



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 01:50 am GMT

PDB ID : 4KKL
Title : Structure of the E148A mutant of CLC-ec1 delta NC construct in 100mM fluoride
Authors : Stockbridge, R.B.; Lim, H.-H.; Miller, C.
Deposited on : 2013-05-06
Resolution : 2.85 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

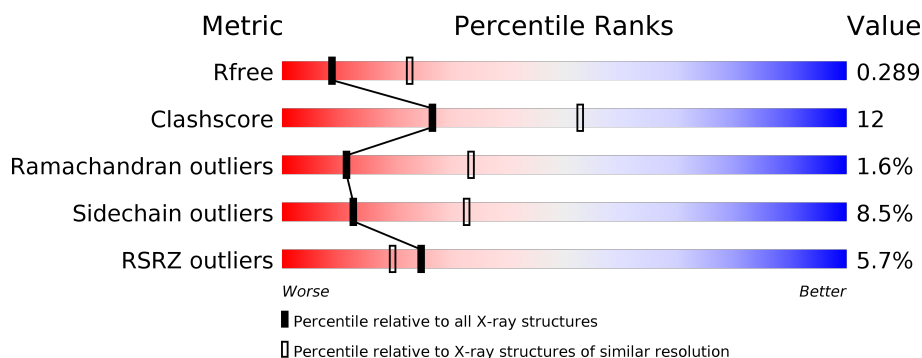
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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}	100719	2469 (2.90-2.82)
Clashscore	112137	2749 (2.90-2.82)
Ramachandran outliers	110173	2687 (2.90-2.82)
Sidechain outliers	110143	2690 (2.90-2.82)
RSRZ outliers	101464	2487 (2.90-2.82)

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	446	<div> <div>4%</div> <div> <div></div> <div>69%</div> <div>26%</div> <div>.</div> </div> </div>
1	B	446	<div> <div>6%</div> <div> <div></div> <div>60%</div> <div>35%</div> <div>.</div> </div> </div>
2	C	222	<div> <div>7%</div> <div> <div></div> <div>75%</div> <div>23%</div> <div>.</div> </div> </div>
2	E	222	<div> <div>5%</div> <div> <div></div> <div>73%</div> <div>24%</div> <div>.</div> </div> </div>
3	D	211	<div> <div>7%</div> <div> <div></div> <div>63%</div> <div>35%</div> <div>.</div> </div> </div>
3	F	211	<div> <div>5%</div> <div> <div></div> <div>74%</div> <div>24%</div> <div>.</div> </div> </div>

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	F	A	501	-	-	-	X
4	F	A	502	-	-	-	X
4	F	B	501	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13219 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
			3329	2188	560	561	20			
1	B	441	Total	C	N	O	S	0	0	0
			3300	2172	553	555	20			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	16	MET	-	INITIATING METHIONINE	UNP P37019
A	148	ALA	GLU	ENGINEERED MUTATION	UNP P37019
A	461	LYS	-	EXPRESSION TAG	UNP P37019
B	16	MET	-	INITIATING METHIONINE	UNP P37019
B	148	ALA	GLU	ENGINEERED MUTATION	UNP P37019
B	461	LYS	-	EXPRESSION TAG	UNP P37019

- Molecule 2 is a protein called Fab, 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, 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			

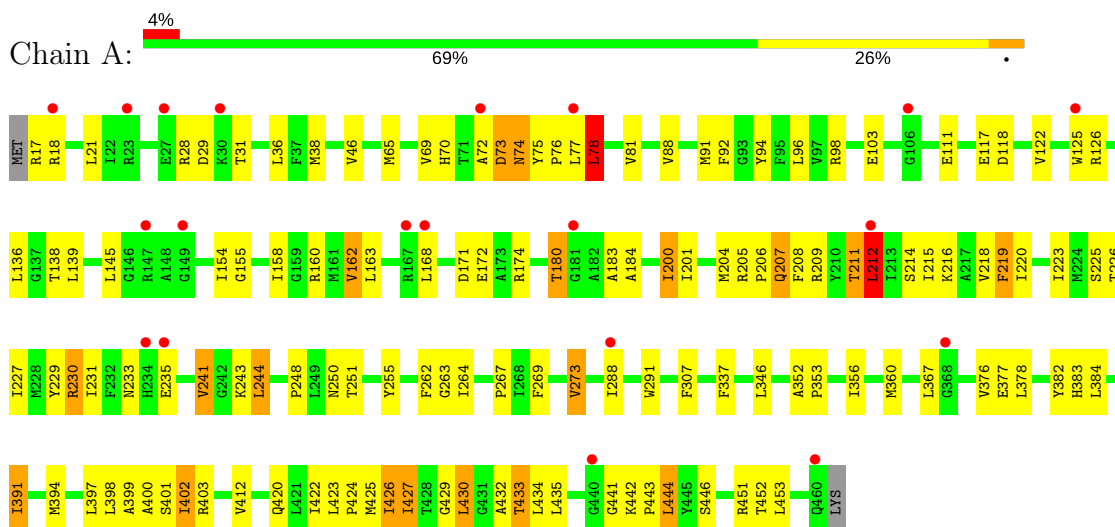
- Molecule 4 is FLUORIDE ION (three-letter code: F) (formula: F).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	2	Total	F	0	0
			2	2		
4	A	2	Total	F	0	0
			2	2		

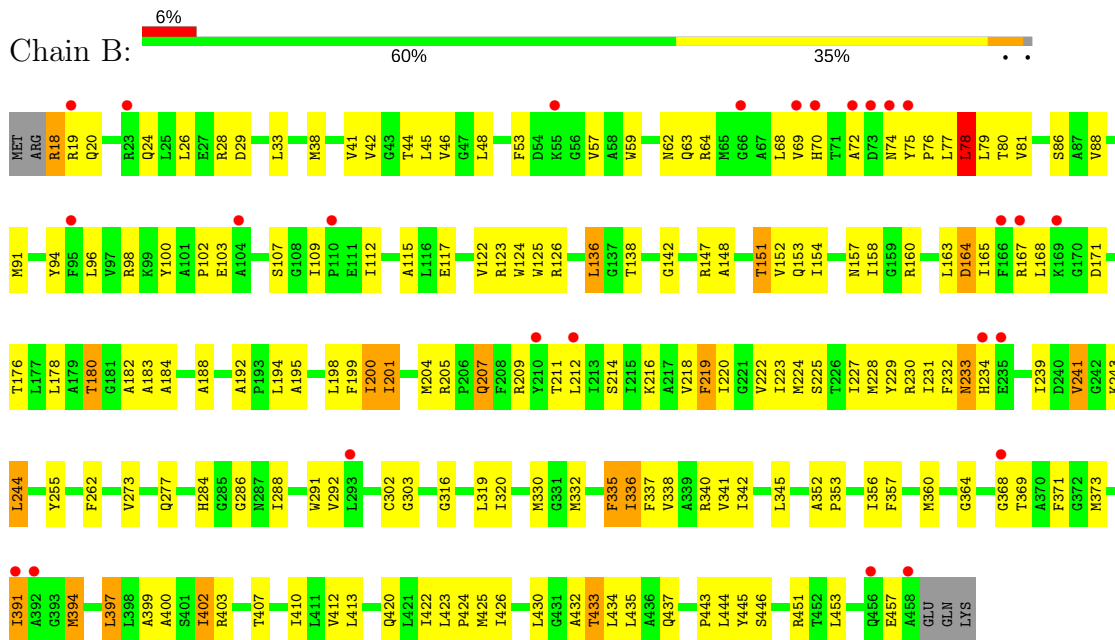
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

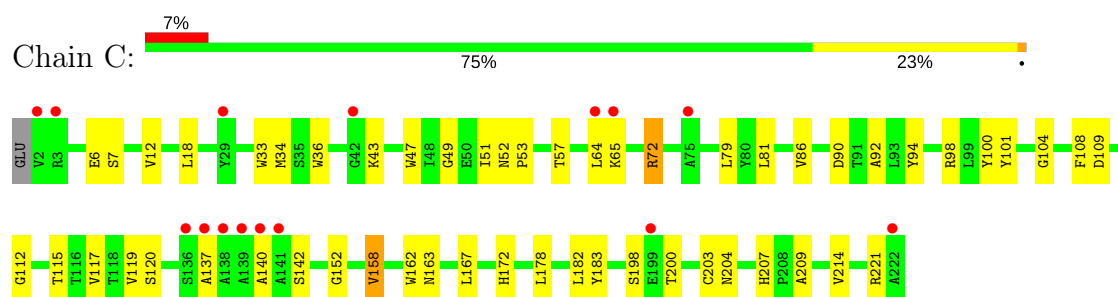
- Molecule 1: H(+)/Cl(-) exchange transporter ClcA



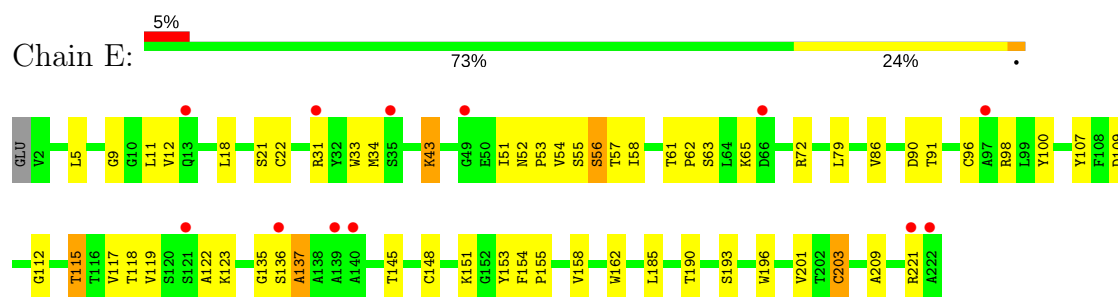
- Molecule 1: H(+)/Cl(-) exchange transporter ClcA



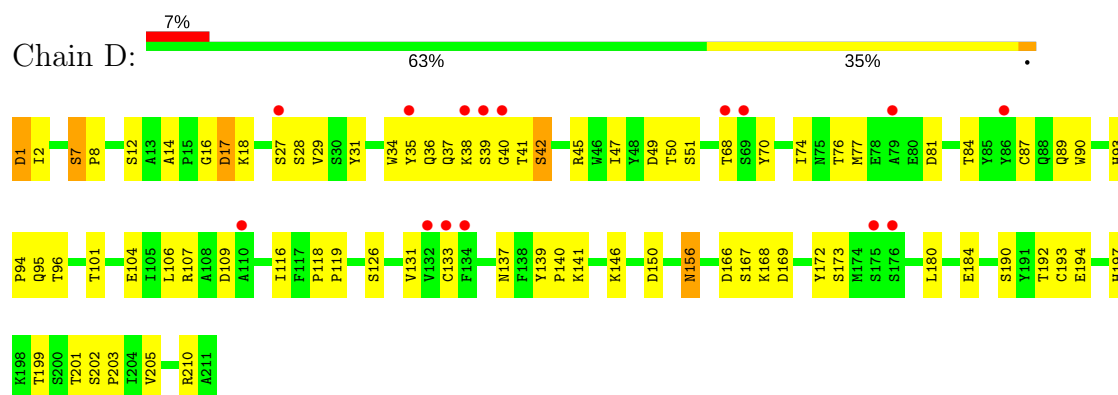
- Molecule 2: Fab, heavy chain



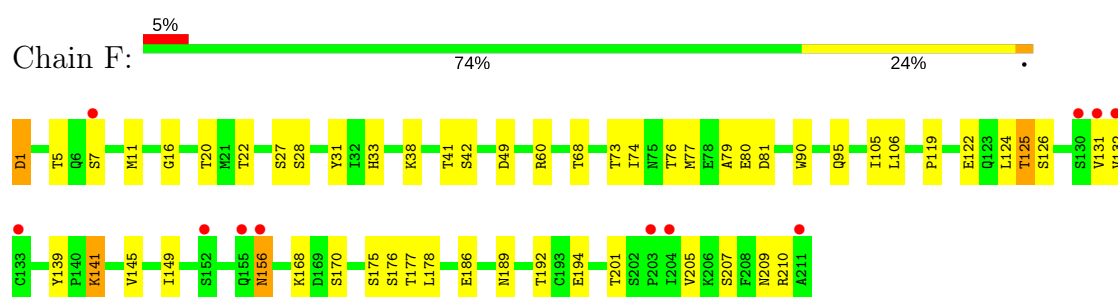
• Molecule 2: Fab, heavy chain



• Molecule 3: Fab, light chain



• Molecule 3: Fab, 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	231.42Å 100.09Å 170.23Å 90.00° 131.82° 90.00°	Depositor
Resolution (Å)	25.00 – 2.85 25.04 – 2.79	Depositor EDS
% Data completeness (in resolution range)	99.0 (25.00-2.85) 99.0 (25.04-2.79)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.74 (at 2.80Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.248 , 0.290 0.232 , 0.289	Depositor DCC
R_{free} test set	3421 reflections (5.06%)	DCC
Wilson B-factor (Å ²)	82.9	Xtriage
Anisotropy	0.425	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 39.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.003 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13219	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.51% 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](#)

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

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	0.41	0/3401	0.61	2/4616 (0.0%)
1	B	0.41	0/3372	0.60	1/4578 (0.0%)
2	C	0.49	0/1721	0.65	0/2355
2	E	0.47	0/1721	0.60	0/2355
3	D	0.45	0/1660	0.62	0/2257
3	F	0.48	0/1660	0.65	0/2257
All	All	0.44	0/13535	0.62	3/18418 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	78	LEU	CA-CB-CG	6.26	129.70	115.30
1	B	78	LEU	CA-CB-CG	5.83	128.72	115.30
1	A	212	LEU	CA-CB-CG	5.08	126.99	115.30

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	3329	0	3483	92	0
1	B	3300	0	3456	115	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	1672	0	1654	26	0
2	E	1672	0	1654	33	0
3	D	1621	0	1546	42	0
3	F	1621	0	1546	31	0
4	A	2	0	0	0	0
4	B	2	0	0	2	0
All	All	13219	0	13339	307	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 12.

All (307) 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:A:219:PHE:HB3	1:B:430:LEU:HD21	1.52	0.90
3:F:95:GLN:N	3:F:95:GLN:OE1	2.08	0.85
3:D:95:GLN:N	3:D:95:GLN:OE1	2.10	0.85
3:D:16:GLY:HA2	3:D:76:THR:HG23	1.59	0.84
3:D:37:GLN:NE2	3:D:41:THR:O	2.17	0.76
1:B:443:PRO:HB2	1:B:446:SER:HB2	1.67	0.76
1:A:18:ARG:HH11	1:B:457:GLU:HB3	1.51	0.75
3:F:1:ASP:OD2	3:F:1:ASP:N	2.20	0.74
3:F:16:GLY:HA2	3:F:76:THR:HG23	1.70	0.73
1:B:147:ARG:O	1:B:151:THR:OG1	2.05	0.72
2:C:7:SER:HA	2:C:115:THR:HG21	1.70	0.72
1:B:117:GLU:OE1	1:B:209:ARG:NH1	2.20	0.71
3:F:60:ARG:NH2	3:F:81:ASP:OD2	2.24	0.71
1:B:243:LYS:HB3	2:E:31:ARG:HH21	1.56	0.70
1:B:200:ILE:HD12	1:B:204:MET:HG3	1.72	0.70
1:B:38:MET:HG3	1:B:168:LEU:HD11	1.71	0.70
2:C:112:GLY:O	3:D:42:SER:OG	2.10	0.70
1:B:18:ARG:O	1:B:20:GLN:N	2.25	0.69
1:B:125:TRP:CD1	1:B:126:ARG:HG3	2.27	0.69
1:A:180:THR:HB	1:A:218:VAL:HA	1.73	0.68
3:D:89:GLN:O	3:D:95:GLN:HB2	1.94	0.68
1:A:398:LEU:O	1:A:402:ILE:HG22	1.93	0.68
1:A:94:TYR:CZ	1:A:352:ALA:HB2	2.29	0.67
1:B:198:LEU:HD13	1:B:201:ILE:HD11	1.77	0.67
1:B:94:TYR:CZ	1:B:352:ALA:HB2	2.30	0.67
3:D:1:ASP:N	3:D:1:ASP:OD2	2.20	0.67
3:D:166:ASP:OD1	3:D:167:SER:N	2.28	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:38:MET:HG3	1:A:168:LEU:HD11	1.78	0.65
1:B:98:ARG:NH2	1:B:102:PRO:HB3	2.11	0.65
1:B:163:LEU:HD12	1:B:168:LEU:HB2	1.77	0.64
3:F:60:ARG:HD2	3:F:81:ASP:OD1	1.97	0.64
1:A:426:ILE:HG22	1:B:223:ILE:HD11	1.80	0.64
1:A:183:ALA:HB2	1:A:200:ILE:HG12	1.79	0.64
1:A:220:ILE:HD11	1:B:434:LEU:HD11	1.80	0.64
1:A:229:TYR:O	1:A:233:ASN:HB2	2.00	0.62
1:A:403:ARG:NH2	1:B:29:ASP:OD1	2.31	0.62
2:C:34:MET:HB3	2:C:79:LEU:HD22	1.82	0.62
1:B:38:MET:O	1:B:42:VAL:HG23	2.00	0.61
2:E:61:THR:O	2:E:63:SER:N	2.33	0.61
1:B:100:TYR:O	1:B:126:ARG:NH1	2.33	0.60
3:F:132:VAL:HG22	3:F:177:THR:HG23	1.84	0.60
2:C:98:ARG:NH1	2:C:109:ASP:OD2	2.34	0.60
2:C:6:GLU:OE1	2:C:6:GLU:N	2.35	0.60
3:F:90:TRP:CG	3:F:95:GLN:HB3	2.37	0.60
1:B:184:ALA:HB1	1:B:225:SER:HB2	1.83	0.59
1:B:229:TYR:O	1:B:233:ASN:HB2	2.01	0.59
3:D:119:PRO:HD3	3:D:131:VAL:HG22	1.83	0.59
1:B:183:ALA:HB2	1:B:200:ILE:HG12	1.83	0.59
3:D:116:ILE:HD12	3:D:193:CYS:HB2	1.85	0.59
3:D:29:VAL:O	3:D:70:TYR:OH	2.17	0.58
1:A:422:ILE:HA	1:A:425:MET:HE3	1.86	0.58
2:E:54:VAL:HG23	2:E:56:SER:HB3	1.86	0.58
1:B:153:GLN:O	1:B:157:ASN:ND2	2.34	0.57
1:B:138:THR:HG21	1:B:353:PRO:HD2	1.86	0.57
1:A:125:TRP:CD1	1:A:126:ARG:HG3	2.40	0.57
1:A:138:THR:HG21	1:A:353:PRO:HD2	1.86	0.57
3:D:49:ASP:O	3:D:51:SER:N	2.33	0.57
3:D:12:SER:HA	3:D:104:GLU:O	2.04	0.56
3:D:194:GLU:HG2	3:D:205:VAL:HG12	1.86	0.56
2:E:145:THR:HG1	2:E:190:THR:HG1	1.51	0.56
1:A:200:ILE:HA	1:A:204:MET:HB2	1.87	0.56
1:B:152:VAL:HG13	1:B:182:ALA:HB1	1.87	0.56
2:E:98:ARG:NH1	2:E:109:ASP:OD2	2.39	0.56
1:A:122:VAL:HB	1:A:160:ARG:HG2	1.87	0.56
2:E:162:TRP:CZ3	2:E:203:CYS:HB3	2.40	0.56
1:A:255:TYR:CG	1:A:424:PRO:HB3	2.41	0.56
1:A:434:LEU:HD11	1:B:220:ILE:HD11	1.87	0.55
3:D:166:ASP:OD1	3:D:168:LYS:N	2.39	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:146:LYS:HB3	3:D:194:GLU:HB2	1.89	0.55
1:A:394:MET:HE2	1:A:412:VAL:HG13	1.88	0.55
1:A:98:ARG:HD2	1:A:291:TRP:CE3	2.42	0.55
2:E:12:VAL:HG11	2:E:18:LEU:HB3	1.89	0.55
1:A:154:ILE:O	1:A:158:ILE:HG12	2.07	0.55
1:A:216:LYS:HD3	1:B:434:LEU:HD23	1.89	0.55
1:A:207:GLN:HG2	1:B:28:ARG:NE	2.21	0.55
1:B:64:ARG:O	1:B:68:LEU:HG	2.06	0.54
1:B:224:MET:O	1:B:228:MET:HG2	2.08	0.54
3:F:38:LYS:NZ	3:F:80:GLU:O	2.40	0.54
1:B:122:VAL:HB	1:B:160:ARG:HG2	1.90	0.54
3:D:41:THR:OG1	3:D:42:SER:N	2.41	0.54
1:A:184:ALA:HB1	1:A:225:SER:HB2	1.90	0.53
1:A:376:VAL:HG22	1:A:384:LEU:HB2	1.89	0.53
1:A:200:ILE:HD12	1:A:204:MET:HG3	1.90	0.53
1:B:241:VAL:HG11	1:B:391:ILE:HD11	1.90	0.53
2:E:9:GLY:H	2:E:115:THR:HG21	1.73	0.53
1:B:200:ILE:HA	1:B:204:MET:HB2	1.91	0.53
2:C:86:VAL:HG13	2:C:90:ASP:HB2	1.91	0.53
1:B:214:SER:O	1:B:218:VAL:HG23	2.08	0.53
1:B:239:ILE:HD13	1:B:320:ILE:HG21	1.91	0.53
1:B:42:VAL:O	1:B:46:VAL:HG23	2.09	0.52
2:C:51:ILE:HD13	2:C:72:ARG:HG2	1.91	0.52
1:B:124:TRP:NE1	1:B:164:ASP:OD1	2.38	0.52
1:A:88:VAL:HA	1:A:91:MET:HE2	1.92	0.52
1:A:426:ILE:O	1:A:429:GLY:N	2.42	0.51
1:B:180:THR:HA	1:B:218:VAL:HG13	1.92	0.51
2:E:18:LEU:HD11	2:E:117:VAL:HG22	1.91	0.51
1:B:400:ALA:HB2	1:B:432:ALA:HB1	1.92	0.51
1:B:41:VAL:O	1:B:45:LEU:HG	2.10	0.51
2:C:53:PRO:HA	2:C:72:ARG:CZ	2.40	0.51
1:B:241:VAL:HG13	1:B:244:LEU:HD21	1.92	0.51
1:B:198:LEU:HG	1:B:410:ILE:HD12	1.93	0.51
2:E:51:ILE:HG13	2:E:58:ILE:HG12	1.93	0.51
1:A:356:ILE:O	1:A:360:MET:HG3	2.11	0.51
1:B:337:PHE:O	1:B:341:VAL:HG23	2.09	0.51
3:D:141:LYS:HB3	3:D:172:TYR:CE1	2.46	0.51
3:D:90:TRP:CG	3:D:95:GLN:HB3	2.46	0.51
3:F:194:GLU:HG2	3:F:205:VAL:HG12	1.92	0.51
1:B:74:ASN:HB3	1:B:77:LEU:HB3	1.93	0.51
3:D:34:TRP:HB2	3:D:47:ILE:HB	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:77:LEU:O	1:A:81:VAL:HG13	2.11	0.50
2:E:22:CYS:HB3	2:E:79:LEU:HB3	1.94	0.50
1:A:31:THR:HB	1:A:36:LEU:HD21	1.93	0.50
1:B:356:ILE:HG22	4:B:502:F:F	2.02	0.50
1:A:118:ASP:CG	1:A:174:ARG:HH21	2.15	0.50
3:D:31:TYR:HB3	3:D:49:ASP:HA	1.93	0.50
2:E:135:GLY:O	2:E:137:ALA:N	2.40	0.50
3:D:14:ALA:O	3:D:17:ASP:HB2	2.12	0.50
3:D:180:LEU:HD22	3:D:184:GLU:HG2	1.94	0.50
2:E:53:PRO:HA	2:E:72:ARG:CZ	2.42	0.49
1:A:92:PHE:O	1:A:96:LEU:HD23	2.12	0.49
1:A:227:ILE:O	1:A:231:ILE:HG12	2.13	0.49
1:B:273:VAL:O	1:B:277:GLN:HG3	2.11	0.49
3:F:95:GLN:CD	3:F:95:GLN:H	2.12	0.49
1:B:188:ALA:HB2	1:B:225:SER:OG	2.12	0.49
1:B:356:ILE:O	1:B:360:MET:HG3	2.13	0.49
1:B:78:LEU:HD13	1:B:79:LEU:HD23	1.95	0.49
2:C:18:LEU:HD11	2:C:117:VAL:HG22	1.93	0.49
1:B:109:ILE:HD12	1:B:445:TYR:CE1	2.48	0.49
2:E:107:TYR:HB3	3:F:33:HIS:CD2	2.47	0.49
1:A:117:GLU:OE1	1:A:209:ARG:NH1	2.46	0.48
2:E:135:GLY:HA2	2:E:221:ARG:HD3	1.96	0.48
2:C:221:ARG:NH2	3:D:118:PRO:HB2	2.27	0.48
1:B:53:PHE:HE2	1:B:147:ARG:HG2	1.77	0.48
3:F:156:ASN:OD1	3:F:156:ASN:N	2.46	0.48
1:A:208:PHE:CE2	1:B:24:GLN:HB3	2.47	0.48
2:C:163:ASN:ND2	2:C:167:LEU:HD22	2.29	0.48
1:B:399:ALA:O	1:B:403:ARG:HA	2.13	0.48
2:E:196:TRP:HD1	2:E:201:VAL:HG23	1.79	0.48
3:D:150:ASP:HA	3:D:190:SER:HB3	1.95	0.48
1:A:377:GLU:HG3	1:A:378:LEU:N	2.29	0.48
1:A:241:VAL:HG11	1:A:391:ILE:HD11	1.96	0.48
1:B:394:MET:HE2	1:B:412:VAL:HG13	1.95	0.48
1:A:216:LYS:NZ	1:B:433:THR:O	2.46	0.48
2:E:86:VAL:HG13	2:E:90:ASP:HB2	1.96	0.48
1:A:443:PRO:HB2	1:A:446:SER:HB2	1.96	0.48
3:F:186:GLU:HA	3:F:210:ARG:CZ	2.44	0.47
1:A:262:PHE:CE2	1:A:367:LEU:HD23	2.50	0.47
1:B:142:GLY:O	1:B:302:CYS:HB3	2.13	0.47
1:A:214:SER:O	1:A:218:VAL:HG23	2.15	0.47
1:B:394:MET:HG2	1:B:412:VAL:HG22	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:53:PHE:O	1:B:57:VAL:HG23	2.15	0.47
3:D:74:ILE:HD13	3:D:81:ASP:OD2	2.15	0.47
3:F:186:GLU:HG2	3:F:210:ARG:NH1	2.30	0.47
3:F:28:SER:HA	3:F:68:THR:HG22	1.96	0.47
1:A:212:LEU:HD12	1:A:212:LEU:H	1.79	0.47
1:A:28:ARG:NE	1:B:207:GLN:HG2	2.30	0.47
1:A:160:ARG:HD2	1:A:160:ARG:HA	1.48	0.47
1:A:337:PHE:CE1	1:A:367:LEU:HB2	2.50	0.47
1:B:335:PHE:O	1:B:338:VAL:N	2.48	0.47
2:C:162:TRP:CZ3	2:C:203:CYS:HB3	2.49	0.47
3:D:36:GLN:HG3	3:D:36:GLN:O	2.13	0.47
1:A:422:ILE:HD11	1:B:194:LEU:HD11	1.98	0.46
1:B:91:MET:HG2	1:B:292:VAL:O	2.15	0.46
3:F:149:ILE:HD11	3:F:178:LEU:HD21	1.97	0.46
3:F:141:LYS:HB2	3:F:141:LYS:HE2	1.64	0.46
1:A:248:PRO:HD2	1:A:251:THR:OG1	2.15	0.46
2:E:153:TYR:CE1	2:E:158:VAL:HG13	2.50	0.46
1:B:397:LEU:HA	1:B:397:LEU:HD23	1.73	0.46
1:B:88:VAL:HA	1:B:91:MET:HE2	1.97	0.46
2:E:34:MET:HB3	2:E:79:LEU:HD22	1.96	0.46
1:A:383:HIS:HD2	2:C:33:TRP:CE3	2.33	0.46
1:B:98:ARG:HH21	1:B:102:PRO:HB3	1.78	0.46
1:A:46:VAL:HG22	1:A:155:GLY:HA2	1.98	0.45
1:A:78:LEU:HD11	1:A:307:PHE:CE2	2.51	0.45
3:F:90:TRP:CD2	3:F:95:GLN:HB3	2.52	0.45
1:A:74:ASN:OD1	1:A:76:PRO:HD2	2.15	0.45
1:B:192:ALA:HB1	1:B:195:ALA:HB3	1.98	0.45
3:D:7:SER:HB3	3:D:8:PRO:HD3	1.98	0.45
2:E:91:THR:OG1	2:E:119:VAL:HG23	2.16	0.45
3:F:7:SER:HB3	3:F:22:THR:HB	1.99	0.45
2:E:155:PRO:HD2	2:E:209:ALA:CB	2.47	0.45
2:E:11:LEU:HA	2:E:11:LEU:HD12	1.79	0.45
1:A:255:TYR:CD2	1:A:424:PRO:HB3	2.52	0.45
1:A:98:ARG:HB3	1:A:288:ILE:HG13	1.99	0.45
1:B:176:THR:O	1:B:180:THR:HG23	2.17	0.45
1:A:117:GLU:HA	1:A:209:ARG:HH22	1.81	0.45
2:C:52:ASN:ND2	2:C:57:THR:HB	2.32	0.45
3:F:145:VAL:HA	3:F:194:GLU:O	2.17	0.45
1:A:139:LEU:CD2	1:A:145:LEU:HB2	2.47	0.45
1:B:59:TRP:O	1:B:62:ASN:HB3	2.17	0.45
3:D:28:SER:HA	3:D:68:THR:HG22	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:158:ILE:O	1:A:162:VAL:HG13	2.17	0.45
1:A:201:ILE:HG21	1:A:215:ILE:HD11	1.98	0.45
3:F:60:ARG:HH21	3:F:81:ASP:CG	2.19	0.45
1:A:399:ALA:O	1:A:433:THR:HG23	2.17	0.45
1:A:73:ASP:OD1	1:A:73:ASP:N	2.34	0.45
1:A:430:LEU:HD21	1:B:219:PHE:HB3	1.99	0.45
1:B:136:LEU:HA	1:B:136:LEU:HD12	1.81	0.44
1:A:230:ARG:NH2	1:B:423:LEU:HD13	2.32	0.44
1:B:75:TYR:HB3	1:B:76:PRO:HD3	1.98	0.44
1:B:86:SER:OG	1:B:303:GLY:HA3	2.16	0.44
3:F:106:LEU:HD23	3:F:139:TYR:OH	2.18	0.44
1:B:422:ILE:HD12	1:B:425:MET:HE3	2.00	0.44
1:B:154:ILE:O	1:B:158:ILE:HG12	2.17	0.44
3:D:35:TYR:CD2	3:D:45:ARG:HA	2.53	0.44
1:A:29:ASP:OD2	1:A:216:LYS:HE3	2.18	0.44
1:B:227:ILE:O	1:B:231:ILE:HG12	2.17	0.44
1:B:332:MET:O	1:B:336:ILE:HG13	2.16	0.44
1:B:369:THR:O	1:B:373:MET:HG3	2.18	0.44
2:C:158:VAL:HG12	2:C:207:HIS:HB2	1.99	0.44
1:A:267:PRO:HG3	1:A:441:GLY:HA3	2.00	0.44
1:B:402:ILE:HD11	1:B:445:TYR:CD1	2.53	0.44
1:B:164:ASP:O	1:B:167:ARG:N	2.39	0.43
1:B:255:TYR:CE2	1:B:424:PRO:HB3	2.52	0.43
2:E:43:LYS:HB3	2:E:43:LYS:HE2	1.78	0.43
1:A:423:LEU:HD13	1:B:230:ARG:NH2	2.33	0.43
1:B:59:TRP:O	1:B:63:GLN:HG2	2.18	0.43
2:C:178:LEU:HD13	2:C:183:TYR:CZ	2.53	0.43
1:B:98:ARG:HB3	1:B:288:ILE:HG13	2.01	0.43
3:F:31:TYR:HB3	3:F:49:ASP:HA	2.00	0.43
1:B:75:TYR:HA	1:B:78:LEU:HD12	2.00	0.43
2:C:101:TYR:HB2	2:C:104:GLY:HA2	2.00	0.43
1:A:433:THR:HB	1:B:216:LYS:HE2	2.01	0.43
1:A:263:GLY:HA3	1:A:435:LEU:HB2	2.00	0.43
2:C:36:TRP:CE2	2:C:81:LEU:HB2	2.54	0.43
3:D:109:ASP:HB3	3:D:199:THR:HG22	2.01	0.43
2:E:196:TRP:CD1	2:E:201:VAL:HG23	2.53	0.43
1:A:160:ARG:O	1:A:163:LEU:HB3	2.18	0.43
1:B:123:ARG:HE	1:B:126:ARG:HD2	1.83	0.43
1:B:341:VAL:O	1:B:345:LEU:HG	2.18	0.43
1:B:402:ILE:HD11	1:B:445:TYR:CE1	2.54	0.43
1:A:400:ALA:HB2	1:A:432:ALA:HB1	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:427:ILE:HG12	1:B:223:ILE:HD13	2.01	0.42
3:F:189:ASN:O	3:F:209:ASN:HA	2.19	0.42
1:B:284:HIS:C	1:B:286:GLY:H	2.23	0.42
3:F:38:LYS:O	3:F:41:THR:HG22	2.20	0.42
3:F:20:THR:HG23	3:F:73:THR:OG1	2.19	0.42
3:D:36:GLN:HA	3:D:84:THR:O	2.19	0.42
1:A:235:GLU:OE1	2:E:100:TYR:HE2	2.03	0.42
1:A:31:THR:N	1:B:437:GLN:HE22	2.18	0.42
1:A:75:TYR:HB3	1:A:76:PRO:HD3	2.01	0.42
1:B:78:LEU:HA	1:B:81:VAL:HG22	2.00	0.42
1:A:206:PRO:HG2	1:A:211:THR:HG21	2.01	0.42
1:A:269:PHE:O	1:A:273:VAL:HG12	2.19	0.42
1:A:430:LEU:HA	1:A:430:LEU:HD12	1.79	0.42
1:B:69:VAL:HA	1:B:72:ALA:HB2	2.01	0.42
2:E:100:TYR:HB3	2:E:107:TYR:CE1	2.55	0.42
3:F:11:MET:HE3	3:F:11:MET:HB2	1.94	0.42
1:A:383:HIS:HD2	2:C:33:TRP:CZ3	2.38	0.42
1:B:255:TYR:CD2	1:B:424:PRO:HB3	2.55	0.42
1:B:316:GLY:O	1:B:319:LEU:HG	2.20	0.42
1:B:243:LYS:HA	1:B:243:LYS:HD3	1.79	0.42
1:B:148:ALA:HB3	4:B:501:F:F	2.09	0.42
1:B:77:LEU:O	1:B:80:THR:HB	2.19	0.42
3:D:2:ILE:O	3:D:96:THR:HG21	2.20	0.42
1:A:139:LEU:HD22	1:A:145:LEU:HB2	2.02	0.42
1:A:163:LEU:HD12	1:A:168:LEU:HB2	2.02	0.42
2:C:152:GLY:HA2	2:C:182:LEU:HB3	2.02	0.42
3:D:93:HIS:CG	3:D:94:PRO:HA	2.55	0.42
2:E:112:GLY:O	3:F:42:SER:OG	2.25	0.42
1:A:21:LEU:HD21	1:B:117:GLU:HG3	2.02	0.41
1:A:226:THR:O	1:A:230:ARG:HG2	2.18	0.41
1:B:435:LEU:HD13	1:B:435:LEU:HA	1.90	0.41
3:D:156:ASN:N	3:D:156:ASN:OD1	2.52	0.41
2:C:108:PHE:HZ	3:D:95:GLN:HE21	1.67	0.41
1:A:243:LYS:HD3	1:A:243:LYS:HA	1.71	0.41
1:B:112:ILE:HD13	1:B:112:ILE:HA	1.94	0.41
1:B:413:LEU:HD12	1:B:413:LEU:HA	1.92	0.41
3:D:107:ARG:HG3	3:D:139:TYR:CD1	2.54	0.41
1:A:122:VAL:HG11	1:A:160:ARG:HB2	2.03	0.41
1:B:48:LEU:HD21	1:B:228:MET:SD	2.60	0.41
2:E:154:PHE:HA	2:E:155:PRO:HA	1.85	0.41
1:A:251:THR:HG22	1:A:255:TYR:CE1	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:98:ARG:HD2	1:B:291:TRP:CD2	2.56	0.41
2:E:51:ILE:HD13	2:E:72:ARG:HG2	2.01	0.41
2:E:52:ASN:ND2	2:E:57:THR:H	2.19	0.41
3:F:124:LEU:C	3:F:126:SER:H	2.24	0.41
2:C:207:HIS:CE1	2:C:209:ALA:HB3	2.56	0.41
2:E:100:TYR:O	2:E:107:TYR:N	2.41	0.41
1:A:250:ASN:ND2	1:A:382:TYR:HE2	2.18	0.41
1:A:442:LYS:HZ2	1:B:26:LEU:HB3	1.86	0.41
2:E:158:VAL:CG2	2:E:185:LEU:HD21	2.51	0.41
3:F:79:ALA:HA	3:F:105:ILE:HD12	2.03	0.41
3:F:119:PRO:HD3	3:F:131:VAL:HG22	2.03	0.41
2:C:47:TRP:CZ2	2:C:49:GLY:HA2	2.55	0.41
2:C:172:HIS:CD2	3:D:173:SER:OG	2.74	0.41
1:A:69:VAL:HA	1:A:72:ALA:HB2	2.02	0.41
3:D:38:LYS:O	3:D:40:GLY:N	2.54	0.41
1:B:115:ALA:HB1	1:B:178:LEU:HD21	2.03	0.40
1:B:262:PHE:HZ	1:B:364:GLY:HA2	1.85	0.40
3:D:90:TRP:CD1	3:D:95:GLN:HB3	2.57	0.40
1:A:223:ILE:HD11	1:B:426:ILE:HG22	2.03	0.40
1:A:226:THR:HG21	1:B:423:LEU:HD11	2.03	0.40
2:C:12:VAL:O	2:C:119:VAL:HA	2.21	0.40
2:C:92:ALA:HB3	2:C:94:TYR:CE2	2.56	0.40
2:E:33:TRP:CH2	2:E:52:ASN:HB3	2.56	0.40
1:B:199:PHE:CD1	1:B:407:THR:HG21	2.56	0.40
1:B:231:ILE:HB	1:B:232:PHE:CD1	2.57	0.40
1:B:44:THR:O	1:B:48:LEU:HG	2.22	0.40
3:D:202:SER:HA	3:D:203:PRO:HD2	1.79	0.40
1:A:160:ARG:NE	1:A:163:LEU:HD23	2.36	0.40
1:A:241:VAL:HG13	1:A:244:LEU:HD21	2.04	0.40
1:B:368:GLY:O	1:B:371:PHE:N	2.55	0.40
3:D:18:LYS:HA	3:D:74:ILE:O	2.21	0.40
1:A:401:SER:O	1:A:444:LEU:HB2	2.21	0.40
1:B:262:PHE:CZ	1:B:364:GLY:HA2	2.56	0.40
3:D:140:PRO:HD2	3:D:197:HIS:HE2	1.86	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	442/446 (99%)	401 (91%)	37 (8%)	4 (1%)	20	49
1	B	439/446 (98%)	400 (91%)	29 (7%)	10 (2%)	7	24
2	C	219/222 (99%)	200 (91%)	17 (8%)	2 (1%)	20	49
2	E	219/222 (99%)	201 (92%)	14 (6%)	4 (2%)	10	31
3	D	209/211 (99%)	187 (90%)	16 (8%)	6 (3%)	5	18
3	F	209/211 (99%)	191 (91%)	17 (8%)	1 (0%)	32	64
All	All	1737/1758 (99%)	1580 (91%)	130 (8%)	27 (2%)	11	33

All (27) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	74	ASN
1	B	19	ARG
2	C	137	ALA
3	D	39	SER
3	D	50	THR
1	A	426	ILE
1	B	335	PHE
2	C	140	ALA
2	E	62	PRO
2	E	136	SER
3	D	169	ASP
2	E	122	ALA
2	E	137	ALA
3	F	125	THR
1	A	427	ILE
1	B	164	ASP
1	B	234	HIS
1	B	357	PHE
3	D	7	SER
3	D	126	SER

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Mol	Chain	Res	Type
1	B	107	SER
1	B	165	ILE
1	B	233	ASN
3	D	137	ASN
1	A	172	GLU
1	B	342	ILE
1	B	336	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	334/336 (99%)	301 (90%)	33 (10%)	9	25
1	B	331/336 (98%)	300 (91%)	31 (9%)	10	27
2	C	181/182 (100%)	169 (93%)	12 (7%)	19	45
2	E	181/182 (100%)	167 (92%)	14 (8%)	15	37
3	D	185/185 (100%)	172 (93%)	13 (7%)	18	42
3	F	185/185 (100%)	169 (91%)	16 (9%)	12	32
All	All	1397/1406 (99%)	1278 (92%)	119 (8%)	12	33

All (119) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	ARG
1	A	65	MET
1	A	70	HIS
1	A	73	ASP
1	A	78	LEU
1	A	103	GLU
1	A	111	GLU
1	A	136	LEU
1	A	162	VAL
1	A	171	ASP
1	A	180	THR

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Mol	Chain	Res	Type
1	A	200	ILE
1	A	205	ARG
1	A	207	GLN
1	A	211	THR
1	A	212	LEU
1	A	219	PHE
1	A	230	ARG
1	A	241	VAL
1	A	244	LEU
1	A	264	ILE
1	A	273	VAL
1	A	346	LEU
1	A	391	ILE
1	A	397	LEU
1	A	402	ILE
1	A	420	GLN
1	A	430	LEU
1	A	433	THR
1	A	444	LEU
1	A	451	ARG
1	A	452	THR
1	A	453	LEU
1	B	18	ARG
1	B	33	LEU
1	B	70	HIS
1	B	78	LEU
1	B	96	LEU
1	B	103	GLU
1	B	136	LEU
1	B	151	THR
1	B	171	ASP
1	B	180	THR
1	B	200	ILE
1	B	201	ILE
1	B	205	ARG
1	B	207	GLN
1	B	211	THR
1	B	212	LEU
1	B	219	PHE
1	B	222	VAL
1	B	241	VAL
1	B	244	LEU

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Mol	Chain	Res	Type
1	B	330	MET
1	B	340	ARG
1	B	391	ILE
1	B	394	MET
1	B	397	LEU
1	B	402	ILE
1	B	420	GLN
1	B	433	THR
1	B	444	LEU
1	B	451	ARG
1	B	453	LEU
2	C	43	LYS
2	C	64	LEU
2	C	65	LYS
2	C	72	ARG
2	C	100	TYR
2	C	120	SER
2	C	142	SER
2	C	158	VAL
2	C	198	SER
2	C	200	THR
2	C	204	ASN
2	C	214	VAL
3	D	1	ASP
3	D	17	ASP
3	D	27	SER
3	D	42	SER
3	D	77	MET
3	D	87	CYS
3	D	101	THR
3	D	106	LEU
3	D	133	CYS
3	D	156	ASN
3	D	192	THR
3	D	201	THR
3	D	210	ARG
2	E	5	LEU
2	E	21	SER
2	E	43	LYS
2	E	55	SER
2	E	56	SER
2	E	65	LYS

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Mol	Chain	Res	Type
2	E	96	CYS
2	E	115	THR
2	E	118	THR
2	E	123	LYS
2	E	148	CYS
2	E	151	LYS
2	E	193	SER
2	E	203	CYS
3	F	1	ASP
3	F	5	THR
3	F	27	SER
3	F	74	ILE
3	F	77	MET
3	F	122	GLU
3	F	125	THR
3	F	141	LYS
3	F	156	ASN
3	F	168	LYS
3	F	170	SER
3	F	175	SER
3	F	176	SER
3	F	192	THR
3	F	201	THR
3	F	207	SER

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

Mol	Chain	Res	Type
1	A	284	HIS
1	B	284	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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 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 [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	444/446 (99%)	0.14	20 (4%) 34 29	42, 62, 91, 125	0
1	B	441/446 (98%)	0.18	26 (5%) 23 18	41, 65, 98, 126	0
2	C	221/222 (99%)	0.03	15 (6%) 18 13	29, 55, 85, 122	0
2	E	221/222 (99%)	-0.06	12 (5%) 26 22	32, 57, 85, 120	0
3	D	211/211 (100%)	0.34	15 (7%) 17 12	38, 65, 89, 108	0
3	F	211/211 (100%)	0.02	11 (5%) 28 23	30, 49, 95, 110	0
All	All	1749/1758 (99%)	0.12	99 (5%) 24 19	29, 60, 92, 126	0

All (99) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	39	SER	5.5
2	C	65	LYS	4.9
2	C	139	ALA	4.8
2	E	222	ALA	4.5
1	A	460	GLN	4.4
1	B	72	ALA	4.4
2	C	136	SER	4.4
1	B	167	ARG	4.4
3	F	156	ASN	4.3
2	C	138	ALA	4.3
2	E	139	ALA	4.2
3	D	133	CYS	4.2
1	B	73	ASP	4.2
3	D	38	LYS	4.0
2	C	137	ALA	3.8
1	A	234	HIS	3.8
1	A	168	LEU	3.7
2	C	140	ALA	3.7
1	B	456	GLN	3.6

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Mol	Chain	Res	Type	RSRZ
2	E	31	ARG	3.5
2	C	141	ALA	3.4
1	B	235	GLU	3.3
3	F	133	CYS	3.3
2	E	136	SER	3.3
3	D	175	SER	3.3
2	C	222	ALA	3.2
1	B	23	ARG	3.2
2	E	35	SER	3.1
2	C	2	VAL	3.1
3	F	155	GLN	3.1
3	D	79	ALA	3.1
1	B	234	HIS	3.0
2	C	29	TYR	3.0
3	F	152	SER	3.0
1	B	66	GLY	3.0
1	A	167	ARG	3.0
1	B	95	PHE	2.9
1	A	212	LEU	2.9
1	B	458	ALA	2.9
2	E	13	GLN	2.9
2	C	75	ALA	2.9
3	D	35	TYR	2.8
3	F	7	SER	2.8
1	A	235	GLU	2.8
2	E	140	ALA	2.8
3	F	131	VAL	2.8
2	C	199	GLU	2.7
1	B	392	ALA	2.7
1	B	74	ASN	2.7
3	D	68	THR	2.6
1	A	125	TRP	2.6
3	D	69	SER	2.6
3	D	132	VAL	2.6
3	D	176	SER	2.5
3	F	130	SER	2.5
1	B	212	LEU	2.5
3	F	203	PRO	2.4
1	A	368	GLY	2.4
1	A	147	ARG	2.4
1	B	368	GLY	2.3
2	E	121	SER	2.3

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Mol	Chain	Res	Type	RSRZ
3	D	86	TYR	2.3
3	F	132	VAL	2.3
1	A	30	LYS	2.3
1	A	72	ALA	2.3
2	E	49	GLY	2.3
1	B	75	TYR	2.2
1	A	27	GLU	2.2
1	A	18	ARG	2.2
1	B	210	TYR	2.2
2	C	42	GLY	2.2
1	A	288	ILE	2.2
1	B	69	VAL	2.2
3	D	110	ALA	2.2
3	D	27	SER	2.2
1	B	19	ARG	2.2
1	A	440	GLY	2.1
1	B	391	ILE	2.1
3	F	204	ILE	2.1
2	E	66	ASP	2.1
1	B	169	LYS	2.1
2	E	97	ALA	2.1
2	E	221	ARG	2.1
3	F	211	ALA	2.1
1	A	149	GLY	2.1
3	D	40	GLY	2.1
1	B	166	PHE	2.1
1	A	106	GLY	2.1
1	A	181	GLY	2.1
1	A	23	ARG	2.1
1	B	110	PRO	2.1
1	A	77	LEU	2.1
1	B	70	HIS	2.1
1	B	104	ALA	2.1
1	B	293	LEU	2.0
2	C	64	LEU	2.0
3	D	134	PHE	2.0
2	C	3	ARG	2.0
1	B	55	LYS	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 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	F	B	501	1/1	0.84	0.68	7.49	64,64,64,64	0
4	F	A	502	1/1	0.81	0.53	4.86	55,55,55,55	0
4	F	A	501	1/1	0.74	0.62	4.43	55,55,55,55	0
4	F	B	502	1/1	0.96	0.17	-0.18	48,48,48,48	0

6.5 Other polymers [i](#)

There are no such residues in this entry.