



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 06:37 pm BST

PDB ID : 1TZH
Title : Crystal Structure of the Fab YADS1 Complexed with h-VEGF
Authors : Fellouse, F.A.; Wiesmann, C.; Sidhu, S.S.
Deposited on : 2004-07-09
Resolution : 2.60 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

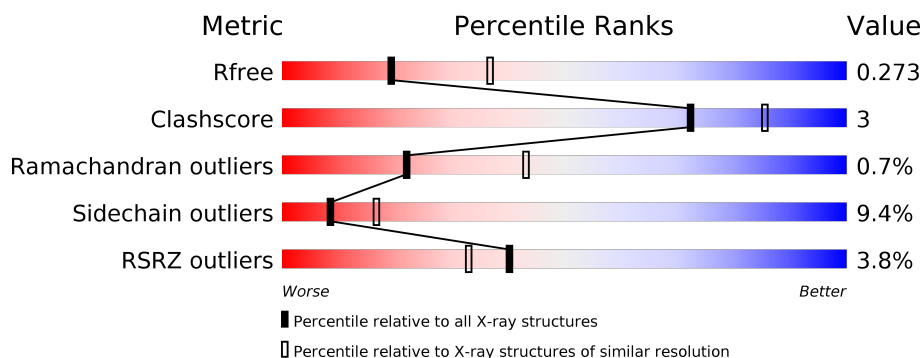
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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	V	102	<div> <div>%</div> <div> <div></div> <div>80%</div> <div>11%</div> <div>8%</div> </div> </div>
1	W	102	<div> <div>4%</div> <div> <div></div> <div>77%</div> <div>13%</div> <div>8%</div> </div> </div>
2	A	213	<div> <div></div> <div> <div></div> <div>78%</div> <div>17%</div> <div>• •</div> </div> </div>
2	L	213	<div> <div>13%</div> <div> <div></div> <div>78%</div> <div>19%</div> <div>• •</div> </div> </div>
3	B	229	<div> <div>%</div> <div> <div></div> <div>82%</div> <div>14%</div> <div>• •</div> </div> </div>
3	H	229	<div> <div>2%</div> <div> <div></div> <div>81%</div> <div>13%</div> <div>• 5%</div> </div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 8104 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 Vascular endothelial growth factor A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	V	94	Total	C	N	O	S	0	0	0
			761	478	128	142	13			
1	W	94	Total	C	N	O	S	0	0	0
			761	478	128	142	13			

- Molecule 2 is a protein called Fab YADS1 Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	210	Total	C	N	O	S	0	0	0
			1599	1001	266	327	5			
2	L	210	Total	C	N	O	S	0	0	0
			1599	1001	266	327	5			

- Molecule 3 is a protein called Fab YADS1 Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	B	222	Total	C	N	O	S	0	0	0
			1653	1040	270	337	6			
3	H	217	Total	C	N	O	S	0	0	0
			1621	1022	264	329	6			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	V	5	Total	O	0	0
			5	5		
4	W	7	Total	O	0	0
			7	7		
4	A	37	Total	O	0	0
			37	37		
4	B	48	Total	O	0	0
			48	48		

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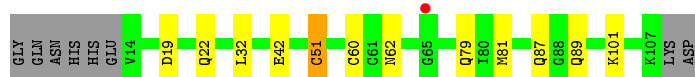
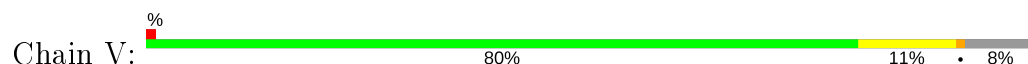
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	8	Total	O	0	0
			8	8		
4	H	5	Total	O	0	0
			5	5		

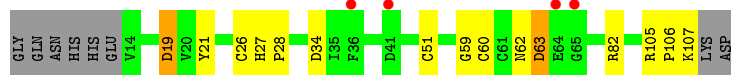
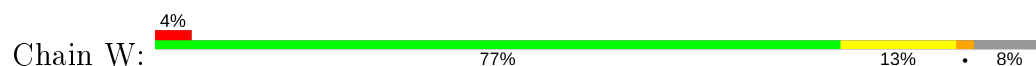
3 Residue-property plots [i](#)

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.

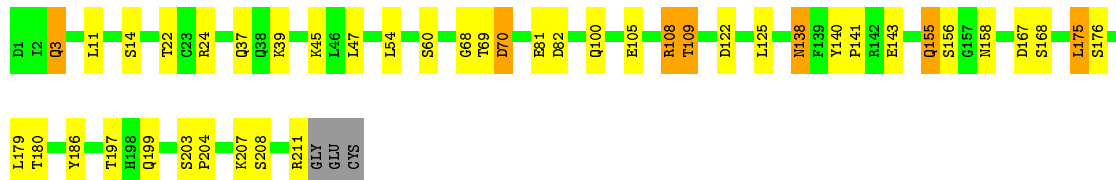
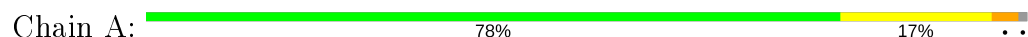
- Molecule 1: Vascular endothelial growth factor A



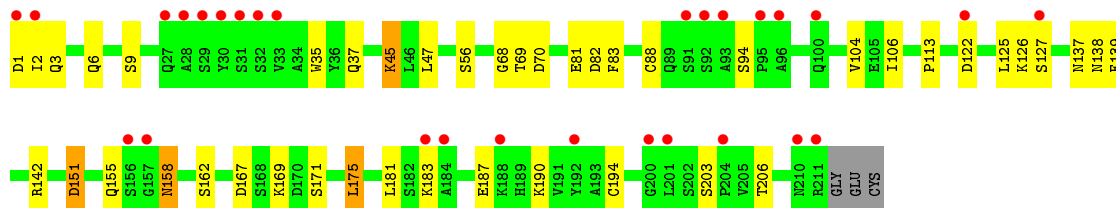
- Molecule 1: Vascular endothelial growth factor A



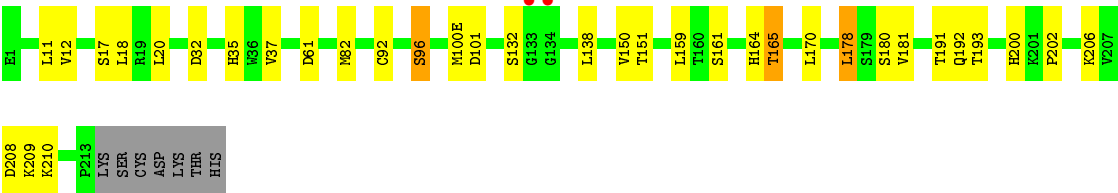
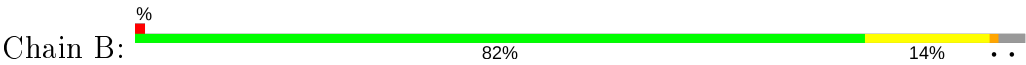
- Molecule 2: Fab YADS1 Light Chain



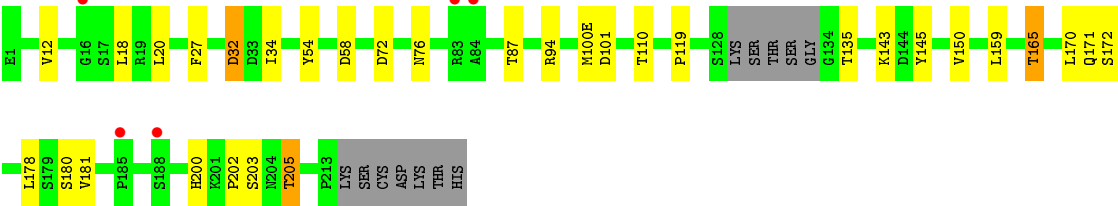
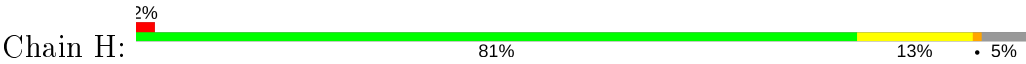
- Molecule 2: Fab YADS1 Light Chain



- Molecule 3: Fab YADS1 Heavy Chain



• Molecule 3: Fab YADS1 Heavy Chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.27Å 76.27Å 112.50Å 90.00° 105.77° 90.00°	Depositor
Resolution (Å)	20.00 – 2.60 29.99 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.8 (20.00-2.60) 99.8 (29.99-2.60)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.50 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.212 , 0.271 0.218 , 0.273	Depositor DCC
R_{free} test set	2108 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	42.0	Xtriage
Anisotropy	0.160	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 47.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8104	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.02% 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	V	0.52	0/779	0.71	1/1050 (0.1%)
1	W	0.56	0/779	0.78	3/1050 (0.3%)
2	A	0.61	0/1634	0.85	5/2220 (0.2%)
2	L	0.44	0/1634	0.72	6/2220 (0.3%)
3	B	0.59	0/1694	0.83	2/2313 (0.1%)
3	H	0.51	0/1661	0.80	4/2268 (0.2%)
All	All	0.54	0/8181	0.79	21/11121 (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
3	H	0	1

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	A	175	LEU	CA-CB-CG	7.90	133.46	115.30
2	L	82	ASP	CB-CG-OD2	6.78	124.40	118.30
1	V	19	ASP	CB-CG-OD2	6.69	124.32	118.30
3	H	58	ASP	CB-CG-OD2	6.52	124.17	118.30
1	W	19	ASP	CB-CG-OD2	6.47	124.12	118.30
3	B	61	ASP	CB-CG-OD2	6.35	124.01	118.30
2	A	167	ASP	CB-CG-OD2	6.20	123.88	118.30
2	L	151	ASP	CB-CG-OD2	6.09	123.78	118.30
2	A	122	ASP	CB-CG-OD2	6.09	123.78	118.30
2	A	70	ASP	CB-CG-OD2	5.90	123.61	118.30
3	H	101	ASP	N-CA-CB	-5.54	100.64	110.60
2	L	1	ASP	CB-CG-OD2	5.53	123.27	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	122	ASP	CB-CG-OD2	5.51	123.25	118.30
2	L	70	ASP	CB-CG-OD2	5.47	123.23	118.30
3	H	72	ASP	CB-CG-OD2	5.47	123.22	118.30
3	H	32	ASP	CB-CG-OD2	5.44	123.19	118.30
2	A	82	ASP	CB-CG-OD2	5.35	123.11	118.30
1	W	63	ASP	CB-CG-OD2	5.32	123.08	118.30
3	B	208	ASP	CB-CG-OD2	5.31	123.08	118.30
2	L	167	ASP	CB-CG-OD2	5.26	123.03	118.30
1	W	34	ASP	CB-CG-OD2	5.21	122.99	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	H	100(E)	MET	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	V	761	0	728	6	0
1	W	761	0	728	6	0
2	A	1599	0	1554	14	0
2	L	1599	0	1554	11	0
3	B	1653	0	1587	12	0
3	H	1621	0	1553	9	0
4	A	37	0	0	2	0
4	B	48	0	0	0	0
4	H	5	0	0	0	0
4	L	8	0	0	1	0
4	V	5	0	0	0	0
4	W	7	0	0	1	0
All	All	8104	0	7704	53	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 3.

All (53) 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)
2:L:45:LYS:NZ	4:L:217:HOH:O	2.08	0.85
2:A:108:ARG:HD3	2:A:109:THR:O	1.79	0.83
1:V:51:CYS:HG	1:W:60:CYS:HG	0.82	0.80
2:A:138:ASN:ND2	3:B:164:HIS:NE2	2.47	0.62
2:A:54:LEU:HD21	2:A:60:SER:HA	1.86	0.57
1:V:81:MET:CE	1:V:89:GLN:HE21	2.17	0.57
2:A:69:THR:HG22	2:A:70:ASP:OD1	2.05	0.56
1:W:19:ASP:HB3	4:W:114:HOH:O	2.06	0.55
3:B:165:THR:HG22	3:B:180:SER:OG	2.07	0.54
1:V:79:GLN:NE2	3:H:54:TYR:O	2.40	0.53
3:H:203:SER:OG	3:H:205:THR:OG1	2.27	0.53
3:B:35:HIS:CB	3:B:100(E):MET:HE2	2.39	0.52
3:B:178:LEU:HD12	3:B:178:LEU:C	2.30	0.52
2:A:140:TYR:CG	2:A:141:PRO:HA	2.45	0.51
3:B:35:HIS:CG	3:B:100(E):MET:HE2	2.46	0.51
3:B:96:SER:HB2	3:B:101:ASP:OD2	2.11	0.51
2:A:45:LYS:NZ	4:A:225:HOH:O	2.44	0.50
2:L:113:PRO:HB3	2:L:139:PHE:HB3	1.94	0.50
3:H:171:GLN:O	3:H:172:SER:HB2	2.12	0.49
2:A:155:GLN:HG3	2:A:179:LEU:HD11	1.95	0.49
2:A:155:GLN:HG3	2:A:158:ASN:HD21	1.78	0.49
2:L:175:LEU:C	2:L:175:LEU:HD12	2.33	0.48
3:H:12:VAL:HG22	3:H:18:LEU:HD22	1.94	0.48
1:V:60:CYS:CB	1:W:51:CYS:HG	2.23	0.48
3:H:200:HIS:CD2	3:H:202:PRO:HD2	2.49	0.48
2:L:158:ASN:C	2:L:158:ASN:HD22	2.17	0.48
1:V:32:LEU:HD13	1:W:59:GLY:HA2	1.97	0.47
1:W:27:HIS:HB2	1:W:28:PRO:HD2	1.96	0.47
2:L:83:PHE:CD1	2:L:104:VAL:HG12	2.50	0.47
2:L:137:ASN:ND2	2:L:138:ASN:OD1	2.49	0.46
2:L:35:TRP:CZ3	2:L:88:CYS:HB3	2.51	0.46
2:A:39:LYS:NZ	2:A:81:GLU:O	2.41	0.45
3:H:165:THR:HB	3:H:180:SER:OG	2.16	0.45
2:L:194:CYS:O	2:L:206:THR:HA	2.16	0.44
3:B:37:VAL:CG2	3:B:100(E):MET:HE1	2.48	0.43
2:A:3:GLN:CG	4:A:242:HOH:O	2.67	0.43
3:B:37:VAL:HG23	3:B:100(E):MET:HE1	2.00	0.43
3:B:12:VAL:HG22	3:B:18:LEU:HD13	2.00	0.43
3:H:119:PRO:HB3	3:H:145:TYR:HB3	2.01	0.43
3:B:193:THR:HG23	3:B:210:LYS:HD2	2.00	0.42
2:L:106:ILE:HG23	2:L:171:SER:HB3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:138:ASN:HD22	2:A:138:ASN:N	2.18	0.42
2:L:83:PHE:HD1	2:L:104:VAL:HG12	1.84	0.42
2:L:37:GLN:HB2	2:L:47:LEU:HD11	2.01	0.42
2:A:37:GLN:HB2	2:A:47:LEU:HD11	2.02	0.42
3:B:200:HIS:CD2	3:B:202:PRO:HD2	2.54	0.42
1:V:81:MET:HE2	1:V:89:GLN:HE21	1.84	0.42
3:H:27:PHE:CZ	3:H:94:ARG:HD2	2.55	0.42
2:A:197:THR:HG22	2:A:204:PRO:HG3	2.02	0.41
1:W:21:TYR:CE1	1:W:62:ASN:ND2	2.88	0.41
3:B:17:SER:HA	3:B:82:MET:O	2.21	0.41
2:A:186:TYR:CE2	2:A:211:ARG:HD2	2.55	0.41
3:H:87:THR:HG23	3:H:110:THR:HA	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	V	92/102 (90%)	88 (96%)	3 (3%)	1 (1%)	14	30
1	W	92/102 (90%)	85 (92%)	4 (4%)	3 (3%)	4	6
2	A	208/213 (98%)	196 (94%)	11 (5%)	1 (0%)	29	52
2	L	208/213 (98%)	196 (94%)	11 (5%)	1 (0%)	29	52
3	B	220/229 (96%)	209 (95%)	11 (5%)	0	100	100
3	H	213/229 (93%)	201 (94%)	11 (5%)	1 (0%)	29	52
All	All	1033/1088 (95%)	975 (94%)	51 (5%)	7 (1%)	22	43

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	L	68	GLY

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Mol	Chain	Res	Type
1	V	42	GLU
1	W	63	ASP
3	H	76	ASN
1	W	26	CYS
1	W	106	PRO
2	A	68	GLY

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	V	89/96 (93%)	84 (94%)	5 (6%)	21	42
1	W	89/96 (93%)	86 (97%)	3 (3%)	37	63
2	A	182/184 (99%)	160 (88%)	22 (12%)	5	9
2	L	182/184 (99%)	158 (87%)	24 (13%)	4	7
3	B	184/191 (96%)	165 (90%)	19 (10%)	7	13
3	H	180/191 (94%)	168 (93%)	12 (7%)	16	33
All	All	906/942 (96%)	821 (91%)	85 (9%)	8	17

All (85) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	V	22	GLN
1	V	51	CYS
1	V	62	ASN
1	V	87	GLN
1	V	101	LYS
1	W	82	ARG
1	W	105	ARG
1	W	107	LYS
2	A	3	GLN
2	A	11	LEU
2	A	14	SER
2	A	22	THR

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Mol	Chain	Res	Type
2	A	24	ARG
2	A	100	GLN
2	A	105	GLU
2	A	108	ARG
2	A	109	THR
2	A	125	LEU
2	A	138	ASN
2	A	143	GLU
2	A	155	GLN
2	A	156	SER
2	A	168	SER
2	A	175	LEU
2	A	176	SER
2	A	180	THR
2	A	199	GLN
2	A	203	SER
2	A	207	LYS
2	A	208	SER
3	B	11	LEU
3	B	20	LEU
3	B	32	ASP
3	B	92	CYS
3	B	96	SER
3	B	132	SER
3	B	138	LEU
3	B	150	VAL
3	B	151	THR
3	B	159	LEU
3	B	161	SER
3	B	165	THR
3	B	170	LEU
3	B	178	LEU
3	B	181	VAL
3	B	191	THR
3	B	192	GLN
3	B	206	LYS
3	B	209	LYS
2	L	2	ILE
2	L	3	GLN
2	L	6	GLN
2	L	9	SER
2	L	45	LYS

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Mol	Chain	Res	Type
2	L	56	SER
2	L	69	THR
2	L	81	GLU
2	L	94	SER
2	L	125	LEU
2	L	126	LYS
2	L	127	SER
2	L	142	ARG
2	L	151	ASP
2	L	155	GLN
2	L	158	ASN
2	L	162	SER
2	L	169	LYS
2	L	175	LEU
2	L	181	LEU
2	L	183	LYS
2	L	187	GLU
2	L	190	LYS
2	L	203	SER
3	H	20	LEU
3	H	32	ASP
3	H	34	ILE
3	H	135	THR
3	H	143	LYS
3	H	150	VAL
3	H	159	LEU
3	H	165	THR
3	H	170	LEU
3	H	178	LEU
3	H	181	VAL
3	H	205	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	V	87	GLN
1	V	89	GLN
1	V	100	ASN
1	W	62	ASN
1	W	87	GLN
1	W	100	ASN
2	A	27	GLN

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Mol	Chain	Res	Type
2	A	137	ASN
2	A	138	ASN
3	B	13	GLN
2	L	89	GLN
2	L	137	ASN
2	L	155	GLN
2	L	158	ASN
2	L	160	GLN
3	H	164	HIS

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	V	94/102 (92%)	-0.20	1 (1%) 80 78	16, 32, 50, 57	0
1	W	94/102 (92%)	0.07	4 (4%) 35 28	14, 33, 50, 57	0
2	A	210/213 (98%)	-0.33	0 100 100	6, 22, 34, 52	0
2	L	210/213 (98%)	0.73	28 (13%) 3 2	17, 30, 45, 65	0
3	B	222/229 (96%)	-0.30	2 (0%) 84 82	10, 19, 32, 48	0
3	H	217/229 (94%)	-0.04	5 (2%) 60 54	10, 23, 33, 47	0
All	All	1047/1088 (96%)	-0.01	40 (3%) 40 33	6, 25, 43, 65	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	92	SER	3.6
2	L	32	SER	3.4
2	L	1	ASP	3.3
2	L	93	ALA	3.3
2	L	184	ALA	3.2
2	L	31	SER	3.1
2	L	201	LEU	3.1
2	L	211	ARG	3.1
2	L	2	ILE	3.0
2	L	28	ALA	2.9
2	L	30	TYR	2.9
2	L	157	GLY	2.9
2	L	200	GLY	2.9
2	L	95	PRO	2.8
1	W	64	GLU	2.8
3	B	133	GLY	2.8
2	L	188	LYS	2.8
2	L	29	SER	2.7
2	L	91	SER	2.7

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Mol	Chain	Res	Type	RSRZ
2	L	127	SER	2.6
2	L	96	ALA	2.6
2	L	33	VAL	2.6
3	H	16	GLY	2.6
3	H	185	PRO	2.5
2	L	183	LYS	2.5
2	L	122	ASP	2.5
2	L	156	SER	2.4
3	H	188	SER	2.3
2	L	27	GLN	2.3
1	W	65	GLY	2.2
2	L	100	GLN	2.2
1	W	36	PHE	2.1
3	H	83	ARG	2.1
1	V	65	GLY	2.1
2	L	204	PRO	2.1
2	L	192	TYR	2.1
2	L	210	ASN	2.1
3	H	84	ALA	2.0
1	W	41	ASP	2.0
3	B	134	GLY	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.