



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 05:24 pm BST

PDB ID : 5WKN
Title : Crystal structure of the parainfluenza virus 5 nucleoprotein-phosphoprotein complex
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Deposited on : 2017-07-25
Resolution : 2.65 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 ⓘ symbol.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
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)	:	2.11

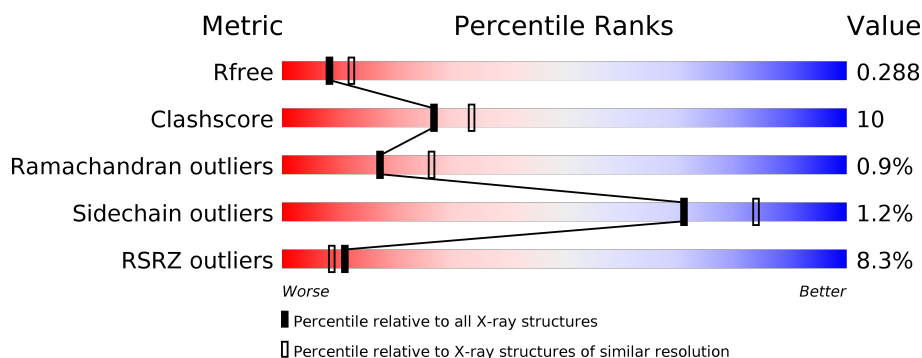
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 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	1332 (2.68-2.64)
Clashscore	141614	1374 (2.68-2.64)
Ramachandran outliers	138981	1349 (2.68-2.64)
Sidechain outliers	138945	1349 (2.68-2.64)
RSRZ outliers	127900	1318 (2.68-2.64)

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 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 $\leq 5\%$. 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	348	<div> <div>6%</div> <div> <div></div> <div>70%</div> <div>21%</div> <div>• 7%</div> </div> </div>
1	B	348	<div> <div>10%</div> <div> <div></div> <div>76%</div> <div>17%</div> <div>• 6%</div> </div> </div>
2	C	49	<div> <div>2%</div> <div> <div></div> <div>31%</div> <div>8%</div> <div>61%</div> </div> </div>
2	D	49	<div> <div>2%</div> <div> <div></div> <div>20%</div> <div>18%</div> <div>61%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5540 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 Nucleoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	323	Total	C	N	O	S	0	0	0
			2588	1658	450	461	19			
1	B	328	Total	C	N	O	S	0	0	0
			2624	1678	455	473	18			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	26	GLY	-	expression tag	UNP Q88435
A	27	SER	-	expression tag	UNP Q88435
A	28	HIS	-	expression tag	UNP Q88435
A	29	MET	-	expression tag	UNP Q88435
A	30	ALA	-	expression tag	UNP Q88435
A	31	SER	-	expression tag	UNP Q88435
A	373	MET	-	expression tag	UNP Q88435
B	26	GLY	-	expression tag	UNP Q88435
B	27	SER	-	expression tag	UNP Q88435
B	28	HIS	-	expression tag	UNP Q88435
B	29	MET	-	expression tag	UNP Q88435
B	30	ALA	-	expression tag	UNP Q88435
B	31	SER	-	expression tag	UNP Q88435
B	373	MET	-	expression tag	UNP Q88435

- Molecule 2 is a protein called Phosphoprotein.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	19	Total	C	N	O	0	0	0
			147	92	22	33			
2	D	19	Total	C	N	O	0	0	0
			146	93	22	31			

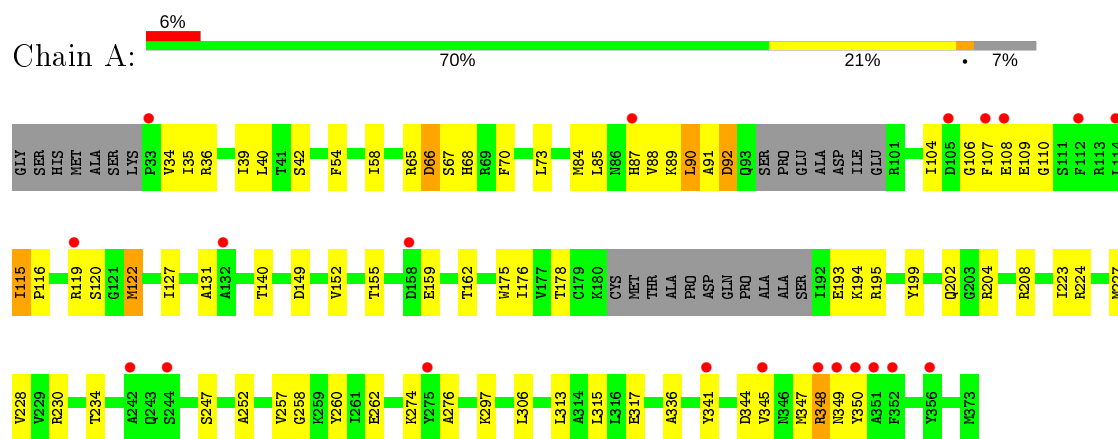
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	17	Total 17	O 17	0	0
3	C	3	Total 3	O 3	0	0
3	B	14	Total 14	O 14	0	0
3	D	1	Total 1	O 1	0	0

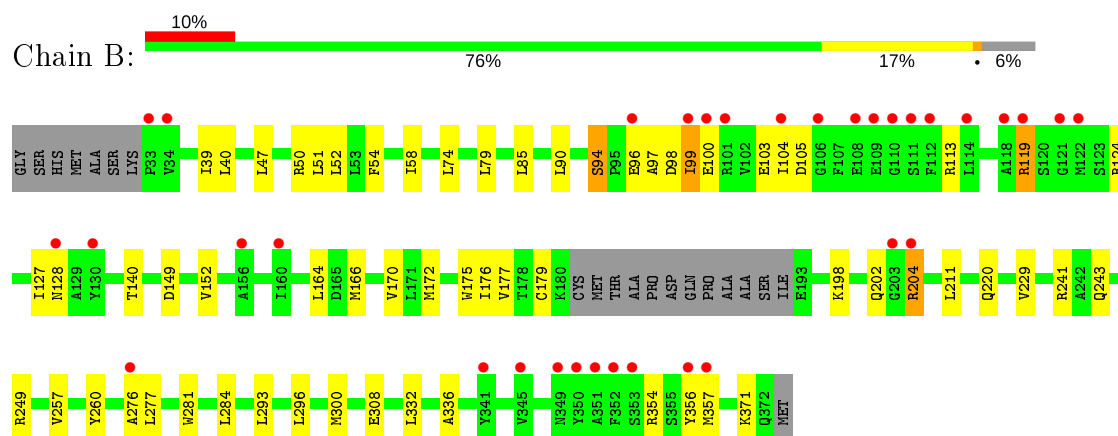
3 Residue-property plots [i](#)

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.

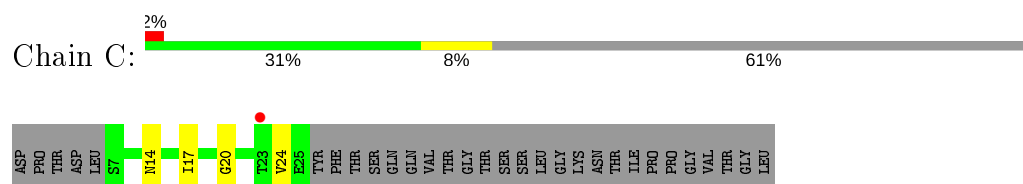
• Molecule 1: Nucleoprotein



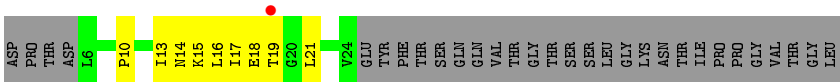
• Molecule 1: Nucleoprotein



• Molecule 2: Phosphoprotein



• Molecule 2: Phosphoprotein



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	79.09 Å 81.06 Å 152.98 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.18 – 2.65 45.50 – 2.65	Depositor EDS
% Data completeness (in resolution range)	99.8 (39.18-2.65) 99.9 (45.50-2.65)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 2.65 Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.251 , 0.288 0.251 , 0.288	Depositor DCC
R_{free} test set	1464 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	66.8	Xtriage
Anisotropy	0.305	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 62.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.015 for k,h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5540	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.69% 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 [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	0.29	0/2640	0.48	0/3562
1	B	0.31	0/2678	0.48	0/3617
2	C	0.29	0/148	0.43	0/200
2	D	0.39	0/147	0.52	0/199
All	All	0.30	0/5613	0.48	0/7578

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	1
1	B	0	3
All	All	0	4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	122	MET	Peptide
1	B	204	ARG	Peptide
1	B	243	GLN	Peptide
1	B	94	SER	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	2588	0	2609	59	0
1	B	2624	0	2634	45	1
2	C	147	0	142	3	0
2	D	146	0	147	12	0
3	A	17	0	0	1	1
3	B	14	0	0	0	0
3	C	3	0	0	0	0
3	D	1	0	0	0	0
All	All	5540	0	5532	113	1

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 10.

All (113) 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:39:ILE:H	1:A:106:GLY:HA2	1.32	0.93
1:B:164:LEU:HD23	1:B:249:ARG:NH2	1.91	0.86
1:A:65:ARG:HD2	1:A:66:ASP:H	1.53	0.73
1:A:65:ARG:HD2	1:A:66:ASP:N	2.04	0.72
1:A:195:ARG:HD3	1:A:260:TYR:HE2	1.54	0.71
1:A:89:LYS:O	1:A:91:ALA:N	2.23	0.71
1:B:354:ARG:NH1	1:B:356:TYR:OH	2.24	0.70
2:D:16:LEU:O	2:D:19:THR:HG22	1.91	0.69
1:A:344:ASP:O	1:A:348:ARG:NH1	2.26	0.68
1:A:109:GLU:OE1	1:A:120:SER:OG	2.08	0.68
1:B:164:LEU:HD23	1:B:249:ARG:HH21	1.59	0.68
1:A:140:THR:O	3:A:401:HOH:O	2.10	0.67
1:B:277:LEU:HB3	2:D:17:ILE:HG23	1.76	0.67
1:B:97:ALA:O	1:B:99:ILE:HD13	1.95	0.66
1:B:90:LEU:HD11	1:B:98:ASP:H	1.62	0.65
1:B:39:ILE:HG21	1:B:85:LEU:HD21	1.77	0.64
1:A:115:ILE:H	1:A:116:PRO:CD	2.11	0.63
1:B:113:ARG:NH1	1:B:128:ASN:OD1	2.31	0.63
1:A:341:TYR:CE2	1:A:348:ARG:HG2	2.35	0.62
1:A:193:GLU:HG2	1:A:194:LYS:H	1.64	0.62
1:B:179:CYS:HB3	1:B:260:TYR:CE2	2.35	0.61
1:B:90:LEU:HA	1:B:94:SER:HB3	1.82	0.61
1:B:179:CYS:HB3	1:B:260:TYR:HE2	1.66	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:277:LEU:HB3	2:D:17:ILE:CG2	2.31	0.61
1:A:88:VAL:HG12	1:A:104:ILE:HG22	1.82	0.61
1:B:52:LEU:HD23	1:B:79:LEU:HD12	1.83	0.60
1:A:341:TYR:CG	1:A:348:ARG:HD2	2.36	0.60
1:B:98:ASP:C	1:B:100:GLU:H	2.05	0.60
1:A:39:ILE:HD12	1:A:85:LEU:HD21	1.84	0.60
1:A:178:THR:HG22	1:A:223:ILE:HD13	1.84	0.59
1:A:116:PRO:HB3	1:A:131:ALA:HB2	1.84	0.59
1:A:149:ASP:HB3	1:A:152:VAL:HG23	1.85	0.59
1:A:274:LYS:HB2	2:C:14:ASN:HD21	1.67	0.58
1:A:104:ILE:HG13	1:A:104:ILE:O	2.02	0.58
1:A:258:GLY:O	1:A:262:GLU:HG3	2.04	0.57
2:D:17:ILE:O	2:D:21:LEU:HD12	2.03	0.57
1:A:195:ARG:HD3	1:A:260:TYR:CE2	2.38	0.56
2:D:13:ILE:C	2:D:15:LYS:H	2.08	0.56
1:A:345:VAL:HA	1:A:348:ARG:HH12	1.69	0.56
1:A:115:ILE:H	1:A:116:PRO:HD2	1.70	0.56
1:A:230:ARG:HH22	1:A:317:GLU:CD	2.10	0.56
1:B:172:MET:O	1:B:176:ILE:HG13	2.06	0.55
1:A:155:THR:OG1	1:A:208:ARG:NH2	2.39	0.55
1:A:40:LEU:HD13	1:A:42:SER:O	2.07	0.55
2:D:17:ILE:O	2:D:21:LEU:CD1	2.56	0.54
1:A:39:ILE:N	1:A:106:GLY:HA2	2.12	0.54
1:A:202:GLN:OE1	1:A:204:ARG:NH1	2.41	0.54
1:A:84:MET:O	1:A:88:VAL:HG23	2.08	0.54
1:B:276:ALA:HB2	1:B:336:ALA:HB2	1.90	0.53
1:B:74:LEU:HD21	1:B:229:VAL:HA	1.90	0.53
1:A:341:TYR:CD2	1:A:348:ARG:HD2	2.43	0.53
1:B:98:ASP:C	1:B:100:GLU:N	2.62	0.52
2:D:18:GLU:CD	2:D:18:GLU:H	2.13	0.52
1:B:281:TRP:HB2	1:B:284:LEU:HD13	1.92	0.52
1:A:35:ILE:HD13	1:A:228:VAL:HG11	1.91	0.51
1:A:224:ARG:O	1:A:230:ARG:NH1	2.44	0.51
1:A:36:ARG:HH22	1:A:122:MET:HG3	1.76	0.50
1:B:300:MET:SD	2:D:18:GLU:OE1	2.70	0.50
1:A:65:ARG:O	1:A:68:HIS:N	2.43	0.50
1:B:170:VAL:HG23	1:B:211:LEU:HD11	1.95	0.49
2:D:13:ILE:HG13	2:D:16:LEU:HD23	1.94	0.49
1:A:40:LEU:HA	1:A:108:GLU:HB3	1.93	0.49
1:B:354:ARG:HB2	1:B:357:MET:HB2	1.95	0.48
1:B:277:LEU:HD13	2:D:17:ILE:HG21	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:124:ARG:HD2	1:B:128:ASN:ND2	2.27	0.48
1:A:306:LEU:HD13	1:A:313:LEU:HD21	1.95	0.47
1:B:293:LEU:HD12	1:B:332:LEU:HD11	1.96	0.47
1:A:347:MET:O	1:A:350:TYR:HB2	2.15	0.47
1:A:348:ARG:HA	1:A:349:ASN:HA	1.61	0.46
1:B:241:ARG:NH2	1:B:308:GLU:OE2	2.46	0.46
1:A:115:ILE:N	1:A:116:PRO:CD	2.77	0.46
1:A:54:PHE:CZ	1:A:58:ILE:HD11	2.50	0.46
1:A:34:VAL:O	1:A:67:SER:OG	2.33	0.46
1:B:198:LYS:O	1:B:202:GLN:HG3	2.16	0.46
1:B:54:PHE:CZ	1:B:58:ILE:HD11	2.51	0.46
1:B:149:ASP:HB3	1:B:152:VAL:HG23	1.97	0.46
1:A:116:PRO:HB2	1:A:127:ILE:CG2	2.46	0.45
1:B:103:GLU:C	1:B:104:ILE:HG13	2.36	0.45
2:C:20:GLY:O	2:C:24:VAL:HG22	2.17	0.45
1:A:176:ILE:HD11	1:A:199:TYR:CD2	2.51	0.45
1:A:87:HIS:ND1	1:A:92:ASP:OD2	2.50	0.45
1:B:40:LEU:HD22	1:B:47:LEU:HD21	1.98	0.44
1:B:177:VAL:HB	1:B:220:GLN:HB2	2.00	0.44
1:B:166:MET:O	1:B:170:VAL:HG12	2.17	0.44
1:A:297:LYS:HG3	2:C:17:ILE:HG23	1.99	0.44
1:A:66:ASP:HB3	1:A:70:PHE:CZ	2.53	0.44
1:B:241:ARG:NH1	1:B:308:GLU:OE2	2.50	0.43
2:D:13:ILE:O	2:D:15:LYS:N	2.51	0.43
1:B:175:TRP:HB3	1:B:257:VAL:HG11	1.99	0.43
1:B:104:ILE:HG22	1:B:105:ASP:N	2.33	0.43
1:B:119:ARG:HA	1:B:119:ARG:HD2	1.83	0.43
1:A:107:PHE:HB3	1:A:120:SER:HB2	2.01	0.43
2:D:18:GLU:CD	2:D:18:GLU:N	2.72	0.43
1:A:108:GLU:HG3	1:A:110:GLY:HA3	2.01	0.43
1:A:109:GLU:N	1:A:110:GLY:CA	2.82	0.42
1:B:113:ARG:HG2	1:B:127:ILE:HG22	2.01	0.42
1:A:276:ALA:HB2	1:A:336:ALA:HB2	2.00	0.42
1:B:99:ILE:CD1	1:B:99:ILE:N	2.82	0.42
1:B:40:LEU:HD21	1:B:51:LEU:HD22	2.02	0.42
1:A:116:PRO:HB2	1:A:127:ILE:HG22	2.01	0.42
1:A:73:LEU:HD21	1:A:223:ILE:HG13	2.01	0.42
1:A:109:GLU:HB2	1:A:110:GLY:HA2	2.00	0.42
1:A:234:THR:HG22	1:A:315:LEU:HD23	2.01	0.42
1:A:39:ILE:HD12	1:A:85:LEU:CD2	2.48	0.41
1:B:103:GLU:O	1:B:104:ILE:HG13	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:227:MET:HA	1:A:230:ARG:NH2	2.36	0.41
1:B:104:ILE:HG22	1:B:105:ASP:H	1.86	0.41
1:B:277:LEU:HD11	1:B:296:LEU:HD13	2.02	0.41
1:B:371:LYS:HB2	1:B:371:LYS:HE3	1.90	0.41
1:A:175:TRP:HB3	1:A:257:VAL:HG11	2.03	0.41
1:B:94:SER:HB2	1:B:96:GLU:O	2.20	0.41
1:A:159:GLU:O	1:A:162:THR:HG22	2.21	0.40
1:A:247:SER:HB2	1:A:252:ALA:HB2	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:140:THR:O	3:A:401:HOH:O[1_455]	2.13	0.07

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	317/348 (91%)	298 (94%)	16 (5%)	3 (1%)	17	26
1	B	324/348 (93%)	307 (95%)	16 (5%)	1 (0%)	41	56
2	C	17/49 (35%)	13 (76%)	4 (24%)	0	100	100
2	D	17/49 (35%)	13 (76%)	2 (12%)	2 (12%)	0	0
All	All	675/794 (85%)	631 (94%)	38 (6%)	6 (1%)	17	26

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	90	LEU
1	B	99	ILE

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Mol	Chain	Res	Type
2	D	14	ASN
1	A	66	ASP
2	D	10	PRO
1	A	115	ILE

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	274/293 (94%)	270 (98%)	4 (2%)	65	80
1	B	278/293 (95%)	275 (99%)	3 (1%)	73	85
2	C	18/44 (41%)	18 (100%)	0	100	100
2	D	18/44 (41%)	18 (100%)	0	100	100
All	All	588/674 (87%)	581 (99%)	7 (1%)	71	84

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	90	LEU
1	A	92	ASP
1	A	119	ARG
1	A	348	ARG
1	B	50	ARG
1	B	119	ARG
1	B	204	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	263	ASN

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	323/348 (92%)	0.59	21 (6%) 18 16	40, 64, 103, 119	1 (0%)
1	B	328/348 (94%)	0.74	34 (10%) 6 4	43, 66, 112, 134	2 (0%)
2	C	19/49 (38%)	0.53	1 (5%) 26 23	70, 85, 97, 109	0
2	D	19/49 (38%)	0.81	1 (5%) 26 23	73, 85, 96, 98	0
All	All	689/794 (86%)	0.66	57 (8%) 11 9	40, 67, 107, 134	3 (0%)

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	110	GLY	15.3
1	B	352	PHE	11.2
1	B	351	ALA	9.8
1	B	99	ILE	7.5
1	B	350	TYR	6.4
1	B	353	SER	6.1
1	B	106	GLY	5.8
1	A	351	ALA	5.5
1	B	118	ALA	5.5
1	B	111	SER	5.0
1	B	122	MET	4.6
1	A	158	ASP	4.3
1	B	109	GLU	4.2
1	A	352	PHE	4.1
1	A	112	PHE	4.0
1	B	356	TYR	3.8
1	B	156	ALA	3.7
1	B	349	ASN	3.7
1	B	341	TYR	3.7
1	B	100	GLU	3.6
1	B	119	ARG	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	345	VAL	3.4
1	A	350	TYR	3.3
1	B	114	LEU	3.3
1	A	244	SER	3.3
1	A	87	HIS	3.2
1	B	34	VAL	3.2
1	B	121	GLY	3.1
1	A	108	GLU	3.1
1	B	160	ILE	3.0
1	A	119	ARG	3.0
1	B	204	ARG	2.9
2	D	19	THR	2.8
1	B	345	VAL	2.8
1	B	276	ALA	2.8
1	B	96	GLU	2.7
1	B	108	GLU	2.7
1	B	101	ARG	2.6
1	A	114	LEU	2.6
1	A	132	ALA	2.6
1	B	130	TYR	2.6
1	B	203	GLY	2.6
1	B	357	MET	2.5
1	B	128	ASN	2.5
1	A	33	PRO	2.5
1	A	356	TYR	2.4
1	B	33	PRO	2.4
1	A	105	ASP	2.3
1	A	107	PHE	2.3
1	B	112	PHE	2.3
1	A	348	ARG	2.2
1	A	349	ASN	2.2
1	A	275	TYR	2.2
1	A	341	TYR	2.1
1	B	104	ILE	2.1
2	C	23	THR	2.1
1	A	242	ALA	2.0

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

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.