



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 29, 2016 – 12:48 PM EDT

PDB ID : 5AOE
Title : Crystal structure of pneumolysin D168A mutant.
Authors : van Pee, K.; Yildiz, O.
Deposited on : 2015-09-10
Resolution : 2.50 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
Mogul : unknown
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027939
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : rb-20027939

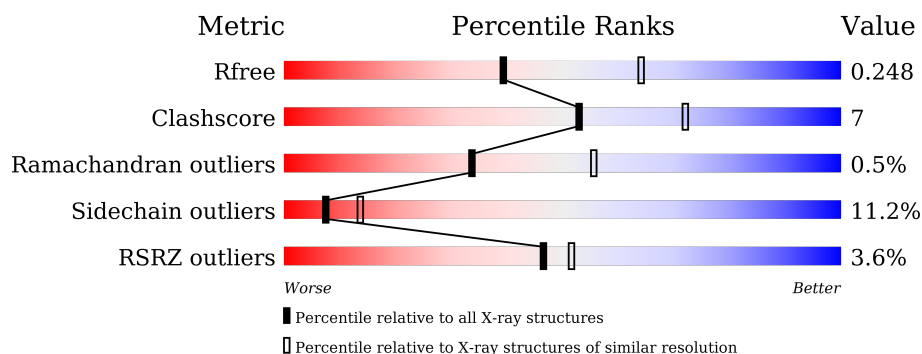
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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	488	<div> <div>3%</div> <div> <div></div> <div>79%</div> <div>17%</div> <div>.</div> </div> </div>
1	B	488	<div> <div>4%</div> <div> <div></div> <div>75%</div> <div>21%</div> <div>.</div> </div> </div>

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 7878 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 PNEUMOLYSIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	488	Total	C	N	O	S	0	0	0
			3862	2432	666	757	7			
1	B	488	Total	C	N	O	S	0	0	0
			3862	2432	666	757	7			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-16	ALA	-	EXPRESSION TAG	UNP Q04IN8
A	-15	HIS	-	EXPRESSION TAG	UNP Q04IN8
A	-14	HIS	-	EXPRESSION TAG	UNP Q04IN8
A	-13	HIS	-	EXPRESSION TAG	UNP Q04IN8
A	-12	HIS	-	EXPRESSION TAG	UNP Q04IN8
A	-11	HIS	-	EXPRESSION TAG	UNP Q04IN8
A	-10	HIS	-	EXPRESSION TAG	UNP Q04IN8
A	-9	SER	-	EXPRESSION TAG	UNP Q04IN8
A	-8	SER	-	EXPRESSION TAG	UNP Q04IN8
A	-7	GLY	-	EXPRESSION TAG	UNP Q04IN8
A	-6	LEU	-	EXPRESSION TAG	UNP Q04IN8
A	-5	VAL	-	EXPRESSION TAG	UNP Q04IN8
A	-4	PRO	-	EXPRESSION TAG	UNP Q04IN8
A	-3	ARG	-	EXPRESSION TAG	UNP Q04IN8
A	-2	GLY	-	EXPRESSION TAG	UNP Q04IN8
A	-1	SER	-	EXPRESSION TAG	UNP Q04IN8
A	0	HIS	-	EXPRESSION TAG	UNP Q04IN8
A	151	GLU	ASP	ENGINEERED MUTATION	UNP Q04IN8
B	-16	ALA	-	EXPRESSION TAG	UNP Q04IN8
B	-15	HIS	-	EXPRESSION TAG	UNP Q04IN8
B	-14	HIS	-	EXPRESSION TAG	UNP Q04IN8
B	-13	HIS	-	EXPRESSION TAG	UNP Q04IN8
B	-12	HIS	-	EXPRESSION TAG	UNP Q04IN8
B	-11	HIS	-	EXPRESSION TAG	UNP Q04IN8
B	-10	HIS	-	EXPRESSION TAG	UNP Q04IN8

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-9	SER	-	EXPRESSION TAG	UNP Q04IN8
B	-8	SER	-	EXPRESSION TAG	UNP Q04IN8
B	-7	GLY	-	EXPRESSION TAG	UNP Q04IN8
B	-6	LEU	-	EXPRESSION TAG	UNP Q04IN8
B	-5	VAL	-	EXPRESSION TAG	UNP Q04IN8
B	-4	PRO	-	EXPRESSION TAG	UNP Q04IN8
B	-3	ARG	-	EXPRESSION TAG	UNP Q04IN8
B	-2	GLY	-	EXPRESSION TAG	UNP Q04IN8
B	-1	SER	-	EXPRESSION TAG	UNP Q04IN8
B	0	HIS	-	EXPRESSION TAG	UNP Q04IN8
B	151	GLU	ASP	ENGINEERED MUTATION	UNP Q04IN8

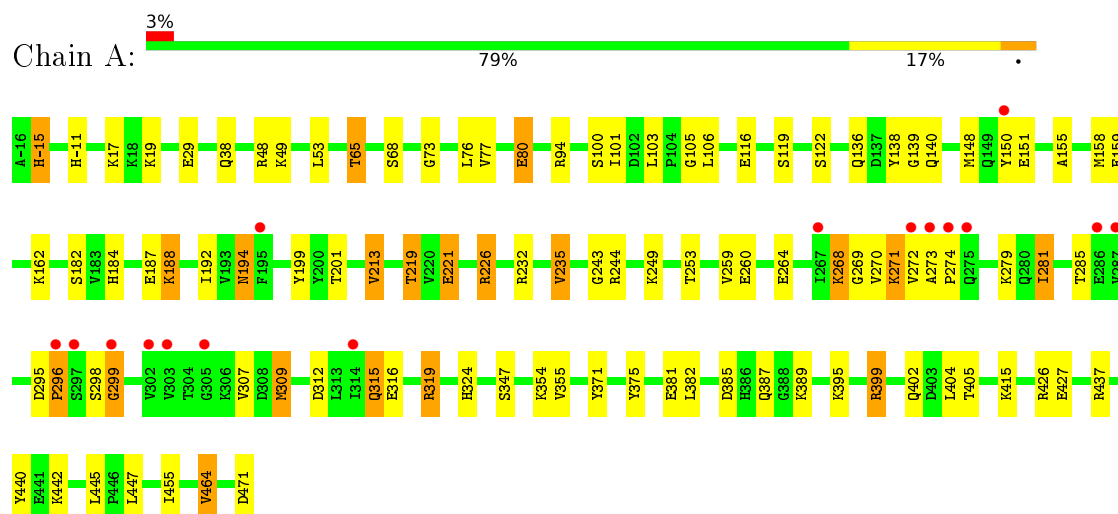
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	80	Total O 80 80	0	0
2	B	74	Total O 74 74	0	0

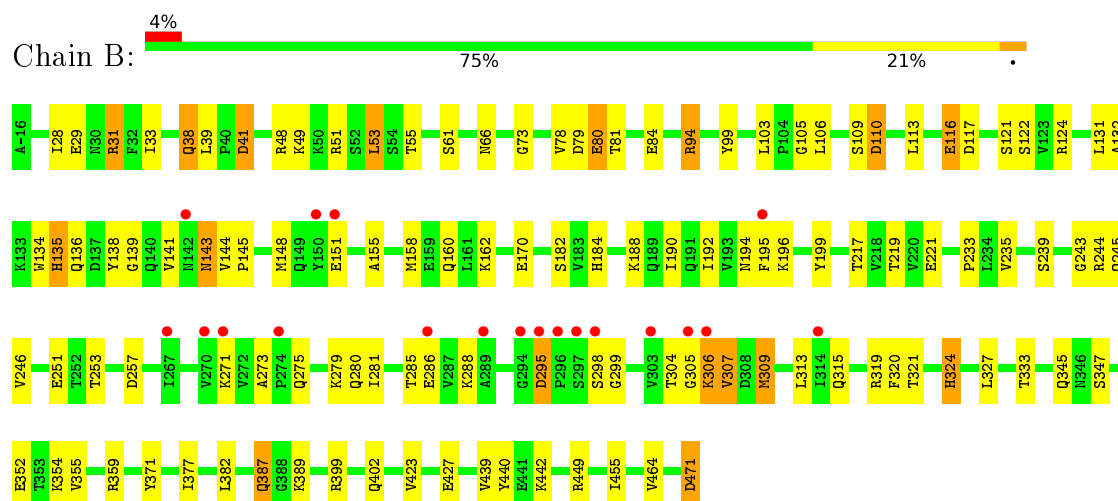
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ($\text{RSRZ} > 2$). 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: PNEUMOLYSIN



• Molecule 1: PNEUMOLYSIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	160.86Å 24.66Å 208.35Å 90.00° 90.26° 90.00°	Depositor
Resolution (Å)	49.49 – 2.50 49.49 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.5 (49.49-2.50) 98.3 (49.49-2.50)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.50 (at 2.51Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.222 , 0.250 0.221 , 0.248	Depositor DCC
R_{free} test set	2960 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	46.6	Xtriage
Anisotropy	0.023	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 34.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.409 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7878	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.27% 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	0.22	0/3943	0.41	0/5359
1	B	0.24	0/3943	0.41	0/5359
All	All	0.23	0/7886	0.41	0/10718

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	3862	0	3793	43	0
1	B	3862	0	3793	58	0
2	A	80	0	0	3	0
2	B	74	0	0	0	0
All	All	7878	0	7586	101	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 7.

All (101) 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:38:GLN:HE21	1:B:38:GLN:HA	1.40	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:73:GLY:O	1:B:94:ARG:NH2	2.16	0.78
1:B:151:GLU:HB2	1:B:194:ASN:HB3	1.69	0.74
1:B:273:ALA:O	1:B:275:GLN:NE2	2.24	0.71
1:A:73:GLY:O	1:A:94:ARG:NH2	2.23	0.71
1:A:295:ASP:HB3	1:A:296:PRO:HD2	1.71	0.71
1:B:321:THR:H	1:B:324:HIS:CE1	2.12	0.68
1:B:105:GLY:HA3	1:B:138:TYR:HE2	1.61	0.66
1:B:116:GLU:OE1	1:B:117:ASP:N	2.27	0.66
1:B:131:LEU:O	1:B:135:HIS:ND1	2.29	0.65
1:B:257:ASP:OD2	1:B:354:LYS:NZ	2.30	0.65
1:A:319:ARG:HE	1:A:324:HIS:HE1	1.43	0.65
1:B:455:ILE:HG22	1:B:464:VAL:HG12	1.79	0.65
1:B:399:ARG:NH1	1:B:402:GLN:OE1	2.30	0.64
1:A:94:ARG:NH1	2:A:2038:HOH:O	2.28	0.63
1:A:48:ARG:NH1	1:A:184:HIS:O	2.32	0.63
1:A:65:THR:HG21	1:A:299:GLY:HA2	1.80	0.63
1:B:324:HIS:O	1:B:324:HIS:ND1	2.31	0.62
1:B:41:ASP:OD1	1:B:41:ASP:N	2.32	0.62
1:A:182:SER:HB3	1:A:188:LYS:HG3	1.81	0.61
1:A:150:TYR:OH	1:A:268:LYS:NZ	2.34	0.61
1:B:143:ASN:OD1	1:B:145:PRO:HG3	1.99	0.61
1:B:320:PHE:HA	1:B:324:HIS:HE1	1.67	0.60
1:A:309:MET:HE3	1:A:309:MET:H	1.66	0.59
1:B:80:GLU:O	1:B:84:GLU:HG2	2.03	0.59
1:B:29:GLU:OE2	1:B:31:ARG:NH1	2.35	0.59
1:A:259:VAL:HG22	1:A:281:ILE:HD12	1.85	0.58
1:A:151:GLU:HB2	1:A:194:ASN:HB2	1.86	0.57
1:B:281:ILE:O	1:B:285:THR:OG1	2.21	0.57
1:B:251:GLU:HG3	1:B:286:GLU:HB2	1.86	0.57
1:B:219:THR:HG22	1:B:221:GLU:H	1.68	0.56
1:B:38:GLN:CA	1:B:38:GLN:HE21	2.17	0.56
1:B:196:LYS:HG2	1:B:245:GLN:HG3	1.88	0.56
1:B:449:ARG:HE	1:B:471:ASP:H	1.53	0.55
1:B:288:LYS:HG3	1:B:304:THR:HG22	1.88	0.55
1:A:159:GLU:OE2	1:A:162:LYS:NZ	2.39	0.55
1:B:103:LEU:HB2	1:B:106:LEU:HD12	1.88	0.55
1:B:79:ASP:OD1	1:B:81:THR:OG1	2.21	0.54
1:B:51:ARG:NE	1:B:352:GLU:OE1	2.40	0.54
1:A:269:GLY:O	1:A:271:LYS:NZ	2.36	0.54
1:B:440:TYR:CZ	1:B:442:LYS:HG3	2.43	0.53
1:B:134:TRP:HE3	1:B:135:HIS:CE1	2.26	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:110:ASP:N	1:B:110:ASP:OD1	2.40	0.53
1:B:48:ARG:NH2	1:B:184:HIS:O	2.37	0.50
1:B:305:GLY:HA3	1:B:309:MET:SD	2.51	0.50
1:A:399:ARG:NH1	1:A:402:GLN:OE1	2.43	0.49
1:B:298:SER:OG	1:B:299:GLY:N	2.45	0.49
1:A:199:TYR:CD2	1:A:243:GLY:HA2	2.47	0.49
1:A:268:LYS:HB3	1:A:270:VAL:HG23	1.95	0.48
1:B:144:VAL:O	1:B:144:VAL:HG23	2.12	0.48
1:B:275:GLN:HB3	1:B:279:LYS:HD3	1.96	0.48
1:A:281:ILE:O	1:A:285:THR:OG1	2.26	0.48
1:B:190:ILE:HG12	1:B:251:GLU:HB3	1.96	0.48
1:B:199:TYR:CD2	1:B:243:GLY:HA2	2.49	0.48
1:B:195:PHE:HB2	1:B:246:VAL:HB	1.95	0.48
1:B:235:VAL:CG1	1:B:333:THR:HB	2.44	0.47
1:A:272:VAL:HB	1:A:274:PRO:HD2	1.97	0.47
1:B:53:LEU:HD12	1:B:160:GLN:HG2	1.95	0.47
1:B:182:SER:HB3	1:B:188:LYS:HG2	1.95	0.47
1:B:135:HIS:HA	1:B:139:GLY:HA3	1.97	0.47
1:A:-15:HIS:O	1:A:-11:HIS:ND1	2.42	0.47
1:A:38:GLN:NE2	1:A:415:LYS:HB3	2.30	0.47
1:B:295:ASP:HA	1:B:324:HIS:CD2	2.50	0.46
1:A:119:SER:H	1:A:122:SER:HB2	1.79	0.46
1:B:321:THR:H	1:B:324:HIS:HE1	1.62	0.46
1:A:38:GLN:HE22	1:A:415:LYS:HB3	1.80	0.46
1:A:105:GLY:HA3	1:A:138:TYR:CD1	2.49	0.46
1:B:306:LYS:HD2	1:B:306:LYS:HA	1.69	0.46
1:A:219:THR:HG23	1:A:221:GLU:H	1.81	0.46
1:A:440:TYR:CZ	1:A:442:LYS:HG3	2.51	0.46
1:A:371:TYR:CG	1:A:427:GLU:HG3	2.50	0.45
1:B:279:LYS:HB3	1:B:279:LYS:NZ	2.32	0.45
1:A:226:ARG:NH1	2:A:2036:HOH:O	2.45	0.45
1:A:381:GLU:HG3	1:A:395:LYS:HG3	1.97	0.45
1:A:77:VAL:HG11	1:A:232:ARG:CZ	2.47	0.44
1:A:94:ARG:NH1	1:A:213:VAL:HG22	2.32	0.44
1:B:78:VAL:HB	1:B:233:PRO:HD2	1.99	0.44
1:A:151:GLU:HB2	1:A:194:ASN:CB	2.47	0.44
1:B:99:TYR:O	1:B:113:LEU:N	2.50	0.44
1:B:155:ALA:HB2	1:B:192:ILE:HD12	2.00	0.44
1:A:309:MET:HB3	1:A:309:MET:HE2	1.81	0.43
1:A:76:LEU:HB2	1:A:235:VAL:HG12	2.00	0.43
1:A:244:ARG:NH2	1:A:316:GLU:OE1	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:272:VAL:HG12	1:A:273:ALA:H	1.84	0.43
1:A:455:ILE:HG22	1:A:464:VAL:HG13	1.99	0.43
1:A:80:GLU:HG3	1:A:80:GLU:H	1.49	0.42
1:B:387:GLN:HB2	1:B:389:LYS:HD2	2.00	0.42
1:A:445:LEU:HD12	1:A:445:LEU:HA	1.87	0.42
1:B:132:ALA:O	1:B:136:GLN:HG2	2.20	0.41
1:B:377:ILE:HG12	1:B:423:VAL:HG13	2.02	0.41
1:A:347:SER:HB3	2:A:2022:HOH:O	2.19	0.41
1:A:103:LEU:HD12	1:A:103:LEU:HA	1.82	0.41
1:A:312:ASP:HA	1:A:315:GLN:HB2	2.02	0.41
1:B:29:GLU:H	1:B:29:GLU:CD	2.24	0.41
1:A:155:ALA:HB2	1:A:192:ILE:HD12	2.02	0.41
1:B:304:THR:O	1:B:304:THR:OG1	2.31	0.41
1:A:375:TYR:CD1	1:A:404:LEU:HD12	2.57	0.40
1:B:162:LYS:HD3	1:B:170:GLU:HB2	2.03	0.40
1:B:244:ARG:NH1	1:B:313:LEU:O	2.43	0.40
1:B:78:VAL:HG21	1:B:233:PRO:HG2	2.03	0.40
1:B:371:TYR:CG	1:B:427:GLU:HG3	2.56	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	486/488 (100%)	467 (96%)	15 (3%)	4 (1%)	24	41
1	B	486/488 (100%)	465 (96%)	20 (4%)	1 (0%)	52	75
All	All	972/976 (100%)	932 (96%)	35 (4%)	5 (0%)	34	55

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	296	PRO
1	B	307	VAL
1	A	-15	HIS
1	A	299	GLY
1	A	139	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	A	429/429 (100%)	378 (88%)	51 (12%)	6	12
1	B	429/429 (100%)	384 (90%)	45 (10%)	8	16
All	All	858/858 (100%)	762 (89%)	96 (11%)	7	14

All (96) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	LYS
1	A	19	LYS
1	A	29	GLU
1	A	49	LYS
1	A	53	LEU
1	A	65	THR
1	A	68	SER
1	A	80	GLU
1	A	100	SER
1	A	101	ILE
1	A	106	LEU
1	A	116	GLU
1	A	136	GLN
1	A	140	GLN
1	A	148	MET
1	A	158	MET
1	A	187	GLU
1	A	188	LYS
1	A	194	ASN

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Mol	Chain	Res	Type
1	A	201	THR
1	A	213	VAL
1	A	219	THR
1	A	221	GLU
1	A	226	ARG
1	A	235	VAL
1	A	249	LYS
1	A	253	THR
1	A	260	GLU
1	A	264	GLU
1	A	268	LYS
1	A	271	LYS
1	A	279	LYS
1	A	281	ILE
1	A	298	SER
1	A	307	VAL
1	A	309	MET
1	A	315	GLN
1	A	319	ARG
1	A	354	LYS
1	A	355	VAL
1	A	382	LEU
1	A	385	ASP
1	A	387	GLN
1	A	389	LYS
1	A	399	ARG
1	A	405	THR
1	A	426	ARG
1	A	437	ARG
1	A	447	LEU
1	A	464	VAL
1	A	471	ASP
1	B	28	ILE
1	B	31	ARG
1	B	33	ILE
1	B	38	GLN
1	B	39	LEU
1	B	41	ASP
1	B	49	LYS
1	B	53	LEU
1	B	55	THR
1	B	61	SER

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Mol	Chain	Res	Type
1	B	66	ASN
1	B	80	GLU
1	B	94	ARG
1	B	109	SER
1	B	110	ASP
1	B	116	GLU
1	B	121	SER
1	B	122	SER
1	B	124	ARG
1	B	135	HIS
1	B	141	VAL
1	B	143	ASN
1	B	148	MET
1	B	158	MET
1	B	217	THR
1	B	239	SER
1	B	253	THR
1	B	271	LYS
1	B	280	GLN
1	B	295	ASP
1	B	306	LYS
1	B	307	VAL
1	B	309	MET
1	B	315	GLN
1	B	319	ARG
1	B	324	HIS
1	B	327	LEU
1	B	345	GLN
1	B	347	SER
1	B	355	VAL
1	B	359	ARG
1	B	382	LEU
1	B	387	GLN
1	B	439	VAL
1	B	471	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	324	HIS
1	B	38	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	488/488 (100%)	0.23	16 (3%) 50 55	27, 67, 131, 188	0
1	B	488/488 (100%)	0.16	19 (3%) 43 48	24, 63, 136, 196	0
All	All	976/976 (100%)	0.20	35 (3%) 46 51	24, 65, 135, 196	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	303	VAL	7.0
1	A	296	PRO	6.1
1	B	297	SER	5.9
1	B	267	ILE	5.5
1	A	275	GLN	5.4
1	B	296	PRO	5.3
1	A	274	PRO	5.2
1	A	273	ALA	4.5
1	B	303	VAL	4.4
1	B	305	GLY	4.4
1	B	295	ASP	4.0
1	B	274	PRO	3.4
1	A	195	PHE	3.4
1	B	150	TYR	3.4
1	A	314	ILE	3.3
1	A	267	ILE	3.3
1	A	272	VAL	3.3
1	A	297	SER	3.3
1	A	299	GLY	3.2
1	B	286	GLU	3.0
1	B	142	ASN	2.8
1	A	150	TYR	2.8
1	B	294	GLY	2.8
1	B	314	ILE	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	270	VAL	2.8
1	A	305	GLY	2.6
1	A	287	VAL	2.5
1	B	289	ALA	2.4
1	A	302	VAL	2.4
1	B	151	GLU	2.4
1	A	286	GLU	2.2
1	B	271	LYS	2.2
1	B	195	PHE	2.1
1	B	306	LYS	2.1
1	B	298	SER	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.