



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 11:54 am GMT

PDB ID : 1ZYM
Title : AMINO TERMINAL DOMAIN OF ENZYME I FROM ESCHERICHIA COLI
Authors : Liao, D.-I.; Davies, D.R.
Deposited on : 1996-05-21
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
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

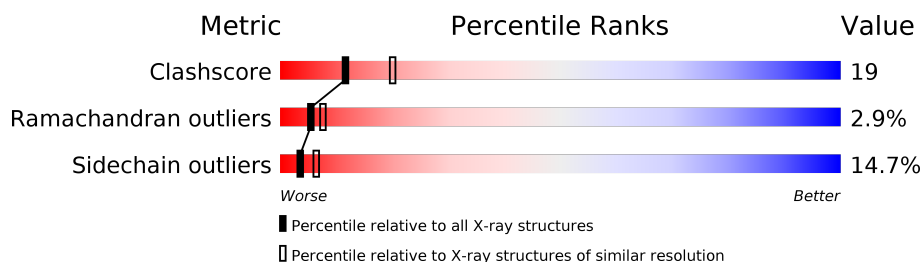
1 Overall quality at a glance

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	112137	4554 (2.50-2.50)
Ramachandran outliers	110173	4463 (2.50-2.50)
Sidechain outliers	110143	4465 (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	258	
1	B	258	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3854 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 ENZYME I.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	247	Total	C	N	O	S	0	0	0
			1891	1181	320	386	4			
1	B	248	Total	C	N	O	S	0	0	0
			1899	1187	321	387	4			

- Molecule 2 is water.

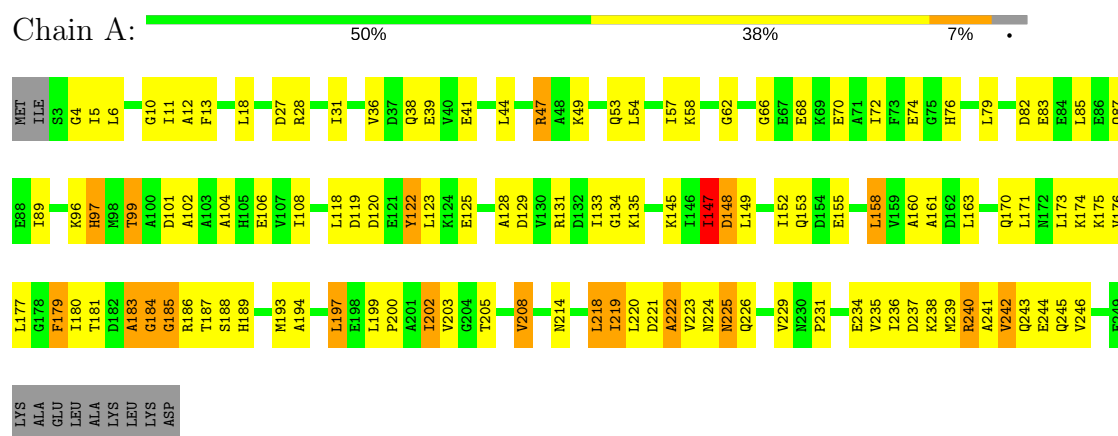
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	30	Total	O	0	0
			30	30		
2	B	34	Total	O	0	0
			34	34		

3 Residue-property plots

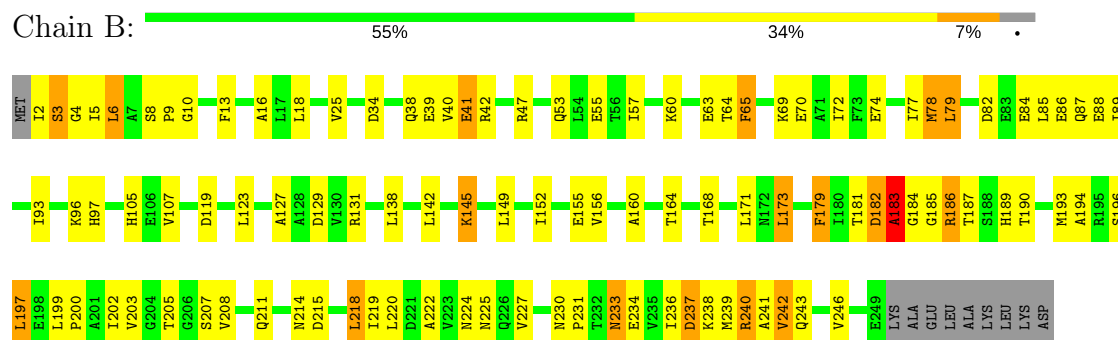
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are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

Note EDS was not executed.

• Molecule 1: ENZYME I



• Molecule 1: ENZYME I



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	47.42Å 75.78Å 170.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.50	Depositor
% Data completeness (in resolution range)	77.5 (10.00-2.50)	Depositor
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.205 , 0.306	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3854	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.59	0/1906	0.89	2/2571 (0.1%)
1	B	0.61	0/1914	0.87	3/2582 (0.1%)
All	All	0.60	0/3820	0.88	5/5153 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	185	GLY	N-CA-C	6.32	128.90	113.10
1	B	185	GLY	N-CA-C	5.76	127.51	113.10
1	B	183	ALA	N-CA-C	5.72	126.43	111.00
1	B	79	LEU	CA-CB-CG	-5.12	103.53	115.30
1	A	184	GLY	N-CA-C	-5.10	100.34	113.10

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	1891	0	1924	78	0
1	B	1899	0	1935	72	0
2	A	30	0	0	1	0
2	B	34	0	0	1	0
All	All	3854	0	3859	149	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 (149) 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:193:MET:SD	1:B:197:LEU:HD21	2.06	0.94
1:B:4:GLY:HA3	1:B:202:ILE:HG23	1.53	0.90
1:A:205:THR:HG21	1:A:208:VAL:HG22	1.57	0.84
1:B:186:ARG:HH11	1:B:186:ARG:HG2	1.41	0.83
1:B:60:LYS:O	1:B:64:THR:HG22	1.79	0.82
1:B:10:GLY:HA3	1:B:200:PRO:HG3	1.61	0.79
1:A:4:GLY:HA3	1:A:225:ASN:HD21	1.51	0.76
1:A:152:ILE:HG22	1:A:175:LYS:HD3	1.69	0.74
1:B:205:THR:HG21	1:B:208:VAL:HB	1.69	0.73
1:B:2:ILE:HG13	1:B:227:VAL:HB	1.71	0.73
1:B:64:THR:HG23	1:B:65:PHE:H	1.54	0.72
1:A:234:GLU:HA	1:A:237:ASP:HB2	1.70	0.72
1:A:185:GLY:HA3	1:A:203:VAL:HG21	1.72	0.72
1:B:160:ALA:O	1:B:181:THR:HA	1.91	0.70
1:A:160:ALA:O	1:A:181:THR:HA	1.90	0.70
1:A:236:ILE:O	1:A:240:ARG:HB2	1.92	0.69
1:A:155:GLU:HB3	1:A:177:LEU:HD21	1.74	0.69
1:A:10:GLY:O	1:A:222:ALA:HB3	1.94	0.68
1:A:155:GLU:HA	1:A:175:LYS:O	1.95	0.67
1:B:85:LEU:O	1:B:89:ILE:HG13	1.93	0.67
1:A:189:HIS:O	1:A:193:MET:HB2	1.96	0.66
1:A:18:LEU:H	1:A:214:ASN:HD22	1.43	0.65
1:A:10:GLY:HA3	1:A:200:PRO:HG3	1.79	0.65
1:B:69:LYS:HA	1:B:72:ILE:HD12	1.79	0.65
1:A:238:LYS:O	1:A:242:VAL:HB	1.97	0.64
1:B:155:GLU:O	1:B:156:VAL:HG13	1.98	0.64
1:A:243:GLN:O	1:A:246:VAL:HG13	2.00	0.62
1:A:120:ASP:HB3	1:A:123:LEU:HB2	1.83	0.61
1:A:27:ASP:O	1:A:102:ALA:HB2	2.01	0.61
1:A:221:ASP:HB3	1:A:226:GLN:H	1.66	0.61
1:A:174:LYS:N	1:A:174:LYS:HD2	2.17	0.60
1:A:238:LYS:NZ	1:A:238:LYS:HB2	2.16	0.60
1:B:47:ARG:NH2	1:B:86:GLU:HG3	2.15	0.60
1:B:10:GLY:O	1:B:222:ALA:HB3	2.02	0.60
1:B:238:LYS:O	1:B:242:VAL:HB	2.03	0.59
1:A:147:ILE:HD13	1:A:147:ILE:H	1.66	0.59
1:A:39:GLU:OE2	1:A:99:THR:HG22	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:ILE:HG12	1:A:97:HIS:O	2.03	0.58
1:B:4:GLY:HA3	1:B:202:ILE:CG2	2.31	0.57
1:B:60:LYS:HA	1:B:63:GLU:OE2	2.05	0.57
1:B:239:MET:O	1:B:241:ALA:N	2.37	0.57
1:A:197:LEU:HD22	1:A:199:LEU:HD13	1.86	0.56
1:A:173:LEU:HD13	1:A:197:LEU:HD11	1.88	0.56
1:B:38:GLN:OE1	1:B:41:GLU:HG2	2.06	0.56
1:A:125:GLU:O	1:A:128:ALA:HB3	2.04	0.56
1:B:5:ILE:HG22	1:B:203:VAL:HG23	1.87	0.56
1:B:16:ALA:HB2	1:B:218:LEU:HD21	1.88	0.56
1:A:185:GLY:HA3	1:A:203:VAL:CG2	2.37	0.55
1:B:197:LEU:HB3	1:B:199:LEU:HD13	1.89	0.55
1:B:8:SER:O	1:B:200:PRO:HA	2.06	0.54
1:A:231:PRO:HB2	1:A:235:VAL:HB	1.90	0.54
1:A:4:GLY:HA3	1:A:225:ASN:ND2	2.22	0.54
1:A:129:ASP:O	1:A:133:ILE:HG13	2.08	0.54
1:B:179:PHE:CD1	1:B:179:PHE:N	2.75	0.53
1:A:47:ARG:HD3	2:A:288:HOH:O	2.09	0.53
1:B:237:ASP:HA	1:B:240:ARG:HB3	1.89	0.53
1:B:189:HIS:O	1:B:193:MET:HB2	2.09	0.53
1:A:194:ALA:HB1	1:A:199:LEU:O	2.08	0.53
1:B:220:LEU:CD2	1:B:222:ALA:HB2	2.39	0.52
1:B:38:GLN:O	1:B:41:GLU:HB3	2.10	0.52
1:B:194:ALA:HB1	1:B:199:LEU:O	2.09	0.52
1:B:186:ARG:NH1	1:B:186:ARG:HG2	2.15	0.52
1:A:239:MET:O	1:A:241:ALA:N	2.42	0.51
1:B:242:VAL:HG12	1:B:243:GLN:N	2.25	0.51
1:A:218:LEU:HB3	1:A:229:VAL:HA	1.92	0.51
1:A:218:LEU:HD23	1:A:229:VAL:HG13	1.93	0.51
1:B:13:PHE:HZ	1:B:239:MET:HB2	1.76	0.51
1:A:89:ILE:HG12	1:A:104:ALA:HA	1.93	0.51
1:A:11:ILE:HA	1:A:220:LEU:O	2.11	0.50
1:A:181:THR:O	1:A:205:THR:HB	2.11	0.50
1:A:173:LEU:HD22	1:A:197:LEU:HD11	1.93	0.50
1:A:220:LEU:CD2	1:A:222:ALA:HB2	2.42	0.49
1:A:49:LYS:O	1:A:53:GLN:HG3	2.11	0.49
1:B:202:ILE:HD11	1:B:225:ASN:HA	1.93	0.49
1:B:18:LEU:H	1:B:214:ASN:HD22	1.60	0.49
1:B:89:ILE:O	1:B:93:ILE:HG13	2.12	0.49
1:A:241:ALA:HA	1:A:244:GLU:HB3	1.95	0.49
1:B:74:GLU:O	1:B:78:MET:HG3	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:145:LYS:HE3	2:B:276:HOH:O	2.13	0.48
1:A:122:TYR:CD1	1:A:122:TYR:C	2.86	0.48
1:A:104:ALA:O	1:A:108:ILE:HG12	2.13	0.48
1:A:158:LEU:HD12	1:A:163:LEU:HD11	1.96	0.48
1:A:183:ALA:HB3	1:A:203:VAL:HB	1.95	0.48
1:B:13:PHE:CZ	1:B:240:ARG:HA	2.49	0.48
1:A:82:ASP:OD2	1:A:85:LEU:HG	2.13	0.48
1:B:197:LEU:H	1:B:197:LEU:HD13	1.79	0.47
1:A:241:ALA:HB1	1:A:245:GLN:OE1	2.15	0.47
1:A:62:GLY:O	1:A:66:GLY:N	2.43	0.47
1:B:168:THR:O	1:B:171:LEU:HB2	2.14	0.47
1:B:193:MET:SD	1:B:197:LEU:CD2	2.92	0.47
1:A:148:ASP:HA	1:A:170:GLN:OE1	2.15	0.47
1:A:4:GLY:HA2	1:A:202:ILE:HG23	1.96	0.47
1:A:12:ALA:O	1:A:219:ILE:HA	2.15	0.46
1:A:179:PHE:CD1	1:A:179:PHE:N	2.83	0.46
1:B:181:THR:O	1:B:205:THR:HB	2.15	0.45
1:B:84:GLU:HG3	1:B:88:GLU:OE2	2.16	0.45
1:B:82:ASP:OD2	1:B:85:LEU:HG	2.17	0.45
1:B:197:LEU:CB	1:B:199:LEU:HD13	2.47	0.45
1:A:44:LEU:HD23	1:A:44:LEU:HA	1.77	0.45
1:B:18:LEU:H	1:B:214:ASN:ND2	2.15	0.45
1:A:145:LYS:HA	1:A:145:LYS:HD2	1.59	0.44
1:A:161:ALA:O	1:A:181:THR:HB	2.18	0.44
1:A:235:VAL:HA	1:A:238:LYS:HZ3	1.82	0.44
1:A:241:ALA:O	1:A:245:GLN:HB2	2.17	0.44
1:B:105:HIS:HB2	1:B:138:LEU:HD21	1.98	0.44
1:A:185:GLY:HA2	1:A:188:SER:OG	2.17	0.44
1:A:13:PHE:HE2	1:A:244:GLU:HG3	1.82	0.44
1:A:57:ILE:O	1:A:58:LYS:C	2.56	0.44
1:B:149:LEU:O	1:B:152:ILE:HG13	2.18	0.44
1:A:38:GLN:O	1:A:41:GLU:HB3	2.18	0.44
1:B:183:ALA:HB1	1:B:184:GLY:H	1.58	0.43
1:B:8:SER:HA	1:B:9:PRO:HD2	1.86	0.43
1:A:186:ARG:NH2	1:B:246:VAL:HG13	2.33	0.43
1:A:99:THR:HB	1:A:101:ASP:OD1	2.18	0.43
1:A:54:LEU:HD23	1:A:54:LEU:HA	1.88	0.43
1:A:58:LYS:HG2	1:A:70:GLU:HG2	1.99	0.43
1:B:138:LEU:O	1:B:142:LEU:HG	2.18	0.43
1:B:65:PHE:CZ	1:B:173:LEU:HD11	2.54	0.43
1:B:39:GLU:O	1:B:40:VAL:C	2.58	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:LEU:HD13	1:A:171:LEU:HD13	2.01	0.42
1:A:180:ILE:HG22	1:A:208:VAL:HG21	2.01	0.42
1:B:47:ARG:HH21	1:B:86:GLU:HG3	1.81	0.42
1:A:11:ILE:HD11	1:A:223:VAL:HG21	2.01	0.42
1:B:69:LYS:HD2	1:B:72:ILE:HD12	2.00	0.42
1:A:197:LEU:HD22	1:A:199:LEU:HD22	2.01	0.42
1:B:197:LEU:CD1	1:B:197:LEU:N	2.82	0.42
1:B:197:LEU:HD13	1:B:197:LEU:N	2.35	0.42
1:A:238:LYS:HB2	1:A:238:LYS:HZ3	1.82	0.41
1:A:131:ARG:O	1:A:135:LYS:HG2	2.20	0.41
1:B:127:ALA:O	1:B:131:ARG:HG3	2.19	0.41
1:B:230:ASN:N	1:B:231:PRO:HD3	2.35	0.41
1:A:234:GLU:O	1:A:238:LYS:N	2.54	0.41
1:B:2:ILE:CG1	1:B:227:VAL:HB	2.45	0.41
1:B:69:LYS:O	1:B:72:ILE:HB	2.21	0.41
1:A:108:ILE:HG21	1:A:134:GLY:HA3	2.03	0.41
1:A:202:ILE:HG12	1:A:225:ASN:ND2	2.36	0.41
1:B:236:ILE:O	1:B:240:ARG:N	2.48	0.41
1:B:60:LYS:O	1:B:63:GLU:HG2	2.20	0.41
1:B:89:ILE:HG12	1:B:107:VAL:HG21	2.03	0.41
1:A:96:LYS:HE2	1:A:96:LYS:HB3	1.83	0.41
1:B:186:ARG:HD3	1:B:186:ARG:HA	1.85	0.41
1:B:182:ASP:OD2	1:B:207:SER:HA	2.20	0.41
1:A:118:LEU:HD12	1:A:123:LEU:CD1	2.50	0.40
1:B:233:ASN:HA	1:B:236:ILE:HD12	2.02	0.40
1:A:149:LEU:HB2	1:A:170:GLN:HG3	2.02	0.40
1:B:53:GLN:O	1:B:57:ILE:HG13	2.22	0.40
1:B:55:GLU:HG2	1:B:77:ILE:HD13	2.02	0.40
1:B:123:LEU:HA	1:B:123:LEU:HD23	1.84	0.40
1:B:196:SER:HB2	1:B:197:LEU:HD13	2.03	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	A	245/258 (95%)	213 (87%)	25 (10%)	7 (3%)	5	7
1	B	246/258 (95%)	216 (88%)	23 (9%)	7 (3%)	6	8
All	All	491/516 (95%)	429 (87%)	48 (10%)	14 (3%)	5	7

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	147	ILE
1	A	183	ALA
1	A	240	ARG
1	B	65	PHE
1	B	183	ALA
1	B	240	ARG
1	B	242	VAL
1	B	3	SER
1	A	36	VAL
1	A	148	ASP
1	A	222	ALA
1	B	173	LEU
1	B	6	LEU
1	A	184	GLY

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	204/213 (96%)	174 (85%)	30 (15%)	3	6
1	B	205/213 (96%)	175 (85%)	30 (15%)	3	6
All	All	409/426 (96%)	349 (85%)	60 (15%)	3	6

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ILE
1	A	6	LEU
1	A	28	ARG
1	A	47	ARG
1	A	68	GLU
1	A	72	ILE
1	A	74	GLU
1	A	76	HIS
1	A	79	LEU
1	A	83	GLU
1	A	87	GLN
1	A	97	HIS
1	A	99	THR
1	A	106	GLU
1	A	119	ASP
1	A	122	TYR
1	A	147	ILE
1	A	153	GLN
1	A	158	LEU
1	A	176	VAL
1	A	179	PHE
1	A	187	THR
1	A	197	LEU
1	A	202	ILE
1	A	208	VAL
1	A	218	LEU
1	A	219	ILE
1	A	224	ASN
1	A	225	ASN
1	A	242	VAL
1	B	3	SER
1	B	6	LEU
1	B	25	VAL
1	B	34	ASP
1	B	41	GLU
1	B	42	ARG
1	B	70	GLU
1	B	78	MET
1	B	79	LEU
1	B	87	GLN
1	B	96	LYS
1	B	97	HIS
1	B	119	ASP

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Mol	Chain	Res	Type
1	B	129	ASP
1	B	145	LYS
1	B	164	THR
1	B	179	PHE
1	B	182	ASP
1	B	186	ARG
1	B	187	THR
1	B	190	THR
1	B	197	LEU
1	B	211	GLN
1	B	215	ASP
1	B	218	LEU
1	B	219	ILE
1	B	224	ASN
1	B	233	ASN
1	B	234	GLU
1	B	237	ASP

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	211	GLN
1	A	214	ASN
1	A	225	ASN
1	A	226	GLN
1	A	233	ASN
1	B	211	GLN
1	B	214	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

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.