



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 10:52 AM UTC

PDB ID : 4KKC / pdb_00004kkc
Title : Structure of the E148A mutant of CLC-ec1 deltaNC construct in 20mM Bromide
Authors : Lim, H.-H.; Miller, C.
Deposited on : 2013-05-05
Resolution : 3.18 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
Xtriage (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

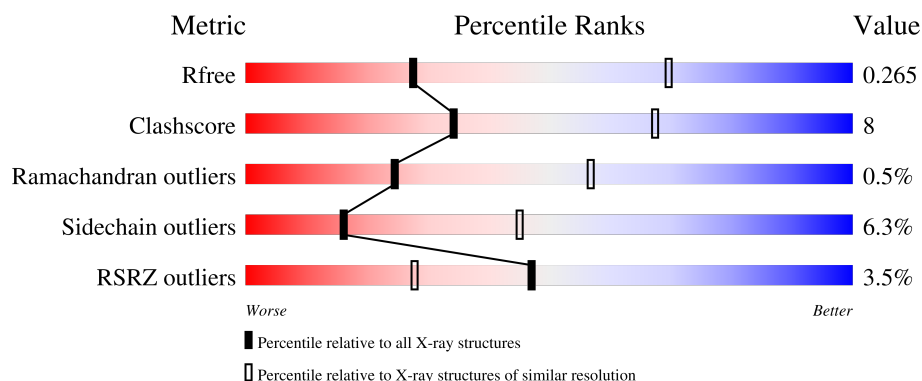
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.18 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	2001 (3.20-3.16)
Clashscore	190562	2119 (3.20-3.16)
Ramachandran outliers	187476	2070 (3.20-3.16)
Sidechain outliers	187428	2069 (3.20-3.16)
RSRZ outliers	180081	2001 (3.20-3.16)

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	446	4% 74% 22% .
1	B	446	4% 71% 25% ..
2	C	222	4% 83% 14% .
2	E	222	3% 80% 17% .
3	D	211	3% 82% 16% .

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Mol	Chain	Length	Quality of chain
3	F	211	% 82% 16%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13215 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	-	expression tag	UNP P37019
A	148	ALA	GLU	engineered mutation	UNP P37019
A	461	LYS	-	expression tag	UNP P37019
B	16	MET	-	expression tag	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			

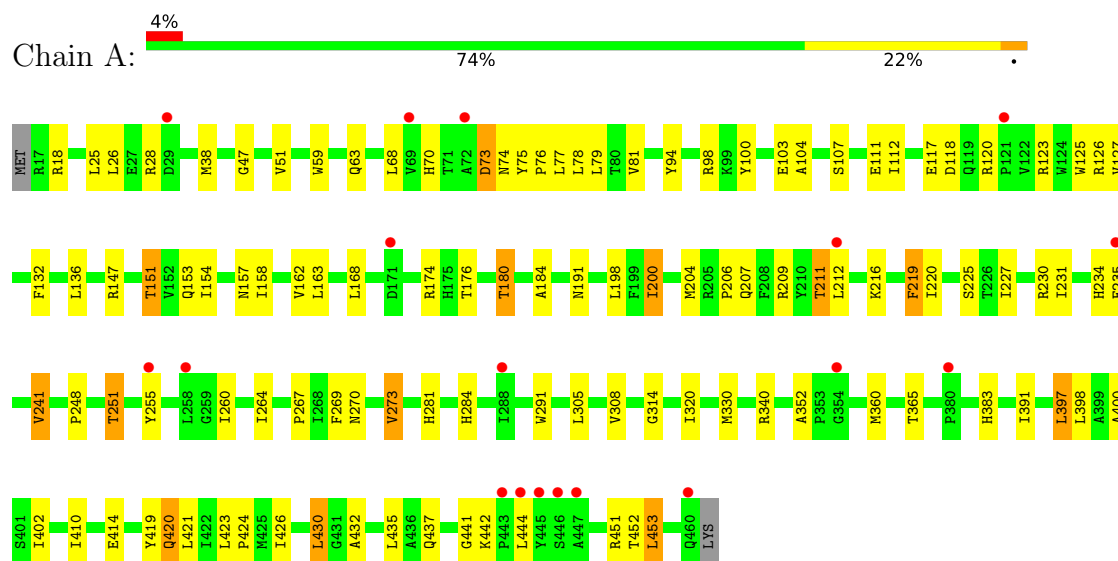
- Molecule 3 is a protein called Fab, light chain.

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			

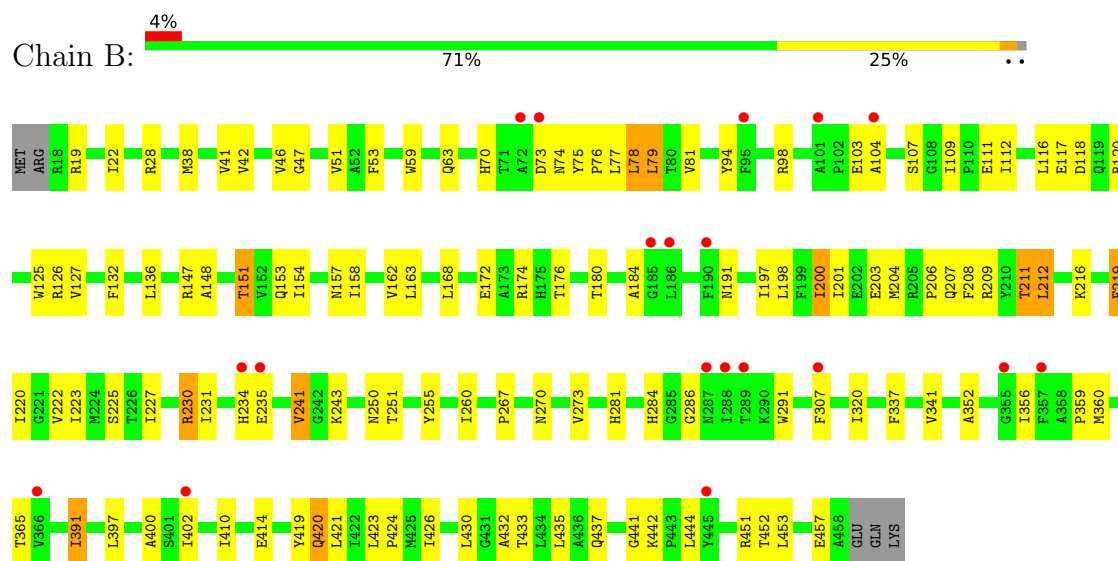
3 Residue-property plots

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.

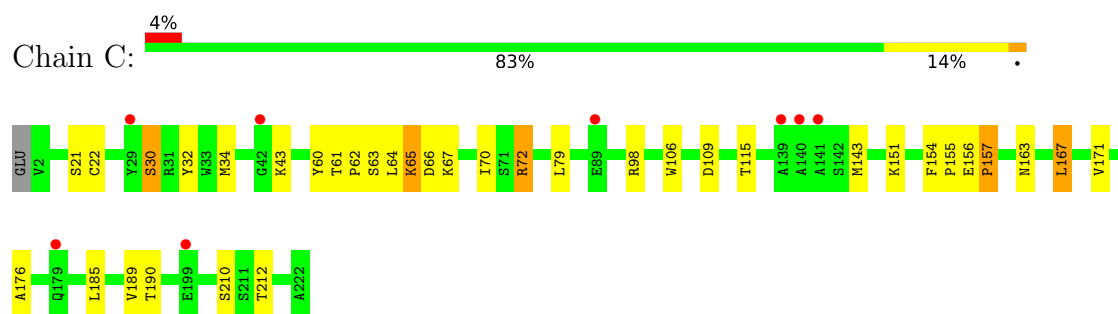
- Molecule 1: H(+)/Cl(-) exchange transporter ClcA



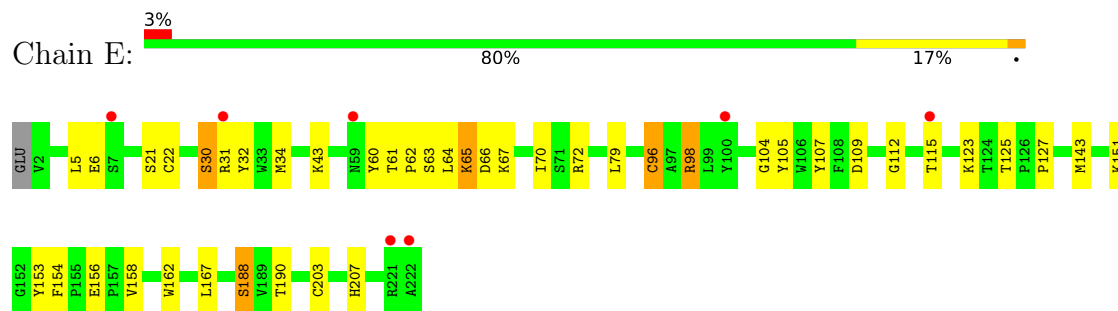
- Molecule 1: H(+)/Cl(-) exchange transporter ClcA



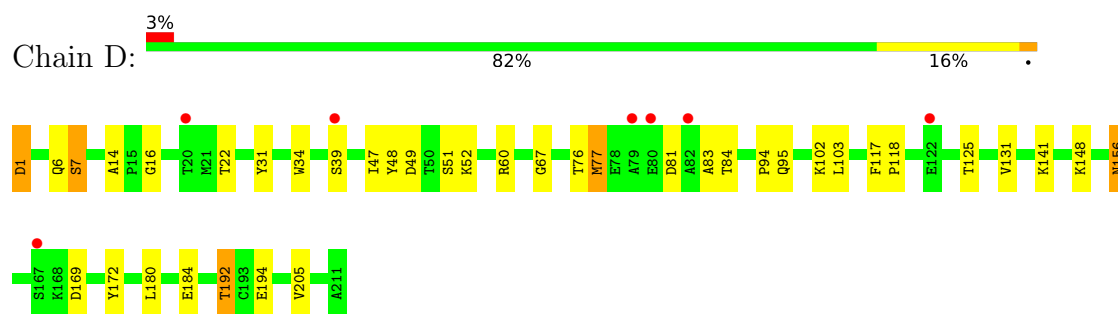
- Molecule 2: Fab, heavy chain



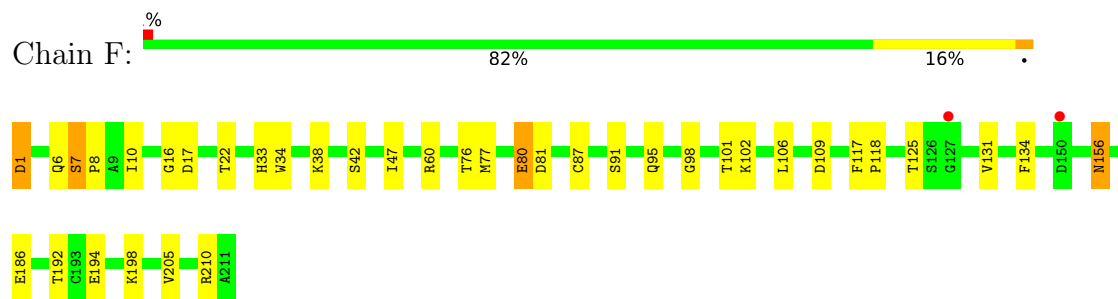
- Molecule 2: Fab, heavy chain



- 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	231.21Å 99.24Å 170.51Å 90.00° 132.00° 90.00°	Depositor
Resolution (Å)	29.53 – 3.18 29.53 – 3.18	Depositor EDS
% Data completeness (in resolution range)	98.3 (29.53-3.18) 98.7 (29.53-3.18)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.96 (at 3.18Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.231 , 0.264 0.230 , 0.265	Depositor DCC
R_{free} test set	2427 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	95.1	Xtriage
Anisotropy	0.505	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 10.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.004 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	13215	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.47% 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	0.43	0/3401	0.76	0/4616
1	B	0.42	0/3372	0.77	0/4578
2	C	0.40	0/1721	0.72	0/2355
2	E	0.41	0/1721	0.75	0/2355
3	D	0.36	0/1660	0.79	4/2257 (0.2%)
3	F	0.41	0/1660	0.77	2/2257 (0.1%)
All	All	0.41	0/13535	0.76	6/18418 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	7	SER	CA-C-N	-7.57	110.95	120.79
3	D	7	SER	C-N-CA	-7.57	110.95	120.79
3	F	7	SER	CA-C-N	-7.50	111.04	120.79
3	F	7	SER	C-N-CA	-7.50	111.04	120.79
3	D	14	ALA	CA-C-N	5.62	126.86	119.84
3	D	14	ALA	C-N-CA	5.62	126.86	119.84

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	72	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3300	0	3456	75	0
2	C	1672	0	1654	17	0
2	E	1672	0	1654	23	0
3	D	1621	0	1546	19	0
3	F	1621	0	1546	21	0
All	All	13215	0	13339	202	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 8.

All (202) 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:18:ARG:HH11	1:B:457:GLU:HB3	1.46	0.79
1:B:147:ARG:O	1:B:151:THR:OG1	2.04	0.76
1:A:200:ILE:HD12	1:A:204:MET:HG3	1.69	0.74
3:D:1:ASP:OD2	3:D:1:ASP:N	2.21	0.73
1:A:241:VAL:HG11	1:A:391:ILE:HD11	1.70	0.72
2:E:98:ARG:NH1	2:E:109:ASP:OD2	2.23	0.71
1:A:216:LYS:NZ	1:B:437:GLN:OE1	2.23	0.71
1:A:147:ARG:O	1:A:151:THR:OG1	2.06	0.71
3:F:1:ASP:OD2	3:F:1:ASP:N	2.23	0.70
1:B:241:VAL:HG11	1:B:391:ILE:HD11	1.74	0.69
1:A:200:ILE:HA	1:A:204:MET:HB2	1.74	0.68
1:B:200:ILE:HD12	1:B:204:MET:HG3	1.74	0.68
1:B:200:ILE:HA	1:B:204:MET:HB2	1.76	0.68
3:F:60:ARG:NH2	3:F:81:ASP:OD2	2.27	0.67
1:A:437:GLN:OE1	1:B:216:LYS:NZ	2.27	0.66
1:B:75:TYR:HA	1:B:78:LEU:HD12	1.77	0.65
3:D:95:GLN:OE1	3:D:95:GLN:N	2.27	0.65
3:F:95:GLN:OE1	3:F:95:GLN:N	2.29	0.64
1:B:117:GLU:OE1	1:B:209:ARG:NH1	2.31	0.63
3:D:194:GLU:HG2	3:D:205:VAL:HG12	1.80	0.61
2:C:22:CYS:HB3	2:C:79:LEU:HB3	1.82	0.61
1:B:206:PRO:HG2	1:B:211:THR:HG21	1.81	0.61
2:C:32:TYR:O	2:C:72:ARG:NH2	2.34	0.61
2:C:98:ARG:NH1	2:C:109:ASP:OD2	2.34	0.59
2:E:61:THR:O	2:E:63:SER:N	2.35	0.59
1:A:111:GLU:OE2	1:A:120:ARG:NE	2.36	0.58
1:A:73:ASP:OD1	1:A:73:ASP:N	2.36	0.58
1:B:38:MET:HG3	1:B:168:LEU:HD11	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:111:GLU:OE2	1:B:120:ARG:NE	2.36	0.58
2:E:64:LEU:HB2	2:E:67:LYS:HB2	1.86	0.58
1:A:260:ILE:HG23	1:A:435:LEU:HG	1.86	0.57
2:C:64:LEU:HB2	2:C:67:LYS:HB2	1.87	0.57
1:A:270:ASN:ND2	1:A:442:LYS:O	2.38	0.56
2:C:163:ASN:HD22	2:C:167:LEU:HD13	1.69	0.56
1:B:163:LEU:HD12	1:B:168:LEU:HB2	1.88	0.56
1:B:94:TYR:CZ	1:B:352:ALA:HB2	2.42	0.55
3:F:38:LYS:NZ	3:F:80:GLU:O	2.36	0.55
3:D:60:ARG:NH2	3:D:81:ASP:OD2	2.39	0.55
2:C:61:THR:O	2:C:63:SER:N	2.40	0.54
1:A:117:GLU:OE1	1:A:209:ARG:NH1	2.40	0.54
3:F:16:GLY:HA2	3:F:76:THR:HG23	1.90	0.54
1:B:198:LEU:HG	1:B:410:ILE:HD12	1.89	0.54
1:A:28:ARG:HE	1:B:207:GLN:HG2	1.73	0.53
1:B:74:ASN:HB3	1:B:77:LEU:HB3	1.91	0.53
1:A:112:ILE:HG13	1:A:153:GLN:HA	1.91	0.52
1:B:125:TRP:CD1	1:B:126:ARG:HG3	2.44	0.52
2:E:34:MET:HB3	2:E:79:LEU:HD22	1.92	0.52
1:B:98:ARG:HD2	1:B:291:TRP:CE3	2.45	0.52
1:A:38:MET:HG3	1:A:168:LEU:HD11	1.91	0.52
1:B:281:HIS:HA	1:B:284:HIS:CE1	2.45	0.52
1:A:206:PRO:HG2	1:A:211:THR:HG21	1.91	0.51
1:A:104:ALA:HB2	1:A:127:VAL:HG13	1.91	0.51
1:A:198:LEU:HG	1:A:410:ILE:HD12	1.92	0.51
3:F:194:GLU:HG2	3:F:205:VAL:HG12	1.92	0.51
1:A:414:GLU:OE1	1:B:419:TYR:OH	2.24	0.51
2:C:34:MET:HB3	2:C:79:LEU:HD22	1.91	0.51
1:A:28:ARG:NH2	1:B:203:GLU:OE1	2.30	0.51
1:A:184:ALA:HB1	1:A:225:SER:HB2	1.92	0.51
1:B:270:ASN:ND2	1:B:442:LYS:O	2.44	0.50
1:B:112:ILE:HG13	1:B:153:GLN:HA	1.94	0.49
3:D:31:TYR:HB3	3:D:49:ASP:HA	1.94	0.49
1:A:227:ILE:O	1:A:231:ILE:HG12	2.13	0.49
1:B:227:ILE:O	1:B:231:ILE:HG12	2.13	0.49
2:E:22:CYS:HB3	2:E:79:LEU:HB3	1.93	0.49
1:A:98:ARG:HD2	1:A:291:TRP:CE3	2.47	0.49
1:B:104:ALA:HB2	1:B:127:VAL:HG13	1.93	0.49
1:A:383:HIS:HD1	2:C:106:TRP:CD1	2.30	0.48
1:A:430:LEU:HD11	1:B:219:PHE:HB3	1.94	0.48
3:D:156:ASN:N	3:D:156:ASN:OD1	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:32:TYR:O	2:E:72:ARG:NH2	2.46	0.48
1:A:267:PRO:HG3	1:A:441:GLY:HA3	1.94	0.48
2:E:125:THR:N	2:E:154:PHE:O	2.39	0.48
3:F:156:ASN:OD1	3:F:156:ASN:N	2.45	0.48
1:A:314:GLY:O	1:A:340:ARG:NH2	2.47	0.48
2:C:30:SER:C	2:C:32:TYR:H	2.21	0.48
1:A:132:PHE:O	1:A:136:LEU:HB2	2.14	0.48
1:A:176:THR:O	1:A:180:THR:HG23	2.14	0.48
1:A:191:ASN:OD1	1:A:230:ARG:NH1	2.47	0.48
2:E:30:SER:C	2:E:32:TYR:H	2.22	0.48
1:A:248:PRO:O	1:A:251:THR:HG22	2.13	0.48
2:C:65:LYS:H	2:C:65:LYS:HG3	1.38	0.47
1:A:281:HIS:HA	1:A:284:HIS:CE1	2.49	0.47
1:B:234:HIS:CD2	1:B:235:GLU:HG2	2.50	0.47
1:A:94:TYR:CZ	1:A:352:ALA:HB2	2.50	0.47
1:B:184:ALA:HB1	1:B:225:SER:HB2	1.95	0.47
2:C:210:SER:OG	2:C:212:THR:OG1	2.31	0.47
1:B:176:THR:O	1:B:180:THR:HG23	2.15	0.47
1:B:118:ASP:CG	1:B:174:ARG:HH21	2.22	0.47
1:A:400:ALA:HB2	1:A:432:ALA:HB1	1.97	0.46
1:A:26:LEU:HD22	1:B:442:LYS:HZ2	1.80	0.46
1:B:73:ASP:N	1:B:73:ASP:OD1	2.48	0.46
1:B:127:VAL:HB	1:B:157:ASN:ND2	2.30	0.46
2:E:127:PRO:HB3	2:E:153:TYR:HB3	1.98	0.46
1:A:216:LYS:O	1:A:220:ILE:HG13	2.16	0.46
1:B:42:VAL:O	1:B:46:VAL:HG23	2.16	0.46
3:D:84:THR:HA	3:D:102:LYS:HA	1.97	0.46
2:E:143:MET:HE3	2:E:190:THR:HG22	1.98	0.46
3:F:109:ASP:OD2	3:F:198:LYS:NZ	2.49	0.46
1:B:400:ALA:HB2	1:B:432:ALA:HB1	1.97	0.45
3:D:1:ASP:HB3	3:D:94:PRO:HD2	1.99	0.45
2:E:65:LYS:H	2:E:65:LYS:HG3	1.40	0.45
2:C:60:TYR:HE2	2:C:70:ILE:HG13	1.81	0.45
3:F:34:TRP:HB2	3:F:47:ILE:HB	1.97	0.45
1:B:53:PHE:HE2	1:B:147:ARG:HG2	1.82	0.45
1:B:216:LYS:O	1:B:220:ILE:HG13	2.16	0.45
2:E:112:GLY:O	3:F:42:SER:OG	2.30	0.45
1:A:360:MET:HE3	1:A:398:LEU:HD23	1.99	0.45
1:A:68:LEU:HB3	1:A:78:LEU:HD11	1.99	0.45
3:F:7:SER:HB3	3:F:8:PRO:HD3	1.98	0.45
1:B:191:ASN:OD1	1:B:230:ARG:NH1	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:6:GLU:HA	2:E:22:CYS:HA	1.97	0.45
1:A:184:ALA:HB1	1:A:225:SER:CB	2.47	0.45
2:E:6:GLU:OE2	2:E:96:CYS:N	2.46	0.45
1:B:59:TRP:O	1:B:63:GLN:HG2	2.18	0.44
1:A:25:LEU:HD23	1:B:208:PHE:HE1	1.81	0.44
3:D:141:LYS:HB3	3:D:172:TYR:CZ	2.52	0.44
2:C:156:GLU:HA	2:C:157:PRO:HA	1.79	0.44
2:E:158:VAL:HG12	2:E:207:HIS:HB2	2.00	0.44
1:B:267:PRO:HG3	1:B:441:GLY:HA3	1.99	0.44
1:A:125:TRP:CD1	1:A:126:ARG:HG3	2.53	0.44
1:B:132:PHE:O	1:B:136:LEU:HB2	2.17	0.44
1:B:284:HIS:C	1:B:286:GLY:H	2.27	0.43
1:A:28:ARG:NE	1:B:207:GLN:HG2	2.33	0.43
1:B:260:ILE:HG23	1:B:435:LEU:HG	1.99	0.43
1:A:397:LEU:HD23	1:A:397:LEU:HA	1.75	0.43
1:B:420:GLN:H	1:B:420:GLN:HG3	1.51	0.43
3:D:117:PHE:HA	3:D:118:PRO:HD3	1.78	0.43
2:C:171:VAL:HG22	2:C:189:VAL:HG23	2.00	0.43
2:E:162:TRP:CZ3	2:E:203:CYS:HB3	2.54	0.43
1:B:250:ASN:OD1	2:E:104:GLY:HA3	2.19	0.43
3:D:6:GLN:HA	3:D:22:THR:O	2.19	0.43
1:B:197:ILE:HG12	1:B:222:VAL:HG21	2.01	0.43
1:B:360:MET:HE1	1:B:402:ILE:HG13	2.00	0.43
1:A:59:TRP:O	1:A:63:GLN:HG2	2.18	0.43
3:D:34:TRP:HB2	3:D:47:ILE:HB	2.01	0.43
3:F:117:PHE:HA	3:F:118:PRO:HD3	1.75	0.43
1:A:163:LEU:HD12	1:A:168:LEU:HB2	2.01	0.43
1:B:78:LEU:HD11	1:B:307:PHE:CE2	2.53	0.43
2:C:154:PHE:HA	2:C:155:PRO:HA	1.85	0.43
3:D:7:SER:HB3	3:D:22:THR:HB	2.00	0.43
3:F:10:ILE:HG12	3:F:102:LYS:HB3	2.01	0.43
2:C:143:MET:HE3	2:C:190:THR:HG22	1.99	0.42
3:D:48:TYR:CE1	3:D:52:LYS:HD2	2.54	0.42
2:C:176:ALA:HB2	2:C:185:LEU:HD23	2.00	0.42
3:D:180:LEU:HD22	3:D:184:GLU:HG2	2.00	0.42
1:B:154:ILE:O	1:B:158:ILE:HG12	2.20	0.42
3:F:6:GLN:HA	3:F:22:THR:O	2.19	0.42
1:A:74:ASN:HB3	1:A:77:LEU:HB3	2.01	0.42
1:A:127:VAL:HB	1:A:157:ASN:ND2	2.34	0.42
1:A:419:TYR:C	1:A:421:LEU:H	2.28	0.42
1:A:255:TYR:CD2	1:A:424:PRO:HB3	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:320:ILE:HG23	1:B:365:THR:HG21	2.01	0.42
1:A:118:ASP:CG	1:A:174:ARG:HH21	2.27	0.42
1:B:78:LEU:HD13	1:B:79:LEU:HD23	2.01	0.42
1:A:426:ILE:HG22	1:B:223:ILE:HD11	2.02	0.42
1:A:430:LEU:HD22	1:B:223:ILE:CD1	2.50	0.42
1:A:154:ILE:O	1:A:158:ILE:HG12	2.20	0.41
1:A:219:PHE:CE2	1:B:426:ILE:HG23	2.55	0.41
1:A:410:ILE:O	1:A:414:GLU:HG3	2.20	0.41
3:F:186:GLU:HG2	3:F:210:ARG:NH1	2.34	0.41
2:E:107:TYR:HB3	3:F:33:HIS:CD2	2.55	0.41
1:A:47:GLY:O	1:A:51:VAL:HG23	2.19	0.41
1:A:234:HIS:CD2	1:A:235:GLU:HG2	2.54	0.41
1:A:420:GLN:H	1:A:420:GLN:HG3	1.57	0.41
1:B:116:LEU:HB3	1:B:206:PRO:HD3	2.01	0.41
1:B:421:LEU:C	1:B:424:PRO:HD2	2.45	0.41
3:F:87:CYS:O	3:F:98:GLY:N	2.52	0.41
1:B:47:GLY:O	1:B:51:VAL:HG23	2.19	0.41
2:E:60:TYR:HE2	2:E:70:ILE:HG13	1.86	0.41
1:A:216:LYS:HE2	1:B:433:THR:HG22	2.03	0.41
1:B:243:LYS:HB3	2:E:31:ARG:HH21	1.85	0.41
3:D:148:LYS:HB2	3:D:192:THR:OG1	2.20	0.41
1:A:305:LEU:HA	1:A:308:VAL:HG22	2.02	0.41
1:A:330:MET:HE3	1:A:330:MET:HA	2.03	0.41
1:A:430:LEU:HD22	1:B:223:ILE:HD12	2.02	0.41
1:B:337:PHE:O	1:B:341:VAL:HG23	2.20	0.41
2:E:188:SER:HB2	3:F:134:PHE:CD2	2.56	0.41
1:A:100:TYR:O	1:A:126:ARG:NH1	2.52	0.41
1:A:207:GLN:HG2	1:B:28:ARG:HE	1.86	0.41
3:F:8:PRO:O	3:F:101:THR:HG23	2.21	0.41
3:F:60:ARG:HH21	3:F:81:ASP:CG	2.27	0.41
1:A:75:TYR:HB3	1:A:76:PRO:HD3	2.02	0.41
1:A:320:ILE:HG23	1:A:365:THR:HG21	2.02	0.41
1:A:421:LEU:C	1:A:424:PRO:HD2	2.46	0.41
1:B:19:ARG:HA	1:B:19:ARG:HD2	1.94	0.41
1:B:75:TYR:HB3	1:B:76:PRO:HD3	2.03	0.41
1:B:356:ILE:C	1:B:359:PRO:HD2	2.46	0.41
3:D:16:GLY:N	3:D:77:MET:O	2.50	0.41
3:D:49:ASP:O	3:D:51:SER:N	2.46	0.41
2:E:105:TYR:CD2	3:F:91:SER:HA	2.56	0.41
1:B:109:ILE:HG23	1:B:204:MET:SD	2.61	0.41
1:B:198:LEU:HD13	1:B:201:ILE:HD11	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:410:ILE:O	1:B:414:GLU:HG3	2.21	0.41
1:A:269:PHE:O	1:A:273:VAL:HG12	2.21	0.40
1:A:419:TYR:OH	1:B:414:GLU:OE1	2.26	0.40
1:B:172:GLU:HG3	1:B:212:LEU:O	2.21	0.40
1:A:123:ARG:HE	1:A:126:ARG:HD2	1.87	0.40
3:D:83:ALA:O	3:D:103:LEU:N	2.52	0.40
2:E:162:TRP:CH2	2:E:203:CYS:HB3	2.56	0.40
1:B:255:TYR:CE2	1:B:424:PRO:HB3	2.57	0.40
1:A:360:MET:HE1	1:A:402:ILE:HG13	2.03	0.40
1:A:453:LEU:HB3	1:B:22:ILE:HG12	2.03	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%)	423 (96%)	18 (4%)	1 (0%)	43	72
1	B	439/446 (98%)	419 (95%)	18 (4%)	2 (0%)	24	57
2	C	219/222 (99%)	200 (91%)	17 (8%)	2 (1%)	14	45
2	E	219/222 (99%)	203 (93%)	15 (7%)	1 (0%)	24	57
3	D	209/211 (99%)	190 (91%)	16 (8%)	3 (1%)	9	36
3	F	209/211 (99%)	192 (92%)	17 (8%)	0	100	100
All	All	1737/1758 (99%)	1627 (94%)	101 (6%)	9 (0%)	24	57

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	E	62	PRO
2	C	62	PRO

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Mol	Chain	Res	Type
3	D	76	THR
1	A	107	SER
1	B	107	SER
1	B	148	ALA
3	D	169	ASP
3	D	67	GLY
2	C	157	PRO

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%)	310 (93%)	24 (7%)	13	40
1	B	331/336 (98%)	306 (92%)	25 (8%)	12	39
2	C	181/182 (100%)	172 (95%)	9 (5%)	22	52
2	E	181/182 (100%)	167 (92%)	14 (8%)	12	38
3	D	185/185 (100%)	178 (96%)	7 (4%)	29	59
3	F	185/185 (100%)	176 (95%)	9 (5%)	22	52
All	All	1397/1406 (99%)	1309 (94%)	88 (6%)	16	45

All (88) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	70	HIS
1	A	73	ASP
1	A	79	LEU
1	A	81	VAL
1	A	103	GLU
1	A	151	THR
1	A	162	VAL
1	A	180	THR
1	A	200	ILE
1	A	211	THR
1	A	212	LEU

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Mol	Chain	Res	Type
1	A	219	PHE
1	A	241	VAL
1	A	251	THR
1	A	264	ILE
1	A	273	VAL
1	A	397	LEU
1	A	420	GLN
1	A	423	LEU
1	A	430	LEU
1	A	444	LEU
1	A	451	ARG
1	A	452	THR
1	A	453	LEU
1	B	41	VAL
1	B	70	HIS
1	B	78	LEU
1	B	79	LEU
1	B	81	VAL
1	B	103	GLU
1	B	151	THR
1	B	162	VAL
1	B	200	ILE
1	B	211	THR
1	B	212	LEU
1	B	219	PHE
1	B	230	ARG
1	B	241	VAL
1	B	251	THR
1	B	273	VAL
1	B	391	ILE
1	B	397	LEU
1	B	420	GLN
1	B	423	LEU
1	B	430	LEU
1	B	444	LEU
1	B	451	ARG
1	B	452	THR
1	B	453	LEU
2	C	21	SER
2	C	30	SER
2	C	43	LYS
2	C	65	LYS

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Mol	Chain	Res	Type
2	C	66	ASP
2	C	72	ARG
2	C	115	THR
2	C	151	LYS
2	C	167	LEU
3	D	1	ASP
3	D	39	SER
3	D	77	MET
3	D	125	THR
3	D	131	VAL
3	D	156	ASN
3	D	192	THR
2	E	5	LEU
2	E	21	SER
2	E	30	SER
2	E	43	LYS
2	E	65	LYS
2	E	66	ASP
2	E	96	CYS
2	E	98	ARG
2	E	115	THR
2	E	123	LYS
2	E	151	LYS
2	E	156	GLU
2	E	167	LEU
2	E	188	SER
3	F	1	ASP
3	F	17	ASP
3	F	77	MET
3	F	80	GLU
3	F	106	LEU
3	F	125	THR
3	F	131	VAL
3	F	156	ASN
3	F	192	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	175	HIS
1	A	284	HIS
1	B	157	ASN

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Mol	Chain	Res	Type
1	B	175	HIS
1	B	284	HIS
2	C	163	ASN
3	D	75	ASN
3	D	209	ASN
3	F	37	GLN
3	F	136	ASN
3	F	209	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no 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	444/446 (99%)	0.08	18 (4%)	41 24	37, 57, 85, 128	0
1	B	441/446 (98%)	0.17	19 (4%)	40 23	37, 61, 96, 128	0
2	C	221/222 (99%)	-0.02	8 (3%)	46 27	29, 56, 95, 126	0
2	E	221/222 (99%)	0.02	7 (3%)	50 30	29, 54, 89, 133	0
3	D	211/211 (100%)	0.31	7 (3%)	49 29	37, 66, 97, 120	0
3	F	211/211 (100%)	-0.00	2 (0%)	81 64	31, 50, 99, 120	0
All	All	1749/1758 (99%)	0.10	61 (3%)	47 28	29, 58, 95, 133	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	79	ALA	8.5
1	A	445	TYR	6.0
3	D	167	SER	5.7
1	A	235	GLU	4.8
1	B	104	ALA	4.7
1	A	447	ALA	4.6
1	B	288	ILE	4.5
1	A	444	LEU	4.4
1	A	460	GLN	4.1
3	D	20	THR	4.0
3	D	82	ALA	3.9
1	A	354	GLY	3.8
1	B	95	PHE	3.8
1	A	171	ASP	3.8
2	E	59	ASN	3.4
1	B	72	ALA	3.3
2	E	31	ARG	3.2
2	E	7	SER	3.2
1	B	357	PHE	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	366	VAL	3.1
1	B	307	PHE	3.1
2	C	199	GLU	3.1
3	F	150	ASP	3.0
1	B	73	ASP	3.0
1	A	69	VAL	2.9
1	A	72	ALA	2.9
3	F	127	GLY	2.9
1	B	186	LEU	2.9
2	C	141	ALA	2.8
2	E	222	ALA	2.8
3	D	80	GLU	2.8
1	B	190	PHE	2.8
1	B	234	HIS	2.7
1	A	258	LEU	2.6
2	C	139	ALA	2.5
1	B	402	ILE	2.5
1	B	355	GLY	2.5
1	A	443	PRO	2.5
2	C	42	GLY	2.4
1	B	101	ALA	2.4
3	D	39	SER	2.4
1	B	445	TYR	2.4
1	A	121	PRO	2.4
3	D	122	GLU	2.4
1	A	288	ILE	2.3
1	A	212	LEU	2.3
2	E	221	ARG	2.3
1	A	380	PRO	2.3
1	A	446	SER	2.2
1	A	29	ASP	2.2
1	B	235	GLU	2.2
1	B	185	GLY	2.2
1	B	289	THR	2.2
2	C	29	TYR	2.2
2	C	89	GLU	2.2
2	C	140	ALA	2.2
1	B	287	ASN	2.1
1	A	255	TYR	2.1
2	E	115	THR	2.1
2	E	100	TYR	2.0
2	C	179	GLN	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 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.