



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2020 – 02:12 PM BST

PDB ID : 6K5I
Title : Crystal structure of the E148D/R147A/F317A mutant CLC-ec1 in the presence of 20 mM NaBr
Authors : Park, K.; Lim, H.H.
Deposited on : 2019-05-29
Resolution : 3.02 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.13.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

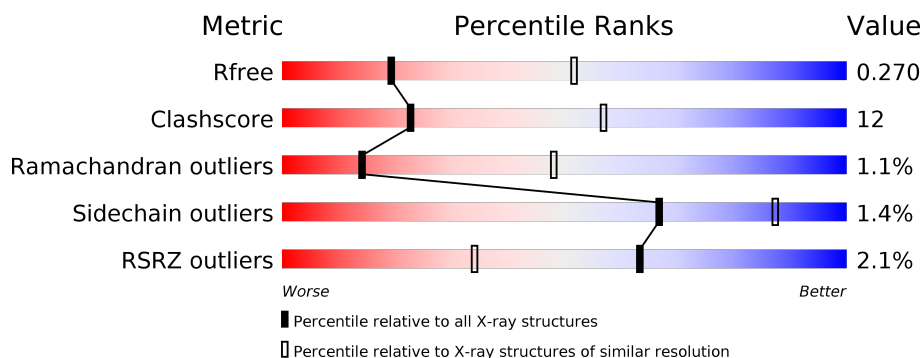
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.02 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}	130704	2399 (3.04-3.00)
Clashscore	141614	2734 (3.04-3.00)
Ramachandran outliers	138981	2640 (3.04-3.00)
Sidechain outliers	138945	2643 (3.04-3.00)
RSRZ outliers	127900	2287 (3.04-3.00)

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 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	473	<div> <div>%</div> <div> <div></div> <div>67%</div> <div>26%</div> <div>6%</div> </div> </div>
1	B	473	<div> <div>2%</div> <div> <div></div> <div>69%</div> <div>24%</div> <div>7%</div> </div> </div>
2	C	222	<div> <div>3%</div> <div> <div></div> <div>75%</div> <div>24%</div> </div> </div>
2	E	222	<div> <div></div> <div> <div></div> <div>76%</div> <div>22%</div> </div> </div>
3	D	211	<div> <div>%</div> <div> <div></div> <div>66%</div> <div>33%</div> </div> </div>
3	F	211	<div> <div>4%</div> <div> <div></div> <div>79%</div> <div>20%</div> </div> </div>

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
			3320	2180	557	563	20			
1	B	442	Total	C	N	O	S	0	0	0
			3302	2170	554	558	20			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	147	ALA	ARG	engineered mutation	UNP E9TIA0
A	148	ASP	GLU	engineered mutation	UNP E9TIA0
A	317	ALA	PHE	engineered mutation	UNP E9TIA0
B	147	ALA	ARG	engineered mutation	UNP E9TIA0
B	148	ASP	GLU	engineered mutation	UNP E9TIA0
B	317	ALA	PHE	engineered mutation	UNP E9TIA0

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	222	Total	C	N	O	S	0	0	0
			1681	1082	275	318	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) (labeled as "Ligand of

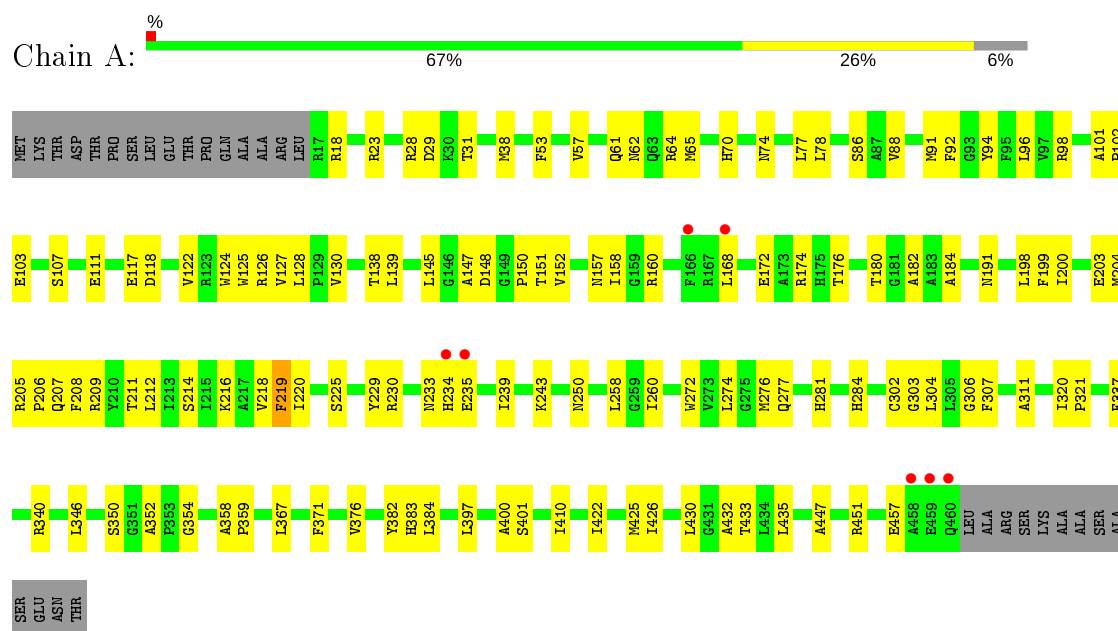
Interest" by author).

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

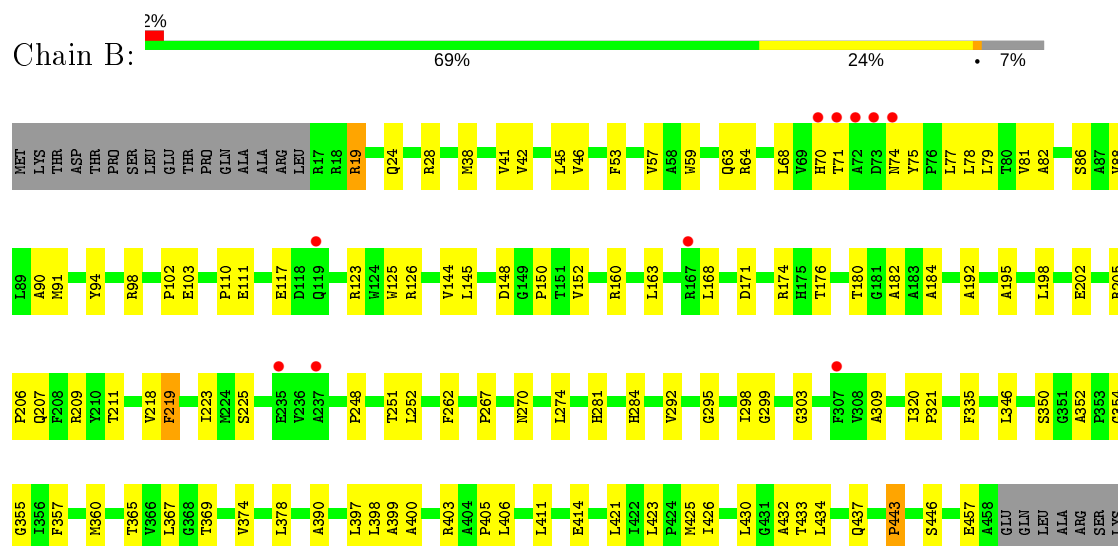
3 Residue-property plots

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

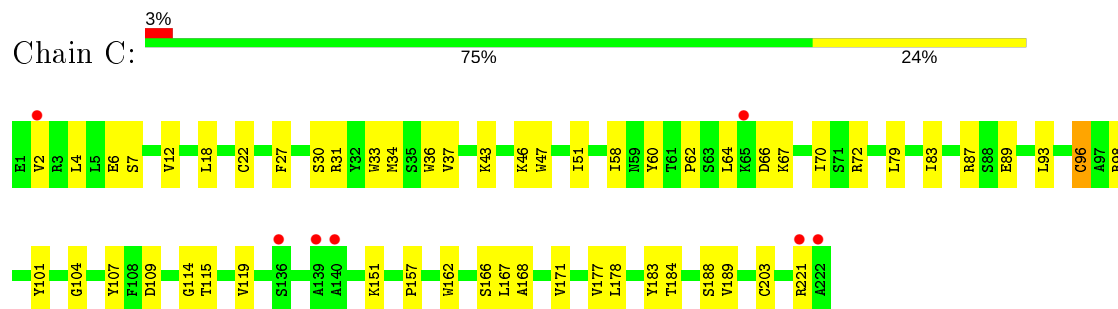


- Molecule 1: H(+)/Cl(-) exchange transporter ClcA

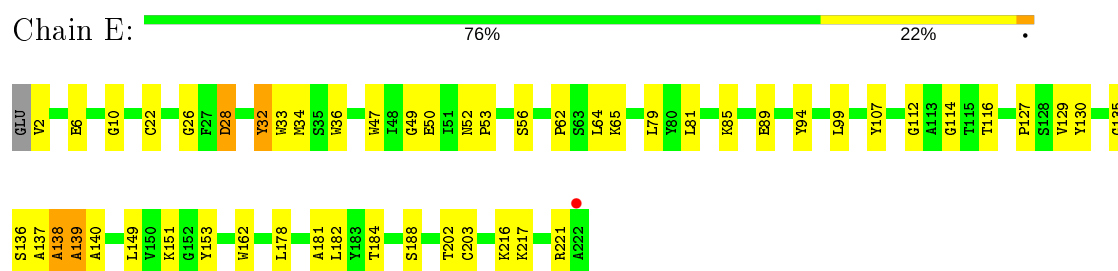


ALA
ALA
SER
ALA
SER
GLU
ASN
THR

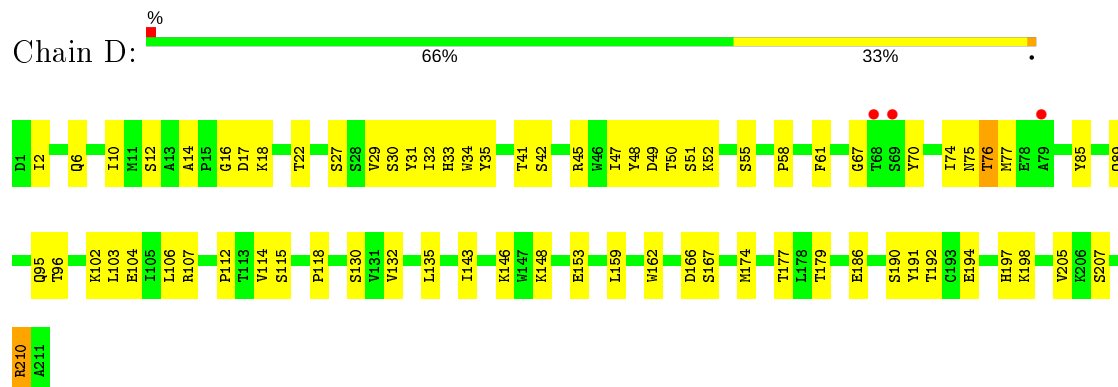
• Molecule 2: Fab fragment, heavy chain



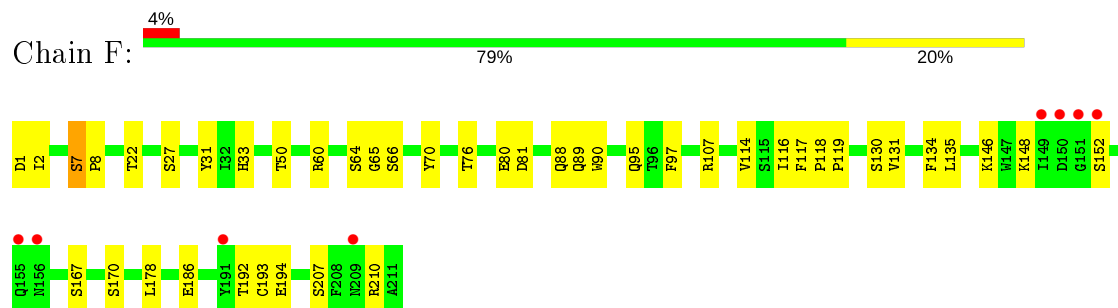
• Molecule 2: Fab fragment, heavy chain



• 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	231.91Å 98.83Å 170.29Å 90.00° 131.88° 90.00°	Depositor
Resolution (Å)	42.89 – 3.02 42.89 – 3.02	Depositor EDS
% Data completeness (in resolution range)	99.5 (42.89-3.02) 99.7 (42.89-3.02)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.49 (at 3.01Å)	Xtriage
Refinement program	PHENIX (1.13 _2998: ???)	Depositor
R, R_{free}	0.210 , 0.264 0.216 , 0.270	Depositor DCC
R_{free} test set	2886 reflections (5.14%)	wwPDB-VP
Wilson B-factor (Å ²)	84.5	Xtriage
Anisotropy	0.509	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 47.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.009 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	13219	wwPDB-VP
Average B, all atoms (Å ²)	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.19% 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: 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	0.46	0/3391	0.62	0/4604
1	B	0.43	0/3373	0.62	0/4580
2	C	0.54	0/1730	0.72	0/2367
2	E	0.53	0/1721	0.70	0/2355
3	D	0.50	0/1660	0.65	0/2257
3	F	0.48	0/1660	0.67	0/2257
All	All	0.48	0/13535	0.65	0/18420

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	3320	0	3470	94	0
1	B	3302	0	3456	90	0
2	C	1681	0	1663	31	0
2	E	1672	0	1654	37	0
3	D	1621	0	1546	53	0
3	F	1621	0	1546	39	0
4	A	1	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	1	0	0	1	0
All	All	13219	0	13335	306	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 (306) 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:D:95:GLN:N	3:D:95:GLN:OE1	1.98	0.95
1:A:239:ILE:HG23	1:A:321:PRO:HG3	1.57	0.87
3:F:95:GLN:OE1	3:F:95:GLN:N	2.10	0.85
3:D:6:GLN:NE2	3:D:85:TYR:O	2.10	0.84
1:A:430:LEU:HD11	1:B:219:PHE:CD1	2.12	0.84
2:E:47:TRP:CD2	3:F:95:GLN:NE2	2.46	0.84
3:F:7:SER:HB3	3:F:22:THR:HB	1.61	0.82
1:A:430:LEU:HD22	1:B:223:ILE:CD1	2.09	0.82
3:D:194:GLU:HG2	3:D:205:VAL:HG12	1.62	0.81
1:A:205:ARG:HH12	1:B:205:ARG:HH12	1.32	0.78
2:E:107:TYR:HB3	3:F:33:HIS:CD2	2.19	0.78
1:B:117:GLU:OE1	1:B:209:ARG:NH1	2.18	0.76
2:C:98:ARG:NH1	2:C:109:ASP:OD2	2.19	0.76
1:A:243:LYS:HB3	2:C:31:ARG:HH21	1.52	0.75
1:B:74:ASN:HB3	1:B:77:LEU:HB3	1.69	0.75
2:C:34:MET:HB3	2:C:79:LEU:HD22	1.68	0.74
1:B:400:ALA:HB2	1:B:432:ALA:HB1	1.67	0.74
3:F:95:GLN:H	3:F:95:GLN:CD	1.90	0.73
1:A:191:ASN:OD1	1:A:230:ARG:NH1	2.21	0.73
1:A:430:LEU:HD22	1:B:223:ILE:HD12	1.70	0.73
3:D:18:LYS:NZ	3:D:75:ASN:OD1	2.21	0.73
3:F:186:GLU:O	3:F:210:ARG:NH2	2.21	0.73
2:E:127:PRO:HB3	2:E:153:TYR:HB3	1.72	0.72
1:A:184:ALA:HB1	1:A:225:SER:HB2	1.73	0.71
3:D:29:VAL:O	3:D:70:TYR:OH	2.06	0.71
1:A:234:HIS:HD1	1:A:235:GLU:HG3	1.56	0.71
1:A:207:GLN:HG2	1:B:28:ARG:HH11	1.55	0.71
2:E:135:GLY:O	2:E:137:ALA:N	2.23	0.71
3:F:95:GLN:CD	3:F:95:GLN:N	2.45	0.70
1:B:123:ARG:HE	1:B:126:ARG:HD2	1.56	0.70
1:A:18:ARG:HH11	1:B:457:GLU:HB3	1.55	0.69
1:A:376:VAL:HG22	1:A:384:LEU:HB2	1.73	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:137:ALA:O	2:E:139:ALA:N	2.24	0.69
2:E:52:ASN:HD21	2:E:56:SER:HB3	1.58	0.68
1:B:38:MET:O	1:B:42:VAL:HG23	1.94	0.68
3:F:90:TRP:CG	3:F:95:GLN:HB3	2.28	0.68
2:E:47:TRP:CE3	3:F:95:GLN:NE2	2.61	0.66
1:A:206:PRO:HD2	1:A:211:THR:HG21	1.78	0.66
1:A:28:ARG:HD2	1:B:207:GLN:HG2	1.78	0.66
1:A:430:LEU:HD22	1:B:223:ILE:HD11	1.77	0.66
1:A:150:PRO:HD3	1:A:354:GLY:HA2	1.78	0.65
1:A:422:ILE:HA	1:A:425:MET:HE3	1.78	0.65
2:C:171:VAL:HG22	2:C:189:VAL:HG23	1.79	0.65
3:D:49:ASP:HB2	3:D:52:LYS:HD2	1.79	0.64
1:A:205:ARG:NH1	1:B:205:ARG:HH12	1.94	0.64
3:D:17:ASP:H	3:D:77:MET:H	1.45	0.64
3:D:197:HIS:CG	3:D:198:LYS:H	2.16	0.63
1:B:125:TRP:CD1	1:B:126:ARG:HG3	2.33	0.63
1:A:53:PHE:O	1:A:57:VAL:HG23	1.99	0.63
2:C:162:TRP:CZ3	2:C:203:CYS:HB3	2.34	0.62
1:A:31:THR:O	1:B:437:GLN:NE2	2.30	0.62
1:B:357:PHE:HE1	1:B:411:LEU:HD22	1.65	0.61
1:B:430:LEU:HA	1:B:433:THR:HG22	1.82	0.60
1:B:206:PRO:HG2	1:B:211:THR:HG21	1.83	0.60
1:A:28:ARG:HH11	1:B:207:GLN:HG2	1.66	0.60
1:A:38:MET:HG3	1:A:168:LEU:HD11	1.83	0.60
2:C:221:ARG:CZ	3:D:118:PRO:HG2	2.32	0.60
3:D:17:ASP:OD1	3:D:18:LYS:N	2.33	0.60
1:B:150:PRO:HD3	1:B:354:GLY:HA2	1.85	0.59
3:D:2:ILE:HD12	3:D:27:SER:HB2	1.84	0.59
3:F:65:GLY:HA3	3:F:70:TYR:HA	1.84	0.59
1:A:138:THR:HG21	1:A:352:ALA:HB1	1.85	0.59
1:A:219:PHE:HB3	1:B:430:LEU:HD21	1.84	0.59
3:D:148:LYS:HB2	3:D:192:THR:OG1	2.03	0.59
3:F:119:PRO:HD3	3:F:131:VAL:HG22	1.86	0.58
1:B:180:THR:HG22	1:B:218:VAL:HG22	1.84	0.58
1:B:163:LEU:HD12	1:B:168:LEU:HB2	1.86	0.58
3:D:2:ILE:O	3:D:96:THR:HG21	2.04	0.57
3:D:146:LYS:NZ	3:D:153:GLU:OE1	2.37	0.57
2:E:162:TRP:CZ3	2:E:203:CYS:HB3	2.39	0.57
3:D:114:VAL:HG22	3:D:135:LEU:HD22	1.86	0.57
2:E:47:TRP:CG	3:F:95:GLN:NE2	2.72	0.57
3:D:162:TRP:CE2	3:D:174:MET:HG3	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:10:ILE:HG23	3:D:102:LYS:HB3	1.86	0.57
1:A:216:LYS:HD3	1:B:434:LEU:HD23	1.87	0.57
1:A:426:ILE:HG23	1:B:219:PHE:CE1	2.40	0.57
1:B:360:MET:HE3	1:B:398:LEU:HD23	1.85	0.57
3:D:48:TYR:CE2	3:D:52:LYS:HD3	2.40	0.57
1:A:117:GLU:OE1	1:A:209:ARG:NH1	2.37	0.56
3:D:31:TYR:HA	3:D:50:THR:OG1	2.05	0.56
2:C:37:VAL:HG22	2:C:47:TRP:HA	1.87	0.56
1:A:124:TRP:HA	1:A:157:ASN:HD22	1.68	0.56
2:C:60:TYR:HE1	2:C:70:ILE:HG13	1.71	0.56
3:D:75:ASN:O	3:D:76:THR:HG22	2.05	0.56
1:A:94:TYR:CZ	1:A:352:ALA:HB2	2.41	0.56
2:E:32:TYR:H	2:E:53:PRO:HG3	1.71	0.56
1:A:219:PHE:CG	1:B:430:LEU:HD11	2.40	0.56
1:B:41:VAL:O	1:B:45:LEU:HG	2.06	0.56
3:D:29:VAL:CG1	3:D:32:ILE:HD11	2.36	0.56
3:D:16:GLY:HA2	3:D:76:THR:OG1	2.06	0.55
1:A:152:VAL:HG13	1:A:182:ALA:HB1	1.87	0.55
1:A:88:VAL:HA	1:A:91:MET:HE2	1.89	0.55
3:D:14:ALA:O	3:D:17:ASP:HB3	2.06	0.55
1:B:38:MET:HG3	1:B:168:LEU:HD11	1.89	0.55
1:B:320:ILE:HG23	1:B:365:THR:HG21	1.88	0.55
1:B:262:PHE:CE2	1:B:367:LEU:HD23	2.42	0.55
1:A:260:ILE:HG23	1:A:435:LEU:HG	1.88	0.55
3:F:148:LYS:HB2	3:F:192:THR:HG23	1.88	0.55
2:C:151:LYS:HA	2:C:184:THR:HG23	1.87	0.55
1:A:18:ARG:NH1	1:B:457:GLU:HB3	2.22	0.54
1:B:176:THR:O	1:B:180:THR:HG23	2.08	0.54
1:A:78:LEU:HD11	1:A:307:PHE:CE2	2.43	0.54
1:A:98:ARG:NH2	1:A:102:PRO:HB3	2.22	0.54
1:A:122:VAL:HB	1:A:160:ARG:HG2	1.90	0.54
3:F:60:ARG:NH2	3:F:81:ASP:OD1	2.40	0.54
1:A:86:SER:OG	1:A:303:GLY:HA3	2.08	0.54
1:A:430:LEU:HD11	1:B:219:PHE:CG	2.43	0.54
1:B:206:PRO:CG	1:B:211:THR:HG21	2.39	0.53
3:D:130:SER:OG	3:D:179:THR:HG23	2.08	0.53
3:D:30:SER:HB3	3:D:31:TYR:CD2	2.43	0.53
1:A:101:ALA:HB3	1:A:130:VAL:HG11	1.89	0.53
2:C:178:LEU:HB2	2:C:183:TYR:CE1	2.43	0.53
1:B:94:TYR:OH	1:B:352:ALA:HB2	2.09	0.53
3:D:132:VAL:HG22	3:D:177:THR:HG23	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:430:LEU:HD21	1:B:219:PHE:HB3	1.91	0.52
3:D:30:SER:HB3	3:D:31:TYR:HD2	1.73	0.52
1:A:243:LYS:HB3	2:C:31:ARG:NH2	2.21	0.52
1:A:447:ALA:O	1:A:451:ARG:HG3	2.10	0.52
2:E:6:GLU:HA	2:E:22:CYS:HA	1.91	0.51
1:B:346:LEU:O	1:B:350:SER:HB3	2.10	0.51
1:B:68:LEU:HD21	1:B:82:ALA:HB2	1.91	0.51
2:E:135:GLY:HA2	2:E:221:ARG:HD2	1.93	0.51
2:E:178:LEU:HD11	2:E:181:ALA:HA	1.93	0.51
2:E:138:ALA:O	2:E:140:ALA:N	2.44	0.51
2:E:28:ASP:O	2:E:32:TYR:HD2	1.93	0.50
1:A:337:PHE:CE1	1:A:367:LEU:HB2	2.46	0.50
3:D:31:TYR:HB3	3:D:49:ASP:HA	1.93	0.50
2:E:47:TRP:HZ2	2:E:50:GLU:HG2	1.76	0.50
2:C:51:ILE:HG13	2:C:58:ILE:HG12	1.93	0.50
1:B:320:ILE:HB	1:B:321:PRO:HD3	1.93	0.50
2:C:2:VAL:HG13	2:C:27:PHE:CD1	2.46	0.50
3:F:192:THR:HB	3:F:207:SER:HB3	1.94	0.50
3:F:90:TRP:CD2	3:F:95:GLN:HB3	2.45	0.50
3:D:34:TRP:HB2	3:D:47:ILE:HB	1.94	0.50
1:A:426:ILE:HG23	1:B:219:PHE:HE1	1.77	0.50
3:D:12:SER:HA	3:D:104:GLU:O	2.11	0.50
1:B:71:THR:O	1:B:78:LEU:HD23	2.12	0.50
1:A:207:GLN:HG2	1:B:28:ARG:NH1	2.26	0.50
1:A:139:LEU:HD22	1:A:145:LEU:HB2	1.95	0.49
1:A:208:PHE:CE1	1:B:24:GLN:HB3	2.48	0.49
1:B:64:ARG:O	1:B:68:LEU:HG	2.12	0.49
1:B:281:HIS:HA	1:B:284:HIS:CE1	2.47	0.49
1:B:94:TYR:CZ	1:B:352:ALA:HB2	2.47	0.49
1:A:457:GLU:OE1	1:B:19:ARG:HD2	2.12	0.49
2:E:89:GLU:N	2:E:89:GLU:OE1	2.46	0.49
1:B:192:ALA:HB1	1:B:195:ALA:HB3	1.95	0.49
1:A:400:ALA:HB2	1:A:432:ALA:HB1	1.94	0.48
1:B:98:ARG:NH2	1:B:102:PRO:HB3	2.27	0.48
2:E:22:CYS:HB3	2:E:79:LEU:HB3	1.95	0.48
3:F:66:SER:HA	3:F:70:TYR:CZ	2.48	0.48
1:A:281:HIS:HA	1:A:284:HIS:CE1	2.48	0.48
2:C:36:TRP:CZ3	2:C:96:CYS:HB2	2.48	0.48
1:B:248:PRO:O	1:B:251:THR:HB	2.14	0.48
2:E:130:TYR:HD2	2:E:149:LEU:HD23	1.79	0.48
1:A:430:LEU:HD11	1:B:219:PHE:CE1	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:74:ILE:HG22	3:D:76:THR:O	2.14	0.48
3:F:7:SER:CB	3:F:8:PRO:HD3	2.43	0.48
3:F:88:GLN:HB2	3:F:97:PHE:CD1	2.49	0.48
1:A:180:THR:HG22	1:A:218:VAL:HG22	1.95	0.47
3:F:31:TYR:HA	3:F:50:THR:OG1	2.14	0.47
1:A:311:ALA:O	1:A:340:ARG:HD2	2.15	0.47
1:A:122:VAL:HG11	1:A:160:ARG:HB2	1.97	0.47
1:A:200:ILE:HA	1:A:204:MET:HB2	1.96	0.47
1:A:92:PHE:O	1:A:96:LEU:HD23	2.14	0.47
3:D:112:PRO:HG3	3:D:143:ILE:HD11	1.96	0.47
3:D:29:VAL:HG12	3:D:32:ILE:HD11	1.96	0.47
1:A:127:VAL:HB	1:A:157:ASN:ND2	2.30	0.47
3:F:148:LYS:HA	3:F:152:SER:O	2.15	0.47
2:C:7:SER:HA	2:C:115:THR:HG21	1.96	0.47
3:D:12:SER:HB3	3:D:106:LEU:HB2	1.97	0.47
3:D:186:GLU:HA	3:D:210:ARG:NH2	2.30	0.46
3:D:6:GLN:HA	3:D:22:THR:O	2.14	0.46
1:A:125:TRP:CD1	1:A:126:ARG:HG3	2.51	0.46
1:A:117:GLU:HA	1:A:209:ARG:HH22	1.79	0.46
1:B:160:ARG:HA	1:B:160:ARG:HD2	1.72	0.46
1:A:214:SER:O	1:A:218:VAL:HG23	2.15	0.46
1:A:229:TYR:O	1:A:233:ASN:HB2	2.15	0.46
1:B:171:ASP:HA	1:B:174:ARG:NH1	2.30	0.46
1:B:219:PHE:O	1:B:223:ILE:HG13	2.16	0.46
1:B:88:VAL:HA	1:B:91:MET:HE2	1.98	0.46
2:C:18:LEU:HD11	2:C:83:ILE:HD12	1.98	0.46
2:E:10:GLY:N	2:E:116:THR:O	2.48	0.46
1:A:198:LEU:HG	1:A:410:ILE:HD12	1.98	0.46
2:E:34:MET:HB3	2:E:79:LEU:HD22	1.98	0.46
1:B:110:PRO:HG2	4:B:501:BR:BR	2.71	0.45
3:F:80:GLU:HA	3:F:167:SER:O	2.15	0.45
1:B:86:SER:OG	1:B:303:GLY:HA3	2.16	0.45
1:A:239:ILE:HD13	1:A:320:ILE:HG21	1.99	0.45
2:E:47:TRP:CZ2	2:E:49:GLY:HA2	2.51	0.45
1:A:147:ALA:O	1:A:151:THR:HG23	2.17	0.45
3:D:197:HIS:CD2	3:D:198:LYS:H	2.33	0.45
1:B:202:GLU:OE2	1:B:405:PRO:HD2	2.17	0.45
3:F:130:SER:HA	3:F:178:LEU:O	2.17	0.45
1:A:358:ALA:HB3	1:A:359:PRO:HD3	1.98	0.45
1:A:302:CYS:O	1:A:306:GLY:N	2.47	0.45
1:A:383:HIS:HD2	2:C:33:TRP:CZ3	2.35	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:93:LEU:HD11	2:C:114:GLY:HA3	1.98	0.45
1:B:91:MET:HG2	1:B:292:VAL:O	2.17	0.45
2:C:166:SER:O	2:C:168:ALA:N	2.49	0.45
2:E:36:TRP:NE1	2:E:81:LEU:HB2	2.31	0.45
3:F:89:GLN:NE2	3:F:95:GLN:HA	2.32	0.45
3:D:49:ASP:O	3:D:51:SER:N	2.49	0.45
2:E:33:TRP:HB2	2:E:99:LEU:HB2	1.99	0.45
3:D:95:GLN:H	3:D:95:GLN:CD	2.02	0.44
1:A:103:GLU:HB2	1:A:111:GLU:CD	2.37	0.44
1:A:430:LEU:HA	1:A:433:THR:HG22	1.99	0.44
2:C:6:GLU:HA	2:C:22:CYS:HA	1.99	0.44
3:D:89:GLN:O	3:D:95:GLN:HB2	2.17	0.44
2:C:47:TRP:CE2	3:D:95:GLN:NE2	2.85	0.44
3:D:35:TYR:CD2	3:D:45:ARG:HA	2.53	0.44
1:B:397:LEU:HA	1:B:397:LEU:HD23	1.77	0.44
1:B:399:ALA:O	1:B:403:ARG:HA	2.18	0.44
1:A:198:LEU:HD11	1:B:198:LEU:HD11	1.99	0.44
2:C:37:VAL:HG13	2:C:46:LYS:O	2.18	0.44
1:A:199:PHE:CE1	1:A:203:GLU:HG2	2.53	0.44
3:D:89:GLN:NE2	3:D:95:GLN:HA	2.33	0.44
3:F:114:VAL:HA	3:F:134:PHE:O	2.18	0.44
3:F:60:ARG:HD2	3:F:81:ASP:OD2	2.18	0.44
1:A:61:GLN:HG2	1:A:64:ARG:HH21	1.82	0.44
2:E:85:LYS:HD3	2:E:85:LYS:HA	1.62	0.44
1:B:94:TYR:CE1	1:B:295:GLY:HA3	2.54	0.43
1:B:421:LEU:O	1:B:425:MET:HG3	2.18	0.43
2:C:30:SER:O	2:C:31:ARG:HB2	2.18	0.43
1:A:128:LEU:HD21	1:A:158:ILE:HG13	2.00	0.43
1:A:346:LEU:O	1:A:350:SER:HB3	2.18	0.43
2:E:221:ARG:NH2	3:F:118:PRO:HD2	2.33	0.43
2:E:151:LYS:HA	2:E:184:THR:HG23	1.99	0.43
3:D:191:TYR:O	3:D:207:SER:HB2	2.18	0.43
3:F:135:LEU:N	3:F:135:LEU:HD12	2.33	0.43
3:F:146:LYS:HB3	3:F:194:GLU:HB2	2.01	0.43
1:A:397:LEU:O	1:A:401:SER:HB2	2.18	0.43
1:B:184:ALA:HB1	1:B:225:SER:HB2	1.99	0.43
1:B:42:VAL:O	1:B:46:VAL:HG23	2.18	0.43
3:F:89:GLN:O	3:F:95:GLN:HB2	2.18	0.43
1:B:150:PRO:CD	1:B:354:GLY:HA2	2.48	0.43
1:A:180:THR:HG22	1:A:218:VAL:HA	2.00	0.43
1:A:250:ASN:ND2	1:A:382:TYR:HE1	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:152:VAL:HG13	1:B:182:ALA:HB1	2.00	0.43
1:B:90:ALA:HB2	1:B:299:GLY:HA3	2.00	0.43
2:C:107:TYR:HB3	3:D:33:HIS:CE1	2.54	0.43
2:E:94:TYR:O	2:E:114:GLY:HA2	2.19	0.43
3:F:7:SER:OG	3:F:8:PRO:CD	2.66	0.43
3:F:88:GLN:HG3	3:F:97:PHE:CE1	2.54	0.43
1:B:148:ASP:OD2	1:B:355:GLY:HA3	2.18	0.43
1:B:369:THR:OG1	1:B:390:ALA:HB2	2.18	0.43
1:B:59:TRP:O	1:B:63:GLN:HG2	2.19	0.43
3:D:197:HIS:CG	3:D:198:LYS:N	2.85	0.43
1:A:220:ILE:HG12	1:B:430:LEU:HD23	2.01	0.42
2:E:6:GLU:OE1	2:E:112:GLY:HA3	2.19	0.42
2:E:178:LEU:HD12	2:E:182:LEU:O	2.19	0.42
1:B:252:LEU:HD11	1:B:423:LEU:HD23	2.01	0.42
3:F:107:ARG:HD2	3:F:170:SER:HB2	2.01	0.42
1:B:374:VAL:O	1:B:378:LEU:HG	2.19	0.42
2:C:64:LEU:HB2	2:C:67:LYS:HB2	2.01	0.42
2:E:64:LEU:HA	2:E:64:LEU:HD23	1.82	0.42
1:A:160:ARG:HD2	1:A:160:ARG:HA	1.78	0.42
3:F:117:PHE:HA	3:F:118:PRO:HD3	1.85	0.42
1:B:68:LEU:HB3	1:B:78:LEU:HD21	2.01	0.42
1:B:75:TYR:O	1:B:79:LEU:HG	2.19	0.42
2:E:33:TRP:CH2	2:E:52:ASN:HB3	2.54	0.42
3:F:7:SER:OG	3:F:8:PRO:HD3	2.20	0.42
1:A:176:THR:O	1:A:180:THR:HG23	2.20	0.42
1:B:77:LEU:O	1:B:81:VAL:HG13	2.20	0.42
2:E:202:THR:HG22	2:E:217:LYS:HA	2.01	0.42
2:E:2:VAL:HA	2:E:26:GLY:HA3	2.02	0.42
1:A:172:GLU:HG3	1:A:212:LEU:HB3	2.02	0.42
2:C:101:TYR:HB2	2:C:104:GLY:HA2	2.01	0.42
2:C:43:LYS:HB3	2:C:43:LYS:HE2	1.52	0.42
3:D:166:ASP:OD1	3:D:167:SER:N	2.52	0.41
1:A:272:TRP:O	1:A:276:MET:HB2	2.20	0.41
1:A:74:ASN:HB3	1:A:77:LEU:HB3	2.01	0.41
2:E:129:VAL:O	2:E:216:LYS:HE3	2.20	0.41
2:E:36:TRP:CE2	2:E:81:LEU:HB2	2.56	0.41
3:D:106:LEU:HD23	3:D:107:ARG:N	2.36	0.41
2:C:4:LEU:HD23	2:C:4:LEU:HA	1.90	0.41
3:D:162:TRP:CD2	3:D:174:MET:HG3	2.56	0.41
1:A:29:ASP:OD1	1:A:216:LYS:HE3	2.20	0.41
1:B:274:LEU:HD23	1:B:274:LEU:HA	1.78	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:406:LEU:HD12	1:B:406:LEU:HA	1.83	0.41
3:F:186:GLU:HA	3:F:210:ARG:CZ	2.51	0.41
3:F:2:ILE:CD1	3:F:27:SER:HB2	2.51	0.41
1:B:53:PHE:O	1:B:57:VAL:HG23	2.21	0.41
2:E:127:PRO:CB	2:E:153:TYR:HB3	2.48	0.41
1:A:107:SER:N	4:A:501:BR:BR	2.98	0.41
1:B:195:ALA:N	1:B:414:GLU:OE2	2.47	0.41
1:B:267:PRO:HA	1:B:270:ASN:HB2	2.03	0.41
1:A:180:THR:HA	1:A:218:VAL:HG13	2.03	0.41
1:A:62:ASN:O	1:A:65:MET:N	2.54	0.41
1:B:145:LEU:HD22	1:B:354:GLY:HA3	2.03	0.41
3:D:58:PRO:HG2	3:D:61:PHE:HE2	1.85	0.41
3:F:116:ILE:HD12	3:F:193:CYS:SG	2.61	0.41
1:A:118:ASP:CG	1:A:174:ARG:HH21	2.23	0.40
1:B:103:GLU:O	1:B:111:GLU:HG3	2.21	0.40
1:B:262:PHE:CZ	1:B:367:LEU:HD23	2.57	0.40
1:B:443:PRO:HB2	1:B:446:SER:HB2	2.03	0.40
1:A:258:LEU:HD13	1:A:371:PHE:CG	2.56	0.40
1:A:94:TYR:OH	1:A:352:ALA:HB2	2.22	0.40
2:C:12:VAL:O	2:C:119:VAL:HA	2.21	0.40
2:C:177:VAL:HG21	3:D:159:LEU:HD13	2.04	0.40
1:A:184:ALA:HB1	1:A:225:SER:CB	2.49	0.40
1:A:219:PHE:CE1	1:B:426:ILE:HG23	2.56	0.40
2:C:87:ARG:HH21	2:C:89:GLU:CD	2.25	0.40
3:D:41:THR:HG22	3:D:42:SER:O	2.21	0.40
3:D:77:MET:SD	3:D:103:LEU:HD21	2.61	0.40
3:F:7:SER:HB2	3:F:8:PRO:HD3	2.04	0.40
1:A:274:LEU:O	1:A:277:GLN:HB2	2.22	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	442/473 (93%)	411 (93%)	31 (7%)	0	100	100
1	B	440/473 (93%)	407 (92%)	28 (6%)	5 (1%)	14	48
2	C	220/222 (99%)	202 (92%)	15 (7%)	3 (1%)	11	41
2	E	219/222 (99%)	193 (88%)	20 (9%)	6 (3%)	5	25
3	D	209/211 (99%)	189 (90%)	17 (8%)	3 (1%)	11	41
3	F	209/211 (99%)	194 (93%)	13 (6%)	2 (1%)	15	50
All	All	1739/1812 (96%)	1596 (92%)	124 (7%)	19 (1%)	14	48

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	E	65	LYS
2	E	138	ALA
3	D	55	SER
2	E	136	SER
3	F	7	SER
2	C	167	LEU
3	D	76	THR
2	E	62	PRO
2	E	139	ALA
2	C	62	PRO
3	F	76	THR
1	B	309	ALA
1	B	335	PHE
2	E	28	ASP
1	B	443	PRO
1	B	144	VAL
3	D	67	GLY
1	B	298	ILE
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	333/356 (94%)	328 (98%)	5 (2%)	65	86
1	B	331/356 (93%)	328 (99%)	3 (1%)	78	92
2	C	182/182 (100%)	178 (98%)	4 (2%)	52	80
2	E	181/182 (100%)	179 (99%)	2 (1%)	73	90
3	D	185/185 (100%)	182 (98%)	3 (2%)	62	86
3	F	185/185 (100%)	183 (99%)	2 (1%)	73	90
All	All	1397/1446 (97%)	1378 (99%)	19 (1%)	67	87

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	ARG
1	A	70	HIS
1	A	148	ASP
1	A	219	PHE
1	A	304	LEU
1	B	19	ARG
1	B	70	HIS
1	B	219	PHE
2	C	66	ASP
2	C	72	ARG
2	C	96	CYS
2	C	188	SER
3	D	115	SER
3	D	190	SER
3	D	210	ARG
2	E	32	TYR
2	E	188	SER
3	F	1	ASP
3	F	64	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	157	ASN
1	B	284	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

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 [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	444/473 (93%)	-0.15	7 (1%)	72 43	55, 78, 107, 143	0
1	B	442/473 (93%)	-0.10	10 (2%)	60 31	56, 79, 116, 141	0
2	C	222/222 (100%)	-0.27	7 (3%)	47 20	41, 70, 110, 138	0
2	E	221/222 (99%)	-0.47	1 (0%)	91 75	48, 72, 105, 140	0
3	D	211/211 (100%)	-0.18	3 (1%)	75 48	56, 80, 100, 113	0
3	F	211/211 (100%)	-0.21	8 (3%)	40 16	48, 64, 111, 129	0
All	All	1751/1812 (96%)	-0.20	36 (2%)	63 34	41, 76, 109, 143	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	222	ALA	4.4
1	A	460	GLN	4.3
1	B	70	HIS	4.1
3	F	156	ASN	3.9
2	C	65	LYS	3.8
1	B	73	ASP	3.7
2	C	140	ALA	3.6
1	B	72	ALA	3.6
1	A	235	GLU	3.6
3	F	151	GLY	3.4
1	B	235	GLU	3.3
3	F	155	GLN	3.3
1	B	71	THR	3.2
2	C	136	SER	3.1
1	A	234	HIS	3.1
1	A	458	ALA	3.0
3	D	68	THR	2.9
1	B	307	PHE	2.8
1	A	459	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	168	LEU	2.6
1	A	166	PHE	2.6
2	C	139	ALA	2.5
2	E	222	ALA	2.5
3	D	79	ALA	2.4
1	B	74	ASN	2.4
3	F	152	SER	2.4
3	F	150	ASP	2.3
2	C	221	ARG	2.3
3	F	149	ILE	2.2
1	B	119	GLN	2.2
3	F	191	TYR	2.2
1	B	237	ALA	2.2
3	F	209	ASN	2.1
3	D	69	SER	2.1
2	C	2	VAL	2.1
1	B	167	ARG	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 ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

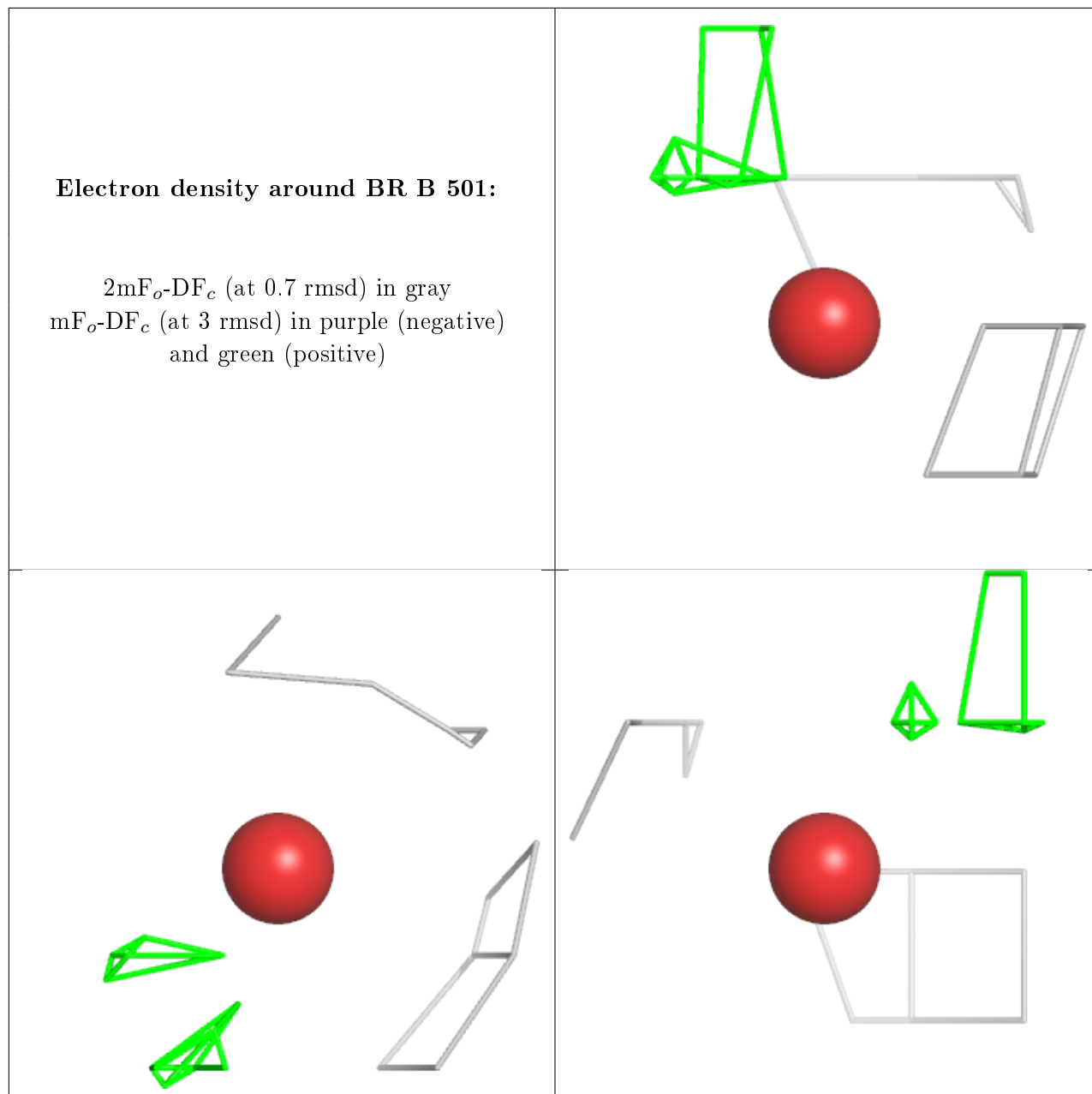
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	BR	B	501	1/1	0.94	0.23	129,129,129,129	0
4	BR	A	501	1/1	0.96	0.24	175,175,175,175	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

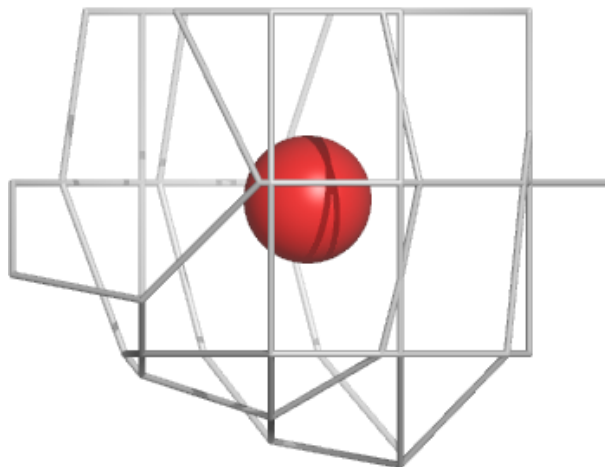
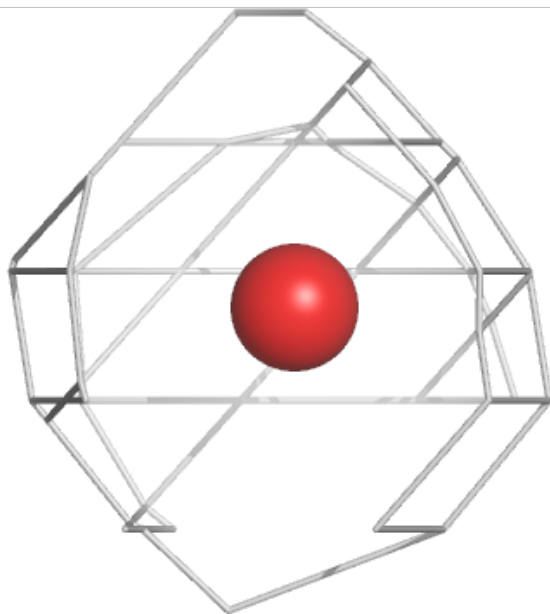
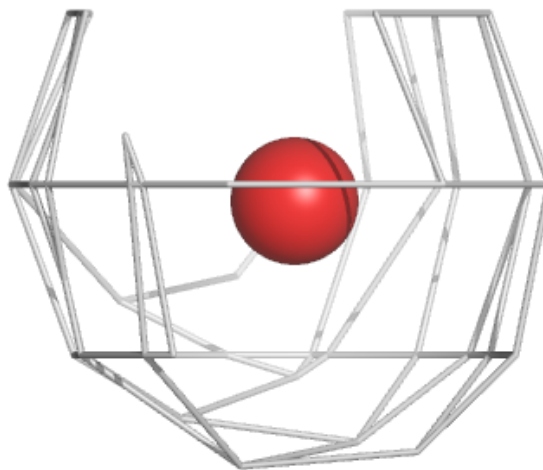
Electron density around BR B 501:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around BR A 501:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.