



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:00 PM GMT

PDB ID : 4E2H
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
<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

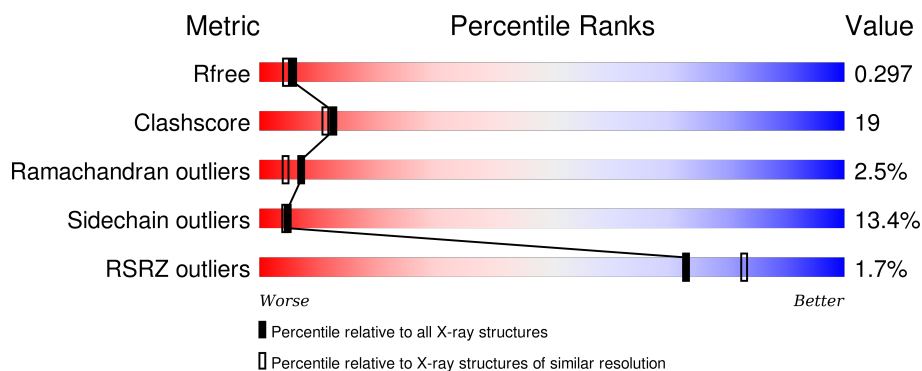
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	91344	1352 (2.38-2.34)
Clashscore	102246	1456 (2.38-2.34)
Ramachandran outliers	100387	1435 (2.38-2.34)
Sidechain outliers	100360	1436 (2.38-2.34)
RSRZ outliers	91569	1358 (2.38-2.34)

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	240	<div> <div>2%</div> <div>55% 30% 5% 10%</div> </div>
1	B	240	<div> <div>2%</div> <div>63% 25% • • 8%</div> </div>
1	C	240	<div> <div>2%</div> <div>53% 26% 5% • 13%</div> </div>

2 Entry composition

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
1	A	216	Total	C	N	O	Se	0	0	0
			1647	1035	285	324	3			
1	C	208	Total	C	N	O	Se	0	0	0
			1578	993	265	317	3			
1	B	222	Total	C	N	O	Se	0	0	0
			1704	1063	296	342	3			

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

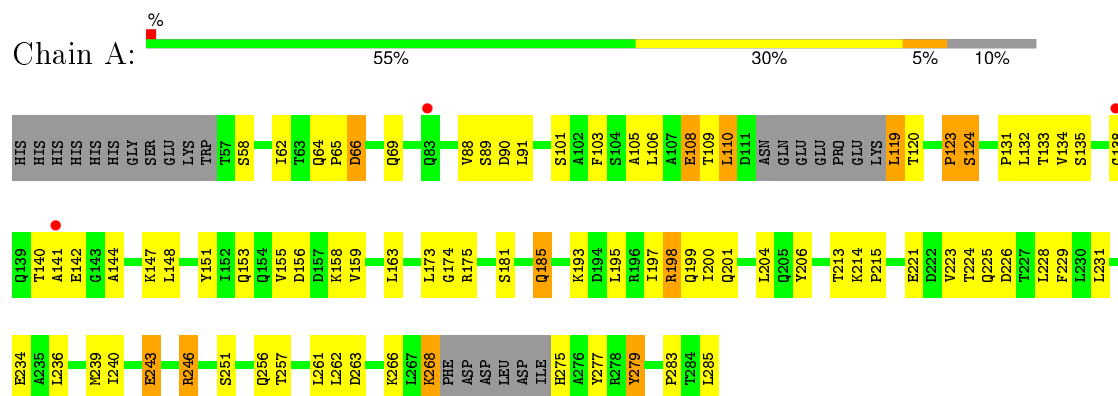
- Molecule 2 is water.

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

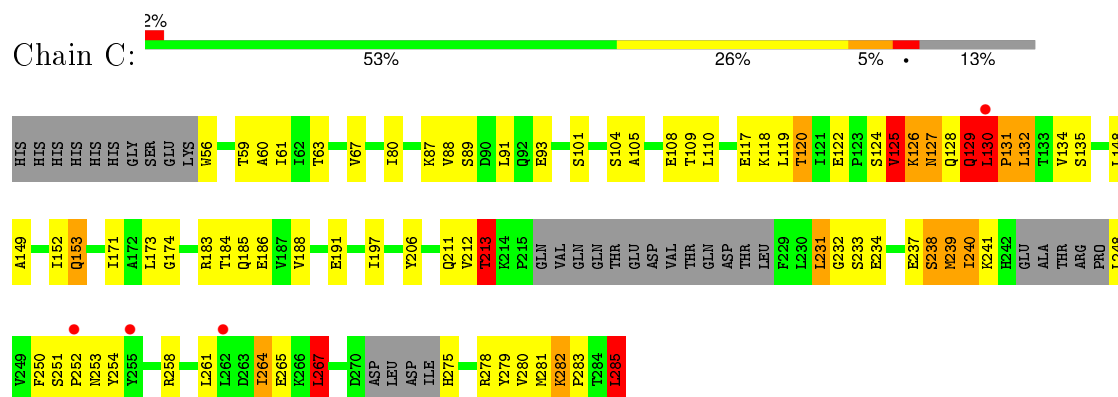
3 Residue-property plots [i](#)

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.

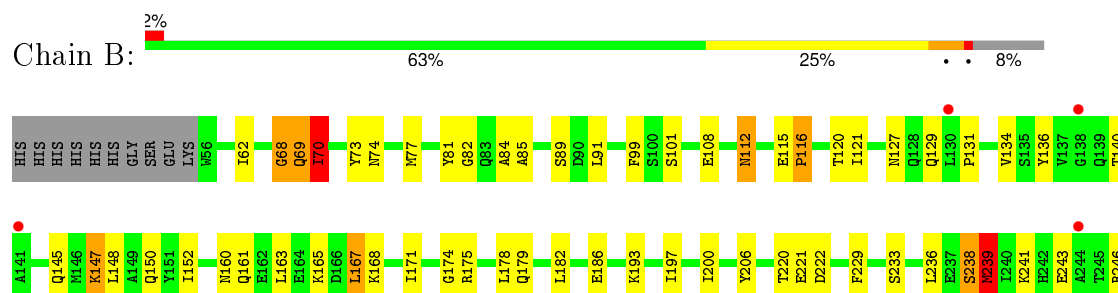
• Molecule 1: Chain length determinant protein



• Molecule 1: Chain length determinant protein



• Molecule 1: Chain length determinant protein



Y254	Y255	Q256	T257	R258
L261	L262	E265	K266	LEU
LYS	PHE	ASP	ASP	LEU
ASP	ASP	LEU	ASP	T274
H275	M281	K282	P283	T284
LEU				

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)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
R, R_{free}	0.245 , 0.302 0.237 , 0.297	Depositor DCC
R_{free} test set	1508 reflections (5.12%)	DCC
Wilson B-factor (Å ²)	33.2	Xtriage
Anisotropy	0.710	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 54.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 29489 reflections	Xtriage
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

Xtriage'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, 0.375 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	0.49	0/1663	0.66	0/2254
1	B	0.48	0/1722	0.66	0/2336
1	C	0.51	0/1595	0.70	4/2161 (0.2%)
All	All	0.49	0/4980	0.67	4/6751 (0.1%)

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 (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	130	LEU	N-CA-C	-6.33	93.92	111.00
1	C	285	LEU	CA-CB-CG	5.47	127.88	115.30
1	C	129	GLN	N-CA-C	-5.41	96.38	111.00
1	C	285	LEU	CB-CG-CD1	5.20	119.85	111.00

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

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Mol	Chain	Res	Type	Group
1	C	213	THR	Peptide
1	C	239	MSE	Peptide

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	1647	0	1618	65	0
1	B	1704	0	1652	55	0
1	C	1578	0	1498	79	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	188	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 (188) 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:O	1:C:285:LEU:HD13	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:B:70:ILE:O	1:B:70:ILE:HG13	1.92	0.69
2:C:324:HOH:O	1:B:239:MSE:SE	2.59	0.69
1:C:129:GLN:CA	1:C:130:LEU:CB	2.70	0.69
1:C:231:LEU:HG	1:C:234:GLU:OE1	1.95	0.67
1:C:80:ILE:HG23	1:C:174:GLY:HA3	1.75	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:C:126:LYS:O	1:C:126:LYS:HG3	1.95	0.64
1:C:240:ILE:O	2:C:313:HOH:O	2.14	0.64
1:A:256:GLN:HG3	1:A:257:THR:N	2.12	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
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:A:156:ASP:C	2:A:322:HOH:O	2.40	0.60
1:B:115:GLU:O	1:B:116:PRO:O	2.19	0.59
1:B:167:LEU:HG	1:B:167:LEU:O	2.02	0.59
1:B:236:LEU:O	1:B:239:MSE:HB3	2.03	0.59
1:B:120:THR:O	1:B:134:VAL:HA	2.03	0.58
1:A:120:THR:O	1:A: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:A:256:GLN:HG3	1:A:257:THR:H	1.68	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:91:LEU:HD23	1:B:91:LEU:O	2.04	0.56
1:A:138:GLY:HA3	1:A:144:ALA:HB2	1.87	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:C:59:THR:O	1:C:283:PRO:HA	2.04	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:C:108:GLU:HG3	1:C:109:THR:N	2.20	0.55
1:A:62:ILE:HG22	1:A:279:TYR:HD1	1.71	0.55
1:C:239:MSE:O	1:C:240:ILE:CB	2.54	0.55
1:A:105:ALA:HA	1:A:108:GLU:OE1	2.07	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
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:B:160:ASN:OD1	1:B:274:ILE:HG12	2.08	0.53
1:C:149:ALA:HB2	1:C:283:PRO:HG2	1.90	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:285:LEU:C	1:C:285:LEU:CD1	2.75	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:B:82:GLY:C	1:B:84:ALA:H	2.11	0.52
1:A:106:LEU:HD21	1:A:151:TYR:HD2	1.74	0.52
1:B:274:ILE:HG22	1:B:275:HIS:HD2	1.73	0.52
1:C:63:THR:HB	1:C:131:PRO:HB3	1.92	0.52
1:C:61:ILE:CG2	1:C:280:VAL:HB	2.37	0.52
1:C:59:THR:HG22	1:C:60:ALA:N	2.25	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: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:O	1:A:246:ARG:HG3	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:136:TYR:CE2	1:B:147:LYS:HG2	2.46	0.51
1:C:237:GLU:O	1:C:238:SER:CB	2.59	0.49
1:C:61:ILE:HG13	1:C:281:MSE:HE2	1.93	0.49
1:A:201:GLN:NE2	2:A:313:HOH:O	2.44	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:N	1:C:127:ASN:OD1	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: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:O	1:C:234:GLU:HG2	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
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: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:B:274:ILE:O	1:B:275:HIS:CB	2.63	0.47
1:C:185:GLN:OE1	1:C:254:TYR:HB2	2.15	0.47
1:A:285:LEU:HD12	1:A:285:LEU:C	2.35	0.47
1:B:62:ILE:O	1:B:62:ILE:HG13	2.15	0.47
1:C:127:ASN:HB2	1:C:130:LEU:HA	1.96	0.47
1:A:123:PRO:O	1:A:124:SER:C	2.53	0.46
1:C:120:THR:HG23	1:C:135:SER:HB2	1.97	0.46
1:B:112:ASN:HD22	1:B:112:ASN:C	2.18	0.46
1:A:204:LEU:HD12	1:A:204:LEU:O	2.15	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:223:VAL:HG11	1:A:231:LEU:HD21	1.98	0.46
1:A:140:THR:HG22	1:A:141:ALA:N	2.29	0.46
1:A:198:ARG:HE	1:A:198:ARG:HB2	1.58	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:B:70:ILE:O	1:B:70:ILE:CG1	2.62	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:200:ILE:HG12	1:B:239:MSE:CE	2.47	0.45
1:C:152:ILE:HD13	1:C:279:TYR:OH	2.16	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:224:THR:C	1:A:225:GLN:O	2.53	0.44
1:A:239:MSE:HG2	1:B:206:TYR:CE2	2.52	0.44
1:A:193:LYS:HE3	1:A:246:ARG:CG	2.48	0.44
1:B:229:PHE:HA	1:B:236:LEU:HD11	1.99	0.44
1:A:214:LYS:HB3	1:A:215:PRO:HD2	2.00	0.44
1:A:140:THR:CG2	1:A:141:ALA:N	2.81	0.44
1:C:282:LYS:HE2	1:C:282:LYS:HB2	1.44	0.44
1:B:70:ILE:HD12	1:B:91:LEU:HD22	2.00	0.44
1:B:167:LEU:O	1:B:171:ILE:HG13	2.18	0.43
1:A:261:LEU:HD12	1:A:261:LEU:HA	1.60	0.43
1:C:173:LEU:HD21	1:B:74:ASN:ND2	2.34	0.43
1:A:155:VAL:O	1:A:159:VAL:HG23	2.19	0.43
1:A:110:LEU:HD12	1:A:110:LEU:HA	1.75	0.43
1:C:211:GLN:O	1:C:213:THR:N	2.50	0.43
1:C:231:LEU:HD23	1:B:222:ASP:HA	2.01	0.43
1:A:103:PHE:HZ	1:A:119:LEU:HD22	1.82	0.43
1:B:200:ILE:HG12	1:B:239:MSE:HE3	1.99	0.43
1:B:81:TYR:HB2	1:B:85:ALA:HB2	2.01	0.43
1:C:120:THR:HG23	1:C:135:SER:N	2.33	0.42
1:A:131:PRO:O	1:A:132:LEU:HD23	2.18	0.42
1:A:277:TYR:CD1	1:A:277:TYR:C	2.92	0.42
1:A:225:GLN:O	1:A:226:ASP:HB2	2.19	0.42
1:C:80:ILE:CG2	1:C:174:GLY:HA3	2.45	0.42
1:A:140:THR:HG22	1:A:142:GLU:H	1.85	0.42
1:A:174:GLY:O	1:A:175:ARG:C	2.57	0.42
1:C:188:VAL:HG22	1:B:255:TYR:HB2	2.00	0.42
1:C:125:VAL:HB	1:C:126:LYS:H	1.49	0.42
1:A:200:ILE:HG22	1:A:240:ILE:HD11	2.02	0.42
1:C:261:LEU:HD12	1:C:261:LEU:HA	1.86	0.41
1:A:275:HIS:CD2	2:A:314:HOH:O	2.72	0.41
1:A:62:ILE:CG2	1:A:279:TYR:CD1	3.03	0.41
1:B:175:ARG:HH11	1:B:175:ARG:HG3	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:99:PHE:CD1	1:B:99:PHE:C	2.94	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:HA	1:A:236:LEU:HD23	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:HG3	1:C:153:GLN:O	2.14	0.40
1:C:128:GLN:O	1:C:129:GLN:CB	2.69	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%)	14	12
1	B	218/240 (91%)	193 (88%)	19 (9%)	6 (3%)	6	3
1	C	200/240 (83%)	171 (86%)	22 (11%)	7 (4%)	4	2
All	All	628/720 (87%)	557 (89%)	55 (9%)	16 (2%)	7	4

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

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Mol	Chain	Res	Type
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 ⓘ

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%)	148 (86%)	25 (14%)	4	3
1	B	180/208 (86%)	160 (89%)	20 (11%)	8	7
1	C	161/208 (77%)	137 (85%)	24 (15%)	4	3
All	All	514/624 (82%)	445 (87%)	69 (13%)	5	4

All (69) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	66	ASP
1	A	88	VAL
1	A	89	SER
1	A	90	ASP
1	A	91	LEU
1	A	101	SER
1	A	108	GLU
1	A	110	LEU
1	A	119	LEU
1	A	133	THR
1	A	153	GLN
1	A	158	LYS
1	A	163	LEU
1	A	173	LEU

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Mol	Chain	Res	Type
1	A	185	GLN
1	A	198	ARG
1	A	206	TYR
1	A	213	THR
1	A	221	GLU
1	A	234	GLU
1	A	246	ARG
1	A	251	SER
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	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	183	ARG
1	C	184	THR
1	C	197	ILE
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

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Mol	Chain	Res	Type
1	B	150	GLN
1	B	161	GLN
1	B	163	LEU
1	B	167	LEU
1	B	186	GLU
1	B	197	ILE
1	B	238	SER
1	B	239	MSE
1	B	241	LYS
1	B	243	GLU
1	B	258	ARG
1	B	261	LEU
1	B	262	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	253	ASN
1	C	275	HIS

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 [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 [i](#)

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

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%)	0.20	3 (1%) 78 87	27, 50, 78, 104	0
1	B	218/240 (90%)	0.26	4 (1%) 71 81	27, 50, 76, 113	0
1	C	204/240 (85%)	0.29	4 (1%) 68 79	18, 51, 85, 110	0
All	All	634/720 (88%)	0.25	11 (1%) 73 83	18, 50, 81, 113	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	255	TYR	4.3
1	C	252	PRO	3.4
1	B	138	GLY	3.3
1	C	262	LEU	3.3
1	B	141	ALA	2.5
1	C	130	LEU	2.5
1	A	83	GLN	2.4
1	A	138	GLY	2.2
1	B	130	LEU	2.2
1	B	244	ALA	2.1
1	A	141	ALA	2.1

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.