



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:23 AM UTC

PDB ID : 3C0M / pdb_00003c0m
Title : Crystal structure of the proaerolysin mutant Y221G
Authors : Pernot, L.; Schiltz, M.; Thurnheer, S.; Burr, S.E.; van der Goot, G.
Deposited on : 2008-01-21
Resolution : 2.88 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

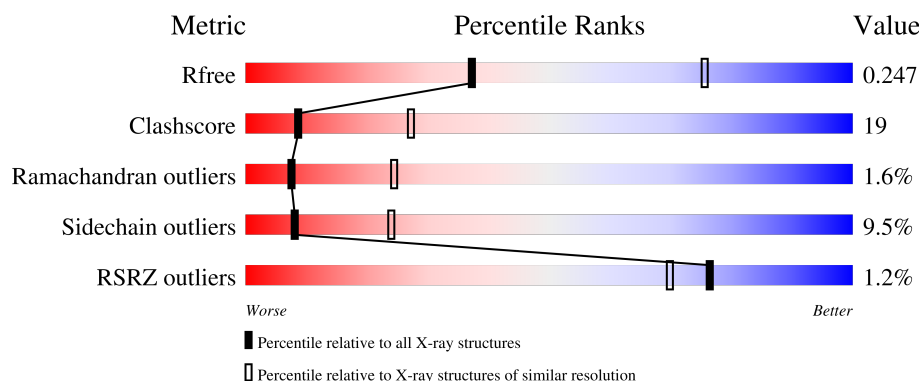
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.88 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}	180053	3557 (2.90-2.86)
Clashscore	190562	3801 (2.90-2.86)
Ramachandran outliers	187476	3699 (2.90-2.86)
Sidechain outliers	187428	3702 (2.90-2.86)
RSRZ outliers	180081	3558 (2.90-2.86)

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 $\leq 5\%$. 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	470	2% 54% 37% 5%
1	B	470	62% 28% 5%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7054 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 Aerolysin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	449	Total	C	N	O	S	0	0	0
			3515	2221	605	680	9			
1	B	449	Total	C	N	O	S	0	0	0
			3515	2221	605	680	9			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	221	GLY	TYR	engineered mutation	UNP P09167
B	221	GLY	TYR	engineered mutation	UNP P09167

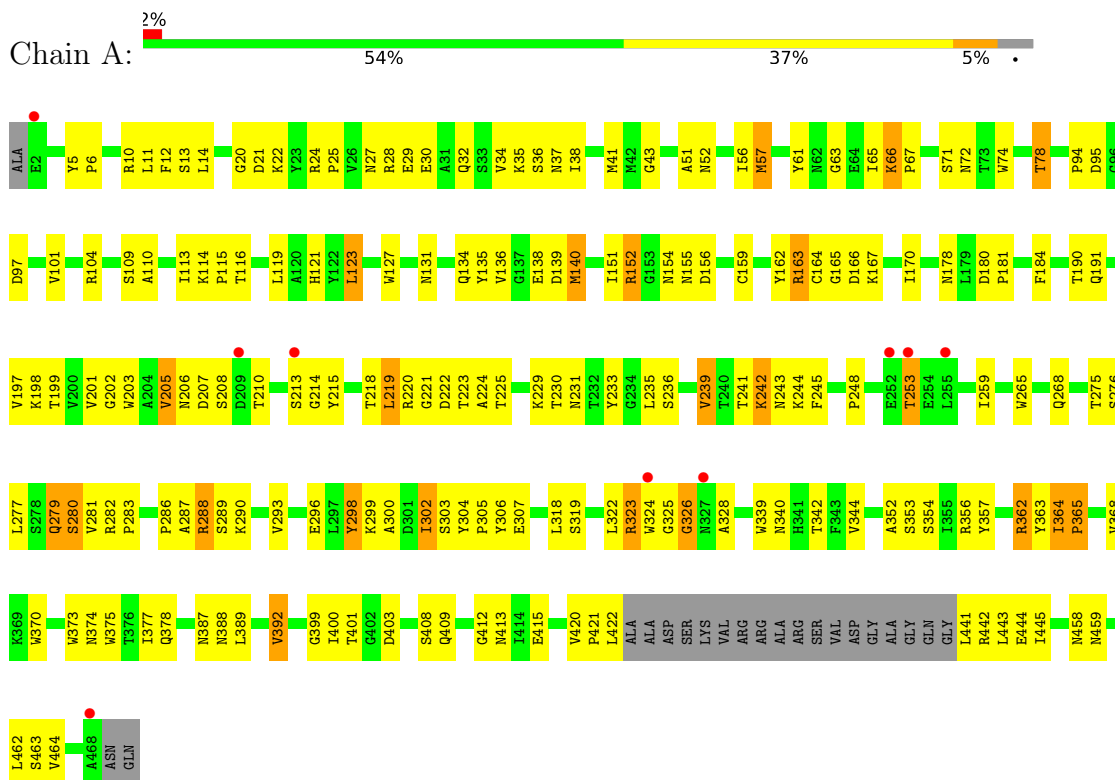
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	8	Total	O	0	0
			8	8		
2	B	16	Total	O	0	0
			16	16		

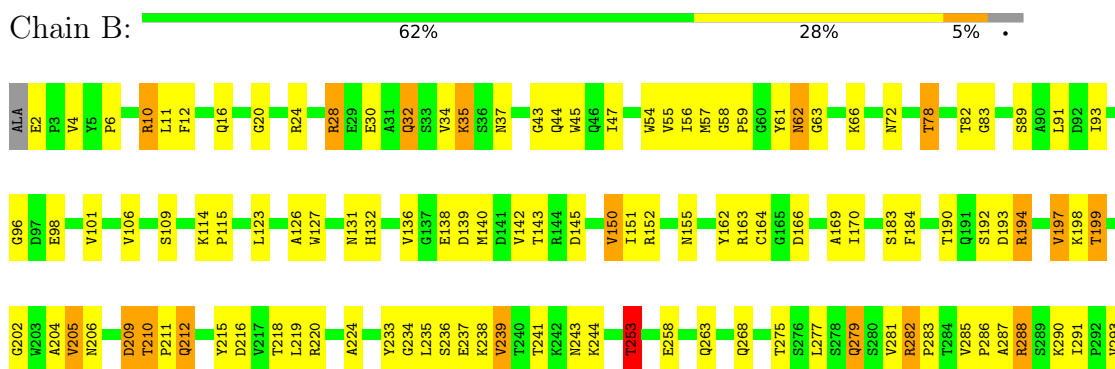
3 Residue-property plots

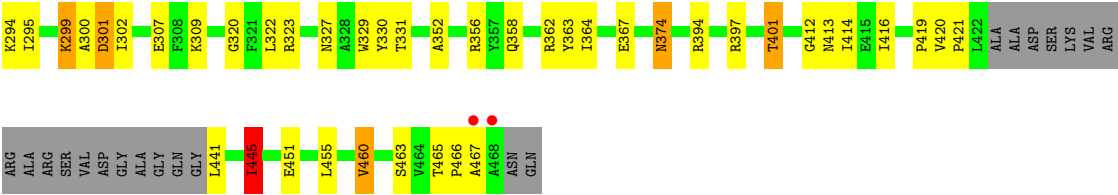
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Aerolysin



• Molecule 1: Aerolysin





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.62Å 92.43Å 169.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.56 – 2.88 44.56 – 2.88	Depositor EDS
% Data completeness (in resolution range)	93.5 (44.56-2.88) 93.4 (44.56-2.88)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.59 (at 2.86Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.199 , 0.252 0.197 , 0.247	Depositor DCC
R_{free} test set	1700 reflections (6.87%)	wwPDB-VP
Wilson B-factor (Å ²)	43.3	Xtriage
Anisotropy	0.240	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 38.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7054	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.26% 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	0.49	0/3612	1.00	10/4930 (0.2%)
1	B	0.52	0/3612	1.04	18/4930 (0.4%)
All	All	0.50	0/7224	1.02	28/9860 (0.3%)

There are no bond length outliers.

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	78	THR	N-CA-C	-7.38	102.93	112.23
1	B	445	ILE	CA-C-N	7.18	127.44	119.90
1	B	445	ILE	C-N-CA	7.18	127.44	119.90
1	B	106	VAL	N-CA-C	6.68	117.44	110.62
1	B	209	ASP	N-CA-C	-6.19	105.56	113.23
1	B	78	THR	N-CA-C	-6.11	105.59	113.16
1	B	412	GLY	N-CA-C	-5.96	103.47	112.41
1	B	401	THR	N-CA-C	5.88	118.12	109.24
1	A	97	ASP	N-CA-C	-5.88	102.62	110.55
1	B	58	GLY	CA-C-N	5.73	125.59	119.28
1	B	58	GLY	C-N-CA	5.73	125.59	119.28
1	A	110	ALA	N-CA-C	5.66	117.45	111.28
1	A	458	ASN	N-CA-C	5.65	118.93	109.95
1	B	184	PHE	N-CA-C	5.51	118.06	110.35
1	A	412	GLY	N-CA-C	-5.42	103.33	112.77
1	B	93	ILE	CA-C-N	-5.38	114.71	120.03
1	B	93	ILE	C-N-CA	-5.38	114.71	120.03
1	A	205	VAL	N-CA-C	5.36	116.20	108.48
1	B	320	GLY	N-CA-C	-5.35	102.41	111.08
1	B	329	TRP	N-CA-C	-5.34	102.37	110.28
1	B	299	LYS	CB-CA-C	5.32	118.01	110.24
1	A	298	TYR	N-CA-C	5.30	118.23	109.85
1	A	52	ASN	N-CA-C	5.29	118.63	111.17
1	B	205	VAL	N-CA-C	5.27	116.17	108.53
1	A	224	ALA	N-CA-C	5.20	117.57	109.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	43	GLY	N-CA-C	-5.17	106.64	112.33
1	A	139	ASP	N-CA-C	-5.14	104.90	110.91
1	B	224	ALA	N-CA-C	5.07	118.28	110.42

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	3515	0	3350	151	0
1	B	3515	0	3350	110	0
2	A	8	0	0	0	0
2	B	16	0	0	2	0
All	All	7054	0	6700	258	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 19.

All (258) 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:445:ILE:HD11	1:A:462:LEU:HD13	1.43	1.00
1:A:202:GLY:HA3	1:A:445:ILE:HG22	1.45	0.95
1:B:47:ILE:HD12	1:B:57:MET:HG2	1.54	0.90
1:A:206:ASN:ND2	1:A:287:ALA:HA	1.87	0.89
1:A:94:PRO:HG2	1:A:104:ARG:NH1	1.87	0.89
1:A:206:ASN:HD22	1:A:287:ALA:HA	1.42	0.84
1:B:322:LEU:HD13	1:B:327:ASN:HD21	1.41	0.84
1:A:445:ILE:HD11	1:A:462:LEU:CD1	2.13	0.78
1:A:279:GLN:OE1	1:A:413:ASN:HA	1.85	0.75
1:B:6:PRO:HG3	1:B:34:VAL:HG11	1.70	0.74
1:A:219:LEU:HB2	1:A:462:LEU:HD21	1.70	0.74
1:A:27:ASN:OD1	1:A:30:GLU:HG3	1.88	0.74
1:A:202:GLY:HA3	1:A:445:ILE:CG2	2.18	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:28:ARG:HD2	1:B:54:TRP:CE2	2.23	0.73
1:A:363:TYR:O	1:A:365:PRO:HD3	1.90	0.72
1:B:322:LEU:HB3	1:B:327:ASN:HD22	1.57	0.70
1:B:282:ARG:NH1	1:B:467:ALA:HB1	2.07	0.70
1:B:101:VAL:HG21	1:B:235:LEU:HD22	1.74	0.69
1:B:114:LYS:HB3	1:B:115:PRO:HD3	1.74	0.69
1:B:293:VAL:HG12	1:B:416:ILE:HG22	1.73	0.69
1:A:362:ARG:HG3	1:A:374:ASN:ND2	2.08	0.69
1:B:28:ARG:HD2	1:B:54:TRP:CD2	2.28	0.69
1:B:6:PRO:HG3	1:B:34:VAL:CG1	2.23	0.68
1:B:57:MET:HE2	1:B:61:TYR:CG	2.29	0.68
1:B:291:ILE:HG22	1:B:419:PRO:HA	1.75	0.67
1:A:362:ARG:HG3	1:A:374:ASN:HD21	1.58	0.67
1:B:219:LEU:HD22	1:B:295:ILE:HD13	1.74	0.67
1:A:276:SER:C	1:A:277:LEU:HD12	2.20	0.67
1:B:11:LEU:HD11	1:B:72:ASN:HD22	1.60	0.67
1:B:132:HIS:HB2	1:B:139:ASP:CG	2.20	0.66
1:B:465:THR:HB	1:B:466:PRO:HD2	1.76	0.66
1:A:57:MET:O	1:A:63:GLY:HA2	1.96	0.66
1:B:322:LEU:HB3	1:B:327:ASN:ND2	2.11	0.65
1:A:94:PRO:O	1:A:104:ARG:NH1	2.30	0.65
1:B:322:LEU:HD13	1:B:327:ASN:ND2	2.11	0.64
1:A:202:GLY:CA	1:A:445:ILE:HG22	2.24	0.64
1:A:43:GLY:HA3	1:B:10:ARG:NH2	2.13	0.63
1:A:219:LEU:HB2	1:A:462:LEU:CD2	2.26	0.63
1:A:389:LEU:HA	1:A:392:VAL:CG1	2.27	0.63
1:A:364:ILE:HD12	1:A:364:ILE:N	2.14	0.63
1:B:281:VAL:HG13	1:B:283:PRO:HD3	1.80	0.63
1:A:207:ASP:OD2	1:A:442:ARG:NH1	2.32	0.62
1:A:191:GLN:H	1:A:303:SER:HB2	1.64	0.62
1:B:47:ILE:CD1	1:B:57:MET:HG2	2.28	0.62
1:A:307:GLU:HG3	1:A:401:THR:HG22	1.81	0.62
1:A:34:VAL:HG12	1:A:38:ILE:HD12	1.81	0.61
1:A:319:SER:HB3	1:A:340:ASN:OD1	1.99	0.61
1:A:259:ILE:HD13	1:A:265:TRP:CZ3	2.35	0.61
1:B:35:LYS:HE2	1:B:63:GLY:O	2.00	0.61
1:A:162:TYR:CD1	1:A:163:ARG:HB2	2.36	0.61
1:B:397:ARG:HD2	2:B:476:HOH:O	2.00	0.61
1:A:56:ILE:HG13	1:A:65:ILE:HG12	1.83	0.60
1:A:441:LEU:O	1:A:464:VAL:HG11	2.01	0.60
1:B:82:THR:HG22	1:B:83:GLY:H	1.67	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:ASN:OD1	1:A:159:CYS:HB3	2.02	0.60
1:A:286:PRO:O	1:A:289:SER:HB2	2.02	0.59
1:A:442:ARG:NH2	1:A:444:GLU:OE2	2.36	0.59
1:B:82:THR:HG22	1:B:83:GLY:N	2.18	0.59
1:A:218:THR:HB	1:A:280:SER:OG	2.02	0.58
1:B:57:MET:O	1:B:63:GLY:HA2	2.04	0.58
1:A:178:ASN:HD21	1:A:242:LYS:HG3	1.67	0.58
1:A:388:ASN:O	1:A:392:VAL:HG12	2.04	0.58
1:B:163:ARG:HG2	1:B:166:ASP:OD2	2.02	0.58
1:B:362:ARG:HD2	1:B:374:ASN:ND2	2.18	0.58
1:A:443:LEU:HG	1:A:464:VAL:CG2	2.33	0.58
1:A:205:VAL:HG22	1:A:290:LYS:HB3	1.86	0.58
1:A:6:PRO:HG3	1:A:34:VAL:HG13	1.86	0.57
1:B:151:ILE:HB	1:B:170:ILE:HB	1.85	0.57
1:A:276:SER:O	1:A:277:LEU:HD12	2.04	0.57
1:A:253:THR:HG21	1:A:302:ILE:HG13	1.86	0.57
1:A:34:VAL:HG12	1:A:34:VAL:O	2.04	0.57
1:B:285:VAL:HG11	1:B:441:LEU:HD11	1.87	0.57
1:A:114:LYS:HE3	1:A:138:GLU:HG3	1.87	0.56
1:A:170:ILE:HD12	1:A:318:LEU:HD22	1.86	0.56
1:A:6:PRO:HG3	1:A:34:VAL:CG1	2.36	0.56
1:B:397:ARG:CD	2:B:476:HOH:O	2.53	0.56
1:A:203:TRP:N	1:A:445:ILE:HG22	2.21	0.56
1:A:373:TRP:O	1:A:377:ILE:HG13	2.06	0.56
1:B:4:VAL:HG23	1:B:34:VAL:HG21	1.88	0.56
1:A:199:THR:HG22	1:A:201:VAL:HG13	1.88	0.56
1:A:208:SER:O	1:A:287:ALA:HB1	2.06	0.56
1:A:344:VAL:HG23	1:A:353:SER:HB2	1.87	0.56
1:B:44:GLN:HE21	1:B:45:TRP:NE1	2.05	0.55
1:A:135:TYR:CE2	1:B:294:LYS:HD3	2.41	0.55
1:A:324:TRP:O	1:A:325:GLY:C	2.49	0.55
1:A:12:PHE:CZ	1:A:20:GLY:HA3	2.42	0.55
1:A:223:THR:HG21	1:A:409:GLN:O	2.07	0.55
1:A:375:TRP:O	1:A:378:GLN:HG2	2.06	0.55
1:A:365:PRO:O	1:A:368:VAL:HG12	2.07	0.54
1:B:12:PHE:CE1	1:B:20:GLY:HA3	2.41	0.54
1:A:322:LEU:HD13	1:A:339:TRP:HB2	1.89	0.54
1:A:399:GLY:C	1:A:400:ILE:HD12	2.32	0.54
1:B:162:TYR:HE1	1:B:163:ARG:HE	1.53	0.54
1:A:287:ALA:C	1:A:288:ARG:HG2	2.33	0.54
1:B:114:LYS:NZ	1:B:138:GLU:HG3	2.22	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:194:ARG:O	1:B:194:ARG:HG3	2.06	0.54
1:A:229:LYS:HA	1:A:403:ASP:O	2.08	0.54
1:A:387:ASN:N	1:A:387:ASN:HD22	2.05	0.54
1:B:279:GLN:HG2	1:B:413:ASN:HA	1.89	0.53
1:B:197:VAL:CG1	1:B:299:LYS:HB3	2.38	0.53
1:A:57:MET:HE2	1:A:61:TYR:CD1	2.44	0.53
1:A:13:SER:O	1:A:14:LEU:HD23	2.09	0.53
1:B:91:LEU:HB2	1:B:394:ARG:HE	1.73	0.53
1:A:151:ILE:HB	1:A:170:ILE:HB	1.89	0.53
1:A:154:ASN:OD1	1:A:156:ASP:HB2	2.09	0.53
1:B:145:ASP:HB2	1:B:150:VAL:CG1	2.40	0.52
1:B:220:ARG:HA	1:B:277:LEU:O	2.08	0.52
1:A:328:ALA:HB2	1:A:375:TRP:CZ3	2.45	0.52
1:A:25:PRO:HB2	1:A:51:ALA:HB2	1.90	0.52
1:A:328:ALA:HB2	1:A:375:TRP:CE3	2.45	0.52
1:A:443:LEU:HG	1:A:464:VAL:HG21	1.90	0.52
1:B:34:VAL:HG12	1:B:34:VAL:O	2.10	0.52
1:B:162:TYR:CD1	1:B:162:TYR:C	2.87	0.52
1:B:205:VAL:HG22	1:B:290:LYS:HB3	1.92	0.52
1:A:389:LEU:HA	1:A:392:VAL:HG13	1.92	0.52
1:A:5:TYR:O	1:A:6:PRO:C	2.53	0.52
1:B:162:TYR:CD1	1:B:163:ARG:HB2	2.44	0.52
1:A:11:LEU:HD11	1:A:72:ASN:HD22	1.75	0.51
1:A:207:ASP:OD1	1:A:422:LEU:HD21	2.10	0.51
1:A:184:PHE:CG	1:A:248:PRO:HG3	2.46	0.51
1:B:451:GLU:O	1:B:455:LEU:HD13	2.11	0.51
1:A:420:VAL:HG13	1:A:421:PRO:HD2	1.92	0.51
1:A:245:PHE:CE1	1:A:259:ILE:HD12	2.46	0.51
1:A:162:TYR:CD1	1:A:162:TYR:C	2.88	0.51
1:B:127:TRP:CD1	1:B:164:CYS:HA	2.46	0.51
1:B:206:ASN:ND2	1:B:212:GLN:H	2.09	0.50
1:B:282:ARG:HD2	1:B:467:ALA:CB	2.41	0.50
1:A:239:VAL:HG22	1:A:265:TRP:HB2	1.92	0.50
1:A:323:ARG:HB2	1:A:326:GLY:O	2.12	0.50
1:A:41:MET:HE2	1:A:74:TRP:HH2	1.76	0.50
1:A:162:TYR:CE1	1:A:163:ARG:HB2	2.47	0.50
1:B:362:ARG:NH1	1:B:363:TYR:CE2	2.79	0.50
1:B:4:VAL:HG11	1:B:30:GLU:HB3	1.94	0.50
1:B:44:GLN:HA	1:B:59:PRO:HG2	1.94	0.49
1:A:127:TRP:CE2	1:A:164:CYS:HA	2.47	0.49
1:A:206:ASN:HD22	1:A:287:ALA:CA	2.21	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:61:TYR:O	1:B:62:ASN:HB2	2.10	0.49
1:B:34:VAL:O	1:B:35:LYS:C	2.56	0.49
1:B:287:ALA:O	1:B:288:ARG:C	2.56	0.49
1:A:230:THR:OG1	1:A:403:ASP:HB2	2.13	0.49
1:A:203:TRP:H	1:A:445:ILE:HG22	1.77	0.48
1:B:47:ILE:HD11	1:B:61:TYR:CE2	2.48	0.48
1:A:222:ASP:HA	1:A:275:THR:O	2.14	0.48
1:A:34:VAL:CG1	1:A:37:ASN:HB2	2.43	0.48
1:A:121:HIS:ND1	1:A:323:ARG:NH2	2.61	0.48
1:A:163:ARG:CG	1:A:166:ASP:OD2	2.62	0.48
1:B:282:ARG:HH11	1:B:467:ALA:HB1	1.78	0.48
1:B:210:THR:CG2	1:B:210:THR:O	2.62	0.48
1:B:291:ILE:HG22	1:B:419:PRO:CA	2.42	0.48
1:B:126:ALA:HB2	1:B:323:ARG:HG2	1.95	0.48
1:A:306:TYR:C	1:A:306:TYR:CD1	2.91	0.48
1:B:234:GLY:O	1:B:235:LEU:C	2.57	0.48
1:A:140:MET:HG3	1:A:152:ARG:O	2.14	0.48
1:A:302:ILE:HG23	1:A:303:SER:N	2.29	0.48
1:B:34:VAL:HG12	1:B:37:ASN:HB2	1.96	0.47
1:A:445:ILE:O	1:A:445:ILE:CG1	2.62	0.47
1:A:131:ASN:OD1	1:A:167:LYS:NZ	2.38	0.47
1:B:307:GLU:HB2	1:B:401:THR:HG22	1.95	0.47
1:B:244:LYS:HD2	1:B:258:GLU:OE1	2.14	0.47
1:A:21:ASP:O	1:A:22:LYS:HB2	2.13	0.47
1:A:225:THR:HG23	1:A:408:SER:HB3	1.97	0.47
1:B:192:SER:HA	1:B:301:ASP:O	2.15	0.47
1:A:205:VAL:HG13	1:A:422:LEU:HD13	1.97	0.47
1:A:400:ILE:HD12	1:A:400:ILE:N	2.30	0.47
1:B:89:SER:O	1:B:394:ARG:CZ	2.63	0.47
1:A:101:VAL:HG21	1:A:235:LEU:HD22	1.96	0.47
1:A:304:TYR:CD1	1:A:304:TYR:C	2.93	0.46
1:B:374:ASN:HD22	1:B:374:ASN:HA	1.53	0.46
1:A:205:VAL:O	1:A:441:LEU:HD12	2.16	0.46
1:A:253:THR:HB	1:A:300:ALA:HB1	1.98	0.46
1:A:307:GLU:HG3	1:A:401:THR:CG2	2.46	0.46
1:B:62:ASN:HD22	1:B:62:ASN:HA	1.58	0.46
1:B:114:LYS:HZ2	1:B:138:GLU:HG3	1.80	0.46
1:B:322:LEU:CD1	1:B:327:ASN:HD21	2.20	0.46
1:A:288:ARG:C	1:A:422:LEU:HB2	2.41	0.46
1:B:2:GLU:O	1:B:2:GLU:HG2	2.14	0.46
1:A:57:MET:HB3	1:A:61:TYR:HB2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:253:THR:HB	1:A:300:ALA:CB	2.46	0.46
1:A:57:MET:HE2	1:A:61:TYR:CG	2.50	0.45
1:A:197:VAL:O	1:A:198:LYS:HB2	2.15	0.45
1:B:34:VAL:CG1	1:B:37:ASN:HB2	2.45	0.45
1:B:364:ILE:HG21	1:B:367:GLU:HG2	1.98	0.45
1:A:198:LYS:HG2	1:A:199:THR:N	2.32	0.45
1:A:221:GLY:C	1:A:277:LEU:HD13	2.42	0.45
1:A:229:LYS:O	1:A:229:LYS:HG3	2.17	0.45
1:A:229:LYS:HE2	1:A:265:TRP:CE2	2.52	0.45
1:B:32:GLN:HE21	1:B:32:GLN:CA	2.30	0.45
1:A:66:LYS:HB2	1:A:67:PRO:CD	2.47	0.44
1:B:234:GLY:O	1:B:237:GLU:HG3	2.16	0.44
1:B:279:GLN:HB3	1:B:414:ILE:HG13	1.99	0.44
1:A:445:ILE:CD1	1:A:462:LEU:HD13	2.30	0.44
1:A:34:VAL:O	1:A:35:LYS:C	2.60	0.44
1:A:445:ILE:O	1:A:445:ILE:HG13	2.18	0.44
1:A:184:PHE:CD1	1:A:248:PRO:HG3	2.52	0.44
1:A:95:ASP:OD1	1:A:233:TYR:HB2	2.18	0.44
1:B:126:ALA:HA	1:B:323:ARG:HE	1.83	0.44
1:A:123:LEU:O	1:A:328:ALA:HB3	2.18	0.43
1:A:155:ASN:O	1:A:156:ASP:C	2.61	0.43
1:A:109:SER:O	1:A:114:LYS:HB2	2.18	0.43
1:B:32:GLN:HA	1:B:32:GLN:NE2	2.33	0.43
1:B:127:TRP:NE1	1:B:164:CYS:HA	2.33	0.43
1:B:352:ALA:O	1:B:358:GLN:NE2	2.51	0.43
1:A:307:GLU:CG	1:A:401:THR:HG22	2.47	0.43
1:B:109:SER:O	1:B:114:LYS:HB2	2.18	0.43
1:B:420:VAL:HG13	1:B:421:PRO:HD2	2.00	0.43
1:A:245:PHE:CZ	1:A:259:ILE:HD12	2.54	0.43
1:A:140:MET:HG2	1:A:151:ILE:HG22	1.99	0.43
1:B:183:SER:OG	1:B:309:LYS:N	2.51	0.43
1:B:330:TYR:CE1	1:B:331:THR:HG23	2.54	0.43
1:A:368:VAL:O	1:A:368:VAL:HG22	2.18	0.43
1:A:364:ILE:N	1:A:364:ILE:CD1	2.82	0.43
1:B:202:GLY:HA3	1:B:445:ILE:HB	2.01	0.43
1:A:352:ALA:HB2	1:A:370:TRP:NE1	2.32	0.43
1:B:198:LYS:HG2	1:B:199:THR:N	2.34	0.43
1:B:206:ASN:HD21	1:B:212:GLN:H	1.65	0.43
1:B:220:ARG:O	1:B:460:VAL:HA	2.18	0.43
1:A:28:ARG:O	1:A:29:GLU:C	2.60	0.43
1:A:210:THR:C	1:A:287:ALA:HB2	2.44	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:354:SER:HB3	1:A:357:TYR:HB3	2.00	0.42
1:B:152:ARG:NH2	1:B:169:ALA:HB2	2.33	0.42
1:A:12:PHE:CE1	1:A:20:GLY:HA3	2.54	0.42
1:A:163:ARG:C	1:A:165:GLY:N	2.76	0.42
1:B:445:ILE:O	1:B:445:ILE:HG12	2.19	0.42
1:A:213:SER:O	1:A:215:TYR:N	2.39	0.42
1:A:259:ILE:HG23	1:A:268:GLN:OE1	2.19	0.42
1:B:96:GLY:HA2	1:B:233:TYR:CG	2.54	0.42
1:B:236:SER:HA	1:B:239:VAL:HG13	2.00	0.42
1:B:253:THR:HB	1:B:300:ALA:CB	2.49	0.42
1:A:109:SER:HA	1:A:113:ILE:HB	2.02	0.42
1:A:116:THR:O	1:A:119:LEU:HB3	2.19	0.42
1:A:180:ASP:HA	1:A:181:PRO:HD2	1.86	0.42
1:A:302:ILE:CG2	1:A:303:SER:N	2.82	0.42
1:B:152:ARG:HH21	1:B:169:ALA:HB2	1.85	0.42
1:B:56:ILE:HG23	1:B:56:ILE:O	2.19	0.41
1:B:132:HIS:HB2	1:B:139:ASP:OD2	2.19	0.41
1:B:162:TYR:CE1	1:B:163:ARG:HB2	2.56	0.41
1:A:296:GLU:OE1	1:A:298:TYR:OH	2.34	0.41
1:B:253:THR:HB	1:B:300:ALA:HB3	2.02	0.41
1:A:304:TYR:HA	1:A:305:PRO:HD3	1.91	0.41
1:B:215:TYR:HA	1:B:467:ALA:H	1.86	0.41
1:A:114:LYS:HB3	1:A:115:PRO:CD	2.50	0.41
1:A:34:VAL:O	1:A:34:VAL:CG1	2.69	0.41
1:B:234:GLY:C	1:B:236:SER:N	2.76	0.41
1:B:263:GLN:HB3	1:B:268:GLN:NE2	2.36	0.41
1:A:207:ASP:HA	1:A:422:LEU:CD2	2.51	0.41
1:A:223:THR:OG1	1:A:409:GLN:HB2	2.21	0.41
1:A:229:LYS:NZ	1:A:231:ASN:HD21	2.19	0.41
1:A:242:LYS:HE3	1:A:242:LYS:HB2	1.86	0.41
1:A:281:VAL:HG12	1:A:282:ARG:N	2.36	0.41
1:B:216:ASP:OD1	1:B:216:ASP:N	2.49	0.41
1:A:415:GLU:OE1	1:B:114:LYS:HE2	2.21	0.41
1:A:239:VAL:HG22	1:A:265:TRP:CB	2.52	0.40
1:B:204:ALA:O	1:B:290:LYS:HA	2.21	0.40
1:B:235:LEU:HD12	1:B:235:LEU:HA	1.88	0.40
1:B:57:MET:SD	1:B:66:LYS:HE2	2.61	0.40
1:B:98:GLU:OE2	1:B:238:LYS:HD2	2.20	0.40
1:A:34:VAL:HG12	1:A:38:ILE:CD1	2.49	0.40
1:A:66:LYS:HB2	1:A:67:PRO:HD2	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	445/470 (95%)	397 (89%)	41 (9%)	7 (2%)	7	24
1	B	445/470 (95%)	406 (91%)	32 (7%)	7 (2%)	7	24
All	All	890/940 (95%)	803 (90%)	73 (8%)	14 (2%)	7	24

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	214	GLY
1	B	190	THR
1	B	253	THR
1	A	323	ARG
1	B	193	ASP
1	A	190	THR
1	A	244	LYS
1	A	326	GLY
1	B	35	LYS
1	A	362	ARG
1	B	28	ARG
1	B	286	PRO
1	A	365	PRO
1	B	211	PRO

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	378/391 (97%)	343 (91%)	35 (9%)	8	25
1	B	378/391 (97%)	341 (90%)	37 (10%)	7	22
All	All	756/782 (97%)	684 (90%)	72 (10%)	8	24

All (72) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	ARG
1	A	24	ARG
1	A	32	GLN
1	A	36	SER
1	A	57	MET
1	A	66	LYS
1	A	71	SER
1	A	78	THR
1	A	123	LEU
1	A	134	GLN
1	A	136	VAL
1	A	140	MET
1	A	152	ARG
1	A	163	ARG
1	A	219	LEU
1	A	220	ARG
1	A	236	SER
1	A	239	VAL
1	A	241	THR
1	A	242	LYS
1	A	243	ASN
1	A	253	THR
1	A	279	GLN
1	A	280	SER
1	A	283	PRO
1	A	288	ARG
1	A	293	VAL
1	A	299	LYS
1	A	302	ILE
1	A	342	THR
1	A	356	ARG
1	A	364	ILE
1	A	392	VAL
1	A	459	ASN
1	A	463	SER

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Mol	Chain	Res	Type
1	B	10	ARG
1	B	16	GLN
1	B	24	ARG
1	B	32	GLN
1	B	55	VAL
1	B	62	ASN
1	B	78	THR
1	B	123	LEU
1	B	131	ASN
1	B	136	VAL
1	B	140	MET
1	B	142	VAL
1	B	143	THR
1	B	150	VAL
1	B	155	ASN
1	B	194	ARG
1	B	197	VAL
1	B	199	THR
1	B	209	ASP
1	B	210	THR
1	B	212	GLN
1	B	218	THR
1	B	239	VAL
1	B	241	THR
1	B	243	ASN
1	B	253	THR
1	B	275	THR
1	B	279	GLN
1	B	282	ARG
1	B	288	ARG
1	B	301	ASP
1	B	302	ILE
1	B	356	ARG
1	B	374	ASN
1	B	445	ILE
1	B	460	VAL
1	B	463	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	GLN

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Mol	Chain	Res	Type
1	A	32	GLN
1	A	62	ASN
1	A	72	ASN
1	A	107	HIS
1	A	178	ASN
1	A	206	ASN
1	A	231	ASN
1	A	243	ASN
1	A	341	HIS
1	A	374	ASN
1	A	379	GLN
1	A	380	ASN
1	A	386	GLN
1	A	387	ASN
1	A	413	ASN
1	A	459	ASN
1	B	32	GLN
1	B	44	GLN
1	B	62	ASN
1	B	72	ASN
1	B	178	ASN
1	B	231	ASN
1	B	279	GLN
1	B	332	HIS
1	B	341	HIS
1	B	374	ASN
1	B	379	GLN
1	B	387	ASN
1	B	388	ASN
1	B	413	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides 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	449/470 (95%)	-0.11	9 (2%) 65 57	15, 35, 64, 85	0
1	B	449/470 (95%)	-0.29	2 (0%) 88 86	11, 29, 61, 84	0
All	All	898/940 (95%)	-0.20	11 (1%) 76 70	11, 33, 64, 85	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	468	ALA	3.7
1	A	253	THR	3.2
1	B	468	ALA	3.0
1	B	467	ALA	3.0
1	A	2	GLU	2.8
1	A	327	ASN	2.8
1	A	213	SER	2.4
1	A	255	LEU	2.3
1	A	209	ASP	2.3
1	A	252	GLU	2.2
1	A	324	TRP	2.1

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 oligosaccharides 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.