



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:18 PM UTC

PDB ID : 1D5S / pdb_00001d5s
Title : CRYSTAL STRUCTURE OF CLEAVED ANTITRYPSIN POLYMER
Authors : Dunstone, M.A.; Dai, W.; Whisstock, J.C.; Rossjohn, J.; Pike, R.N.; Feil, S.C.; Le Bonneic, B.F.; Parker, M.W.; Bottomley, S.P.
Deposited on : 1999-10-11
Resolution : 3.00 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

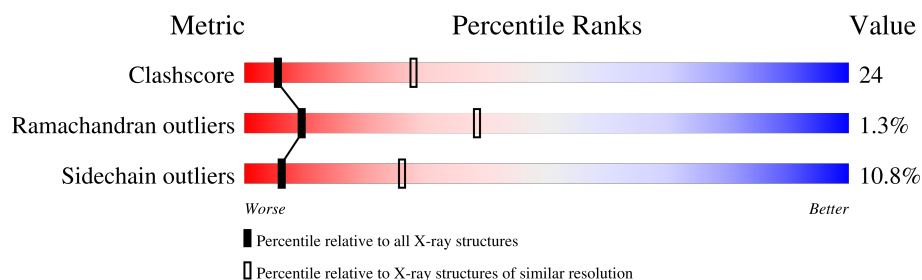
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)

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 $\leq 5\%$

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	334	
2	B	41	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2984 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 P1-ARG ANTITRYPSIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	334	Total	C	N	O	S	0	0	0
			2653	1702	436	508	7			

- Molecule 2 is a protein called P1-ARG ANTITRYPSIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	41	Total	C	N	O	S	0	0	0
			331	218	54	57	2			

There is a discrepancy between the modelled and reference sequences:

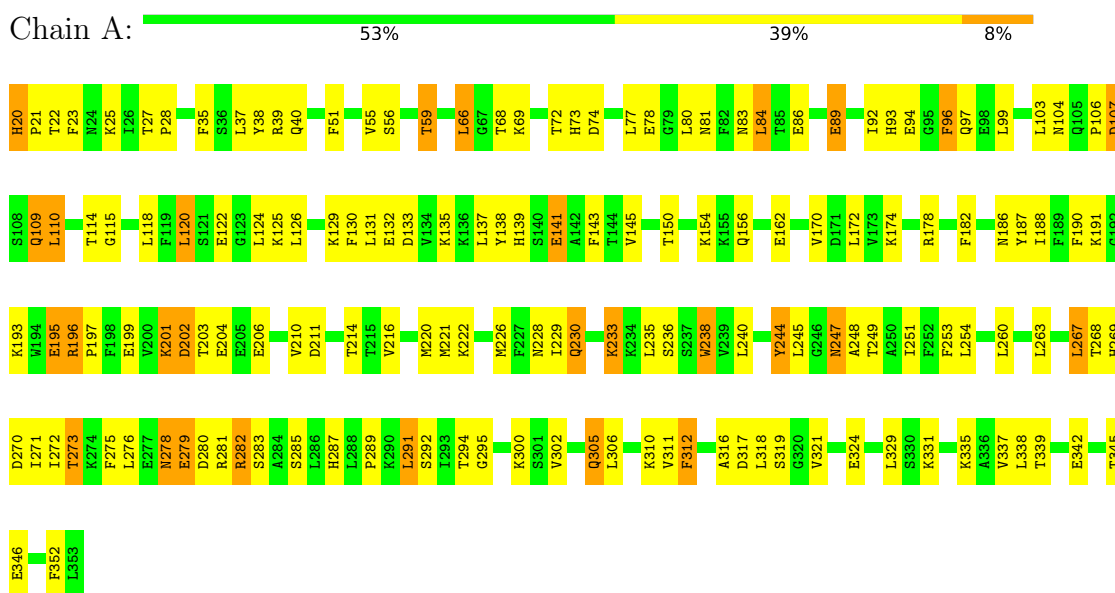
Chain	Residue	Modelled	Actual	Comment	Reference
B	358	ARG	MET	engineered mutation	UNP P01009

3 Residue-property plots [i](#)

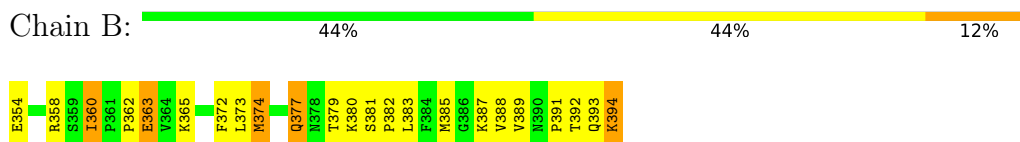
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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: P1-ARG ANTITRYPSIN



• Molecule 2: P1-ARG ANTITRYPSIN



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	110.17Å 110.17Å 76.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.00	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-3.00)	Depositor
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.208 , 0.262	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2984	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

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	0.53	0/2706	0.98	8/3655 (0.2%)
2	B	0.62	0/340	1.21	4/458 (0.9%)
All	All	0.54	0/3046	1.01	12/4113 (0.3%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	393	GLN	N-CA-C	-8.79	96.32	110.20
1	A	195	GLU	N-CA-C	-8.77	101.67	111.14
1	A	244	TYR	N-CA-C	-6.89	100.25	110.46
1	A	122	GLU	N-CA-C	6.63	118.17	111.07
2	B	377	GLN	N-CA-C	5.79	118.06	111.11
1	A	238	TRP	N-CA-C	-5.69	99.46	108.73
2	B	381	SER	CA-C-N	5.66	126.92	119.84
2	B	381	SER	C-N-CA	5.66	126.92	119.84
1	A	20	HIS	CA-C-N	5.54	125.47	119.76
1	A	20	HIS	C-N-CA	5.54	125.47	119.76
1	A	312	PHE	N-CA-C	-5.16	107.50	112.97
1	A	110	LEU	N-CA-C	5.06	117.03	108.02

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	2653	0	2651	133	1
2	B	331	0	348	27	1
All	All	2984	0	2999	142	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 24.

All (142) 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:B:394:LYS:H	2:B:394:LYS:HE3	1.14	1.07
1:A:201:LYS:H	1:A:201:LYS:HD2	1.27	0.97
1:A:162:GLU:HG3	1:A:170:VAL:HG12	1.48	0.95
2:B:394:LYS:HE3	2:B:394:LYS:N	1.83	0.91
1:A:104:ASN:O	1:A:106:PRO:HD3	1.76	0.86
1:A:99:LEU:HD12	2:B:379:THR:HG21	1.58	0.84
1:A:55:VAL:O	1:A:59:THR:HG23	1.78	0.81
1:A:294:THR:HG22	1:A:337:VAL:HG13	1.62	0.81
1:A:195:GLU:O	1:A:197:PRO:HD3	1.79	0.81
2:B:360:ILE:H	2:B:360:ILE:HD13	1.43	0.81
1:A:114:THR:HG22	1:A:188:ILE:HG13	1.66	0.78
1:A:68:THR:HG22	1:A:318:LEU:HD23	1.66	0.78
1:A:125:LYS:O	1:A:125:LYS:HG3	1.88	0.74
1:A:66:LEU:HD13	1:A:93:HIS:CE1	2.23	0.73
1:A:72:THR:HG22	1:A:310:LYS:HG2	1.69	0.73
1:A:251:ILE:HD11	1:A:276:LEU:HD11	1.71	0.72
1:A:69:LYS:HD2	1:A:319:SER:HB2	1.70	0.72
1:A:229:ILE:O	1:A:230:GLN:HB3	1.90	0.70
1:A:226:MET:HG3	1:A:281:ARG:HB3	1.72	0.70
1:A:83:ASN:HD22	1:A:86:GLU:HG3	1.59	0.69
1:A:20:HIS:N	1:A:21:PRO:HD2	2.08	0.68
1:A:228:ASN:ND2	1:A:279:GLU:HA	2.09	0.67
1:A:141:GLU:HG2	1:A:143:PHE:CZ	2.30	0.67
1:A:107:ASP:HB2	1:A:247:ASN:ND2	2.11	0.66
1:A:104:ASN:C	1:A:106:PRO:HD3	2.21	0.66
1:A:93:HIS:HB3	1:A:137:LEU:HD23	1.78	0.66
1:A:253:PHE:HB3	1:A:263:LEU:HD21	1.78	0.66
1:A:23:PHE:CD1	2:B:379:THR:HG22	2.33	0.64
1:A:55:VAL:HG23	2:B:382:PRO:O	1.98	0.63
1:A:83:ASN:HD22	1:A:86:GLU:CG	2.10	0.63
1:A:72:THR:HG21	1:A:316:ALA:HA	1.81	0.62
1:A:72:THR:HG21	1:A:317:ASP:H	1.63	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:300:LYS:HE3	1:A:312:PHE:HB3	1.82	0.61
1:A:22:THR:HG21	1:A:94:GLU:OE2	2.01	0.61
1:A:206:GLU:OE2	1:A:222:LYS:HE2	1.99	0.61
1:A:251:ILE:HD12	2:B:373:LEU:HD11	1.81	0.61
1:A:260:LEU:HD21	2:B:387:LYS:HE3	1.82	0.60
1:A:201:LYS:HD2	1:A:201:LYS:N	2.08	0.60
1:A:263:LEU:O	1:A:267:LEU:HD12	2.01	0.60
1:A:59:THR:HG22	1:A:96:PHE:HE1	1.66	0.60
1:A:110:LEU:HD12	1:A:191:LYS:O	2.01	0.59
1:A:201:LYS:H	1:A:201:LYS:CD	2.07	0.59
1:A:216:VAL:HG11	2:B:392:THR:HG23	1.85	0.58
1:A:278:ASN:ND2	1:A:279:GLU:N	2.51	0.58
1:A:295:GLY:O	1:A:335:LYS:HA	2.03	0.58
1:A:199:GLU:HB2	1:A:202:ASP:OD1	2.03	0.58
1:A:72:THR:HG22	1:A:310:LYS:CG	2.34	0.57
2:B:360:ILE:HD13	2:B:360:ILE:N	2.16	0.57
1:A:272:ILE:O	1:A:276:LEU:HD13	2.05	0.57
1:A:78:GLU:HG3	1:A:84:LEU:HD12	1.87	0.56
1:A:107:ASP:HB2	1:A:247:ASN:HD21	1.71	0.56
1:A:226:MET:HG2	1:A:281:ARG:HD3	1.86	0.56
1:A:40:GLN:OE1	1:A:305:GLN:HG2	2.06	0.55
1:A:89:GLU:HA	1:A:92:ILE:HD12	1.88	0.55
1:A:211:ASP:HB3	1:A:214:THR:H	1.72	0.55
1:A:83:ASN:HB3	1:A:86:GLU:HG3	1.88	0.54
1:A:133:ASP:O	1:A:137:LEU:HB2	2.08	0.54
1:A:110:LEU:HD23	1:A:247:ASN:OD1	2.07	0.54
1:A:268:THR:HB	1:A:271:ILE:HG13	1.89	0.54
1:A:291:LEU:HD22	2:B:391:PRO:HB3	1.90	0.54
1:A:131:LEU:O	1:A:135:LYS:HB2	2.07	0.54
1:A:51:PHE:CZ	1:A:338:LEU:HB2	2.43	0.53
1:A:193:LYS:O	1:A:244:TYR:HA	2.09	0.53
1:A:202:ASP:O	1:A:203:THR:C	2.50	0.53
1:A:291:LEU:CD1	2:B:388:VAL:HG21	2.38	0.53
1:A:124:LEU:HD23	1:A:126:LEU:HD21	1.90	0.52
1:A:35:PHE:O	1:A:39:ARG:HG3	2.09	0.52
1:A:40:GLN:HB3	1:A:302:VAL:HG13	1.91	0.52
1:A:226:MET:HB3	1:A:281:ARG:HD3	1.92	0.52
1:A:202:ASP:OD1	1:A:202:ASP:N	2.43	0.51
1:A:226:MET:CG	1:A:281:ARG:HB3	2.39	0.51
1:A:278:ASN:ND2	1:A:279:GLU:H	2.09	0.51
1:A:228:ASN:HD21	1:A:279:GLU:HA	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:PHE:C	1:A:25:LYS:H	2.19	0.51
1:A:268:THR:HG22	1:A:269:HIS:N	2.25	0.51
2:B:373:LEU:HD23	2:B:385:MET:HE3	1.92	0.51
1:A:244:TYR:HE2	1:A:345:THR:HB	1.76	0.51
1:A:120:LEU:HD22	1:A:126:LEU:HD11	1.91	0.51
1:A:99:LEU:O	1:A:103:LEU:HG	2.11	0.51
1:A:72:THR:HG21	1:A:317:ASP:N	2.26	0.50
1:A:115:GLY:HA3	1:A:187:TYR:CE2	2.47	0.50
1:A:253:PHE:CB	1:A:263:LEU:HD21	2.39	0.50
1:A:238:TRP:HB2	1:A:254:LEU:HB3	1.93	0.50
1:A:270:ASP:HA	1:A:273:THR:OG1	2.12	0.50
1:A:110:LEU:HD11	1:A:190:PHE:CE1	2.47	0.50
1:A:120:LEU:HD13	1:A:126:LEU:HD11	1.95	0.49
1:A:235:LEU:O	1:A:236:SER:HB2	2.10	0.49
1:A:77:LEU:HB3	1:A:84:LEU:HD21	1.93	0.49
1:A:154:LYS:HD2	1:A:174:LYS:O	2.12	0.49
1:A:118:LEU:HD13	1:A:120:LEU:HD21	1.95	0.49
1:A:78:GLU:HG3	1:A:84:LEU:CD1	2.42	0.49
1:A:216:VAL:HG11	2:B:392:THR:CG2	2.43	0.49
1:A:204:GLU:HB3	1:A:222:LYS:HE3	1.95	0.49
1:A:283:SER:O	2:B:362:PRO:HD2	2.12	0.48
1:A:247:ASN:HA	2:B:377:GLN:NE2	2.28	0.48
1:A:28:PRO:HB3	1:A:269:HIS:CD2	2.49	0.48
1:A:182:PHE:CE1	1:A:321:VAL:HG11	2.49	0.48
1:A:278:ASN:ND2	1:A:280:ASP:H	2.11	0.48
1:A:35:PHE:HD1	2:B:385:MET:HE1	1.79	0.47
1:A:35:PHE:CD1	2:B:385:MET:HE1	2.48	0.47
1:A:68:THR:HG22	1:A:318:LEU:CD2	2.42	0.47
1:A:80:LEU:O	1:A:81:ASN:HB2	2.13	0.47
1:A:210:VAL:HG13	2:B:389:VAL:O	2.15	0.47
1:A:279:GLU:O	1:A:280:ASP:C	2.57	0.47
1:A:124:LEU:CD2	1:A:126:LEU:HD21	2.44	0.46
1:A:226:MET:CG	1:A:281:ARG:HD3	2.45	0.46
1:A:268:THR:O	1:A:272:ILE:HD13	2.16	0.46
1:A:114:THR:HG22	1:A:188:ILE:CG1	2.42	0.46
1:A:138:TYR:O	1:A:139:HIS:C	2.59	0.46
1:A:195:GLU:C	1:A:197:PRO:HD3	2.41	0.45
1:A:118:LEU:HD23	1:A:118:LEU:HA	1.77	0.45
1:A:38:TYR:OH	2:B:387:LYS:HD2	2.17	0.45
2:B:379:THR:O	2:B:380:LYS:HB2	2.17	0.45
2:B:372:PHE:HE2	2:B:374:MET:HE1	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:LYS:O	1:A:130:PHE:C	2.59	0.44
1:A:22:THR:O	1:A:25:LYS:HB3	2.18	0.44
1:A:73:HIS:HD2	1:A:74:ASP:OD2	2.01	0.44
1:A:289:PRO:HG2	1:A:291:LEU:HD21	2.00	0.44
1:A:99:LEU:HD12	2:B:379:THR:CG2	2.40	0.44
1:A:285:SER:HB3	2:B:363:GLU:CG	2.47	0.44
1:A:244:TYR:CE2	1:A:345:THR:HB	2.53	0.43
1:A:23:PHE:C	1:A:25:LYS:N	2.76	0.43
1:A:271:ILE:HG22	1:A:275:PHE:HE2	1.83	0.43
1:A:196:ARG:HA	1:A:196:ARG:HD2	1.35	0.43
1:A:221:MET:O	1:A:287:HIS:HA	2.19	0.43
1:A:244:TYR:HD1	1:A:248:ALA:O	2.01	0.43
2:B:363:GLU:OE2	2:B:365:LYS:HE3	2.19	0.42
1:A:172:LEU:HD21	1:A:352:PHE:HB2	2.01	0.42
1:A:206:GLU:HB2	1:A:220:MET:HB2	2.01	0.42
1:A:282:ARG:HB2	1:A:282:ARG:HH11	1.84	0.42
1:A:346:GLU:O	1:A:346:GLU:HG3	2.19	0.42
1:A:233:LYS:O	1:A:233:LYS:HG3	2.12	0.42
2:B:360:ILE:H	2:B:360:ILE:CD1	2.14	0.42
1:A:27:THR:N	1:A:28:PRO:HD2	2.34	0.42
1:A:106:PRO:HB2	1:A:110:LEU:O	2.20	0.42
1:A:56:SER:OG	1:A:186:ASN:ND2	2.53	0.41
1:A:294:THR:CG2	1:A:337:VAL:HG13	2.42	0.41
1:A:37:LEU:HD12	1:A:306:LEU:HD11	2.01	0.41
1:A:74:ASP:O	1:A:78:GLU:HB2	2.21	0.41
1:A:109:GLN:HB3	1:A:245:LEU:HB2	2.03	0.41
1:A:338:LEU:HD22	2:B:372:PHE:HZ	1.86	0.41
1:A:129:LYS:O	1:A:132:GLU:HG3	2.21	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:A:331:LYS:NZ	2:B:354:GLU:OE2[5_555]	1.99	0.21

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	A	332/334 (99%)	292 (88%)	35 (10%)	5 (2%)	8	35
2	B	39/41 (95%)	38 (97%)	1 (3%)	0	100	100
All	All	371/375 (99%)	330 (89%)	36 (10%)	5 (1%)	9	38

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	84	LEU
1	A	230	GLN
1	A	89	GLU
1	A	107	ASP
1	A	150	THR

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	293/293 (100%)	263 (90%)	30 (10%)	7	29
2	B	39/39 (100%)	33 (85%)	6 (15%)	2	13
All	All	332/332 (100%)	296 (89%)	36 (11%)	6	26

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	THR

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Mol	Chain	Res	Type
1	A	66	LEU
1	A	96	PHE
1	A	97	GLN
1	A	109	GLN
1	A	120	LEU
1	A	141	GLU
1	A	145	VAL
1	A	156	GLN
1	A	178	ARG
1	A	196	ARG
1	A	201	LYS
1	A	202	ASP
1	A	233	LYS
1	A	240	LEU
1	A	247	ASN
1	A	249	THR
1	A	267	LEU
1	A	273	THR
1	A	278	ASN
1	A	279	GLU
1	A	282	ARG
1	A	291	LEU
1	A	292	SER
1	A	305	GLN
1	A	311	VAL
1	A	324	GLU
1	A	329	LEU
1	A	339	THR
1	A	342	GLU
2	B	358	ARG
2	B	360	ILE
2	B	363	GLU
2	B	374	MET
2	B	383	LEU
2	B	394	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	ASN
1	A	73	HIS
1	A	83	ASN

Continued on next page...

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Mol	Chain	Res	Type
1	A	97	GLN
1	A	139	HIS
1	A	146	ASN
1	A	186	ASN
1	A	212	GLN
1	A	265	ASN
1	A	278	ASN
2	B	377	GLN
2	B	378	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 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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.