



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:00 AM UTC

PDB ID : 4KI5 / pdb_00004ki5
Title : Crystal structure of human factor VIII C2 domain in a ternary complex with murine inhibitory antibodies 3E6 and G99
Authors : Walter, J.D.; Meeks, S.L.; Healey, J.F.; Lollar, P.; Spiegel, P.C.
Deposited on : 2013-05-01
Resolution : 2.47 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

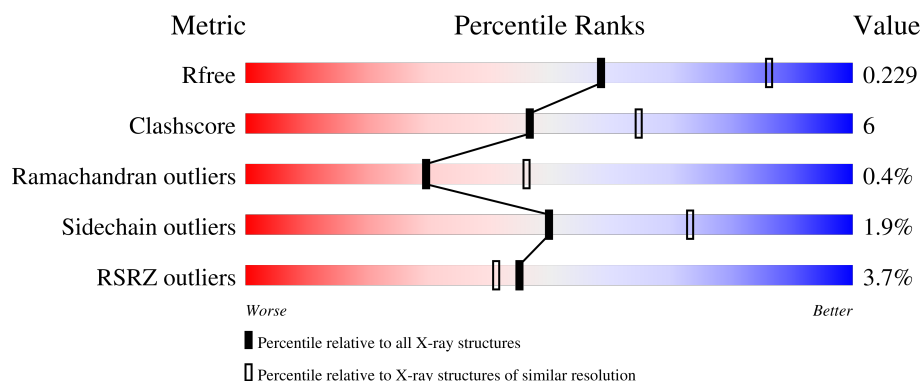
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.47 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	7589 (2.50-2.46)
Clashscore	190562	8295 (2.50-2.46)
Ramachandran outliers	187476	8164 (2.50-2.46)
Sidechain outliers	187428	8166 (2.50-2.46)
RSRZ outliers	180081	7593 (2.50-2.46)

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	C	219	3% 89% 9% .
2	D	213	2% 84% 15% .
3	E	224	2% 87% 12% .
4	F	214	5% 78% 19% ..
5	M	183	5% 75% 9% 16%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8074 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 MURINE MONOCLONAL 3E6 FAB HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	218	Total	C	N	O	S	0	0	0
			1647	1045	270	326	6			

- Molecule 2 is a protein called MURINE MONOCLONAL 3E6 FAB LIGHT CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	211	Total	C	N	O	S	0	0	0
			1620	1009	269	333	9			

- Molecule 3 is a protein called MURINE MONOCLONAL G99 FAB HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	224	Total	C	N	O	S	0	0	0
			1685	1065	272	341	7			

- Molecule 4 is a protein called MURINE MONOCLONAL G99 FAB LIGHT CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	F	209	Total	C	N	O	S	0	0	0
			1624	1011	271	335	7			

- Molecule 5 is a protein called Coagulation factor VIII.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	M	154	Total	C	N	O	S	0	0	0
			1231	785	214	224	8			

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
M	2150	MET	-	expression tag	UNP P00451
M	2151	GLY	-	expression tag	UNP P00451

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Chain	Residue	Modelled	Actual	Comment	Reference
M	2152	SER	-	expression tag	UNP P00451
M	2153	SER	-	expression tag	UNP P00451
M	2154	HIS	-	expression tag	UNP P00451
M	2155	HIS	-	expression tag	UNP P00451
M	2156	HIS	-	expression tag	UNP P00451
M	2157	HIS	-	expression tag	UNP P00451
M	2158	HIS	-	expression tag	UNP P00451
M	2159	HIS	-	expression tag	UNP P00451
M	2160	SER	-	expression tag	UNP P00451
M	2161	SER	-	expression tag	UNP P00451
M	2162	GLY	-	expression tag	UNP P00451
M	2163	LEU	-	expression tag	UNP P00451
M	2164	VAL	-	expression tag	UNP P00451
M	2165	PRO	-	expression tag	UNP P00451
M	2166	ARG	-	expression tag	UNP P00451
M	2167	GLY	-	expression tag	UNP P00451
M	2168	SER	-	expression tag	UNP P00451
M	2169	HIS	-	expression tag	UNP P00451
M	2170	MET	-	expression tag	UNP P00451

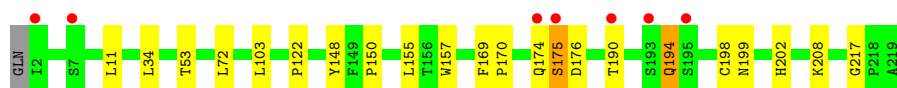
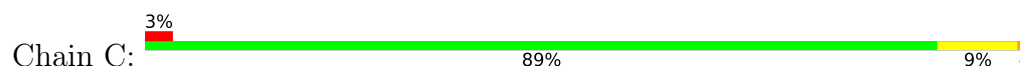
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	C	76	Total O 76 76	0	0
6	D	54	Total O 54 54	0	0
6	E	46	Total O 46 46	0	0
6	F	40	Total O 40 40	0	0
6	M	51	Total O 51 51	0	0

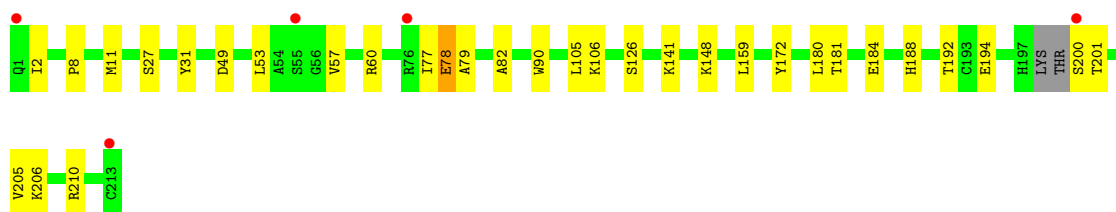
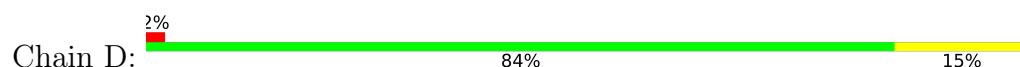
3 Residue-property plots [i](#)

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.

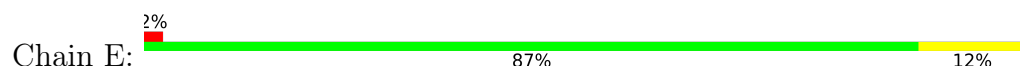
- Molecule 1: MURINE MONOCLONAL 3E6 FAB HEAVY CHAIN



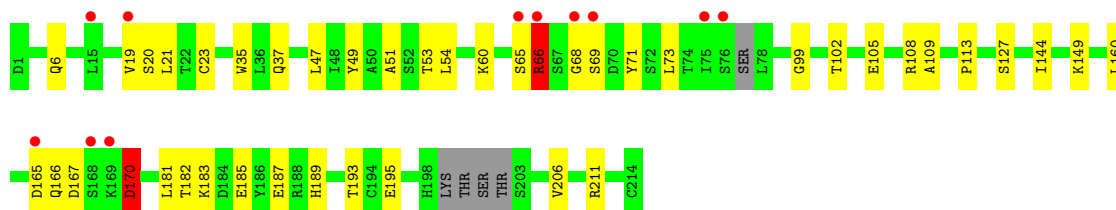
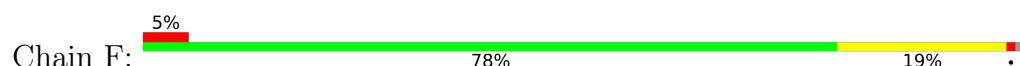
- Molecule 2: MURINE MONOCLONAL 3E6 FAB LIGHT CHAIN



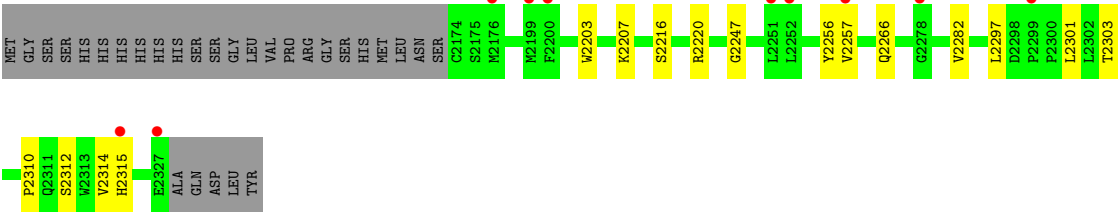
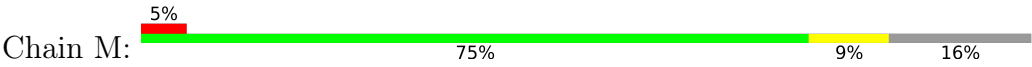
- Molecule 3: MURINE MONOCLONAL G99 FAB HEAVY CHAIN



- Molecule 4: MURINE MONOCLONAL G99 FAB LIGHT CHAIN



- Molecule 5: Coagulation factor VIII



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	54.25Å 71.16Å 277.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	56.43 – 2.47 56.43 – 2.47	Depositor EDS
% Data completeness (in resolution range)	99.8 (56.43-2.47) 99.8 (56.43-2.47)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.80 (at 2.48Å)	Xtriage
Refinement program	REFMAC, PHENIX 1.8_1069	Depositor
R, R_{free}	0.188 , 0.238 (Not available) , 0.229	Depositor DCC
R_{free} test set	2000 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	35.3	Xtriage
Anisotropy	0.663	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 42.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8074	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.36% 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	C	0.40	0/1690	0.73	3/2309 (0.1%)
2	D	0.38	0/1658	0.69	0/2253
3	E	0.43	0/1731	0.72	2/2365 (0.1%)
4	F	0.43	0/1657	0.74	0/2244
5	M	0.38	0/1263	0.69	0/1713
All	All	0.41	0/7999	0.72	5/10884 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	208	HIS	CA-C-N	5.60	125.06	119.24
3	E	208	HIS	C-N-CA	5.60	125.06	119.24
1	C	202	HIS	CA-C-N	5.59	125.05	119.24
1	C	202	HIS	C-N-CA	5.59	125.05	119.24
1	C	176	ASP	N-CA-C	-5.10	107.03	114.12

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	C	1647	0	1611	13	0
2	D	1620	0	1541	19	1
3	E	1685	0	1619	23	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	F	1624	0	1551	31	1
5	M	1231	0	1218	8	0
6	C	76	0	0	1	0
6	D	54	0	0	0	0
6	E	46	0	0	0	0
6	F	40	0	0	1	0
6	M	51	0	0	1	0
All	All	8074	0	7540	87	1

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 6.

All (87) 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:E:178:LEU:CD2	4:F:160:LEU:HB3	1.97	0.95
3:E:178:LEU:HD21	4:F:160:LEU:HB3	1.49	0.94
1:C:174:GLN:HG3	2:D:159:LEU:CD1	1.98	0.94
1:C:174:GLN:HG3	2:D:159:LEU:HD12	1.49	0.93
4:F:66:ARG:HD2	4:F:71:TYR:CE1	2.13	0.82
1:C:194:GLN:NE2	1:C:194:GLN:HA	1.95	0.79
4:F:108:ARG:NH1	4:F:109:ALA:O	2.18	0.76
2:D:8:PRO:HD2	2:D:11:MET:HE2	1.69	0.75
3:E:128:PRO:HB3	3:E:154:TYR:HB3	1.71	0.73
3:E:154:TYR:CE1	3:E:159:VAL:HG23	2.25	0.72
4:F:65:SER:O	4:F:66:ARG:HB3	1.92	0.69
4:F:195:GLU:HG2	4:F:206:VAL:HG12	1.74	0.69
2:D:194:GLU:HG2	2:D:205:VAL:HG12	1.75	0.69
2:D:2:ILE:HG12	2:D:27:SER:HB3	1.73	0.69
1:C:174:GLN:HG3	2:D:159:LEU:HD11	1.76	0.67
3:E:13:LYS:HG3	3:E:122:SER:HA	1.77	0.65
3:E:201:SER:OG	3:E:218:LYS:HE3	1.98	0.64
4:F:37:GLN:HB2	4:F:47:LEU:HD11	1.81	0.62
4:F:19:VAL:HG12	4:F:20:SER:N	2.15	0.61
5:M:2257:VAL:HG13	5:M:2312:SER:O	2.01	0.61
4:F:21:LEU:HD22	4:F:102:THR:HG21	1.83	0.60
1:C:208:LYS:NZ	6:C:361:HOH:O	2.36	0.59
1:C:53:THR:HA	1:C:72:LEU:HD21	1.85	0.57
2:D:31:TYR:HB3	2:D:49:ASP:HA	1.86	0.56
4:F:165:ASP:OD1	4:F:166:GLN:N	2.36	0.56
2:D:180:LEU:HB3	2:D:184:GLU:HG3	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:178:LEU:HD23	4:F:160:LEU:HB3	1.87	0.55
3:E:195:SER:O	3:E:199:SER:OG	2.25	0.55
3:E:38:LYS:HB2	3:E:48:ILE:HD11	1.90	0.54
3:E:154:TYR:CE1	3:E:159:VAL:CG2	2.91	0.53
4:F:66:ARG:HD2	4:F:71:TYR:CZ	2.44	0.53
2:D:200:SER:OG	2:D:201:THR:N	2.43	0.51
3:E:147:LEU:HD12	3:E:202:ILE:HD13	1.92	0.51
3:E:222:ARG:C	3:E:223:GLY:O	2.51	0.51
4:F:19:VAL:CG1	4:F:20:SER:N	2.74	0.51
4:F:66:ARG:O	4:F:66:ARG:HG3	2.09	0.50
2:D:82:ALA:HB2	2:D:105:LEU:HD13	1.93	0.50
3:E:222:ARG:O	3:E:223:GLY:O	2.30	0.50
5:M:2282:VAL:O	6:M:2448:HOH:O	2.20	0.49
2:D:60:ARG:CZ	2:D:78:GLU:HG3	2.44	0.47
3:E:178:LEU:HD21	4:F:160:LEU:CB	2.35	0.47
4:F:51:ALA:O	6:F:307:HOH:O	2.20	0.47
4:F:113:PRO:HG3	4:F:144:ILE:HD11	1.96	0.47
3:E:197:TRP:CG	3:E:198:PRO:HA	2.49	0.47
2:D:141:LYS:HG3	2:D:172:TYR:CG	2.49	0.47
1:C:190:THR:O	1:C:194:GLN:HB2	2.15	0.46
4:F:35:TRP:CD2	4:F:73:LEU:HB2	2.51	0.46
4:F:35:TRP:CE2	4:F:73:LEU:HB2	2.51	0.46
5:M:2257:VAL:HG21	5:M:2310:PRO:HB3	1.98	0.46
5:M:2314:VAL:O	5:M:2315:HIS:HB2	2.15	0.46
4:F:181:LEU:HB3	4:F:185:GLU:HG3	1.97	0.45
4:F:6:GLN:HG3	4:F:99:GLY:HA3	1.98	0.45
4:F:182:THR:OG1	4:F:185:GLU:HG2	2.17	0.45
3:E:197:TRP:CD1	3:E:198:PRO:HA	2.52	0.45
4:F:54:LEU:HD11	4:F:60:LYS:HA	1.98	0.45
5:M:2203:TRP:CD2	5:M:2220:ARG:HD3	2.52	0.45
1:C:122:PRO:HB3	1:C:148:TYR:HB3	1.99	0.44
5:M:2207:LYS:HD3	5:M:2216:SER:O	2.17	0.44
3:E:67:LYS:HE2	3:E:84:SER:O	2.18	0.44
4:F:183:LYS:HE3	4:F:183:LYS:HB3	1.81	0.44
2:D:181:THR:OG1	2:D:184:GLU:HG2	2.18	0.43
2:D:206:LYS:HD3	2:D:206:LYS:HA	1.83	0.43
3:E:7:SER:HG	3:E:21:SER:HG	1.62	0.43
4:F:6:GLN:HG2	4:F:23:CYS:SG	2.59	0.43
1:C:11:LEU:HD22	1:C:150:PRO:HD3	2.01	0.42
2:D:79:ALA:HA	2:D:105:LEU:HD22	2.00	0.42
4:F:167:ASP:OD2	4:F:170:ASP:HB3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:7:SER:OG	3:E:21:SER:OG	2.29	0.42
2:D:148:LYS:HB2	2:D:192:THR:HB	2.02	0.42
5:M:2266:GLN:HE22	5:M:2303:THR:HA	1.84	0.42
4:F:68:GLY:HA2	4:F:69:SER:HA	1.52	0.41
3:E:40:ARG:HB2	3:E:43:HIS:HB2	2.01	0.41
1:C:174:GLN:N	1:C:175:SER:HA	2.34	0.41
3:E:196:THR:O	3:E:200:GLN:HB2	2.21	0.41
3:E:161:LEU:HA	3:E:205:ASN:O	2.21	0.41
3:E:197:TRP:CD1	3:E:202:ILE:HD12	2.55	0.41
5:M:2247:GLY:HA3	5:M:2256:TYR:HB3	2.03	0.41
4:F:183:LYS:O	4:F:187:GLU:HG3	2.21	0.41
2:D:53:LEU:HG	2:D:57:VAL:HB	2.02	0.41
2:D:159:LEU:HD23	2:D:159:LEU:HA	1.91	0.41
4:F:49:TYR:CZ	4:F:53:THR:HG21	2.56	0.41
4:F:189:HIS:O	4:F:211:ARG:HD3	2.21	0.41
4:F:149:LYS:HB2	4:F:193:THR:HB	2.04	0.40
2:D:188:HIS:O	2:D:210:ARG:HD3	2.21	0.40
1:C:155:LEU:HA	1:C:199:ASN:O	2.21	0.40
1:C:169:PHE:HA	1:C:170:PRO:HD3	1.93	0.40
1:C:157:TRP:CZ3	1:C:198:CYS:HB3	2.56	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:126:SER:OG	4:F:127:SER:O[4_535]	2.01	0.19

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	C	216/219 (99%)	207 (96%)	8 (4%)	1 (0%)	24 40

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	207/213 (97%)	201 (97%)	6 (3%)	0	100	100
3	E	222/224 (99%)	216 (97%)	5 (2%)	1 (0%)	24	40
4	F	203/214 (95%)	194 (96%)	7 (3%)	2 (1%)	12	22
5	M	152/183 (83%)	145 (95%)	7 (5%)	0	100	100
All	All	1000/1053 (95%)	963 (96%)	33 (3%)	4 (0%)	30	46

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	F	170	ASP
4	F	66	ARG
1	C	217	GLY
3	E	223	GLY

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	C	185/186 (100%)	181 (98%)	4 (2%)	45	70
2	D	186/188 (99%)	182 (98%)	4 (2%)	45	70
3	E	191/191 (100%)	187 (98%)	4 (2%)	47	71
4	F	186/191 (97%)	183 (98%)	3 (2%)	55	77
5	M	139/164 (85%)	137 (99%)	2 (1%)	59	79
All	All	887/920 (96%)	870 (98%)	17 (2%)	50	73

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	34	LEU
1	C	103	LEU
1	C	175	SER
1	C	194	GLN

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Mol	Chain	Res	Type
2	D	77	ILE
2	D	78	GLU
2	D	90	TRP
2	D	106	LYS
3	E	19	LYS
3	E	67	LYS
3	E	200	GLN
3	E	201	SER
4	F	66	ARG
4	F	105	GLU
4	F	170	ASP
5	M	2297	LEU
5	M	2301	LEU

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

Mol	Chain	Res	Type
1	C	194	GLN
3	E	200	GLN

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	C	218/219 (99%)	0.35	7 (3%) 50 47	36, 48, 63, 72	0
2	D	211/213 (99%)	0.20	5 (2%) 59 56	34, 47, 62, 74	0
3	E	224/224 (100%)	0.51	5 (2%) 62 59	39, 52, 66, 77	0
4	F	209/214 (97%)	0.47	11 (5%) 32 28	35, 49, 68, 77	0
5	M	154/183 (84%)	0.53	10 (6%) 25 22	42, 54, 72, 81	0
All	All	1016/1053 (96%)	0.41	38 (3%) 45 41	34, 50, 66, 81	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
5	M	2252	LEU	4.6
2	D	1	GLN	4.5
5	M	2278	GLY	4.3
4	F	19	VAL	4.0
3	E	1	GLN	3.4
5	M	2251	LEU	3.3
1	C	195	SER	3.1
1	C	175	SER	3.1
1	C	174	GLN	3.0
3	E	178	LEU	2.9
3	E	13	LYS	2.9
5	M	2257	VAL	2.8
4	F	76	SER	2.7
4	F	68	GLY	2.7
1	C	2	ILE	2.7
4	F	66	ARG	2.6
4	F	165	ASP	2.5
5	M	2200	PHE	2.5
2	D	200	SER	2.5
4	F	69	SER	2.5

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Mol	Chain	Res	Type	RSRZ
5	M	2199	MET	2.4
4	F	65	SER	2.4
3	E	224	PRO	2.4
4	F	169	LYS	2.4
2	D	213	CYS	2.3
5	M	2299	PRO	2.3
2	D	55	SER	2.2
1	C	193	SER	2.2
5	M	2327	GLU	2.2
2	D	76	ARG	2.2
5	M	2315	HIS	2.2
1	C	190	THR	2.2
3	E	17	SER	2.1
4	F	168	SER	2.1
5	M	2176	MET	2.1
4	F	15	LEU	2.0
1	C	7	SER	2.0
4	F	75	ILE	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.