:A:108:ARG:NH2	1:A:363:LYS:HD3	2.06	0.70
1:B:50:THR:CG2	1:B:51:ASP:H	2.05	0.70
1:A:321:ARG:HG3	1:A:321:ARG:HH11	1.58	0.68
1:B:265:ASP:O	1:B:283:ARG:HD3	1.92	0.68
1:B:329:ILE:HD12	1:B:345:VAL:HG21	1.75	0.68
1:B:245:VAL:CG2	1:B:250:LYS:HB3	2.24	0.68
1:B:81:ASN:CB	1:B:82:PRO:CD	2.70	0.67
1:B:27:LEU:CD2	1:B:217:LYS:HG3	2.19	0.67
1:B:69:GLY:HA2	1:B:83:LYS:CD	2.25	0.67
1:B:69:GLY:CA	1:B:83:LYS:CD	2.72	0.66
1:B:246:ASP:CB	1:B:247:LYS:HE2	2.26	0.66
1:B:206:LEU:CD1	1:B:239:ILE:HD13	2.26	0.65
1:A:69:GLY:C	1:A:83:LYS:HB2	2.17	0.64
1:B:172:LYS:HZ2	1:B:172:LYS:HB3	1.61	0.64
1:B:68:VAL:HA	1:B:83:LYS:O	1.97	0.63
1:B:262:ARG:NH1	1:B:262:ARG:CG	2.44	0.63
1:A:222:LYS:NZ	1:B:28:GLN:OE1	2.31	0.62
1:A:246:ASP:O	1:A:247:LYS:C	2.33	0.62
1:A:15:LEU:HD12	1:A:32:ASP:OD2	1.99	0.62
1:B:252:VAL:HG12	1:B:257:LEU:HD12	1.82	0.61
1:B:82:PRO:CG	1:B:83:LYS:N	2.39	0.61
1:A:62:TYR:CE2	1:A:91:MET:HE3	2.35	0.61
1:B:247:LYS:O	1:B:248:LYS:HB2	2.00	0.61
1:A:54:GLU:HA	1:A:183:ARG:NH2	2.16	0.60
1:A:313:LYS:NZ	1:B:310:GLN:HE22	1.99	0.60
1:A:348:THR:O	1:A:349:LYS:HD2	2.01	0.60
1:B:119:LYS:HA	1:B:127:MET:HE1	1.84	0.60
1:B:119:LYS:HD2	1:B:127:MET:CE	2.31	0.60
1:A:23:PHE:CE1	1:A:86:LEU:HG	2.36	0.60
1:B:283:ARG:HH11	1:B:283:ARG:CG	1.94	0.60
1:B:267:SER:HB2	1:B:283:ARG:HE	1.67	0.60
1:A:253:THR:C	1:A:255:ASP:N	2.56	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:SER:OG	1:A:37:ILE:HG12	2.03	0.59
1:A:43:ILE:HG21	1:A:103:LEU:HD23	1.85	0.59
1:A:321:ARG:HD3	1:A:321:ARG:N	2.19	0.58
1:B:226:LYS:HD3	1:B:231:TYR:CE1	2.39	0.58
1:B:47:ASP:HA	1:B:54:GLU:OE1	2.03	0.58
1:A:15:LEU:N	1:A:16:PRO:CD	2.67	0.57
1:B:247:LYS:H	1:B:247:LYS:CE	2.18	0.57
1:B:181:VAL:HG22	1:B:182:PRO:HD2	1.86	0.57
1:A:227:SER:OG	1:A:323:ARG:NH2	2.37	0.57
1:A:31:GLY:O	1:A:35:LYS:HA	2.05	0.57
1:B:119:LYS:CD	1:B:127:MET:CE	2.81	0.57
1:B:246:ASP:HB3	1:B:247:LYS:HE2	1.86	0.57
1:A:321:ARG:HH11	1:A:321:ARG:CG	2.17	0.56
1:B:81:ASN:OD1	1:B:82:PRO:HD3	1.99	0.56
1:B:246:ASP:HB2	1:B:247:LYS:HE2	1.88	0.55
1:A:248:LYS:CE	1:A:248:LYS:HA	2.31	0.55
1:A:253:THR:O	1:A:255:ASP:N	2.39	0.55
1:A:158:LYS:HA	1:A:313:LYS:HE3	1.87	0.55
1:A:313:LYS:NZ	1:B:310:GLN:NE2	2.55	0.55
1:A:313:LYS:HZ1	1:B:310:GLN:HE22	1.54	0.55
1:B:58:PHE:HZ	1:B:183:ARG:HD2	1.72	0.55
1:B:95:GLY:HA2	1:B:274:LEU:HD11	1.89	0.54
1:A:273:VAL:HG23	1:A:273:VAL:O	2.07	0.54
1:A:202:TYR:CE2	1:A:244:THR:HG22	2.43	0.54
1:B:261:ILE:CD1	1:B:329:ILE:HD11	2.37	0.54
1:B:127:MET:HA	1:B:127:MET:CE	2.38	0.54
1:B:248:LYS:O	1:B:249:GLY:C	2.44	0.53
1:B:249:GLY:C	1:B:250:LYS:O	2.46	0.52
1:B:228:ASP:OD1	1:B:228:ASP:N	2.29	0.51
1:A:62:TYR:HE2	1:A:91:MET:HE3	1.74	0.51
1:A:194:LEU:HB3	1:B:122:THR:O	2.11	0.51
1:A:15:LEU:HG	1:A:33:GLY:N	2.26	0.50
1:B:146:ASN:O	1:B:147:LYS:HB2	2.12	0.50
1:B:47:ASP:CG	1:B:54:GLU:HG2	2.32	0.49
1:B:192:ALA:O	1:B:196:VAL:HG23	2.12	0.49
1:B:47:ASP:CA	1:B:54:GLU:CD	2.57	0.49
1:A:191:THR:HG23	1:B:123:ASP:OD1	2.13	0.49
1:A:37:ILE:CG1	1:A:38:ALA:N	2.75	0.49
1:B:201:LEU:O	1:B:332:GLY:HA3	2.12	0.49
1:B:206:LEU:HD12	1:B:239:ILE:CD1	2.42	0.48
1:A:119:LYS:HE3	1:A:163:ILE:HB	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:15:LEU:HG	1:A:33:GLY:H	1.79	0.48
1:A:170:GLU:CD	2:A:495:HOH:O	2.51	0.48
1:A:262:ARG:HH11	1:A:347:SER:CB	2.26	0.48
1:B:81:ASN:ND2	1:B:82:PRO:N	2.62	0.48
1:B:117:CYS:HA	1:B:128:VAL:O	2.14	0.47
1:A:256:LYS:HA	1:A:256:LYS:HD3	1.44	0.47
1:B:119:LYS:HD2	1:B:127:MET:HE1	1.96	0.47
1:A:323:ARG:NH1	2:A:453:HOH:O	2.43	0.47
1:A:254:PHE:HA	1:A:257:LEU:HD22	1.97	0.47
1:B:226:LYS:HD2	1:B:227:SER:H	1.79	0.47
1:B:252:VAL:CG1	1:B:257:LEU:HD12	2.44	0.47
1:B:59:ILE:HG12	1:B:178:LEU:HD22	1.97	0.47
1:A:56:SER:OG	1:A:183:ARG:NH1	2.44	0.46
1:B:245:VAL:HG22	1:B:250:LYS:CB	2.45	0.46
1:B:352:LYS:HG2	2:B:475:HOH:O	2.14	0.46
1:B:329:ILE:HD12	1:B:345:VAL:CG2	2.44	0.46
1:B:245:VAL:CG2	1:B:250:LYS:CB	2.93	0.46
1:A:23:PHE:CG	1:A:24:PRO:HA	2.51	0.46
1:A:139:GLN:O	1:A:145:ALA:HB2	2.16	0.46
1:B:203:ASN:OD1	1:B:238:HIS:ND1	2.48	0.46
1:A:114:VAL:HG12	1:A:132:VAL:HB	1.98	0.46
1:B:245:VAL:HG23	1:B:250:LYS:HB3	1.98	0.45
1:B:172:LYS:HB3	1:B:172:LYS:HZ3	1.77	0.45
1:B:69:GLY:H	1:B:83:LYS:CA	2.28	0.45
1:B:189:ILE:HD11	1:B:300:ALA:CB	2.47	0.45
1:A:256:LYS:HE3	1:A:262:ARG:HH22	1.82	0.45
1:B:15:LEU:HD13	1:B:19:ASN:HD22	1.82	0.45
1:A:138:LEU:HA	1:A:138:LEU:HD23	1.66	0.44
1:A:15:LEU:HG	1:A:32:ASP:O	2.18	0.44
1:B:32:ASP:HB3	1:B:33:GLY:HA2	1.99	0.44
1:B:44:GLN:HB2	2:B:424:HOH:O	2.16	0.44
1:B:81:ASN:CB	1:B:82:PRO:HD2	2.47	0.44
1:B:63:GLY:HA3	1:B:172:LYS:O	2.18	0.44
1:A:248:LYS:HE3	1:A:248:LYS:CA	2.32	0.44
1:A:45:ARG:NH1	1:A:47:ASP:OD1	2.51	0.43
1:B:113:MET:HB2	2:B:453:HOH:O	2.18	0.43
1:B:54:GLU:OE2	1:B:182:PRO:HD3	2.18	0.43
1:A:313:LYS:HZ3	1:B:310:GLN:NE2	2.16	0.43
1:B:50:THR:CG2	1:B:51:ASP:N	2.53	0.43
1:B:215:ASP:OD2	1:B:217:LYS:HG2	2.18	0.43
1:A:241:LEU:HD22	1:A:293:PHE:CZ	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:360:ASN:HB2	2:A:462:HOH:O	2.19	0.43
1:B:32:ASP:CB	1:B:33:GLY:HA2	2.49	0.43
1:A:256:LYS:HE3	1:A:259:ARG:HB2	2.00	0.42
1:B:81:ASN:HB3	1:B:82:PRO:HD2	2.02	0.42
1:A:199:SER:HB2	2:A:439:HOH:O	2.20	0.42
1:A:32:ASP:HA	1:A:33:GLY:HA2	1.79	0.42
1:B:59:ILE:HG21	1:B:94:LEU:HD12	2.02	0.42
1:B:310:GLN:NE2	2:B:465:HOH:O	2.44	0.42
1:B:202:TYR:CE2	1:B:257:LEU:HD21	2.55	0.42
1:A:62:TYR:CE2	1:A:91:MET:CE	3.01	0.42
1:A:256:LYS:HD2	1:A:259:ARG:N	2.27	0.42
1:B:133:GLN:HB3	1:B:133:GLN:HE21	1.71	0.42
1:B:329:ILE:CD1	1:B:345:VAL:HG21	2.48	0.41
1:B:127:MET:HE2	1:B:127:MET:CA	2.49	0.41
1:B:353:ARG:HE	1:B:353:ARG:HB2	1.44	0.41
1:A:69:GLY:H	1:A:84:HIS:HD2	1.69	0.41

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	321/364 (88%)	306 (95%)	15 (5%)	0	100	100
1	B	350/364 (96%)	337 (96%)	12 (3%)	1 (0%)	44	49
All	All	671/728 (92%)	643 (96%)	27 (4%)	1 (0%)	55	63

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	82	PRO

5.3.2 Protein sidechains ⓘ

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	280/317 (88%)	254 (91%)	26 (9%)	10	10
1	B	305/317 (96%)	269 (88%)	36 (12%)	6	5
All	All	585/634 (92%)	523 (89%)	62 (11%)	8	7

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	ILE
1	A	68	VAL
1	A	83	LYS
1	A	87	LEU
1	A	92	LEU
1	A	94	LEU
1	A	154	VAL
1	A	186	VAL
1	A	199	SER
1	A	212	VAL
1	A	227	SER
1	A	241	LEU
1	A	247	LYS
1	A	248	LYS
1	A	254	PHE
1	A	256	LYS
1	A	257	LEU
1	A	259	ARG
1	A	264	LEU
1	A	280	VAL
1	A	283	ARG
1	A	286	ARG
1	A	315	LEU
1	A	321	ARG
1	A	340	THR
1	A	349	LYS
1	B	37	ILE

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Mol	Chain	Res	Type
1	B	47	ASP
1	B	49	TRP
1	B	51	ASP
1	B	52	SER
1	B	53	LYS
1	B	81	ASN
1	B	82	PRO
1	B	91	MET
1	B	100	VAL
1	B	127	MET
1	B	133	GLN
1	B	172	LYS
1	B	183	ARG
1	B	184	LYS
1	B	186	VAL
1	B	218	SER
1	B	227	SER
1	B	228	ASP
1	B	229	SER
1	B	239	ILE
1	B	241	LEU
1	B	245	VAL
1	B	246	ASP
1	B	247	LYS
1	B	248	LYS
1	B	251	LYS
1	B	256	LYS
1	B	262	ARG
1	B	283	ARG
1	B	288	ARG
1	B	296	SER
1	B	335	ARG
1	B	353	ARG
1	B	363	LYS
1	B	364	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	84	HIS
1	A	156	HIS
1	B	19	ASN

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Mol	Chain	Res	Type
1	B	81	ASN
1	B	133	GLN
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 carbohydrates 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

6.1 Protein, DNA and RNA chains

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%)	-0.17	15 (4%) 33 32	14, 25, 62, 96	0
1	B	354/364 (97%)	-0.17	11 (3%) 49 47	13, 25, 53, 83	0
All	All	683/728 (93%)	-0.17	26 (3%) 41 39	13, 25, 56, 96	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	248	LYS	5.5
1	B	83	LYS	5.1
1	A	34	LYS	4.2
1	A	54	GLU	4.1
1	A	69	GLY	4.1
1	A	35	LYS	3.4
1	B	248	LYS	3.2
1	A	66	PHE	3.1
1	B	249	GLY	3.1
1	B	74	THR	3.0
1	A	32	ASP	2.9
1	B	288	ARG	2.7
1	A	254	PHE	2.7
1	B	32	ASP	2.5
1	B	52	SER	2.5
1	A	183	ARG	2.5
1	A	246	ASP	2.4
1	B	51	ASP	2.4
1	A	229	SER	2.3
1	A	33	GLY	2.3
1	A	247	LYS	2.3
1	A	228	ASP	2.3
1	B	364	LYS	2.1
1	B	183	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	82	PRO	2.0
1	A	37	ILE	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 carbohydrates 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.