



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:31 PM GMT

PDB ID : 1PFG
Title : Strategy to design inhibitors: Structure of a complex of Proteinase K with a designed octapeptide inhibitor N-Ac-Pro-Ala-Pro-Phe-DAla-Ala-Ala-Ala-NH₂ at 2.5Å resolution
Authors : Saxena, A.K.; Singh, T.P.; Peters, K.; Fittkau, S.; Betzel, C.
Deposited on : 2003-05-27
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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

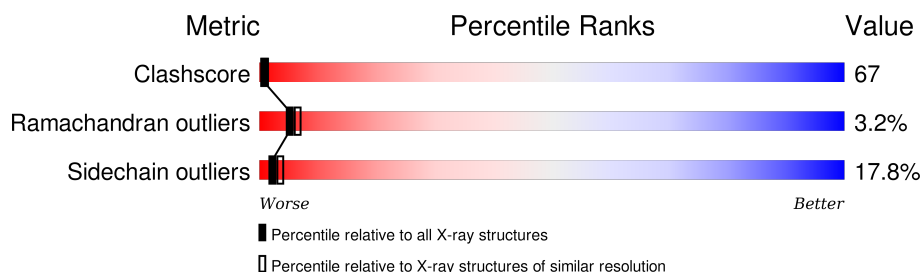
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

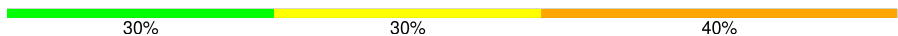
Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	279	 35% 47% 15% •
2	B	10	 30% 30% 40%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2277 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 Proteinase K.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	279	Total	C	N	O	S	0	0	0
			2017	1242	352	413	10			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	85	VAL	ALA	CONFLICT	UNP P06873

- Molecule 2 is a protein called N-Ac-PAPFAAAA-NH₂.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	10	Total	C	N	O	0	0	1
			54	36	9	9			

- Molecule 3 is water.

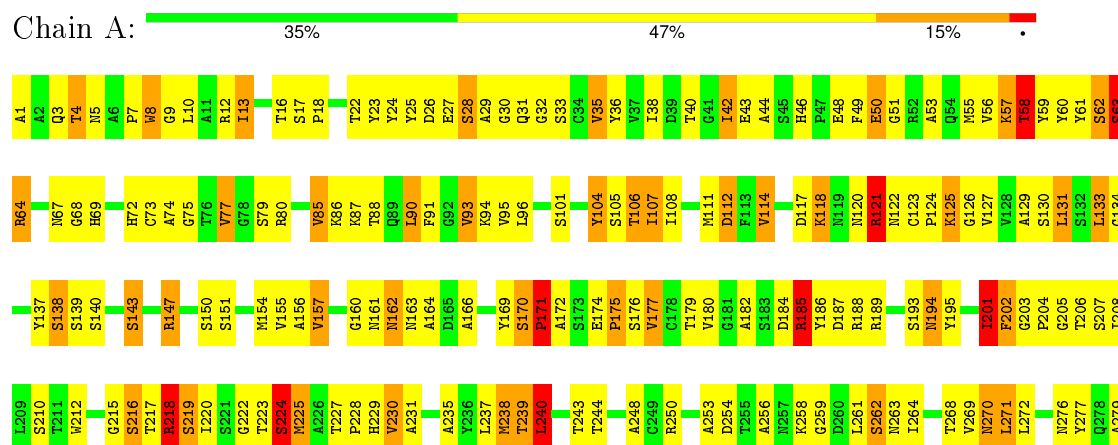
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	196	Total	O	0	0
			196	196		
3	B	10	Total	O	0	0
			10	10		

3 Residue-property plots

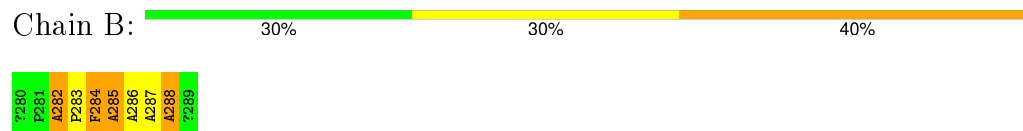
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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 ($RSRZ > 2$). 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.

Note EDS was not executed.

• Molecule 1: Proteinase K



• Molecule 2: N-Ac-PAPFAAAA-NH2



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	68.00 Å 68.00 Å 107.70 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.50	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.50)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.167 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2277	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ACE, NH2

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	1.26	6/2056 (0.3%)	1.87	46/2796 (1.6%)
2	B	2.71	1/52 (1.9%)	2.48	4/71 (5.6%)
All	All	1.32	7/2108 (0.3%)	1.89	50/2867 (1.7%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	285	ALA	N-CA	-17.11	1.12	1.46
1	A	238	MET	CG-SD	-6.13	1.65	1.81
1	A	125	LYS	CB-CG	-5.83	1.36	1.52
1	A	123	CYS	CB-SG	-5.49	1.72	1.81
1	A	202	PHE	CB-CG	-5.36	1.42	1.51
1	A	35	VAL	CA-CB	-5.13	1.44	1.54
1	A	268	THR	C-N	-5.05	1.22	1.34

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	80	ARG	NE-CZ-NH1	-19.01	110.80	120.30
1	A	121	ARG	NE-CZ-NH1	17.91	129.25	120.30
1	A	171	PRO	CA-N-CD	-17.36	87.19	111.50
1	A	170	SER	C-N-CD	-16.43	84.45	120.60
1	A	121	ARG	NE-CZ-NH2	-16.28	112.16	120.30
1	A	80	ARG	NE-CZ-NH2	12.65	126.63	120.30
1	A	171	PRO	N-CA-CB	12.35	118.12	103.30
1	A	171	PRO	N-CD-CG	11.99	121.19	103.20
1	A	170	SER	C-N-CA	11.65	170.94	122.00
2	B	285	ALA	N-CA-CB	-10.15	95.89	110.10
1	A	185	ARG	NE-CZ-NH2	10.00	125.30	120.30
1	A	147	ARG	NE-CZ-NH2	9.65	125.12	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	185	ARG	NE-CZ-NH1	-9.15	115.72	120.30
1	A	147	ARG	NE-CZ-NH1	-8.66	115.97	120.30
1	A	133	LEU	CA-CB-CG	8.54	134.95	115.30
1	A	112	ASP	CB-CG-OD2	8.30	125.77	118.30
1	A	90	LEU	CB-CG-CD2	8.15	124.86	111.00
1	A	188	ARG	NE-CZ-NH1	-7.95	116.32	120.30
1	A	175	PRO	N-CD-CG	-7.83	91.46	103.20
1	A	125	LYS	CB-CA-C	-7.38	95.64	110.40
1	A	240	LEU	CA-CB-CG	7.37	132.25	115.30
1	A	58	THR	CB-CA-C	-7.26	91.99	111.60
1	A	117	ASP	CB-CG-OD2	7.26	124.84	118.30
1	A	174	GLU	N-CA-C	-7.16	91.66	111.00
2	B	284	PHE	CB-CG-CD2	7.06	125.74	120.80
1	A	224	SER	N-CA-CB	6.72	120.58	110.50
1	A	90	LEU	CB-CG-CD1	-6.66	99.67	111.00
1	A	63	SER	CB-CA-C	6.63	122.70	110.10
1	A	218	ARG	NE-CZ-NH1	6.59	123.60	120.30
1	A	106	THR	N-CA-CB	6.59	122.82	110.30
1	A	85	VAL	CB-CA-C	-6.57	98.91	111.40
1	A	157	VAL	N-CA-CB	-6.46	97.29	111.50
1	A	224	SER	CB-CA-C	-6.38	97.98	110.10
2	B	288	ALA	N-CA-C	6.37	128.21	111.00
1	A	93	VAL	N-CA-C	-6.32	93.95	111.00
1	A	80	ARG	CB-CA-C	-6.08	98.24	110.40
1	A	133	LEU	CB-CG-CD2	6.08	121.33	111.00
1	A	230	VAL	CB-CA-C	-6.02	99.96	111.40
1	A	117	ASP	CB-CA-C	-6.01	98.38	110.40
1	A	254	ASP	CB-CG-OD2	5.91	123.62	118.30
1	A	219	SER	CB-CA-C	-5.85	98.98	110.10
1	A	240	LEU	CB-CG-CD1	5.77	120.82	111.00
1	A	121	ARG	CD-NE-CZ	5.60	131.44	123.60
1	A	63	SER	N-CA-CB	-5.58	102.14	110.50
1	A	218	ARG	NE-CZ-NH2	-5.43	117.59	120.30
1	A	131	LEU	CB-CG-CD1	-5.37	101.86	111.00
1	A	271	LEU	CB-CG-CD1	-5.34	101.92	111.00
2	B	284	PHE	CB-CG-CD1	-5.34	117.06	120.80
1	A	80	ARG	CD-NE-CZ	-5.24	116.26	123.60
1	A	129	ALA	CB-CA-C	5.18	117.88	110.10

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	2017	0	1916	261	3
2	B	54	0	50	31	0
3	A	196	0	0	35	5
3	B	10	0	0	1	0
All	All	2277	0	1966	271	5

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 67.

All (271) 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:224:SER:CB	1:A:225:MET:HE3	1.61	1.30
1:A:126:GLY:HA3	1:A:238:MET:CE	1.60	1.28
1:A:85:VAL:CG1	1:A:85:VAL:O	1.73	1.23
1:A:220:ILE:HD12	2:B:286:ALA:HB2	1.23	1.18
1:A:269:VAL:CG1	1:A:271:LEU:HD12	1.75	1.16
1:A:224:SER:HB3	1:A:225:MET:CE	1.76	1.15
1:A:85:VAL:HG11	1:A:88:THR:HB	1.29	1.13
1:A:126:GLY:HA3	1:A:238:MET:HE3	1.19	1.12
1:A:224:SER:HB2	1:A:225:MET:HE3	1.27	1.09
1:A:269:VAL:HG11	1:A:271:LEU:HD12	1.37	1.04
1:A:164:ALA:H	1:A:194:ASN:ND2	1.54	1.03
1:A:13:ILE:CD1	1:A:204:PRO:HG2	1.88	1.02
1:A:230:VAL:HB	3:A:359:HOH:O	1.58	1.02
1:A:224:SER:CB	1:A:225:MET:CE	2.34	1.02
1:A:137:TYR:HA	1:A:170:SER:OG	1.59	1.02
2:B:283:PRO:O	2:B:284:PHE:HD1	1.43	1.01
1:A:9:GLY:O	1:A:13:ILE:HD12	1.61	0.99
1:A:180:VAL:HG21	1:A:230:VAL:HG21	1.44	0.99
1:A:220:ILE:HD12	2:B:286:ALA:CB	1.92	0.98
1:A:13:ILE:HD11	1:A:204:PRO:HG2	1.48	0.95
1:A:176:SER:HA	3:A:415:HOH:O	1.66	0.95
1:A:126:GLY:CA	1:A:238:MET:CE	2.44	0.95
1:A:125:LYS:HG3	1:A:239:THR:HB	1.50	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:85:VAL:HG11	1:A:88:THR:CB	1.96	0.94
1:A:29:ALA:HB3	1:A:87:LYS:HD2	1.50	0.94
1:A:224:SER:HB3	1:A:225:MET:HE3	1.41	0.94
1:A:16:THR:HG22	3:A:421:HOH:O	1.70	0.92
1:A:261:LEU:H	1:A:270:ASN:HD21	1.18	0.91
1:A:269:VAL:HG13	1:A:271:LEU:HD12	1.48	0.91
1:A:164:ALA:H	1:A:194:ASN:HD22	1.20	0.89
1:A:220:ILE:CD1	2:B:286:ALA:HB2	2.02	0.88
1:A:13:ILE:HD11	1:A:204:PRO:CG	2.06	0.85
1:A:225:MET:O	1:A:229:HIS:HD2	1.59	0.85
1:A:207:SER:HB2	3:A:376:HOH:O	1.77	0.84
2:B:283:PRO:C	2:B:284:PHE:HD1	1.80	0.84
1:A:85:VAL:CG1	1:A:88:THR:HB	2.08	0.83
1:A:175:PRO:HG2	1:A:176:SER:N	1.92	0.83
1:A:207:SER:HB3	3:A:405:HOH:O	1.76	0.83
2:B:283:PRO:O	2:B:284:PHE:CD1	2.32	0.83
1:A:85:VAL:HG12	1:A:85:VAL:O	0.96	0.82
1:A:13:ILE:HD11	1:A:204:PRO:CD	2.09	0.82
1:A:163:ASN:HA	1:A:193:SER:O	1.79	0.82
3:A:601:HOH:O	2:B:282:ALA:HB2	1.79	0.81
1:A:220:ILE:CD1	2:B:286:ALA:CB	2.58	0.81
2:B:283:PRO:C	2:B:284:PHE:CD1	2.56	0.79
1:A:73:CYS:SG	3:A:516:HOH:O	2.41	0.78
1:A:127:VAL:N	1:A:238:MET:HE1	1.98	0.78
1:A:13:ILE:HD11	1:A:204:PRO:HD2	1.65	0.78
1:A:95:VAL:HB	1:A:107:ILE:HG22	1.64	0.78
1:A:32:GLY:HA3	1:A:125:LYS:HG2	1.66	0.78
1:A:9:GLY:O	1:A:13:ILE:CD1	2.32	0.77
1:A:85:VAL:HG11	1:A:88:THR:CG2	2.14	0.77
1:A:13:ILE:HD12	1:A:204:PRO:HG2	1.67	0.77
1:A:225:MET:O	1:A:229:HIS:CD2	2.38	0.76
1:A:177:VAL:HG23	3:A:513:HOH:O	1.87	0.74
1:A:67:ASN:HD21	2:B:288:ALA:HB3	1.52	0.74
1:A:161:ASN:HD22	2:B:285:ALA:HB2	1.53	0.73
1:A:262:SER:HB3	3:A:436:HOH:O	1.89	0.72
1:A:175:PRO:HG2	1:A:176:SER:H	1.51	0.72
1:A:13:ILE:CD1	1:A:204:PRO:CG	2.61	0.72
1:A:107:ILE:C	1:A:107:ILE:HD12	2.10	0.72
1:A:147:ARG:HG3	3:A:423:HOH:O	1.88	0.72
1:A:95:VAL:HG23	1:A:96:LEU:HG	1.72	0.71
1:A:46:HIS:CD2	1:A:48:GLU:HB2	2.26	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:GLY:CA	1:A:238:MET:HE1	2.22	0.69
1:A:269:VAL:HG13	1:A:271:LEU:CD1	2.20	0.69
1:A:107:ILE:HD13	1:A:111:MET:CE	2.23	0.69
1:A:126:GLY:HA3	1:A:238:MET:HE1	1.70	0.69
1:A:73:CYS:O	1:A:77:VAL:HG12	1.92	0.68
1:A:224:SER:HB3	1:A:225:MET:HE2	1.73	0.68
2:B:282:ALA:HB3	2:B:283:PRO:HD3	1.75	0.68
1:A:107:ILE:HG12	3:A:531:HOH:O	1.92	0.67
1:A:134:GLY:C	2:B:284:PHE:CE2	2.68	0.67
1:A:262:SER:O	1:A:263:ASN:CB	2.43	0.67
1:A:180:VAL:CG2	1:A:230:VAL:HG21	2.24	0.66
1:A:227:THR:O	1:A:230:VAL:HG22	1.95	0.66
1:A:69:HIS:CD2	3:A:516:HOH:O	2.49	0.66
1:A:225:MET:CE	3:A:517:HOH:O	2.43	0.66
1:A:169:TYR:HB3	3:A:532:HOH:O	1.95	0.65
1:A:276:ASN:CG	3:A:420:HOH:O	2.34	0.65
1:A:155:VAL:HG12	1:A:157:VAL:HG23	1.79	0.65
1:A:30:GLY:C	1:A:239:THR:HG21	2.18	0.64
1:A:157:VAL:HG21	1:A:177:VAL:HG21	1.79	0.64
1:A:216:SER:O	1:A:217:THR:HG22	1.98	0.64
1:A:243:THR:HG21	1:A:248:ALA:HA	1.79	0.64
1:A:25:TYR:HD1	1:A:26:ASP:O	1.80	0.64
1:A:32:GLY:CA	1:A:125:LYS:HG2	2.28	0.63
1:A:180:VAL:HG21	1:A:230:VAL:CG2	2.23	0.63
1:A:164:ALA:N	1:A:194:ASN:HD22	1.95	0.63
1:A:49:PHE:O	1:A:51:GLY:N	2.31	0.63
1:A:107:ILE:HD13	1:A:111:MET:HE2	1.80	0.63
1:A:155:VAL:HG12	1:A:157:VAL:CG2	2.30	0.62
1:A:224:SER:HB2	1:A:225:MET:CE	2.14	0.62
1:A:224:SER:N	2:B:284:PHE:O	2.33	0.62
1:A:201:ILE:HD11	1:A:256:ALA:HB2	1.81	0.62
1:A:67:ASN:ND2	2:B:288:ALA:HB3	2.14	0.62
1:A:126:GLY:C	1:A:127:VAL:HG23	2.18	0.62
1:A:127:VAL:H	1:A:238:MET:HE1	1.64	0.62
1:A:137:TYR:CA	1:A:170:SER:OG	2.43	0.62
1:A:227:THR:O	1:A:230:VAL:CG2	2.47	0.61
1:A:107:ILE:CD1	1:A:111:MET:HE3	2.31	0.61
1:A:57:LYS:HD2	1:A:59:TYR:CE2	2.36	0.60
1:A:277:TYR:CZ	1:A:279:ALA:HA	2.36	0.60
1:A:29:ALA:HB3	1:A:87:LYS:CD	2.29	0.60
1:A:60:TYR:CD2	1:A:60:TYR:N	2.69	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:ARG:HB2	3:A:400:HOH:O	2.01	0.60
1:A:220:ILE:CD1	2:B:286:ALA:HB1	2.32	0.58
1:A:208:ILE:HD13	3:A:320:HOH:O	2.04	0.58
1:A:261:LEU:HD21	1:A:272:LEU:HD13	1.86	0.57
1:A:227:THR:HB	1:A:228:PRO:CD	2.35	0.57
1:A:74:ALA:O	1:A:77:VAL:HG13	2.03	0.57
1:A:107:ILE:CD1	1:A:111:MET:CE	2.82	0.57
1:A:35:VAL:HG11	1:A:77:VAL:HG21	1.86	0.57
1:A:224:SER:CA	2:B:284:PHE:O	2.52	0.57
1:A:46:HIS:HD2	1:A:48:GLU:H	1.53	0.57
1:A:163:ASN:CG	1:A:163:ASN:O	2.43	0.56
1:A:58:THR:HG21	1:A:62:SER:O	2.06	0.56
1:A:68:GLY:HA3	1:A:212:TRP:CZ2	2.41	0.56
1:A:67:ASN:O	2:B:288:ALA:HB1	2.06	0.56
1:A:161:ASN:ND2	2:B:285:ALA:HB2	2.20	0.55
1:A:4:THR:HG22	1:A:5:ASN:OD1	2.07	0.55
1:A:67:ASN:ND2	2:B:288:ALA:CB	2.70	0.55
1:A:118:LYS:NZ	1:A:151:SER:O	2.40	0.55
1:A:126:GLY:C	1:A:238:MET:HE1	2.27	0.55
1:A:58:THR:HG22	1:A:59:TYR:H	1.72	0.55
1:A:58:THR:HG23	1:A:94:LYS:HB3	1.89	0.54
1:A:230:VAL:HG23	1:A:231:ALA:N	2.22	0.54
1:A:1:ALA:O	1:A:24:TYR:HA	2.07	0.54
1:A:186:TYR:O	1:A:187:ASP:HB2	2.06	0.54
1:A:225:MET:HE2	3:A:517:HOH:O	2.07	0.54
1:A:1:ALA:N	1:A:27:GLU:OE1	2.33	0.53
1:A:75:GLY:HA2	1:A:79:SER:HB3	1.90	0.53
1:A:77:VAL:O	1:A:85:VAL:HB	2.08	0.53
1:A:25:TYR:CD1	1:A:26:ASP:O	2.61	0.53
1:A:139:SER:O	1:A:143:SER:HB2	2.08	0.53
1:A:57:LYS:HG3	1:A:93:VAL:HG13	1.90	0.53
1:A:56:VAL:HG21	1:A:91:PHE:CD2	2.44	0.53
1:A:49:PHE:O	1:A:50:GLU:C	2.48	0.52
1:A:228:PRO:HA	1:A:231:ALA:HB3	1.91	0.52
1:A:262:SER:O	1:A:263:ASN:HB2	2.09	0.52
1:A:4:THR:O	1:A:5:ASN:C	2.46	0.52
1:A:59:TYR:C	1:A:60:TYR:CD2	2.82	0.52
1:A:69:HIS:O	1:A:72:HIS:HB3	2.10	0.52
1:A:61:TYR:CE1	3:A:345:HOH:O	2.53	0.52
1:A:131:LEU:HD12	1:A:131:LEU:N	2.25	0.51
1:A:126:GLY:C	1:A:127:VAL:CG2	2.78	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:32:GLY:C	1:A:125:LYS:HG2	2.30	0.51
1:A:3:GLN:HE22	1:A:86:LYS:HZ1	1.57	0.51
1:A:139:SER:HB2	3:A:474:HOH:O	2.11	0.51
1:A:161:ASN:HD21	2:B:285:ALA:N	2.07	0.51
1:A:46:HIS:HE1	1:A:216:SER:O	1.94	0.50
1:A:138:SER:N	1:A:170:SER:OG	2.44	0.50
1:A:201:ILE:HD11	1:A:256:ALA:CB	2.41	0.50
1:A:269:VAL:HG13	1:A:271:LEU:CG	2.42	0.50
1:A:3:GLN:O	1:A:22:THR:HA	2.12	0.50
1:A:208:ILE:O	1:A:219:SER:HA	2.10	0.50
1:A:237:LEU:HA	1:A:240:LEU:HB2	1.93	0.50
1:A:38:ILE:HD11	1:A:114:VAL:HG11	1.93	0.50
1:A:107:ILE:HD13	1:A:111:MET:HE3	1.90	0.50
1:A:28:SER:O	1:A:31:GLN:HG2	2.12	0.50
1:A:107:ILE:HD12	1:A:108:ILE:N	2.27	0.50
1:A:42:ILE:O	1:A:44:ALA:N	2.45	0.50
1:A:13:ILE:HD12	1:A:204:PRO:CG	2.35	0.49
1:A:220:ILE:HD13	2:B:286:ALA:HB1	1.95	0.49
2:B:284:PHE:N	2:B:284:PHE:CD1	2.76	0.49
1:A:69:HIS:CG	3:A:516:HOH:O	2.65	0.49
1:A:164:ALA:N	1:A:194:ASN:ND2	2.39	0.49
1:A:185:ARG:NH1	3:A:380:HOH:O	2.44	0.49
1:A:216:SER:C	1:A:217:THR:CG2	2.82	0.49
1:A:126:GLY:C	1:A:238:MET:CE	2.81	0.49
1:A:216:SER:O	1:A:217:THR:CG2	2.61	0.48
1:A:138:SER:H	1:A:170:SER:CB	2.26	0.48
1:A:185:ARG:NH2	3:A:380:HOH:O	2.46	0.48
1:A:243:THR:CG2	1:A:248:ALA:HA	2.42	0.48
1:A:64:ARG:CD	3:A:327:HOH:O	2.62	0.48
1:A:7:PRO:O	1:A:10:LEU:HB2	2.14	0.48
1:A:182:ALA:HB1	1:A:205:GLY:HA3	1.96	0.48
1:A:224:SER:HA	2:B:284:PHE:O	2.13	0.48
1:A:134:GLY:O	2:B:284:PHE:CE2	2.67	0.47
1:A:85:VAL:CG1	1:A:88:THR:H	2.27	0.47
1:A:157:VAL:CG1	3:A:513:HOH:O	2.62	0.47
1:A:171:PRO:O	1:A:172:ALA:C	2.52	0.47
1:A:73:CYS:O	1:A:74:ALA:C	2.50	0.47
1:A:122:ASN:C	1:A:124:PRO:HD3	2.35	0.47
1:A:57:LYS:HE3	1:A:57:LYS:HB2	1.55	0.47
1:A:12:ARG:HG3	1:A:12:ARG:O	2.14	0.46
1:A:201:ILE:HG13	1:A:202:PHE:H	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:LYS:C	1:A:87:LYS:HG3	2.35	0.46
1:A:120:ASN:C	1:A:121:ARG:HG2	2.34	0.46
1:A:130:SER:HB2	1:A:231:ALA:HB1	1.96	0.46
1:A:69:HIS:CE1	3:A:516:HOH:O	2.68	0.46
2:B:282:ALA:HB3	2:B:283:PRO:CD	2.45	0.46
1:A:55:MET:CE	1:A:63:SER:O	2.64	0.46
1:A:46:HIS:CE1	1:A:216:SER:O	2.68	0.46
1:A:189:ARG:NH2	3:A:386:HOH:O	2.45	0.45
1:A:64:ARG:HD2	3:A:327:HOH:O	2.15	0.45
1:A:8:TRP:C	1:A:8:TRP:CE3	2.89	0.45
1:A:157:VAL:HG11	3:A:513:HOH:O	2.16	0.45
1:A:270:ASN:HD22	1:A:270:ASN:C	2.19	0.45
1:A:25:TYR:HE1	3:A:445:HOH:O	2.00	0.45
1:A:243:THR:HA	3:A:344:HOH:O	2.16	0.45
1:A:201:ILE:CD1	1:A:253:ALA:HA	2.47	0.45
1:A:164:ALA:O	1:A:194:ASN:HA	2.17	0.45
1:A:36:TYR:CD1	1:A:114:VAL:HB	2.52	0.45
1:A:7:PRO:HD2	1:A:10:LEU:HD12	1.98	0.45
1:A:224:SER:OG	2:B:284:PHE:C	2.56	0.44
1:A:46:HIS:HD2	1:A:48:GLU:HB2	1.80	0.44
1:A:264:ILE:HB	3:A:437:HOH:O	2.17	0.44
1:A:17:SER:O	1:A:17:SER:OG	2.31	0.44
1:A:72:HIS:CD2	1:A:210:SER:HB3	2.52	0.44
1:A:57:LYS:C	1:A:57:LYS:HD3	2.38	0.44
1:A:30:GLY:HA3	1:A:85:VAL:HG22	2.00	0.44
1:A:107:ILE:O	1:A:108:ILE:C	2.56	0.43
1:A:157:VAL:O	1:A:180:VAL:N	2.47	0.43
1:A:161:ASN:ND2	2:B:285:ALA:CB	2.81	0.43
1:A:216:SER:HB2	1:A:217:THR:H	1.59	0.43
1:A:107:ILE:CD1	1:A:107:ILE:C	2.85	0.43
1:A:218:ARG:NH2	3:A:530:HOH:O	2.50	0.43
1:A:224:SER:O	1:A:228:PRO:CD	2.66	0.43
1:A:182:ALA:HA	1:A:203:GLY:O	2.18	0.43
1:A:157:VAL:O	1:A:179:THR:HA	2.18	0.43
1:A:187:ASP:O	1:A:261:LEU:HA	2.18	0.43
1:A:225:MET:N	1:A:225:MET:CE	2.81	0.43
1:A:3:GLN:N	1:A:23:TYR:O	2.47	0.43
2:B:284:PHE:HZ	3:B:506:HOH:O	1.98	0.43
1:A:27:GLU:C	1:A:29:ALA:N	2.72	0.43
1:A:55:MET:HE1	1:A:63:SER:O	2.19	0.43
1:A:64:ARG:HG2	1:A:64:ARG:HH11	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:258:LYS:HA	1:A:270:ASN:O	2.19	0.42
1:A:31:GLN:HB3	1:A:87:LYS:HB3	2.01	0.42
1:A:235:ALA:O	1:A:239:THR:HG23	2.18	0.42
1:A:3:GLN:NE2	1:A:86:LYS:HE3	2.35	0.42
1:A:85:VAL:CG1	1:A:88:THR:HG22	2.49	0.42
1:A:49:PHE:C	1:A:51:GLY:N	2.73	0.42
1:A:125:LYS:O	1:A:238:MET:HB3	2.19	0.42
1:A:46:HIS:CE1	1:A:215:GLY:HA2	2.54	0.42
1:A:189:ARG:NH1	1:A:193:SER:O	2.52	0.42
1:A:58:THR:CG2	1:A:94:LYS:HG2	2.49	0.42
1:A:156:ALA:C	1:A:157:VAL:HG23	2.40	0.41
1:A:227:THR:O	1:A:230:VAL:HG23	2.19	0.41
1:A:202:PHE:HB2	1:A:271:LEU:O	2.20	0.41
1:A:56:VAL:CG2	1:A:91:PHE:CD2	3.03	0.41
1:A:18:PRO:HD3	1:A:187:ASP:OD1	2.21	0.41
1:A:25:TYR:CD2	1:A:25:TYR:N	2.86	0.41
1:A:189:ARG:HB3	1:A:263:ASN:HB3	2.02	0.41
1:A:161:ASN:ND2	2:B:285:ALA:N	2.68	0.41
1:A:227:THR:HB	1:A:228:PRO:HD3	2.01	0.41
1:A:8:TRP:HZ2	1:A:206:THR:O	2.03	0.41
1:A:195:TYR:C	1:A:195:TYR:CD2	2.93	0.41
1:A:224:SER:C	1:A:225:MET:CE	2.89	0.41
1:A:162:ASN:N	1:A:162:ASN:ND2	2.68	0.41
1:A:184:ASP:C	1:A:184:ASP:OD1	2.58	0.41
1:A:104:TYR:O	1:A:108:ILE:N	2.47	0.41
1:A:130:SER:HB2	1:A:231:ALA:CB	2.51	0.41
1:A:32:GLY:HA3	1:A:125:LYS:CG	2.43	0.41
1:A:261:LEU:N	1:A:270:ASN:HD21	2.00	0.41
1:A:258:LYS:HG2	1:A:271:LEU:HD23	2.02	0.41
1:A:53:ALA:HA	1:A:90:LEU:O	2.21	0.41
1:A:40:THR:O	1:A:94:LYS:HD2	2.21	0.41
1:A:30:GLY:O	1:A:239:THR:HG21	2.20	0.41
3:A:530:HOH:O	2:B:287:ALA:HB3	2.21	0.41
1:A:162:ASN:HD22	1:A:162:ASN:H	1.69	0.41
1:A:160:GLY:O	1:A:223:THR:HG21	2.20	0.41
1:A:210:SER:O	1:A:217:THR:HB	2.21	0.41
1:A:220:ILE:HD13	1:A:220:ILE:HG21	1.73	0.40
1:A:256:ALA:HB1	1:A:272:LEU:O	2.20	0.40
1:A:25:TYR:CZ	1:A:86:LYS:HD2	2.56	0.40
1:A:58:THR:HG22	1:A:59:TYR:N	2.34	0.40
1:A:222:GLY:C	1:A:224:SER:N	2.75	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:ALA:HB1	3:A:459:HOH:O	2.21	0.40
1:A:33:SER:HG	1:A:33:SER:H	1.58	0.40

All (5) 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)
3:A:457:HOH:O	3:A:457:HOH:O[7_465]	0.69	1.51
3:A:476:HOH:O	3:A:543:HOH:O[6_455]	0.81	1.39
1:A:112:ASP:OD1	3:A:339:HOH:O[7_465]	1.77	0.43
1:A:216:SER:CB	3:A:511:HOH:O[6_555]	1.80	0.40
1:A:150:SER:OG	3:A:423:HOH:O[7_465]	1.93	0.27

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	277/279 (99%)	244 (88%)	25 (9%)	8 (3%)	6	8
2	B	6/10 (60%)	4 (67%)	1 (17%)	1 (17%)	0	0
All	All	283/289 (98%)	248 (88%)	26 (9%)	9 (3%)	5	6

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	171	PRO
1	A	50	GLU
2	B	282	ALA
1	A	42	ILE
1	A	43	GLU
1	A	166	ALA
1	A	201	ILE

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Mol	Chain	Res	Type
1	A	259	GLY
1	A	177	VAL

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	211/214 (99%)	173 (82%)	38 (18%)	2	3
2	B	3/3 (100%)	3 (100%)	0	100	100
All	All	214/217 (99%)	176 (82%)	38 (18%)	2	3

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	THR
1	A	8	TRP
1	A	13	ILE
1	A	28	SER
1	A	57	LYS
1	A	58	THR
1	A	62	SER
1	A	63	SER
1	A	64	ARG
1	A	77	VAL
1	A	101	SER
1	A	104	TYR
1	A	105	SER
1	A	106	THR
1	A	107	ILE
1	A	114	VAL
1	A	118	LYS
1	A	121	ARG
1	A	133	LEU
1	A	138	SER
1	A	140	SER

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Mol	Chain	Res	Type
1	A	143	SER
1	A	154	MET
1	A	162	ASN
1	A	171	PRO
1	A	185	ARG
1	A	194	ASN
1	A	201	ILE
1	A	216	SER
1	A	218	ARG
1	A	224	SER
1	A	225	MET
1	A	239	THR
1	A	240	LEU
1	A	244	THR
1	A	250	ARG
1	A	262	SER
1	A	270	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	46	HIS
1	A	162	ASN
1	A	194	ASN
1	A	229	HIS
1	A	270	ASN
1	A	276	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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.