



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:37 AM GMT

PDB ID : 2HTK
Title : Structure of the Escherichia coli ClC chloride channel Y445A mutant and Fab complex
Authors : Accardi, A.; Lobet, S.; Williams, C.; Miller, C.; Dutzler, R.
Deposited on : 2006-07-26
Resolution : 3.41 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

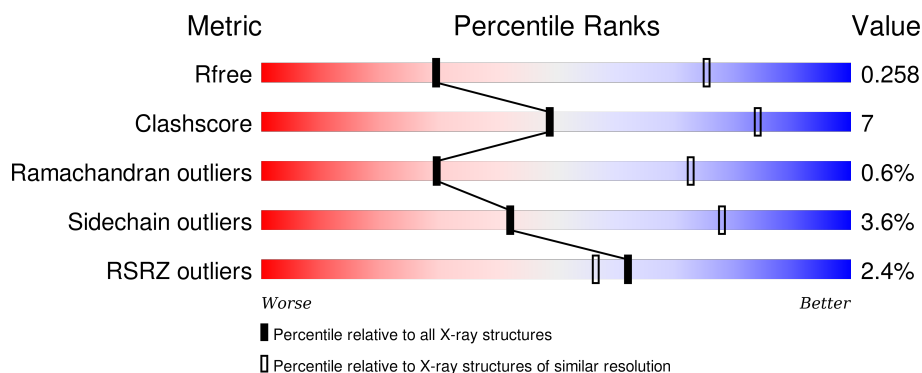
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.41 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	1049 (3.52-3.32)
Clashscore	102246	1032 (3.50-3.34)
Ramachandran outliers	100387	1002 (3.50-3.34)
Sidechain outliers	100360	1003 (3.50-3.34)
RSRZ outliers	91569	1054 (3.52-3.32)

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	473	<div> <div>75%</div> <div>17%</div> <div>• 6%</div> </div>
1	B	473	<div> <div>74%</div> <div>18%</div> <div>• 7%</div> </div>
2	C	221	<div> <div>3%</div> <div>89%</div> <div>11%</div> </div>
2	E	221	<div> <div>4%</div> <div>82%</div> <div>17%</div> <div>•</div> </div>
3	D	211	<div> <div>2%</div> <div>85%</div> <div>14%</div> </div>

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Mol	Chain	Length	Quality of chain
3	F	211	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	BR	A	474	-	-	X	X
4	BR	B	474	-	-	X	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13211 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 H(+)/Cl(-) exchange transporter clcA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	444	Total	C	N	O	S	0	0	0
			3326	2184	560	562	20			
1	B	441	Total	C	N	O	S	0	0	0
			3297	2168	553	556	20			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	445	ALA	TYR	ENGINEERED	UNP P37019
B	445	ALA	TYR	ENGINEERED	UNP P37019

- Molecule 2 is a protein called Fab fragment, Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	221	Total	C	N	O	S	0	0	0
			1672	1077	274	315	6			
2	E	221	Total	C	N	O	S	0	0	0
			1672	1077	274	315	6			

- Molecule 3 is a protein called Fab fragment, Light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	211	Total	C	N	O	S	0	0	0
			1621	1008	271	334	8			
3	F	211	Total	C	N	O	S	0	0	0
			1621	1008	271	334	8			

- Molecule 4 is BROMIDE ION (three-letter code: BR) (formula: Br).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Br	0	0
			1	1		

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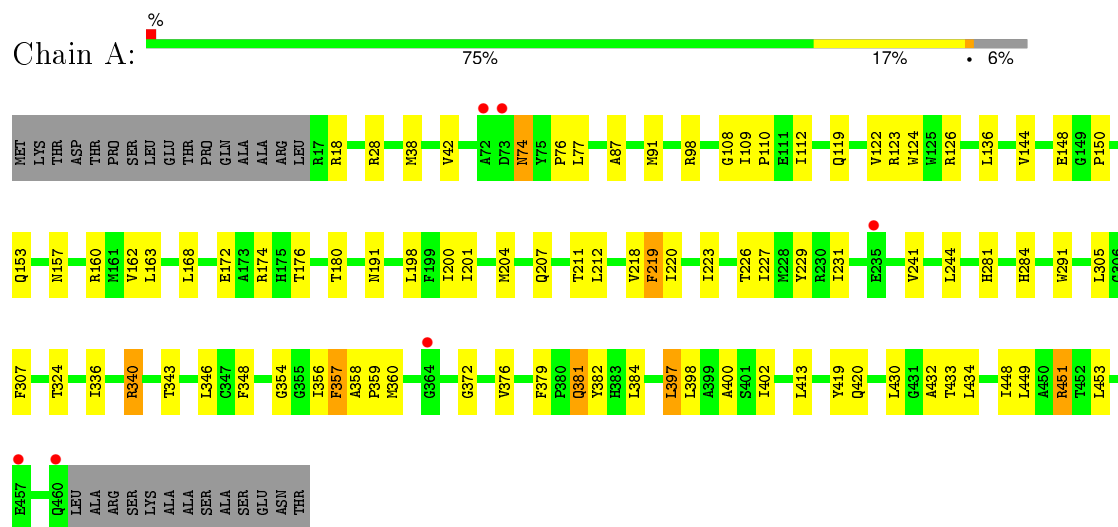
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Br	0	0
			1	1		

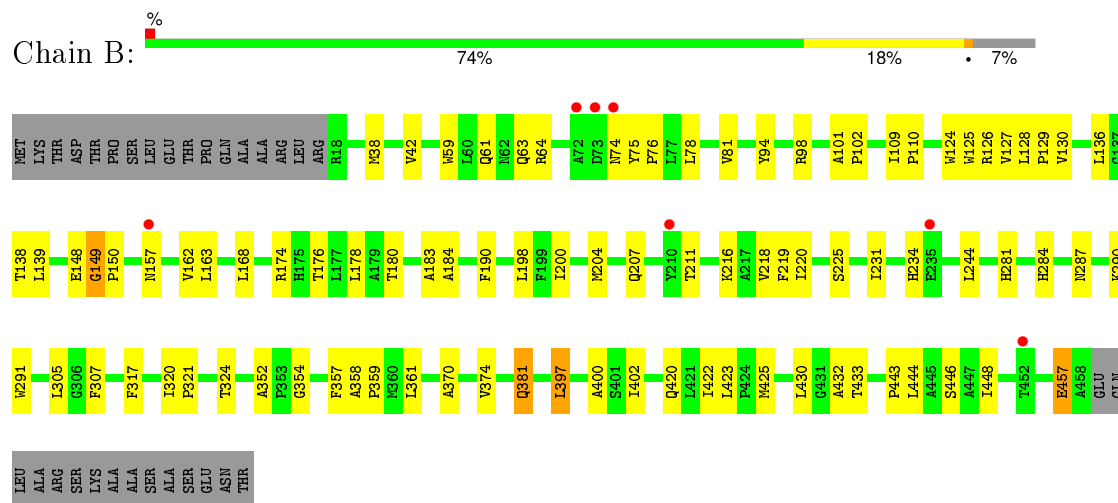
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

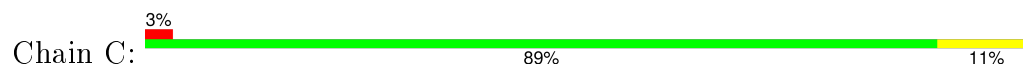
- Molecule 1: H(+)/Cl(-) exchange transporter clcA



- Molecule 1: H(+)/Cl(-) exchange transporter clcA

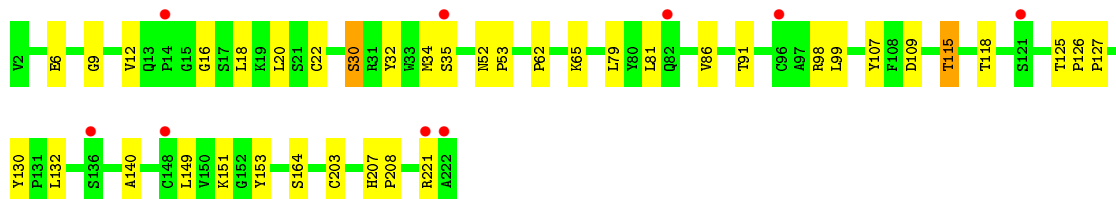
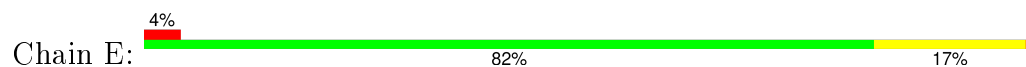


- Molecule 2: Fab fragment, Heavy chain

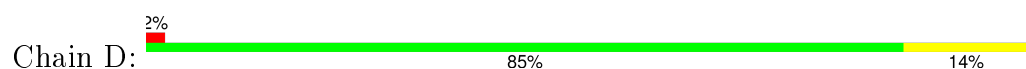




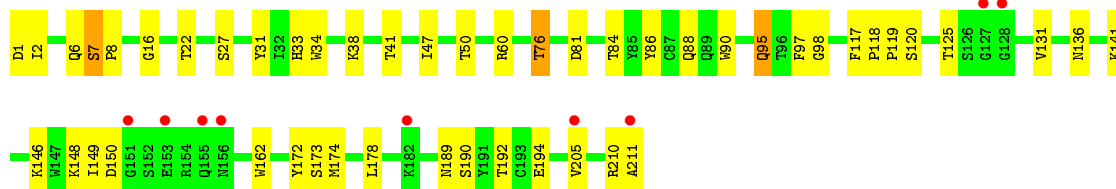
- Molecule 2: Fab fragment, Heavy chain



- Molecule 3: Fab fragment, Light chain



- Molecule 3: Fab fragment, Light chain



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	232.07Å 98.14Å 171.07Å 90.00° 131.56° 90.00°	Depositor
Resolution (Å)	40.00 – 3.41 19.96 – 3.40	Depositor EDS
% Data completeness (in resolution range)	96.9 (40.00-3.41) 97.0 (19.96-3.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.20 (at 3.44Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.266 , 0.276 0.246 , 0.258	Depositor DCC
R_{free} test set	1921 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	113.4	Xtriage
Anisotropy	0.467	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 69.0	EDS
Estimated twinning fraction	0.000 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 39582 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13211	wwPDB-VP
Average B, all atoms (Å ²)	101.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.94% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BR

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	0.35	0/3397	0.50	0/4610
1	B	0.35	0/3368	0.49	0/4572
2	C	0.35	0/1721	0.52	0/2355
2	E	0.35	0/1721	0.51	0/2355
3	D	0.34	0/1660	0.49	0/2257
3	F	0.33	0/1660	0.50	0/2257
All	All	0.35	0/13527	0.50	0/18406

There are no bond length outliers.

There are no bond angle outliers.

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	3326	0	3480	62	0
1	B	3297	0	3453	56	0
2	C	1672	0	1654	11	0
2	E	1672	0	1654	20	0
3	D	1621	0	1546	18	0
3	F	1621	0	1546	36	0
4	A	1	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	1	0	0	2	0
All	All	13211	0	13333	190	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 7.

All (190) 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:F:7:SER:HB3	3:F:8:PRO:HD3	1.42	0.98
1:A:381:GLN:HE21	1:A:381:GLN:H	0.95	0.91
1:B:381:GLN:HE21	1:B:381:GLN:H	1.15	0.90
1:A:148:GLU:HG2	1:A:357:PHE:HB3	1.63	0.80
1:B:148:GLU:CD	1:B:148:GLU:H	1.85	0.79
3:D:194:GLU:HG2	3:D:205:VAL:HG12	1.64	0.79
1:A:381:GLN:N	1:A:381:GLN:HE21	1.79	0.76
1:B:150:PRO:HD3	1:B:354:GLY:HA2	1.69	0.74
3:D:148:LYS:HB2	3:D:192:THR:OG1	1.87	0.74
1:B:400:ALA:HB2	1:B:432:ALA:HB1	1.70	0.72
1:B:78:LEU:HA	1:B:81:VAL:HG22	1.71	0.71
3:F:7:SER:HB2	3:F:22:THR:HB	1.73	0.70
3:F:146:LYS:HB3	3:F:194:GLU:HB2	1.73	0.69
3:F:16:GLY:HA2	3:F:76:THR:HG23	1.73	0.68
3:F:6:GLN:HE22	3:F:86:TYR:HA	1.59	0.67
1:A:200:ILE:HA	1:A:204:MET:HB2	1.76	0.67
1:A:123:ARG:HE	1:A:126:ARG:HD2	1.58	0.67
2:E:12:VAL:HG11	2:E:18:LEU:HB3	1.76	0.67
2:C:34:MET:HB3	2:C:79:LEU:HD22	1.76	0.67
1:A:150:PRO:HD3	1:A:354:GLY:HA2	1.76	0.67
1:B:38:MET:O	1:B:42:VAL:HG23	1.95	0.67
1:A:74:ASN:HD22	1:A:76:PRO:HD2	1.60	0.66
3:D:7:SER:HB3	3:D:8:PRO:HD3	1.76	0.66
1:B:381:GLN:N	1:B:381:GLN:HE21	1.91	0.66
1:A:124:TRP:HA	1:A:157:ASN:HD22	1.61	0.66
1:A:176:THR:O	1:A:180:THR:HG23	1.95	0.66
1:A:381:GLN:NE2	1:A:381:GLN:H	1.80	0.65
1:A:400:ALA:HB2	1:A:432:ALA:HB1	1.79	0.64
1:B:200:ILE:HA	1:B:204:MET:HB2	1.79	0.63
1:A:148:GLU:H	1:A:148:GLU:CD	2.02	0.63
3:F:7:SER:HB3	3:F:8:PRO:CD	2.22	0.63
3:D:6:GLN:HE22	3:D:86:TYR:HA	1.63	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:VAL:HG21	1:A:343:THR:HB	1.80	0.62
1:B:42:VAL:HG22	1:B:162:VAL:HG21	1.81	0.62
1:B:176:THR:O	1:B:180:THR:HG23	2.00	0.62
3:D:95:GLN:H	3:D:95:GLN:CD	2.04	0.61
2:C:127:PRO:HB3	2:C:153:TYR:HB3	1.84	0.59
2:E:91:THR:HG23	2:E:118:THR:HA	1.86	0.58
1:A:430:LEU:HD21	1:B:220:ILE:HG12	1.84	0.58
1:A:38:MET:HG3	1:A:168:LEU:HD11	1.86	0.57
1:B:180:THR:HG22	1:B:218:VAL:HA	1.87	0.57
3:F:194:GLU:HG2	3:F:205:VAL:HG12	1.86	0.57
3:D:95:GLN:N	3:D:95:GLN:CD	2.58	0.56
1:B:281:HIS:HA	1:B:284:HIS:CE1	2.41	0.56
1:B:320:ILE:HB	1:B:321:PRO:HD3	1.87	0.56
1:A:148:GLU:HG2	1:A:357:PHE:CB	2.35	0.56
1:A:74:ASN:ND2	1:A:76:PRO:HD2	2.21	0.55
1:A:360:MET:HG2	1:A:397:LEU:HD12	1.87	0.55
3:F:189:ASN:HD21	3:F:211:ALA:H	1.54	0.55
2:E:6:GLU:HA	2:E:22:CYS:HA	1.89	0.55
2:E:130:TYR:HB3	3:F:120:SER:OG	2.06	0.55
3:F:95:GLN:H	3:F:95:GLN:CD	2.10	0.55
1:A:112:ILE:HD11	1:A:153:GLN:HG3	1.90	0.54
1:A:227:ILE:O	1:A:231:ILE:HG12	2.07	0.54
3:D:1:ASP:HB3	3:D:94:PRO:HD2	1.88	0.54
1:B:98:ARG:HD2	1:B:291:TRP:CE3	2.42	0.54
3:F:7:SER:CB	3:F:8:PRO:HD3	2.27	0.53
2:E:98:ARG:NH1	2:E:109:ASP:OD2	2.41	0.53
1:B:75:TYR:HB3	1:B:76:PRO:HD3	1.91	0.53
2:E:132:LEU:HD11	2:E:149:LEU:HB2	1.92	0.52
1:A:241:VAL:HG11	1:A:324:THR:HG21	1.93	0.52
1:A:110:PRO:HG2	4:A:474:BR:BR	2.65	0.51
1:A:74:ASN:HB3	1:A:77:LEU:HB3	1.91	0.51
1:B:94:TYR:CZ	1:B:352:ALA:HB2	2.45	0.51
1:A:356:ILE:HG23	1:A:360:MET:CE	2.40	0.51
2:E:34:MET:HB3	2:E:79:LEU:HD22	1.93	0.51
1:A:18:ARG:HH11	1:B:457:GLU:HB3	1.76	0.51
2:E:52:ASN:HB2	2:E:53:PRO:CD	2.41	0.51
3:F:60:ARG:NH2	3:F:81:ASP:OD2	2.44	0.51
1:A:398:LEU:HD22	1:A:402:ILE:HD12	1.93	0.51
2:E:127:PRO:HB3	2:E:153:TYR:HB3	1.93	0.51
1:B:61:GLN:HG2	1:B:64:ARG:HH21	1.76	0.50
3:F:119:PRO:HD3	3:F:131:VAL:HG22	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:148:GLU:CG	1:A:357:PHE:HB3	2.39	0.50
3:F:189:ASN:HD22	3:F:210:ARG:HB2	1.77	0.50
1:B:101:ALA:HB3	1:B:130:VAL:HG11	1.94	0.50
3:F:88:GLN:HB2	3:F:97:PHE:CD1	2.47	0.49
1:A:220:ILE:HG12	1:B:430:LEU:HD21	1.92	0.49
1:B:124:TRP:HA	1:B:157:ASN:HD22	1.77	0.49
3:D:16:GLY:HA2	3:D:76:THR:HG23	1.94	0.49
1:B:422:ILE:HA	1:B:425:MET:HE2	1.93	0.49
2:C:91:THR:HG23	2:C:118:THR:HA	1.94	0.49
2:E:107:TYR:HB3	3:F:33:HIS:CD2	2.47	0.49
3:D:189:ASN:HD21	3:D:211:ALA:H	1.61	0.49
1:A:160:ARG:HH12	1:A:174:ARG:HD2	1.78	0.48
3:F:95:GLN:N	3:F:95:GLN:OE1	2.47	0.48
1:A:172:GLU:HG3	1:A:212:LEU:O	2.13	0.48
1:B:127:VAL:HB	1:B:157:ASN:ND2	2.29	0.48
2:E:35:SER:HB2	2:E:99:LEU:HD11	1.95	0.48
3:F:90:TRP:CZ2	3:F:95:GLN:NE2	2.81	0.48
3:D:124:LEU:C	3:D:126:SER:H	2.17	0.48
1:B:138:THR:HG21	1:B:352:ALA:HB1	1.95	0.48
1:A:336:ILE:O	1:A:340:ARG:HG3	2.14	0.47
3:F:6:GLN:HE21	3:F:98:GLY:HA3	1.79	0.47
3:F:95:GLN:N	3:F:95:GLN:CD	2.67	0.47
1:A:376:VAL:HG22	1:A:384:LEU:HB2	1.96	0.47
1:A:98:ARG:HD2	1:A:291:TRP:CE3	2.50	0.47
1:A:122:VAL:HB	1:A:160:ARG:HG2	1.97	0.47
3:F:34:TRP:HB2	3:F:47:ILE:HB	1.96	0.47
3:F:150:ASP:HA	3:F:190:SER:HB3	1.97	0.47
3:D:7:SER:CB	3:D:8:PRO:HD3	2.45	0.46
3:F:141:LYS:HB3	3:F:172:TYR:CD1	2.50	0.46
1:A:191:ASN:HB2	1:A:229:TYR:CE2	2.50	0.46
3:D:2:ILE:HD12	3:D:27:SER:HB2	1.97	0.46
2:E:32:TYR:CE2	2:E:98:ARG:HD3	2.50	0.46
1:B:234:HIS:H	1:B:234:HIS:CD2	2.34	0.46
1:B:125:TRP:CD1	1:B:126:ARG:HG3	2.51	0.46
1:A:305:LEU:C	1:A:307:PHE:H	2.20	0.46
1:A:348:PHE:CD1	1:A:356:ILE:HB	2.50	0.45
3:F:148:LYS:HB2	3:F:192:THR:OG1	2.15	0.45
3:F:149:ILE:HD11	3:F:178:LEU:HD21	1.97	0.45
3:F:7:SER:CB	3:F:22:THR:HB	2.42	0.45
1:A:379:PHE:HB3	1:A:382:TYR:CD1	2.52	0.45
3:F:90:TRP:CG	3:F:95:GLN:HB3	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:110:PRO:HG2	4:B:474:BR:BR	2.72	0.45
1:B:448:ILE:HG21	4:B:474:BR:BR	2.72	0.45
1:B:190:PHE:HE2	1:B:317:PHE:HZ	1.65	0.45
2:E:30:SER:C	2:E:32:TYR:H	2.20	0.45
2:C:69:ILE:HB	2:C:82:GLN:HB2	1.97	0.45
2:C:134:PRO:HD2	2:C:196:TRP:CH2	2.52	0.45
1:A:358:ALA:HB3	1:A:359:PRO:HD3	1.98	0.44
1:B:443:PRO:HB2	1:B:446:SER:HB2	1.98	0.44
1:A:451:ARG:HB3	1:A:451:ARG:HH11	1.83	0.44
1:A:42:VAL:HG22	1:A:162:VAL:HG21	1.99	0.44
2:E:6:GLU:HB3	2:E:22:CYS:HB2	2.00	0.44
2:C:52:ASN:HB2	2:C:53:PRO:HD2	2.00	0.44
1:A:28:ARG:HD3	1:B:443:PRO:HG2	2.00	0.44
3:F:162:TRP:CD1	3:F:174:MET:HG3	2.52	0.44
1:A:180:THR:HG22	1:A:218:VAL:HA	1.98	0.44
1:A:356:ILE:HG23	1:A:360:MET:HE2	1.98	0.44
1:A:281:HIS:HA	1:A:284:HIS:CE1	2.53	0.44
1:A:223:ILE:CD1	1:B:430:LEU:HD22	2.48	0.44
2:C:195:SER:O	2:C:199:GLU:HB3	2.17	0.44
2:C:29:TYR:OH	2:C:79:LEU:HB2	2.18	0.44
1:B:128:LEU:HB2	1:B:129:PRO:CD	2.48	0.44
2:E:221:ARG:HH22	3:F:120:SER:HA	1.83	0.43
1:A:219:PHE:HB3	1:B:430:LEU:CD1	2.48	0.43
2:C:30:SER:C	2:C:32:TYR:H	2.20	0.43
3:F:117:PHE:HA	3:F:118:PRO:HD3	1.86	0.43
1:B:287:ASN:HD22	1:B:290:LYS:HG3	1.83	0.43
1:B:59:TRP:O	1:B:63:GLN:HG2	2.17	0.43
2:E:9:GLY:H	2:E:115:THR:HG21	1.84	0.43
1:A:112:ILE:HG13	1:A:153:GLN:HA	2.01	0.43
3:D:150:ASP:HA	3:D:190:SER:HB3	1.99	0.43
1:A:413:LEU:HD11	1:A:419:TYR:HA	2.00	0.43
3:D:162:TRP:CD1	3:D:174:MET:HG3	2.53	0.43
1:B:358:ALA:HB3	1:B:359:PRO:HD3	2.00	0.43
1:B:148:GLU:O	1:B:149:GLY:C	2.56	0.43
1:A:109:ILE:N	1:A:110:PRO:CD	2.82	0.43
3:F:31:TYR:HA	3:F:50:THR:OG1	2.19	0.43
3:F:2:ILE:HD12	3:F:27:SER:HB2	2.00	0.43
3:D:90:TRP:CG	3:D:95:GLN:HB3	2.54	0.43
1:A:434:LEU:HD23	1:B:216:LYS:HD3	2.00	0.43
1:A:449:LEU:O	1:A:453:LEU:HB2	2.19	0.43
1:A:198:LEU:HD11	1:B:198:LEU:HD21	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:20:LEU:HD12	2:E:81:LEU:HD23	2.01	0.42
1:B:184:ALA:HB1	1:B:225:SER:HB3	2.00	0.42
3:F:38:LYS:O	3:F:41:THR:HG22	2.19	0.42
1:B:163:LEU:HD12	1:B:168:LEU:HB2	2.01	0.42
1:A:223:ILE:HD12	1:B:430:LEU:HD22	2.01	0.42
1:B:98:ARG:NH2	1:B:102:PRO:HB3	2.34	0.42
1:B:38:MET:HG3	1:B:168:LEU:HD11	2.01	0.42
1:A:38:MET:O	1:A:42:VAL:HG23	2.19	0.42
1:A:372:GLY:O	1:A:376:VAL:HG23	2.18	0.42
1:A:119:GLN:HB3	1:A:453:LEU:HD11	2.01	0.42
1:B:381:GLN:NE2	1:B:381:GLN:H	1.98	0.42
2:C:127:PRO:CB	2:C:153:TYR:HB3	2.50	0.42
3:F:90:TRP:CD2	3:F:95:GLN:HB3	2.55	0.42
1:B:358:ALA:HA	1:B:361:LEU:HD12	2.01	0.42
1:A:87:ALA:O	1:A:91:MET:HG3	2.20	0.41
1:A:226:THR:HG21	1:B:423:LEU:HD11	2.02	0.41
3:F:136:ASN:HD22	3:F:173:SER:HB3	1.85	0.41
2:E:16:GLY:O	2:E:86:VAL:HG23	2.20	0.41
2:E:125:THR:HA	2:E:126:PRO:HD2	1.90	0.41
3:D:6:GLN:HE21	3:D:98:GLY:HA3	1.85	0.41
1:B:109:ILE:N	1:B:110:PRO:CD	2.83	0.41
3:D:148:LYS:HB2	3:D:192:THR:HG1	1.83	0.41
1:B:183:ALA:HB2	1:B:200:ILE:HD13	2.03	0.41
3:F:189:ASN:ND2	3:F:210:ARG:HB2	2.34	0.41
1:A:108:GLY:O	1:A:112:ILE:HG12	2.20	0.41
2:E:207:HIS:HA	2:E:208:PRO:HD2	1.93	0.41
2:C:16:GLY:O	2:C:86:VAL:HG23	2.20	0.41
1:A:448:ILE:HG21	4:A:474:BR:BR	2.76	0.41
1:B:397:LEU:HA	1:B:397:LEU:HD22	1.95	0.41
1:B:305:LEU:C	1:B:307:PHE:H	2.24	0.41
3:D:95:GLN:N	3:D:95:GLN:OE1	2.54	0.40
1:B:231:ILE:CG2	1:B:231:ILE:O	2.69	0.40
1:A:163:LEU:HD12	1:A:168:LEU:HB2	2.03	0.40
1:B:370:ALA:O	1:B:374:VAL:HG23	2.21	0.40
1:B:174:ARG:O	1:B:178:LEU:HB2	2.21	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	442/473 (93%)	415 (94%)	26 (6%)	1 (0%)	52	87
1	B	439/473 (93%)	408 (93%)	30 (7%)	1 (0%)	52	87
2	C	219/221 (99%)	202 (92%)	16 (7%)	1 (0%)	34	75
2	E	219/221 (99%)	197 (90%)	18 (8%)	4 (2%)	11	51
3	D	209/211 (99%)	187 (90%)	21 (10%)	1 (0%)	34	75
3	F	209/211 (99%)	191 (91%)	16 (8%)	2 (1%)	19	64
All	All	1737/1810 (96%)	1600 (92%)	127 (7%)	10 (1%)	30	73

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	E	65	LYS
2	E	62	PRO
2	E	140	ALA
2	C	62	PRO
3	F	7	SER
3	F	76	THR
1	B	149	GLY
3	D	125	THR
2	E	164	SER
1	A	201	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	334/357 (94%)	320 (96%)	14 (4%)	36	74
1	B	331/357 (93%)	315 (95%)	16 (5%)	31	71
2	C	181/181 (100%)	175 (97%)	6 (3%)	45	80
2	E	181/181 (100%)	177 (98%)	4 (2%)	60	86
3	D	185/185 (100%)	179 (97%)	6 (3%)	46	80
3	F	185/185 (100%)	181 (98%)	4 (2%)	60	86
All	All	1397/1446 (97%)	1347 (96%)	50 (4%)	42	78

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	74	ASN
1	A	136	LEU
1	A	207	GLN
1	A	211	THR
1	A	219	PHE
1	A	244	LEU
1	A	340	ARG
1	A	346	LEU
1	A	357	PHE
1	A	381	GLN
1	A	397	LEU
1	A	420	GLN
1	A	433	THR
1	A	451	ARG
1	B	74	ASN
1	B	136	LEU
1	B	139	LEU
1	B	207	GLN
1	B	211	THR
1	B	219	PHE
1	B	244	LEU
1	B	324	THR
1	B	357	PHE
1	B	381	GLN
1	B	397	LEU
1	B	402	ILE
1	B	420	GLN
1	B	433	THR
1	B	444	LEU
1	B	457	GLU

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Mol	Chain	Res	Type
2	C	98	ARG
2	C	115	THR
2	C	118	THR
2	C	151	LYS
2	C	188	SER
2	C	203	CYS
3	D	5	THR
3	D	22	THR
3	D	46	TRP
3	D	59	VAL
3	D	95	GLN
3	D	199	THR
2	E	30	SER
2	E	115	THR
2	E	151	LYS
2	E	203	CYS
3	F	1	ASP
3	F	84	THR
3	F	95	GLN
3	F	125	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	ASN
1	A	119	GLN
1	A	153	GLN
1	A	157	ASN
1	A	207	GLN
1	A	270	ASN
1	A	277	GLN
1	A	287	ASN
1	A	327	ASN
1	A	381	GLN
1	A	420	GLN
1	A	437	GLN
1	B	62	ASN
1	B	74	ASN
1	B	157	ASN
1	B	207	GLN
1	B	234	HIS
1	B	270	ASN

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Mol	Chain	Res	Type
1	B	277	GLN
1	B	287	ASN
1	B	327	ASN
1	B	381	GLN
1	B	420	GLN
1	B	437	GLN
3	D	6	GLN
3	D	36	GLN
3	D	136	ASN
3	D	189	ASN
2	E	172	HIS
3	F	6	GLN
3	F	136	ASN
3	F	137	ASN
3	F	189	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 ⓘ

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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

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	444/473 (93%)	-0.23	6 (1%) 78 72	40, 96, 146, 200	0
1	B	441/473 (93%)	-0.19	7 (1%) 74 69	44, 97, 154, 200	0
2	C	221/221 (100%)	-0.05	6 (2%) 58 52	53, 90, 146, 197	0
2	E	221/221 (100%)	0.03	9 (4%) 41 36	48, 99, 157, 200	0
3	D	211/211 (100%)	0.05	5 (2%) 62 57	48, 109, 157, 179	0
3	F	211/211 (100%)	0.07	9 (4%) 39 34	40, 89, 147, 181	0
All	All	1749/1810 (96%)	-0.09	42 (2%) 62 57	40, 97, 154, 200	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	73	ASP	7.2
2	C	139	ALA	5.6
2	E	35	SER	4.9
1	A	235	GLU	4.7
3	F	127	GLY	4.6
1	B	235	GLU	3.8
2	C	137	ALA	3.6
2	C	136	SER	3.4
3	D	68	THR	3.2
2	C	140	ALA	3.2
3	F	155	GLN	3.2
3	F	205	VAL	3.2
3	F	151	GLY	3.2
2	E	136	SER	3.1
2	C	138	ALA	3.0
2	E	148	CYS	2.9
2	E	82	GLN	2.9
1	B	210	TYR	2.8
1	A	73	ASP	2.7

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Mol	Chain	Res	Type	RSRZ
3	D	153	GLU	2.7
3	F	128	GLY	2.7
3	D	87	CYS	2.7
2	C	222	ALA	2.7
3	D	39	SER	2.6
3	D	1	ASP	2.6
1	A	364	GLY	2.5
1	A	457	GLU	2.5
2	E	14	PRO	2.5
1	B	72	ALA	2.5
2	E	121	SER	2.5
3	F	211	ALA	2.5
2	E	222	ALA	2.4
1	B	74	ASN	2.4
2	E	96	CYS	2.3
1	A	72	ALA	2.3
3	F	182	LYS	2.2
2	E	221	ARG	2.2
1	B	157	ASN	2.2
1	A	460	GLN	2.1
3	F	153	GLU	2.0
3	F	156	ASN	2.0
1	B	452	THR	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	BR	B	474	1/1	0.92	0.58	10.67	56,56,56,56	0
4	BR	A	474	1/1	0.88	0.62	10.57	56,56,56,56	0

6.5 Other polymers [i](#)

There are no such residues in this entry.