



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:44 PM UTC

PDB ID : 4E2H / pdb_00004e2h
Title : Crystal structure of the periplasmic domain of Shigella flexneri WzzB
Authors : Kalynych, S.; Yao, D.; Magee, J.D.; Cygler, M.
Deposited on : 2012-03-08
Resolution : 2.36 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
Mogul : 2022.3.0, CSD as543be (2022)
Xtrriage (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

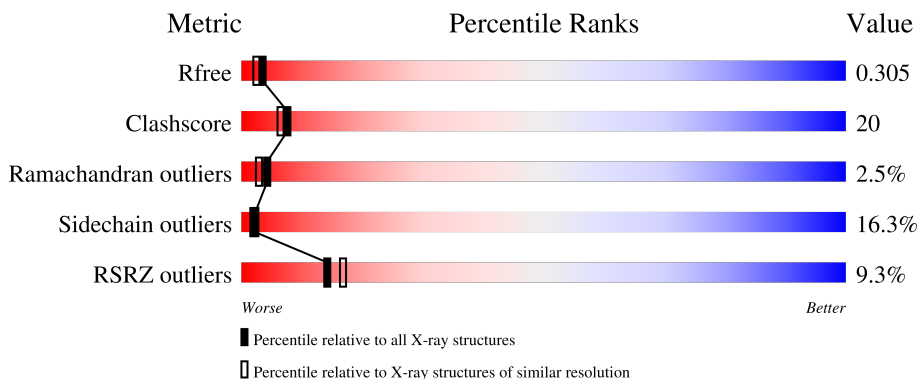
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.36 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	240	 8% 52% 30% 7% 10%
1	B	240	 8% 59% 27% 5% 8%
1	C	240	 10% 48% 29% 8% 13%

2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 5007 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 Chain length determinant protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	Se			
1	A	216	1647	1035	285	324	3	0	0	0
1	C	208	1578	993	265	317	3	0	0	0
1	B	222	1704	1063	296	342	3	0	0	0

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	46	HIS	-	expression tag	UNP P37792
A	47	HIS	-	expression tag	UNP P37792
A	48	HIS	-	expression tag	UNP P37792
A	49	HIS	-	expression tag	UNP P37792
A	50	HIS	-	expression tag	UNP P37792
A	51	HIS	-	expression tag	UNP P37792
A	52	GLY	-	expression tag	UNP P37792
A	53	SER	-	expression tag	UNP P37792
C	46	HIS	-	expression tag	UNP P37792
C	47	HIS	-	expression tag	UNP P37792
C	48	HIS	-	expression tag	UNP P37792
C	49	HIS	-	expression tag	UNP P37792
C	50	HIS	-	expression tag	UNP P37792
C	51	HIS	-	expression tag	UNP P37792
C	52	GLY	-	expression tag	UNP P37792
C	53	SER	-	expression tag	UNP P37792
B	46	HIS	-	expression tag	UNP P37792
B	47	HIS	-	expression tag	UNP P37792
B	48	HIS	-	expression tag	UNP P37792
B	49	HIS	-	expression tag	UNP P37792
B	50	HIS	-	expression tag	UNP P37792
B	51	HIS	-	expression tag	UNP P37792
B	52	GLY	-	expression tag	UNP P37792

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Chain	Residue	Modelled	Actual	Comment	Reference
B	53	SER	-	expression tag	UNP P37792

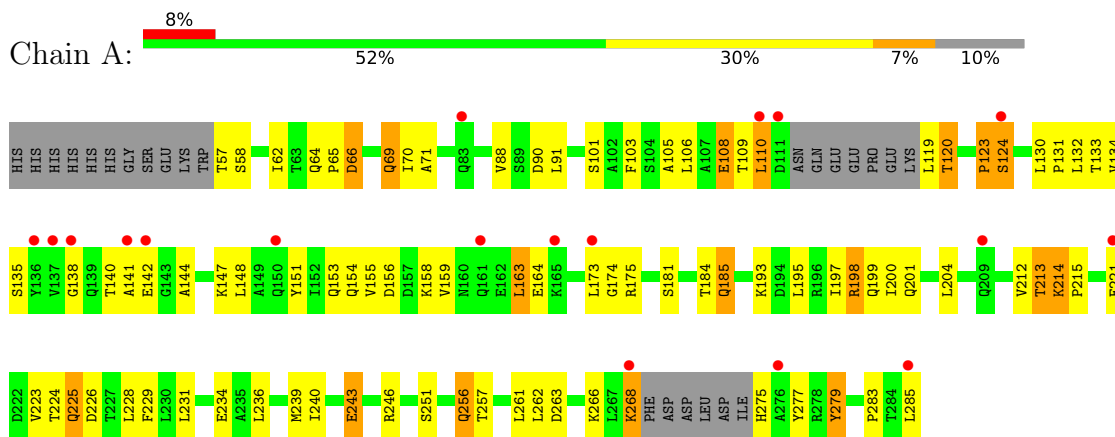
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	26	Total O 26 26	0	0
2	C	31	Total O 31 31	0	0
2	B	21	Total O 21 21	0	0

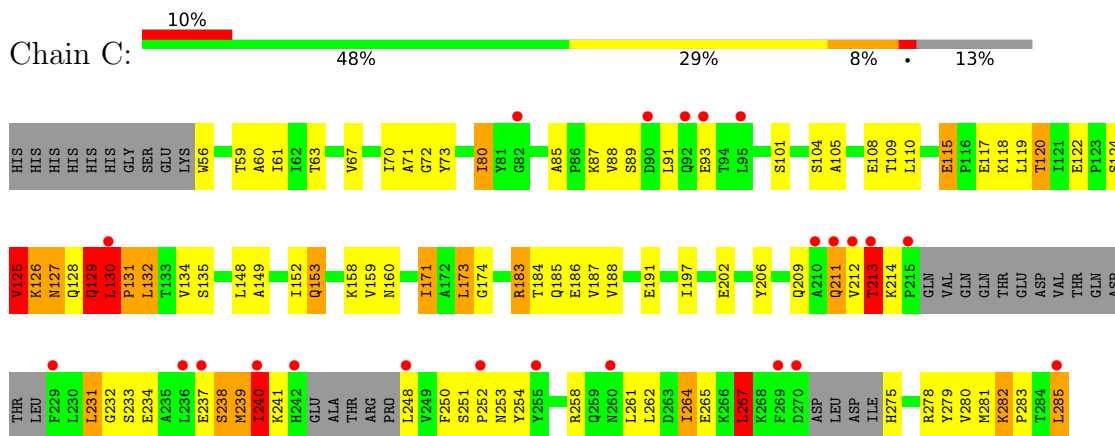
3 Residue-property plots [i](#)

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 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.

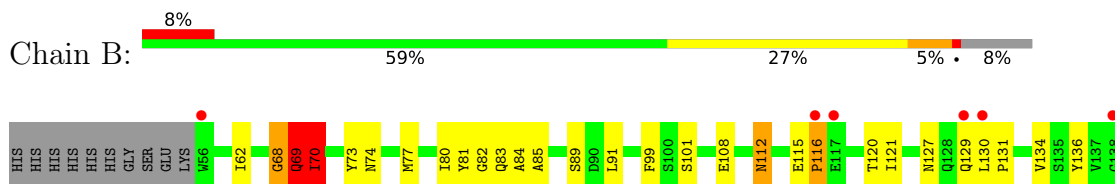
- Molecule 1: Chain length determinant protein

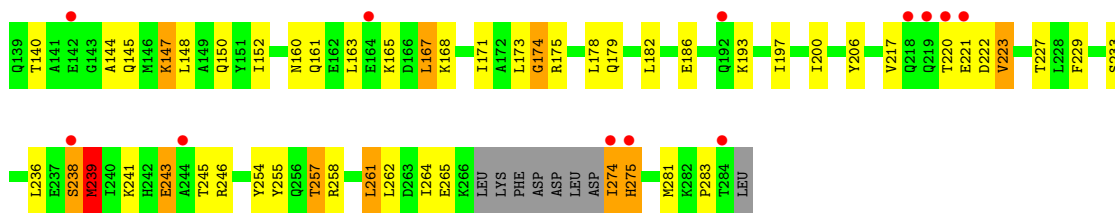


- Molecule 1: Chain length determinant protein



- Molecule 1: Chain length determinant protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	81.51Å 62.62Å 90.74Å 90.00° 94.67° 90.00°	Depositor
Resolution (Å)	49.60 – 2.36 49.60 – 2.36	Depositor EDS
% Data completeness (in resolution range)	78.0 (49.60-2.36) 78.0 (49.60-2.36)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.63 (at 2.37Å)	Xtrriage
Refinement program	PHENIX 1.7_650	Depositor
R, R_{free}	0.245 , 0.302 0.246 , 0.305	Depositor DCC
R_{free} test set	1509 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	33.2	Xtrriage
Anisotropy	0.710	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 49.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5007	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.70% 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 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	0.60	0/1663	1.00	5/2254 (0.2%)
1	B	0.60	0/1722	1.03	9/2336 (0.4%)
1	C	0.60	0/1595	1.05	12/2161 (0.6%)
All	All	0.60	0/4980	1.03	26/6751 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	2
1	C	0	4
All	All	0	6

There are no bond length outliers.

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	225	GLN	N-CA-C	-8.17	99.79	110.33
1	B	144	ALA	N-CA-C	-8.06	103.91	114.31
1	C	183	ARG	N-CA-C	-7.54	103.17	112.38
1	A	184	THR	N-CA-C	-7.21	106.40	114.62
1	C	129	GLN	N-CA-C	-6.77	96.38	110.80
1	B	83	GLN	N-CA-C	-6.75	105.11	113.15
1	C	130	LEU	N-CA-C	-6.10	93.92	111.00
1	B	69	GLN	N-CA-C	-6.01	98.00	110.80
1	C	211	GLN	N-CA-C	-5.62	104.58	112.12
1	C	85	ALA	CA-C-N	5.48	125.43	119.78
1	C	85	ALA	C-N-CA	5.48	125.43	119.78
1	B	227	THR	N-CA-C	5.38	119.32	112.87
1	B	80	ILE	N-CA-C	5.33	115.95	110.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	174	GLY	N-CA-C	-5.29	107.76	114.16
1	C	262	LEU	N-CA-C	-5.28	107.50	114.04
1	A	69	GLN	N-CA-C	-5.25	106.83	114.12
1	B	223	VAL	N-CA-C	5.24	115.67	108.12
1	B	130	LEU	N-CA-C	5.24	121.38	109.81
1	C	240	ILE	N-CA-C	-5.23	98.45	109.34
1	A	120	THR	N-CA-C	5.21	117.57	108.76
1	B	217	VAL	N-CA-C	5.15	115.32	108.11
1	A	213	THR	N-CA-C	-5.14	107.18	113.50
1	C	214	LYS	N-CA-C	5.08	119.70	108.27
1	C	115	GLU	CA-C-N	5.07	125.23	119.90
1	C	115	GLU	C-N-CA	5.07	125.23	119.90
1	C	80	ILE	CB-CA-C	-5.04	105.27	112.22

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	129	GLN	Peptide
1	B	68	GLY	Peptide
1	C	129	GLN	Peptide
1	C	130	LEU	Peptide
1	C	213	THR	Peptide
1	C	239	MSE	Peptide

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	1647	0	1618	67	0
1	B	1704	0	1652	58	0
1	C	1578	0	1498	83	0
2	A	26	0	0	9	0
2	B	21	0	0	1	0
2	C	31	0	0	8	0
All	All	5007	0	4768	196	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 20.

All (196) 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:C:285:LEU:HD13	1:C:285:LEU:O	1.58	1.02
1:B:148:LEU:HD23	1:B:283:PRO:HB3	1.43	0.99
1:C:238:SER:O	1:C:239:MSE:HB3	1.77	0.82
1:C:285:LEU:O	1:C:285:LEU:CD1	2.28	0.81
1:C:241:LYS:HA	2:C:313:HOH:O	1.80	0.80
1:A:148:LEU:HD23	1:A:283:PRO:HB3	1.63	0.79
1:C:120:THR:CG2	1:C:135:SER:HB2	2.14	0.78
1:A:147:LYS:O	1:A:151:TYR:HD1	1.67	0.77
1:C:264:ILE:HG12	1:C:265:GLU:HG2	1.66	0.76
1:C:129:GLN:N	1:C:130:LEU:CB	2.49	0.76
1:C:61:ILE:HG22	1:C:280:VAL:HB	1.69	0.75
1:C:231:LEU:HD12	1:C:233:SER:H	1.52	0.75
1:C:129:GLN:H	1:C:130:LEU:CB	2.00	0.75
1:C:239:MSE:HG3	1:C:240:ILE:H	1.54	0.72
1:C:275:HIS:N	2:C:301:HOH:O	2.22	0.72
1:A:268:LYS:O	2:A:319:HOH:O	2.07	0.72
1:A:151:TYR:O	1:A:155:VAL:HG23	1.89	0.71
2:C:324:HOH:O	1:B:239:MSE:SE	2.59	0.69
1:B:70:ILE:HG13	1:B:70:ILE:O	1.92	0.69
1:C:129:GLN:CA	1:C:130:LEU:CB	2.70	0.69
1:C:80:ILE:HG23	1:C:174:GLY:HA3	1.75	0.67
1:C:231:LEU:HG	1:C:234:GLU:OE1	1.95	0.67
1:C:212:VAL:HG22	1:C:239:MSE:HE3	1.77	0.67
1:B:68:GLY:O	1:B:69:GLN:HB2	1.95	0.66
1:C:119:LEU:HD12	1:C:135:SER:O	1.95	0.66
1:C:206:TYR:OH	2:C:307:HOH:O	2.10	0.65
1:C:108:GLU:OE2	1:B:281:MSE:CB	2.45	0.65
1:A:123:PRO:O	1:A:124:SER:O	2.15	0.65
1:B:274:ILE:O	1:B:275:HIS:HB2	1.97	0.65
1:A:156:ASP:C	2:A:322:HOH:O	2.40	0.65
1:C:126:LYS:O	1:C:126:LYS:HG3	1.95	0.64
1:A:256:GLN:HG3	1:A:257:THR:N	2.12	0.64
1:C:240:ILE:O	2:C:313:HOH:O	2.14	0.64
1:C:105:ALA:O	1:C:108:GLU:HG2	1.98	0.64
1:C:239:MSE:HG3	1:C:240:ILE:N	2.09	0.63
1:C:128:GLN:HG3	1:B:127:ASN:O	1.99	0.63
1:A:231:LEU:HD13	1:A:239:MSE:HE1	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:PRO:HG3	1:B:101:SER:HB2	1.81	0.62
1:C:278:ARG:HD3	2:C:318:HOH:O	2.00	0.62
1:C:120:THR:HG21	1:C:135:SER:HB2	1.83	0.61
1:A:228:LEU:HB2	2:A:323:HOH:O	1.99	0.61
1:C:231:LEU:CD1	1:C:233:SER:H	2.14	0.60
1:A:195:LEU:O	1:A:199:GLN:HG3	2.01	0.60
1:B:115:GLU:O	1:B:116:PRO:O	2.19	0.59
1:B:236:LEU:O	1:B:239:MSE:HB3	2.03	0.59
1:A:120:THR:O	1:A:134:VAL:HA	2.03	0.58
1:B:120:THR:O	1:B:134:VAL:HA	2.03	0.58
1:B:69:GLN:HG2	1:B:275:HIS:ND1	2.18	0.58
1:C:56:TRP:N	2:C:303:HOH:O	2.36	0.58
1:B:73:TYR:OH	1:B:77:MSE:HE3	2.02	0.58
1:C:285:LEU:CD1	1:C:285:LEU:C	2.75	0.58
1:A:256:GLN:HG3	1:A:257:THR:H	1.68	0.58
1:B:82:GLY:C	1:B:84:ALA:H	2.11	0.58
1:B:73:TYR:CD2	1:B:91:LEU:HD21	2.39	0.57
1:C:171:ILE:HD11	1:C:267:LEU:HD23	1.85	0.57
1:B:68:GLY:O	1:B:69:GLN:CB	2.53	0.57
1:C:234:GLU:HG3	1:B:220:THR:O	2.03	0.57
1:A:159:VAL:N	2:A:322:HOH:O	2.37	0.57
1:C:191:GLU:O	2:C:305:HOH:O	2.17	0.57
1:A:138:GLY:HA3	1:A:144:ALA:HB2	1.87	0.56
1:B:91:LEU:O	1:B:91:LEU:HD23	2.04	0.56
1:C:59:THR:O	1:C:283:PRO:HA	2.04	0.56
1:B:73:TYR:CZ	1:B:77:MSE:HE3	2.40	0.56
1:B:238:SER:O	1:B:239:MSE:HB2	2.05	0.56
1:A:197:ILE:HG13	1:A:243:GLU:HG3	1.87	0.56
1:A:263:ASP:O	1:A:266:LYS:HG3	2.06	0.56
1:A:103:PHE:CZ	1:A:119:LEU:HD22	2.40	0.56
1:A:229:PHE:HA	1:A:236:LEU:HD11	1.87	0.56
1:A:62:ILE:HG22	1:A:279:TYR:HD1	1.71	0.55
1:C:108:GLU:HG3	1:C:109:THR:N	2.20	0.55
1:B:167:LEU:O	1:B:167:LEU:HG	2.02	0.55
1:A:105:ALA:HA	1:A:108:GLU:OE1	2.07	0.55
1:C:239:MSE:O	1:C:240:ILE:CB	2.54	0.55
1:C:120:THR:O	1:C:134:VAL:HA	2.07	0.55
1:C:231:LEU:CD2	1:B:222:ASP:HA	2.37	0.55
1:C:80:ILE:HD11	1:C:171:ILE:HG13	1.88	0.54
1:C:120:THR:HG23	1:C:135:SER:H	1.73	0.54
1:B:145:GLN:HG2	1:B:145:GLN:O	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:64:GLN:HB2	1:A:65:PRO:HD2	1.88	0.54
1:A:105:ALA:O	1:A:108:GLU:OE1	2.26	0.54
1:C:149:ALA:HB2	1:C:283:PRO:HG2	1.90	0.53
1:B:160:ASN:OD1	1:B:274:ILE:HG12	2.08	0.53
1:B:179:GLN:HG3	1:B:261:LEU:HD11	1.90	0.53
1:A:140:THR:HG22	1:A:142:GLU:N	2.23	0.53
1:C:101:SER:OG	1:B:131:PRO:HB3	2.09	0.53
1:A:193:LYS:HE3	1:A:246:ARG:HG3	1.91	0.53
1:C:127:ASN:HB2	1:C:130:LEU:CA	2.38	0.52
1:A:106:LEU:HD21	1:A:151:TYR:HD2	1.74	0.52
1:C:63:THR:HB	1:C:131:PRO:HB3	1.92	0.52
1:B:274:ILE:HG22	1:B:275:HIS:HD2	1.73	0.52
1:C:59:THR:HG22	1:C:60:ALA:N	2.25	0.52
1:C:61:ILE:CG2	1:C:280:VAL:HB	2.37	0.52
1:A:62:ILE:HG22	1:A:279:TYR:CD1	2.44	0.52
1:A:181:SER:O	1:A:185:GLN:HG3	2.10	0.52
1:A:285:LEU:HD12	1:A:285:LEU:C	2.35	0.52
1:C:108:GLU:HG3	1:C:109:THR:HG23	1.91	0.51
1:B:182:LEU:HD21	1:B:257:THR:HG22	1.92	0.51
1:A:246:ARG:HG3	1:A:246:ARG:O	2.10	0.51
1:A:123:PRO:O	1:A:124:SER:C	2.53	0.51
1:C:186:GLU:HA	1:C:250:PHE:HE1	1.76	0.51
1:C:110:LEU:HB3	1:C:117:GLU:HB2	1.92	0.51
1:C:251:SER:OG	1:C:252:PRO:HD2	2.10	0.51
1:B:112:ASN:C	1:B:112:ASN:HD22	2.18	0.51
1:B:136:TYR:CE2	1:B:147:LYS:HG2	2.46	0.51
1:A:201:GLN:NE2	2:A:313:HOH:O	2.44	0.49
1:C:61:ILE:HG13	1:C:281:MSE:HE2	1.93	0.49
1:C:237:GLU:O	1:C:238:SER:CB	2.59	0.49
1:A:123:PRO:HA	1:A:132:LEU:HD23	1.94	0.49
1:A:120:THR:OG1	1:A:135:SER:HB2	2.12	0.49
1:C:124:SER:HB2	1:C:131:PRO:HD2	1.94	0.49
1:C:127:ASN:OD1	1:C:127:ASN:N	2.44	0.49
1:C:173:LEU:HD21	1:B:74:ASN:HD22	1.77	0.49
1:B:221:GLU:O	1:B:222:ASP:HB3	2.13	0.49
1:A:224:THR:C	1:A:225:GLN:O	2.53	0.49
1:C:261:LEU:HA	1:C:264:ILE:HG23	1.95	0.49
1:B:182:LEU:HD22	1:B:254:TYR:CE1	2.47	0.49
1:C:234:GLU:HG2	1:C:234:GLU:O	2.12	0.48
1:B:69:GLN:HG2	1:B:275:HIS:CE1	2.48	0.48
1:C:148:LEU:HD23	1:C:283:PRO:HB3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:231:LEU:HD12	1:C:232:GLY:N	2.29	0.47
1:B:193:LYS:HE3	1:B:246:ARG:HB3	1.95	0.47
1:B:73:TYR:HD2	1:B:91:LEU:HD21	1.78	0.47
1:A:277:TYR:CD1	1:A:277:TYR:C	2.92	0.47
1:A:197:ILE:CD1	1:A:243:GLU:HG3	2.44	0.47
1:C:129:GLN:HA	1:C:130:LEU:CB	2.45	0.47
1:C:185:GLN:OE1	1:C:254:TYR:HB2	2.15	0.47
1:B:62:ILE:O	1:B:62:ILE:HG13	2.15	0.47
1:B:274:ILE:O	1:B:275:HIS:CB	2.63	0.47
1:A:174:GLY:O	1:A:175:ARG:C	2.57	0.47
1:C:127:ASN:HB2	1:C:130:LEU:HA	1.96	0.47
1:A:204:LEU:HD12	1:A:204:LEU:O	2.15	0.46
1:C:120:THR:HG23	1:C:135:SER:HB2	1.97	0.46
1:A:105:ALA:O	1:A:109:THR:HG23	2.14	0.46
1:A:58:SER:O	1:A:135:SER:HA	2.16	0.46
1:A:140:THR:HG22	1:A:141:ALA:N	2.29	0.46
1:A:223:VAL:HG11	1:A:231:LEU:HD21	1.98	0.46
1:A:198:ARG:HE	1:A:198:ARG:HB2	1.58	0.46
1:B:99:PHE:C	1:B:99:PHE:CD1	2.94	0.46
1:B:174:GLY:O	1:B:178:LEU:HG	2.16	0.46
1:A:119:LEU:N	2:A:307:HOH:O	2.49	0.45
1:B:134:VAL:HG21	1:B:152:ILE:HD11	1.98	0.45
1:C:152:ILE:HD13	1:C:279:TYR:OH	2.16	0.45
1:B:70:ILE:O	1:B:70:ILE:CG1	2.62	0.45
1:B:200:ILE:HG12	1:B:239:MSE:CE	2.47	0.45
1:C:124:SER:CB	1:C:131:PRO:HD2	2.47	0.45
1:B:165:LYS:HD2	1:B:168:LYS:HE2	1.99	0.45
1:C:67:VAL:HG12	1:C:88:VAL:HG11	1.99	0.45
1:C:186:GLU:HA	1:C:250:PHE:CE1	2.52	0.44
1:A:193:LYS:HE3	1:A:246:ARG:CG	2.48	0.44
1:A:239:MSE:HG2	1:B:206:TYR:CE2	2.52	0.44
1:A:214:LYS:HB3	1:A:215:PRO:HD2	2.00	0.44
1:B:229:PHE:HA	1:B:236:LEU:HD11	1.99	0.44
1:A:140:THR:CG2	1:A:141:ALA:N	2.81	0.44
1:C:282:LYS:HB2	1:C:282:LYS:HE2	1.44	0.44
1:B:70:ILE:HD12	1:B:91:LEU:HD22	2.00	0.44
1:C:202:GLU:CD	1:B:245:THR:HG22	2.42	0.43
1:A:261:LEU:HD12	1:A:261:LEU:HA	1.60	0.43
1:B:82:GLY:C	1:B:84:ALA:N	2.76	0.43
1:B:167:LEU:O	1:B:171:ILE:HG13	2.18	0.43
1:C:173:LEU:HD21	1:B:74:ASN:ND2	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:LEU:HA	1:A:110:LEU:HD12	1.75	0.43
1:A:155:VAL:O	1:A:159:VAL:HG23	2.19	0.43
1:C:72:GLY:O	1:C:73:TYR:C	2.61	0.43
1:C:211:GLN:O	1:C:213:THR:N	2.50	0.43
1:A:103:PHE:HZ	1:A:119:LEU:HD22	1.82	0.43
1:C:231:LEU:HD23	1:B:222:ASP:HA	2.01	0.43
1:B:81:TYR:HB2	1:B:85:ALA:HB2	2.01	0.43
1:B:200:ILE:HG12	1:B:239:MSE:HE3	1.99	0.43
1:A:131:PRO:O	1:A:132:LEU:HD23	2.18	0.42
1:C:120:THR:HG23	1:C:135:SER:N	2.33	0.42
1:B:243:GLU:CD	1:B:243:GLU:C	2.86	0.42
1:A:225:GLN:O	1:A:226:ASP:HB2	2.19	0.42
1:A:140:THR:HG22	1:A:142:GLU:H	1.85	0.42
1:C:80:ILE:CG2	1:C:174:GLY:HA3	2.45	0.42
1:C:188:VAL:HG22	1:B:255:TYR:HB2	2.00	0.42
1:A:200:ILE:HG22	1:A:240:ILE:HD11	2.02	0.42
1:C:70:ILE:O	1:C:71:ALA:C	2.63	0.42
1:C:125:VAL:HB	1:C:126:LYS:H	1.49	0.42
1:C:231:LEU:O	1:C:232:GLY:C	2.61	0.42
1:A:275:HIS:CD2	2:A:314:HOH:O	2.72	0.41
1:C:261:LEU:HA	1:C:261:LEU:HD12	1.86	0.41
1:A:62:ILE:CG2	1:A:279:TYR:CD1	3.03	0.41
1:A:163:LEU:O	1:A:164:GLU:C	2.62	0.41
1:B:175:ARG:HG3	1:B:175:ARG:HH11	1.85	0.41
1:A:66:ASP:O	1:A:69:GLN:HB2	2.19	0.41
1:B:233:SER:HB2	2:B:302:HOH:O	2.19	0.41
1:C:129:GLN:CB	1:C:132:LEU:HD13	2.50	0.41
1:A:197:ILE:CG1	1:A:243:GLU:HG3	2.51	0.41
1:C:91:LEU:HD12	1:C:91:LEU:HA	1.92	0.41
1:A:236:LEU:HD23	1:A:236:LEU:HA	1.89	0.40
1:A:275:HIS:CA	2:A:318:HOH:O	2.68	0.40
1:A:275:HIS:N	2:A:318:HOH:O	2.53	0.40
1:C:153:GLN:O	1:C:153:GLN:HG3	2.14	0.40
1:C:128:GLN:O	1:C:129:GLN:CB	2.69	0.40
1:A:70:ILE:O	1:A:71:ALA:C	2.65	0.40
1:A:132:LEU:HD23	1:A:132:LEU:HA	1.88	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	210/240 (88%)	193 (92%)	14 (7%)	3 (1%)	9	7
1	B	218/240 (91%)	193 (88%)	19 (9%)	6 (3%)	4	2
1	C	200/240 (83%)	171 (86%)	22 (11%)	7 (4%)	3	1
All	All	628/720 (87%)	557 (89%)	55 (9%)	16 (2%)	4	3

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	124	SER
1	C	129	GLN
1	C	130	LEU
1	C	131	PRO
1	C	238	SER
1	B	116	PRO
1	B	239	MSE
1	B	275	HIS
1	C	267	LEU
1	B	69	GLN
1	A	123	PRO
1	C	125	VAL
1	B	265	GLU
1	A	243	GLU
1	C	240	ILE
1	B	70	ILE

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	173/208 (83%)	146 (84%)	27 (16%)	2	2
1	B	180/208 (86%)	155 (86%)	25 (14%)	3	3
1	C	161/208 (77%)	129 (80%)	32 (20%)	1	1
All	All	514/624 (82%)	430 (84%)	84 (16%)	2	2

All (84) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	THR
1	A	66	ASP
1	A	88	VAL
1	A	90	ASP
1	A	91	LEU
1	A	101	SER
1	A	108	GLU
1	A	110	LEU
1	A	130	LEU
1	A	133	THR
1	A	153	GLN
1	A	154	GLN
1	A	158	LYS
1	A	163	LEU
1	A	173	LEU
1	A	185	GLN
1	A	198	ARG
1	A	212	VAL
1	A	213	THR
1	A	214	LYS
1	A	221	GLU
1	A	234	GLU
1	A	251	SER
1	A	256	GLN
1	A	262	LEU
1	A	268	LYS
1	A	279	TYR
1	C	87	LYS
1	C	89	SER
1	C	93	GLU
1	C	104	SER
1	C	115	GLU

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Mol	Chain	Res	Type
1	C	118	LYS
1	C	120	THR
1	C	122	GLU
1	C	125	VAL
1	C	126	LYS
1	C	127	ASN
1	C	132	LEU
1	C	153	GLN
1	C	158	LYS
1	C	159	VAL
1	C	160	ASN
1	C	171	ILE
1	C	173	LEU
1	C	183	ARG
1	C	184	THR
1	C	187	VAL
1	C	197	ILE
1	C	209	GLN
1	C	213	THR
1	C	231	LEU
1	C	248	LEU
1	C	253	ASN
1	C	258	ARG
1	C	264	ILE
1	C	267	LEU
1	C	282	LYS
1	C	285	LEU
1	B	70	ILE
1	B	89	SER
1	B	108	GLU
1	B	112	ASN
1	B	121	ILE
1	B	140	THR
1	B	147	LYS
1	B	150	GLN
1	B	161	GLN
1	B	163	LEU
1	B	167	LEU
1	B	173	LEU
1	B	186	GLU
1	B	197	ILE
1	B	223	VAL

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Mol	Chain	Res	Type
1	B	238	SER
1	B	239	MSE
1	B	241	LYS
1	B	243	GLU
1	B	257	THR
1	B	258	ARG
1	B	261	LEU
1	B	262	LEU
1	B	264	ILE
1	B	274	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	177	ASN
1	A	192	GLN
1	A	219	GLN
1	A	253	ASN
1	C	74	ASN
1	C	78	ASN
1	C	179	GLN
1	C	205	GLN
1	C	209	GLN
1	C	253	ASN
1	C	259	GLN
1	C	275	HIS
1	B	74	ASN
1	B	113	GLN
1	B	185	GLN
1	B	209	GLN
1	B	253	ASN
1	B	259	GLN

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

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	212/240 (88%)	0.59	18 (8%) 16 18	27, 50, 78, 104	0
1	B	218/240 (90%)	0.66	18 (8%) 17 19	27, 50, 76, 113	0
1	C	204/240 (85%)	0.77	23 (11%) 10 11	18, 51, 85, 110	0
All	All	634/720 (88%)	0.67	59 (9%) 14 17	18, 50, 81, 113	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	130	LEU	6.0
1	B	56	TRP	5.8
1	B	221	GLU	5.6
1	A	141	ALA	5.6
1	C	269	PHE	5.4
1	A	285	LEU	5.1
1	C	210	ALA	4.5
1	C	229	PHE	4.5
1	C	130	LEU	4.3
1	B	164	GLU	4.1
1	B	274	ILE	4.0
1	C	215	PRO	3.9
1	A	142	GLU	3.5
1	C	270	ASP	3.4
1	C	248	LEU	3.4
1	C	240	ILE	3.3
1	C	242	HIS	3.3
1	B	220	THR	3.3
1	B	116	PRO	3.1
1	B	284	THR	3.0
1	A	138	GLY	2.8
1	C	255	TYR	2.8
1	C	236	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	212	VAL	2.7
1	A	268	LYS	2.7
1	A	124	SER	2.6
1	C	285	LEU	2.6
1	C	260	ASN	2.6
1	C	211	GLN	2.5
1	A	111	ASP	2.5
1	A	165	LYS	2.5
1	B	218	GLN	2.4
1	B	117	GLU	2.4
1	B	138	GLY	2.4
1	C	213	THR	2.3
1	A	136	TYR	2.3
1	C	82	GLY	2.3
1	B	244	ALA	2.3
1	A	83	GLN	2.3
1	C	90	ASP	2.3
1	B	142	GLU	2.2
1	B	219	GLN	2.2
1	B	238	SER	2.2
1	A	276	ALA	2.2
1	A	173	LEU	2.2
1	C	92	GLN	2.2
1	C	95	LEU	2.1
1	A	137	VAL	2.1
1	A	161	GLN	2.1
1	B	129	GLN	2.1
1	A	150	GLN	2.1
1	A	110	LEU	2.1
1	A	221	GLU	2.1
1	C	93	GLU	2.1
1	C	252	PRO	2.1
1	C	237	GLU	2.1
1	A	209	GLN	2.1
1	B	192	GLN	2.0
1	B	275	HIS	2.0

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

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.