



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:32 PM GMT

PDB ID : 4CHA
Title : STRUCTURE OF ALPHA-CHYMOTRYPSIN REFINED AT 1.68
ANGSTROMS RESOLUTION
Authors : Tsukada, H.; Blow, D.M.
Deposited on : 1984-11-26
Resolution : 1.68 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) : 1.9-1692
EDS : rb-20026688
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) : trunk26865

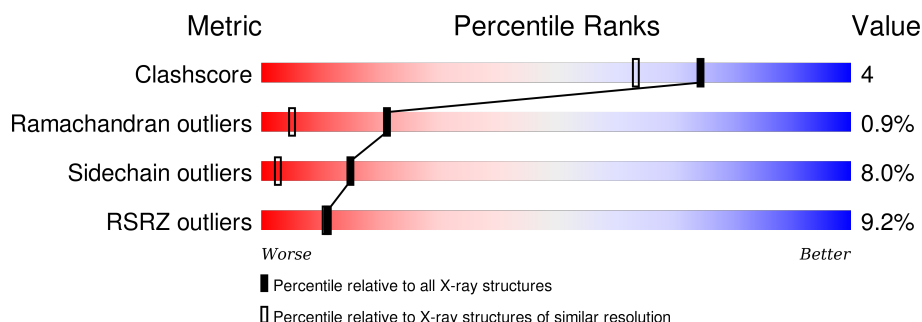
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.68 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	5317 (1.70-1.66)
Ramachandran outliers	100387	5225 (1.70-1.66)
Sidechain outliers	100360	5224 (1.70-1.66)
RSRZ outliers	91569	4813 (1.70-1.66)

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	13	<div> <div>23%</div> <div>62%</div> <div>23%</div> <div>15%</div> </div>
1	E	13	<div> <div>15%</div> <div>54%</div> <div>23%</div> <div>23%</div> </div>
2	B	131	<div> <div>11%</div> <div>79%</div> <div>18%</div> <div>•</div> </div>
2	F	131	<div> <div>8%</div> <div>74%</div> <div>21%</div> <div>5% •</div> </div>
3	C	97	<div> <div>7%</div> <div>78%</div> <div>18%</div> <div>• •</div> </div>
3	G	97	<div> <div>8%</div> <div>81%</div> <div>13%</div> <div>• •</div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3591 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 ALPHA-CHYMOTRYPSIN A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	11	Total	C	N	O	S	0	0	0
			74	48	12	13	1			
1	E	10	Total	C	N	O	S	0	0	0
			68	45	11	11	1			

- Molecule 2 is a protein called ALPHA-CHYMOTRYPSIN A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	131	Total	C	N	O	S	0	0	0
			980	618	162	196	4			
2	F	131	Total	C	N	O	S	0	0	0
			980	618	162	196	4			

- Molecule 3 is a protein called ALPHA-CHYMOTRYPSIN A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	97	Total	C	N	O	S	0	0	0
			702	436	123	136	7			
3	G	97	Total	C	N	O	S	0	0	0
			702	436	123	136	7			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	O	0	0
			1	1		
4	B	23	Total	O	0	0
			23	23		
4	C	18	Total	O	0	0
			18	18		
4	F	21	Total	O	0	0
			21	21		

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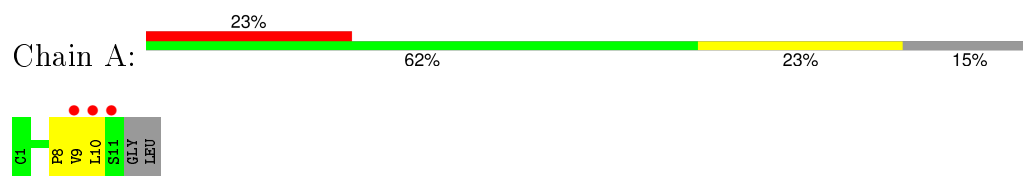
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	G	22	Total	O	0	0
			22	22		

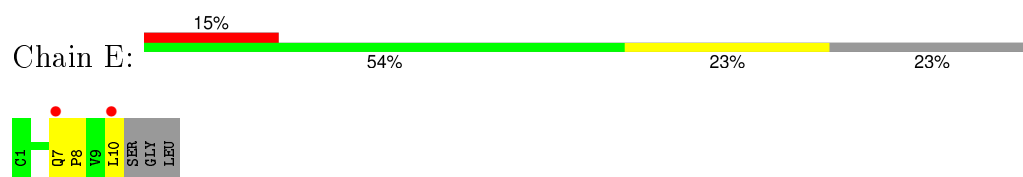
3 Residue-property plots [i](#)

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.

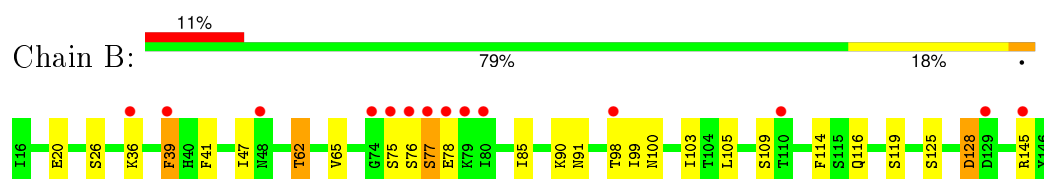
- Molecule 1: ALPHA-CHYMOTRYPSIN A



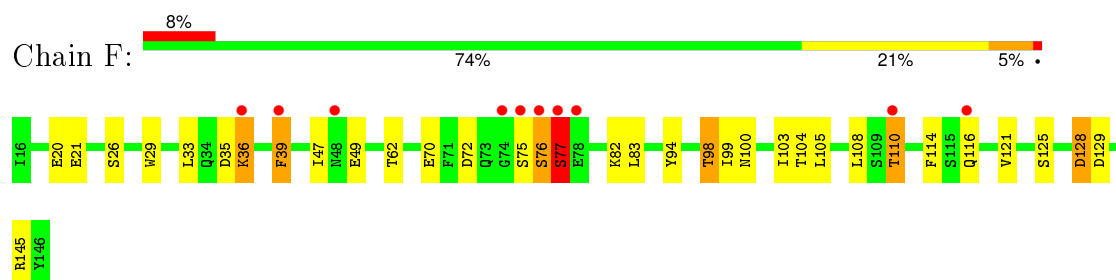
- Molecule 1: ALPHA-CHYMOTRYPSIN A



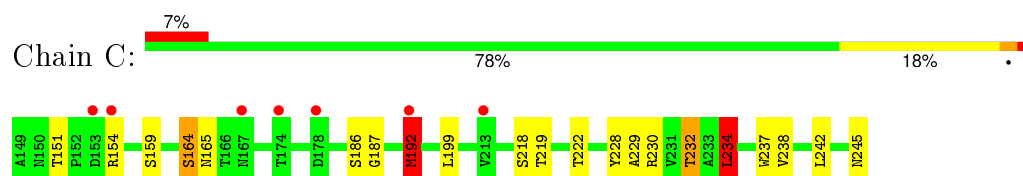
- Molecule 2: ALPHA-CHYMOTRYPSIN A



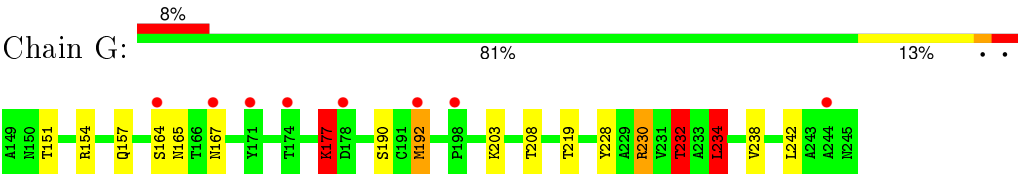
- Molecule 2: ALPHA-CHYMOTRYPSIN A



- Molecule 3: ALPHA-CHYMOTRYPSIN A



- Molecule 3: ALPHA-CHYMOTRYPSIN A



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	49.23 Å 67.39 Å 65.99 Å 90.00° 101.80° 90.00°	Depositor
Resolution (Å)	(Not available) – 1.68 29.87 – 1.69	Depositor EDS
% Data completeness (in resolution range)	(Not available) ((Not available)-1.68) 94.4 (29.87-1.69)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.93 (at 1.70 Å)	Xtriage
Refinement program	unknown	Depositor
R, R_{free}	(Not available) , (Not available) 0.239 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	15.2	Xtriage
Anisotropy	0.273	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 46.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 44397 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3591	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.18% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.71	0/75	1.79	2/103 (1.9%)
1	E	0.78	0/69	1.54	0/95
2	B	0.91	0/1000	1.79	14/1361 (1.0%)
2	F	0.92	0/1000	1.91	24/1361 (1.8%)
3	C	0.94	1/715 (0.1%)	1.69	12/973 (1.2%)
3	G	0.93	0/715	1.72	14/973 (1.4%)
All	All	0.92	1/3574 (0.0%)	1.79	66/4866 (1.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	245	ASN	C-OXT	5.16	1.33	1.23

All (66) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	128	ASP	CB-CG-OD1	23.73	139.66	118.30
2	B	128	ASP	CB-CG-OD1	15.12	131.91	118.30
2	B	114	PHE	CB-CG-CD1	11.31	128.71	120.80
2	F	145	ARG	NE-CZ-NH2	10.08	125.34	120.30
2	B	125	SER	N-CA-CB	9.08	124.12	110.50
2	B	114	PHE	CB-CG-CD2	-8.92	114.56	120.80
2	F	114	PHE	CB-CG-CD1	8.83	126.98	120.80
3	G	232	THR	N-CA-CB	-8.54	94.08	110.30
3	G	154	ARG	NE-CZ-NH2	8.36	124.48	120.30
2	F	128	ASP	OD1-CG-OD2	-8.01	108.07	123.30
2	B	65	VAL	CA-CB-CG1	7.98	122.88	110.90
3	C	154	ARG	NE-CZ-NH2	7.96	124.28	120.30
2	F	125	SER	N-CA-CB	7.82	122.24	110.50
2	F	72	ASP	CB-CG-OD1	7.80	125.32	118.30
2	F	128	ASP	CA-CB-CG	7.74	130.42	113.40
3	G	228	TYR	CB-CG-CD2	-7.67	116.40	121.00
2	B	47	ILE	C-N-CA	7.27	139.88	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	114	PHE	CB-CG-CD2	-7.21	115.75	120.80
3	G	234	LEU	CA-CB-CG	6.95	131.29	115.30
2	F	145	ARG	CA-CB-CG	6.94	128.67	113.40
3	G	192	MET	CA-CB-CG	-6.91	101.55	113.30
2	F	128	ASP	CB-CG-OD2	-6.85	112.13	118.30
3	C	232	THR	N-CA-CB	-6.84	97.31	110.30
2	B	128	ASP	CA-CB-CG	6.78	128.31	113.40
3	G	230	ARG	NE-CZ-NH2	6.70	123.65	120.30
3	C	186	SER	N-CA-CB	-6.62	100.57	110.50
2	B	145	ARG	CD-NE-CZ	6.60	132.84	123.60
3	C	219	THR	N-CA-CB	-6.55	97.86	110.30
3	C	192	MET	CA-CB-CG	-6.42	102.38	113.30
1	A	9	VAL	CA-CB-CG1	6.34	120.41	110.90
2	F	47	ILE	C-N-CA	6.33	137.53	121.70
3	C	164	SER	N-CA-CB	6.24	119.86	110.50
3	C	229	ALA	N-CA-CB	6.20	118.78	110.10
3	G	177	LYS	CA-CB-CG	6.15	126.92	113.40
2	B	76	SER	C-N-CA	6.13	137.03	121.70
3	G	219	THR	CA-CB-CG2	6.10	120.94	112.40
2	F	20	GLU	OE1-CD-OE2	5.98	130.47	123.30
3	G	190	SER	N-CA-CB	5.97	119.46	110.50
2	B	41	PHE	CB-CG-CD1	-5.92	116.65	120.80
3	G	219	THR	N-CA-CB	-5.88	99.12	110.30
3	G	232	THR	CB-CA-C	5.87	127.45	111.60
3	C	218	SER	C-N-CA	5.86	136.35	121.70
2	F	105	LEU	CA-CB-CG	5.85	128.76	115.30
2	B	91	ASN	CA-CB-CG	5.82	126.21	113.40
2	F	104	THR	CA-CB-CG2	5.82	120.55	112.40
3	G	234	LEU	N-CA-CB	-5.82	98.77	110.40
2	F	35	ASP	CB-CG-OD2	5.72	123.44	118.30
2	F	76	SER	C-N-CA	5.70	135.94	121.70
2	B	91	ASN	CB-CA-C	5.69	121.78	110.40
2	F	77	SER	C-N-CA	5.67	135.88	121.70
3	C	228	TYR	CB-CG-CD2	-5.67	117.60	121.00
3	G	154	ARG	NH1-CZ-NH2	-5.58	113.27	119.40
2	B	62	THR	CA-CB-CG2	5.52	120.13	112.40
3	C	199	LEU	N-CA-CB	5.44	121.27	110.40
1	A	10	LEU	CA-CB-CG	5.42	127.78	115.30
2	F	94	TYR	CB-CG-CD2	-5.38	117.77	121.00
2	F	33	LEU	O-C-N	5.30	131.18	122.70
2	F	110	THR	CA-CB-CG2	5.26	119.76	112.40
3	C	228	TYR	CB-CG-CD1	5.24	124.15	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	145	ARG	NH1-CZ-NH2	-5.22	113.66	119.40
3	C	234	LEU	CA-CB-CG	5.22	127.30	115.30
2	B	128	ASP	OD1-CG-OD2	-5.20	113.43	123.30
3	G	208	THR	O-C-N	5.14	130.92	122.70
2	F	98	THR	CA-CB-CG2	5.10	119.55	112.40
2	F	21	GLU	CG-CD-OE1	5.07	128.44	118.30
2	F	70	GLU	CA-CB-CG	5.04	124.49	113.40

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	74	0	81	1	0
1	E	68	0	76	3	0
2	B	980	0	950	7	0
2	F	980	0	951	9	1
3	C	702	0	698	10	0
3	G	702	0	698	11	1
4	A	1	0	0	0	0
4	B	23	0	0	0	0
4	C	18	0	0	0	0
4	F	21	0	0	0	0
4	G	22	0	0	1	0
All	All	3591	0	3454	28	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 4.

All (28) 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)
3:C:192:MET:HG2	3:G:192:MET:HG2	1.66	0.78
2:F:98:THR:HG22	2:F:100:ASN:HB2	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:39:PHE:HE2	3:G:151:THR:HG21	1.64	0.61
3:G:165:ASN:HD21	3:G:230:ARG:HH11	1.48	0.60
2:B:98:THR:HG22	2:B:100:ASN:HB2	1.85	0.58
2:F:103:ILE:HG21	3:G:234:LEU:HG	1.88	0.55
2:F:83:LEU:HD13	2:F:108:LEU:HD13	1.89	0.53
3:C:151:THR:HG21	2:F:39:PHE:CZ	2.44	0.52
2:F:36:LYS:HE2	2:F:36:LYS:HA	1.95	0.48
2:B:105:LEU:HD23	3:C:237:TRP:CZ3	2.49	0.47
2:B:39:PHE:CE2	3:G:151:THR:HG21	2.47	0.46
3:C:151:THR:HG21	2:F:39:PHE:HZ	1.80	0.45
2:F:29:TRP:CG	2:F:121:VAL:HB	2.52	0.45
2:F:98:THR:HG23	3:G:177:LYS:HE2	1.98	0.45
1:E:10:LEU:HD22	3:G:157:GLN:HE21	1.83	0.44
3:C:238:VAL:O	3:C:242:LEU:HG	2.18	0.44
2:B:62:THR:HG23	2:B:85:ILE:O	2.18	0.43
3:G:238:VAL:O	3:G:242:LEU:HG	2.18	0.43
1:E:7:GLN:HA	1:E:8:PRO:HD3	1.92	0.43
3:G:232:THR:HB	4:G:570:HOH:O	2.16	0.43
3:C:165:ASN:HD21	3:C:230:ARG:HH11	1.66	0.43
3:C:187:GLY:C	3:C:222:THR:HB	2.40	0.42
3:C:165:ASN:ND2	3:C:230:ARG:HH11	2.17	0.42
3:G:164:SER:HB2	3:G:167:ASN:OD1	2.20	0.42
3:C:192:MET:HG2	3:G:192:MET:CG	2.42	0.41
2:B:103:ILE:HG21	3:C:234:LEU:HG	2.01	0.41
1:E:8:PRO:HA	2:F:26:SER:HB2	2.02	0.41
1:A:8:PRO:HA	2:B:26:SER:HB2	2.03	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)
2:F:76:SER:CB	3:G:167:ASN:ND2[1_455]	2.16	0.04

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	9/13 (69%)	9 (100%)	0	0	100	100
1	E	8/13 (62%)	8 (100%)	0	0	100	100
2	B	129/131 (98%)	125 (97%)	2 (2%)	2 (2%)	12	1
2	F	129/131 (98%)	125 (97%)	2 (2%)	2 (2%)	12	1
3	C	95/97 (98%)	91 (96%)	4 (4%)	0	100	100
3	G	95/97 (98%)	90 (95%)	5 (5%)	0	100	100
All	All	465/482 (96%)	448 (96%)	13 (3%)	4 (1%)	21	5

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	77	SER
2	B	77	SER
2	B	99	ILE
2	F	99	ILE

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	9/10 (90%)	9 (100%)	0	100	100
1	E	8/10 (80%)	8 (100%)	0	100	100
2	B	109/109 (100%)	98 (90%)	11 (10%)	9	1
2	F	109/109 (100%)	98 (90%)	11 (10%)	9	1
3	C	77/77 (100%)	72 (94%)	5 (6%)	21	5
3	G	77/77 (100%)	73 (95%)	4 (5%)	29	8
All	All	389/392 (99%)	358 (92%)	31 (8%)	15	3

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	20	GLU
2	B	36	LYS
2	B	39	PHE
2	B	75	SER
2	B	77	SER
2	B	78	GLU
2	B	90	LYS
2	B	109	SER
2	B	116	GLN
2	B	119	SER
2	B	128	ASP
3	C	159	SER
3	C	164	SER
3	C	192	MET
3	C	232	THR
3	C	234	LEU
2	F	36	LYS
2	F	39	PHE
2	F	49	GLU
2	F	62	THR
2	F	75	SER
2	F	77	SER
2	F	82	LYS
2	F	110	THR
2	F	116	GLN
2	F	128	ASP
2	F	129	ASP
3	G	177	LYS
3	G	203	LYS
3	G	232	THR
3	G	234	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	18	ASN
2	B	34	GLN
2	B	48	ASN
3	C	165	ASN
3	G	165	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 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	11/13 (84%)	1.81	3 (27%) 1 1	10, 14, 29, 30	0
1	E	10/13 (76%)	1.11	2 (20%) 1 1	9, 12, 19, 22	0
2	B	131/131 (100%)	0.90	14 (10%) 8 7	5, 13, 27, 36	0
2	F	131/131 (100%)	0.85	10 (7%) 17 17	6, 14, 26, 35	0
3	C	97/97 (100%)	0.64	7 (7%) 18 18	5, 12, 20, 21	0
3	G	97/97 (100%)	0.61	8 (8%) 14 14	5, 12, 22, 23	0
All	All	477/482 (98%)	0.80	44 (9%) 11 11	5, 13, 23, 36	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	74	GLY	9.9
2	F	76	SER	8.1
1	E	10	LEU	7.2
1	A	11	SER	6.8
2	B	76	SER	6.2
1	A	10	LEU	6.0
2	B	77	SER	6.0
2	F	39	PHE	4.8
2	F	77	SER	4.2
2	B	79	LYS	4.2
2	F	74	GLY	4.2
2	F	75	SER	4.2
2	B	36	LYS	4.0
2	F	110	THR	3.8
2	B	110	THR	3.5
2	B	39	PHE	3.5
1	A	9	VAL	3.5
2	B	75	SER	3.1
2	F	36	LYS	3.1

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Mol	Chain	Res	Type	RSRZ
2	B	78	GLU	3.0
3	G	244	ALA	3.0
3	G	171	TYR	3.0
3	G	167	ASN	2.9
3	C	192	MET	2.8
3	C	153	ASP	2.8
3	C	174	THR	2.7
3	C	167	ASN	2.6
3	C	154	ARG	2.5
2	F	78	GLU	2.4
3	G	178	ASP	2.4
3	C	178	ASP	2.4
2	F	48	ASN	2.3
3	G	164	SER	2.3
2	B	129	ASP	2.3
2	B	98	THR	2.2
1	E	7	GLN	2.2
2	F	116	GLN	2.2
2	B	145	ARG	2.2
3	G	198	PRO	2.2
2	B	48	ASN	2.2
3	G	174	THR	2.1
2	B	80	ILE	2.1
3	C	213	VAL	2.0
3	G	192	MET	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.