



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:19 AM UTC

PDB ID : 1XX9 / pdb_00001xx9
Title : Crystal Structure of the FXIa Catalytic Domain in Complex with EcotinM84R
Authors : Jin, L.; Pandey, P.; Babine, R.E.; Gorga, J.C.; Seidl, K.J.; Gelfand, E.;
Weaver, D.T.; Abdel-Meguid, S.S.; Strickler, J.E.
Deposited on : 2004-11-04
Resolution : 2.20 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
Mogul : 2022.3.0, CSD as543be (2022)
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

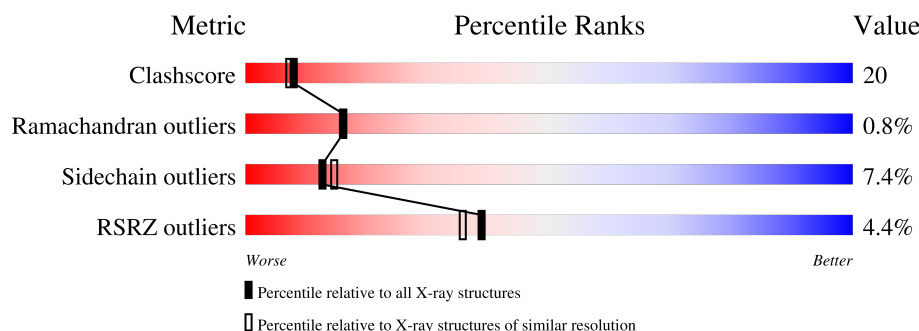
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	238	5% 57% 36% 6%
1	B	238	5% 61% 31% 5%
2	C	142	4% 64% 31%
2	D	142	3% 62% 34%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6244 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 Coagulation factor XI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	237	Total	C	N	O	S	0	1	0
			1882	1189	333	349	11			
1	B	233	Total	C	N	O	S	0	2	0
			1854	1171	325	347	11			

- Molecule 2 is a protein called Ecotin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	140	Total	C	N	O	S	0	0	0
			1120	714	188	213	5			
2	D	139	Total	C	N	O	S	0	0	0
			1114	711	187	211	5			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	84	ARG	MET	engineered mutation	UNP P23827
D	84	ARG	MET	engineered mutation	UNP P23827

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		

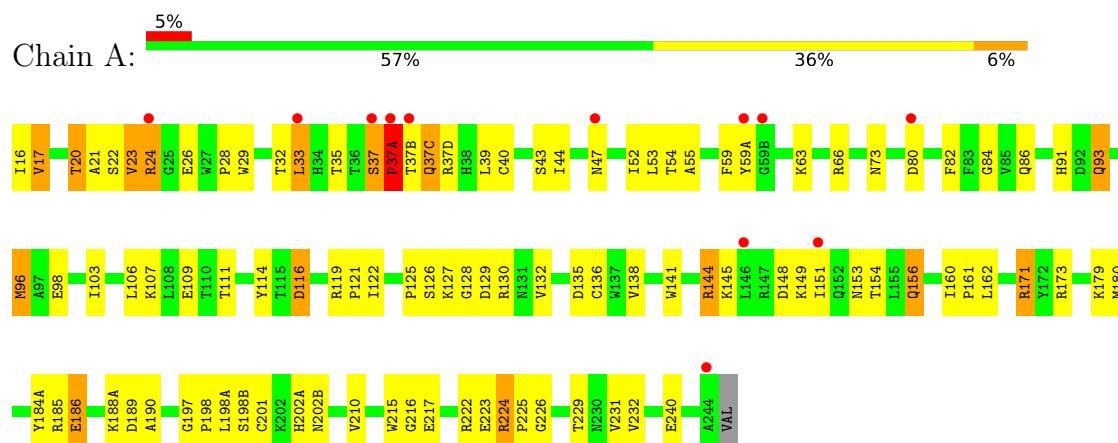
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	87	Total	O	0	0
			87	87		
4	B	81	Total	O	0	0
			81	81		
4	C	51	Total	O	0	0
			51	51		
4	D	41	Total	O	0	0
			41	41		

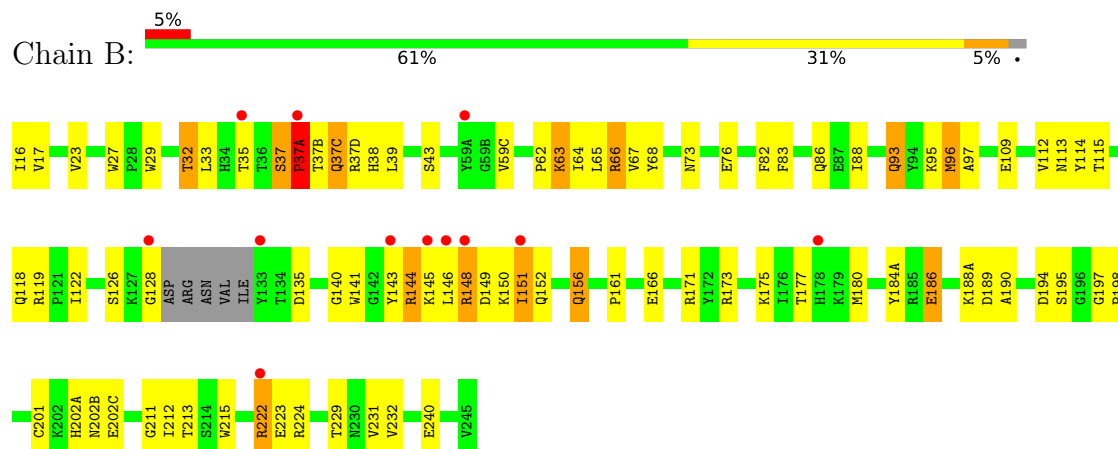
3 Residue-property plots

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.

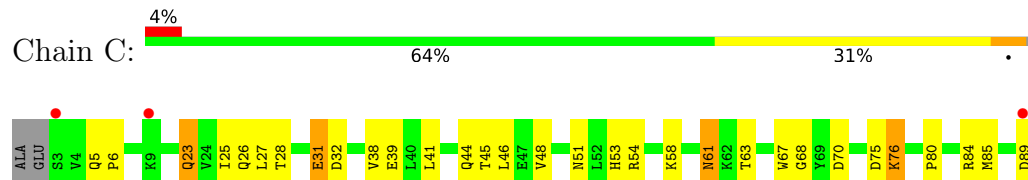
• Molecule 1: Coagulation factor XI



• Molecule 1: Coagulation factor XI

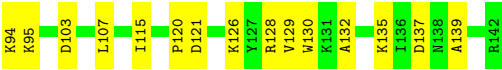


• Molecule 2: Ecotin





● Molecule 2: Ecotin



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	44.62Å 92.67Å 186.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.63 – 2.20 24.63 – 2.20	Depositor EDS
% Data completeness (in resolution range)	82.7 (24.63-2.20) 82.7 (24.63-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.45 (at 2.20Å)	Xtriage
Refinement program	CNX 2002	Depositor
R, R_{free}	0.235 , 0.285 0.237 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	31.7	Xtriage
Anisotropy	0.698	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 33.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6244	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.37% 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

5.1 Standard geometry

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

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	0.52	0/1932	0.94	7/2619 (0.3%)
1	B	0.45	0/1908	0.90	5/2583 (0.2%)
2	C	0.43	0/1142	0.94	3/1543 (0.2%)
2	D	0.42	0/1136	0.93	1/1535 (0.1%)
All	All	0.46	0/6118	0.93	16/8280 (0.2%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	231	VAL	N-CA-C	7.42	119.08	110.62
1	A	224	ARG	CA-C-N	6.52	126.49	119.78
1	A	224	ARG	C-N-CA	6.52	126.49	119.78
1	B	144	ARG	N-CA-C	-6.47	105.13	113.16
1	A	226	GLY	N-CA-C	-6.45	103.82	112.14
1	A	215	TRP	N-CA-C	6.14	116.67	108.38
2	C	135	LYS	N-CA-C	6.00	118.70	110.24
2	C	76	LYS	N-CA-C	-5.90	105.38	112.58
1	A	190	ALA	N-CA-C	-5.75	102.78	110.55
2	D	139	ALA	N-CA-C	-5.68	102.88	110.55
1	B	27	TRP	CA-C-N	5.37	125.04	119.56
1	B	27	TRP	C-N-CA	5.37	125.04	119.56
1	B	215	TRP	N-CA-C	5.34	115.59	108.38
1	A	144	ARG	N-CA-C	-5.22	106.94	113.72
2	C	89	ASP	N-CA-C	5.18	117.21	110.53
1	A	231	VAL	N-CA-C	5.09	116.42	110.62

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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	1882	0	1845	94	0
1	B	1854	0	1810	74	0
2	C	1120	0	1134	45	0
2	D	1114	0	1129	44	0
3	A	14	0	13	1	0
4	A	87	0	0	7	0
4	B	81	0	0	8	0
4	C	51	0	0	2	0
4	D	41	0	0	2	0
All	All	6244	0	5931	235	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 20.

All (235) 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)
2:C:132:ALA:HB1	2:D:128:ARG:HD3	1.36	1.07
2:C:103:ASP:HB3	2:D:103:ASP:HB3	1.36	1.06
1:B:177:THR:H	1:B:180:MET:HE3	1.22	1.01
1:B:113:ASN:HA	4:B:324:HOH:O	1.64	0.97
2:C:63:THR:HG23	4:C:171:HOH:O	1.67	0.95
2:C:61:ASN:C	2:C:61:ASN:HD22	1.77	0.92
2:C:51:ASN:HD21	2:C:85:MET:H	1.17	0.90
1:A:240:GLU:HG3	2:D:67:TRP:HH2	1.38	0.88
1:A:47:ASN:OD1	1:A:111:THR:HB	1.76	0.86
1:B:151:ILE:HB	4:B:259:HOH:O	1.74	0.85
1:B:144:ARG:HD2	1:B:152:GLN:CD	2.03	0.83
1:B:109:GLU:HG2	4:B:290:HOH:O	1.79	0.83
1:A:232:VAL:HG23	4:A:254:HOH:O	1.80	0.81
1:A:229:THR:HA	4:A:305:HOH:O	1.81	0.81
1:A:138:VAL:HG12	1:A:198(A):LEU:HD13	1.63	0.81
1:B:73:ASN:O	1:B:76:GLU:HG2	1.80	0.80
1:B:184(A):TYR:HB3	1:B:186[A]:GLU:OE1	1.83	0.79
1:A:184(A):TYR:HE2	1:A:188(A):LYS:HD2	1.48	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:184(A):TYR:CE2	1:A:188(A):LYS:HD2	2.19	0.76
1:B:33:LEU:HD23	1:B:67:VAL:HG22	1.67	0.76
1:A:240:GLU:HG3	2:D:67:TRP:CH2	2.22	0.75
1:B:149:ASP:CG	1:B:150:LYS:H	1.93	0.75
2:C:122:ASN:HD22	2:C:122:ASN:H	1.32	0.74
1:B:43:SER:OG	1:B:198:PRO:HB3	1.87	0.74
1:B:240:GLU:HG3	2:C:67:TRP:HH2	1.53	0.73
2:C:61:ASN:C	2:C:61:ASN:ND2	2.45	0.73
1:B:177:THR:N	1:B:180:MET:HE3	2.01	0.73
1:A:198(A):LEU:O	1:A:210:VAL:HG22	1.90	0.72
1:A:24:ARG:HG2	4:A:268:HOH:O	1.87	0.71
1:A:185:ARG:HH12	1:A:222:ARG:NH2	1.89	0.70
1:B:177:THR:H	1:B:180:MET:CE	2.02	0.70
1:A:63:LYS:HE2	1:A:63:LYS:HA	1.75	0.69
1:A:43:SER:OG	1:A:198:PRO:HB3	1.93	0.69
1:B:33:LEU:CD2	1:B:67:VAL:HG22	2.23	0.67
2:D:88:PRO:O	2:D:89:ASP:HB3	1.94	0.66
1:B:63:LYS:HE3	1:B:63:LYS:O	1.95	0.66
2:C:45:THR:C	2:C:46:LEU:HD12	2.22	0.65
1:B:37:SER:HA	1:B:37(A):PRO:C	2.22	0.65
1:A:23[B]:VAL:HG22	1:A:26:GLU:OE1	1.97	0.64
1:A:197:GLY:HA3	4:A:294:HOH:O	1.97	0.63
2:D:48:VAL:O	2:D:92:LYS:HA	1.99	0.63
2:C:23:GLN:HG2	2:C:125:VAL:HB	1.81	0.63
2:D:25:ILE:HB	2:D:115:ILE:HB	1.80	0.63
1:B:202(B):ASN:O	1:B:202(C):GLU:HB2	1.99	0.62
2:C:130:TRP:CZ2	2:D:132:ALA:HB2	2.34	0.62
1:A:179:LYS:NZ	2:D:63:THR:HG21	2.15	0.62
1:A:129:ASP:O	1:A:132:VAL:HG12	2.00	0.61
1:A:96:MET:HE2	1:A:98:GLU:OE1	2.00	0.61
1:A:144:ARG:HG2	1:A:144:ARG:HH11	1.66	0.61
2:C:127:TYR:CZ	2:D:137:ASP:HB2	2.36	0.60
1:B:35:THR:CG2	1:B:39:LEU:HD22	2.31	0.60
2:C:51:ASN:ND2	2:C:85:MET:H	1.92	0.60
1:B:93:GLN:HE22	2:C:70:ASP:H	1.48	0.60
1:A:73:ASN:HD22	1:A:154:THR:HA	1.67	0.60
1:B:143:TYR:HB3	1:B:145:LYS:O	2.02	0.59
2:C:122:ASN:H	2:C:122:ASN:ND2	2.00	0.59
1:B:202(A):HIS:O	1:B:202(B):ASN:HB2	2.03	0.59
2:C:51:ASN:HD21	2:C:85:MET:HB2	1.67	0.59
1:A:91:HIS:HE1	2:D:67:TRP:O	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:54:ARG:O	2:C:80:PRO:HA	2.04	0.57
1:B:149:ASP:CG	1:B:150:LYS:N	2.63	0.57
2:D:6:PRO:HG2	2:D:9:LYS:HD2	1.84	0.57
1:A:84:GLY:HA3	1:A:109:GLU:OE1	2.04	0.56
2:C:130:TRP:CE2	2:D:132:ALA:HB2	2.41	0.56
2:C:51:ASN:HD21	2:C:85:MET:N	1.95	0.56
2:C:51:ASN:ND2	2:C:85:MET:HB2	2.20	0.56
1:A:33:LEU:HD21	1:A:40:CYS:CB	2.35	0.56
1:A:222:ARG:HG2	1:A:223:GLU:HG3	1.87	0.56
1:B:95:LYS:HD2	2:C:108:ARG:HD2	1.88	0.56
1:A:73:ASN:HA	1:A:153:ASN:O	2.05	0.56
1:B:188(A):LYS:O	1:B:189:ASP:HB2	2.06	0.55
1:B:189:ASP:CG	1:B:190:ALA:H	2.14	0.55
2:D:135:LYS:HE3	2:D:137:ASP:OD1	2.07	0.55
2:D:5:GLN:N	2:D:6:PRO:HD2	2.20	0.55
1:B:232:VAL:HB	4:B:306:HOH:O	2.07	0.55
1:A:91:HIS:CD2	1:A:93:GLN:HB2	2.42	0.55
1:A:224:ARG:NE	4:A:291:HOH:O	2.37	0.55
1:A:91:HIS:CD2	1:A:93:GLN:H	2.25	0.55
1:B:35:THR:O	1:B:37(C):GLN:HA	2.07	0.54
1:A:93:GLN:HE22	2:D:70:ASP:CG	2.15	0.54
1:B:175:LYS:O	1:B:180:MET:HE1	2.07	0.54
1:B:114:TYR:HA	1:B:118:GLN:O	2.08	0.54
2:C:132:ALA:HB2	2:D:130:TRP:CE2	2.42	0.53
1:A:22:SER:HA	4:A:270:HOH:O	2.07	0.53
1:A:28:PRO:HB2	1:A:119:ARG:H	1.73	0.53
1:A:141:TRP:O	1:A:151:ILE:HG13	2.08	0.53
1:A:145:LYS:O	1:A:148:ASP:HB2	2.08	0.53
1:B:16:ILE:HB	1:B:156:GLN:HB3	1.91	0.53
1:B:16:ILE:HD12	1:B:194:ASP:CG	2.33	0.53
2:C:97:VAL:HG23	2:C:97:VAL:O	2.08	0.53
1:B:171:ARG:NH1	1:B:223:GLU:OE2	2.40	0.53
2:C:45:THR:O	2:C:46:LEU:HD12	2.08	0.53
1:A:138:VAL:HG11	1:A:160:ILE:HD12	1.90	0.53
1:A:129:ASP:HB3	1:A:132:VAL:CG1	2.39	0.52
1:A:148:ASP:CG	1:A:149:LYS:H	2.17	0.52
3:A:1:NAG:O7	3:A:1:NAG:H3	2.09	0.52
2:D:15:GLN:HE21	2:D:15:GLN:HA	1.75	0.52
1:B:195:SER:HA	1:B:213:THR:HG22	1.92	0.51
2:C:44:GLN:O	2:C:97:VAL:HG22	2.10	0.51
1:A:33:LEU:HD21	1:A:40:CYS:HB2	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:98:GLU:O	1:A:180:MET:HE1	2.10	0.51
1:B:161:PRO:HD3	1:B:184(A):TYR:CZ	2.45	0.51
1:A:33:LEU:CD1	1:A:39:LEU:HD23	2.41	0.51
1:A:136:CYS:HA	1:A:198(B):SER:O	2.10	0.51
1:A:129:ASP:HB3	1:A:132:VAL:HG11	1.92	0.51
1:B:35:THR:HG21	1:B:39:LEU:HD22	1.92	0.51
1:B:65:LEU:C	1:B:66:ARG:HG2	2.35	0.51
2:C:122:ASN:HD22	2:C:122:ASN:N	1.96	0.51
2:D:77:VAL:HG21	2:D:120:PRO:HG2	1.92	0.51
1:A:186:GLU:CD	1:A:186:GLU:H	2.15	0.51
1:B:37:SER:HA	1:B:37(A):PRO:O	2.11	0.51
1:B:189:ASP:CG	1:B:190:ALA:N	2.69	0.51
1:A:91:HIS:HD2	1:A:93:GLN:H	1.57	0.50
1:B:186[B]:GLU:HG3	4:B:282:HOH:O	2.11	0.50
1:A:171:ARG:HG3	1:A:223:GLU:O	2.11	0.50
2:D:72:TYR:CD1	2:D:72:TYR:N	2.80	0.50
1:A:23[A]:VAL:HG13	1:A:24:ARG:N	2.25	0.50
1:A:73:ASN:HD21	1:A:154:THR:HG23	1.77	0.50
1:B:115:THR:O	1:B:119:ARG:HG3	2.12	0.50
1:A:125:PRO:CB	1:A:232:VAL:HG22	2.42	0.50
1:A:179:LYS:HZ2	2:D:63:THR:HG21	1.75	0.50
1:A:86:GLN:HB3	1:A:107:LYS:O	2.12	0.50
1:B:126:SER:C	1:B:128:GLY:H	2.19	0.50
1:B:222:ARG:HH11	1:B:222:ARG:HG2	1.77	0.50
2:C:25:ILE:HB	2:C:115:ILE:HB	1.93	0.50
1:A:173:ARG:O	2:C:58:LYS:HE3	2.12	0.50
1:B:126:SER:C	1:B:128:GLY:N	2.70	0.50
1:A:16:ILE:HB	1:A:156:GLN:HB3	1.94	0.49
1:B:38:HIS:HB3	1:B:151:ILE:HD12	1.93	0.49
1:A:73:ASN:ND2	1:A:154:THR:HA	2.25	0.49
1:B:140:GLY:HA3	1:B:194:ASP:OD1	2.12	0.49
2:C:48:VAL:HB	2:C:53:HIS:CD2	2.48	0.49
1:A:201:CYS:SG	1:A:210:VAL:HG21	2.52	0.49
2:D:5:GLN:H	2:D:6:PRO:HD2	1.77	0.49
1:A:138:VAL:HG12	1:A:198(A):LEU:CD1	2.40	0.49
1:B:240:GLU:HG3	2:C:67:TRP:CH2	2.42	0.49
1:B:195:SER:C	1:B:197:GLY:H	2.21	0.49
1:A:216:GLY:HA3	2:C:84:ARG:NH2	2.28	0.48
1:B:62:PRO:HB3	1:B:88:ILE:HG13	1.95	0.48
1:A:66:ARG:HG2	1:A:82:PHE:CE1	2.49	0.48
1:B:211:GLY:HA2	1:B:229:THR:O	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:68:GLY:C	4:C:171:HOH:O	2.56	0.48
1:B:146:LEU:HD12	1:B:148:ARG:H	1.78	0.48
1:B:146:LEU:O	1:B:148:ARG:HB2	2.12	0.48
1:B:93:GLN:HE22	2:C:70:ASP:N	2.12	0.47
1:A:44:ILE:O	1:A:121:PRO:HA	2.13	0.47
1:A:185:ARG:HG3	1:A:185:ARG:HH11	1.80	0.47
1:B:43:SER:HG	1:B:198:PRO:HB3	1.79	0.47
2:C:31:GLU:HG3	2:C:32:ASP:N	2.29	0.47
1:A:130:ARG:NH1	2:D:65:GLU:HG3	2.30	0.47
1:B:59(C):VAL:HG13	1:B:64:ILE:HD11	1.96	0.47
1:B:113:ASN:OD1	4:B:325:HOH:O	2.20	0.47
2:C:54:ARG:HG3	2:C:98:THR:O	2.15	0.47
2:D:27:LEU:HD11	2:D:38:VAL:HG21	1.96	0.46
1:A:161:PRO:HD3	1:A:184(A):TYR:CZ	2.50	0.46
1:A:59:PHE:CZ	1:A:106:LEU:HD21	2.50	0.46
1:A:17:VAL:HG13	1:A:189:ASP:O	2.15	0.46
1:A:17:VAL:HB	1:A:144:ARG:O	2.15	0.46
1:A:223:GLU:O	1:A:224:ARG:HG3	2.15	0.46
1:B:29:TRP:CG	1:B:122:ILE:HB	2.50	0.46
1:B:186[B]:GLU:OE1	1:B:222:ARG:NH1	2.46	0.46
1:A:29:TRP:CG	1:A:122:ILE:HB	2.51	0.46
1:A:135:ASP:O	1:A:201:CYS:HA	2.16	0.46
2:D:19:GLY:HA2	2:D:121:ASP:OD2	2.16	0.45
1:A:138:VAL:CG1	1:A:160:ILE:HD12	2.46	0.45
1:A:114:TYR:CZ	1:A:121:PRO:HG3	2.52	0.45
1:B:68:TYR:CE2	1:B:82:PHE:HB3	2.52	0.45
2:C:39:GLU:HB2	2:C:106:MET:SD	2.57	0.45
1:B:83:PHE:CZ	1:B:112:VAL:HG22	2.52	0.45
2:D:16:ALA:HA	4:D:174:HOH:O	2.16	0.45
1:B:171:ARG:HG2	1:B:223:GLU:HB3	1.99	0.45
2:C:26:GLN:HE21	2:C:26:GLN:HB3	1.58	0.45
2:C:48:VAL:HB	2:C:53:HIS:NE2	2.31	0.45
2:C:75:ASP:O	2:C:76:LYS:HB2	2.17	0.44
2:D:7:LEU:HG	2:D:13:TYR:CD2	2.52	0.44
2:D:49:ASP:OD1	2:D:53:HIS:HE1	2.00	0.44
1:B:223:GLU:O	1:B:224:ARG:NH1	2.38	0.44
2:C:132:ALA:HB2	2:D:130:TRP:CZ2	2.52	0.44
2:D:5:GLN:HB2	2:D:6:PRO:CD	2.48	0.44
1:B:135:ASP:O	1:B:201:CYS:HA	2.18	0.44
1:A:33:LEU:HG	1:A:33:LEU:O	2.18	0.44
1:B:194:ASP:O	1:B:195:SER:C	2.61	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:59(A):TYR:CD2	1:A:59(A):TYR:C	2.96	0.43
1:A:144:ARG:HG2	1:A:144:ARG:NH1	2.33	0.43
2:D:53:HIS:CD2	2:D:82:SER:HB3	2.53	0.43
2:D:4:VAL:HG13	2:D:5:GLN:N	2.33	0.43
1:A:202(A):HIS:O	1:A:202(B):ASN:HB3	2.19	0.43
2:D:11:ALA:HB1	2:D:12:PRO:HD2	2.00	0.43
2:D:45:THR:HG21	2:D:94:LYS:HE2	2.00	0.43
1:B:126:SER:O	1:B:128:GLY:N	2.52	0.43
1:A:184(A):TYR:O	1:A:185:ARG:C	2.60	0.43
1:A:24:ARG:HH22	1:A:116:ASP:HB3	1.84	0.43
1:A:28:PRO:HB2	1:A:119:ARG:N	2.34	0.43
1:B:86:GLN:HE21	1:B:86:GLN:HB2	1.63	0.43
2:D:5:GLN:O	2:D:6:PRO:C	2.62	0.43
1:A:127:LYS:HB3	1:A:127:LYS:HE2	1.83	0.42
1:B:212:ILE:HB	1:B:229:THR:HB	2.01	0.42
1:A:125:PRO:HB2	1:A:232:VAL:HG22	2.02	0.42
2:D:79:SER:HA	2:D:80:PRO:HD3	1.93	0.42
1:A:173:ARG:CG	1:A:173:ARG:HH11	2.32	0.42
1:A:188(A):LYS:O	1:A:189:ASP:HB2	2.18	0.42
1:A:20:THR:CG2	1:A:21:ALA:N	2.79	0.42
1:A:80:ASP:O	1:A:80:ASP:CG	2.63	0.42
1:B:66:ARG:HD2	1:B:82:PHE:CE1	2.55	0.42
1:A:53:LEU:HD11	1:A:103:ILE:HD11	2.02	0.42
1:B:32:THR:HG23	1:B:68:TYR:HB2	2.01	0.42
2:D:33:GLU:OE1	4:D:155:HOH:O	2.22	0.42
1:A:171:ARG:HG2	1:A:225:PRO:HD3	2.02	0.42
1:B:202(B):ASN:O	1:B:202(C):GLU:CB	2.68	0.42
1:A:52:ILE:HB	1:A:106:LEU:HB2	2.01	0.42
2:D:21:LYS:HE2	2:D:121:ASP:O	2.19	0.42
1:A:37:SER:O	1:A:37:SER:OG	2.32	0.41
2:D:4:VAL:HG13	2:D:5:GLN:HG2	2.02	0.41
1:A:37:SER:HA	1:A:37(A):PRO:HA	1.72	0.41
2:D:8:GLU:HA	2:D:11:ALA:O	2.20	0.41
1:A:217:GLU:HG2	4:A:308:HOH:O	2.20	0.41
1:B:96:MET:O	1:B:97:ALA:C	2.63	0.41
2:C:5:GLN:HA	2:C:6:PRO:HD2	1.97	0.41
1:A:35:THR:O	1:A:37(C):GLN:HA	2.20	0.41
2:C:41:LEU:HB2	2:C:126:LYS:HB2	2.01	0.41
2:C:133:GLU:HG3	2:D:129:VAL:O	2.21	0.41
2:D:15:GLN:HA	2:D:15:GLN:NE2	2.35	0.41
1:B:114:TYR:N	4:B:324:HOH:O	2.44	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:146:LEU:O	1:B:148:ARG:CB	2.68	0.41
1:A:54:THR:OG1	1:A:55:ALA:N	2.54	0.41
2:C:27:LEU:HD11	2:C:38:VAL:HG21	2.01	0.41
1:B:32:THR:HB	1:B:141:TRP:CZ3	2.56	0.41
2:C:138:ASN:OD1	2:D:126:LYS:HG2	2.21	0.41
1:A:35:THR:O	1:A:35:THR:HG23	2.21	0.41
1:A:82:PHE:CD1	1:A:82:PHE:C	2.98	0.41
1:A:144:ARG:C	1:A:145:LYS:HG3	2.46	0.40
1:B:113:ASN:CA	4:B:324:HOH:O	2.44	0.40
2:D:95:LYS:HD2	2:D:95:LYS:HA	1.77	0.40
1:A:32:THR:O	1:A:32:THR:HG23	2.20	0.40
1:A:126:SER:C	1:A:128:GLY:N	2.77	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	236/238 (99%)	213 (90%)	22 (9%)	1 (0%)	30	34
1	B	231/238 (97%)	207 (90%)	22 (10%)	2 (1%)	14	14
2	C	138/142 (97%)	132 (96%)	6 (4%)	0	100	100
2	D	137/142 (96%)	127 (93%)	7 (5%)	3 (2%)	5	3
All	All	742/760 (98%)	679 (92%)	57 (8%)	6 (1%)	16	16

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	37(A)	PRO
1	B	222	ARG
2	D	6	PRO

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Mol	Chain	Res	Type
2	D	89	ASP
2	D	92	LYS
1	B	37(A)	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	204/204 (100%)	186 (91%)	18 (9%)	9	10
1	B	201/204 (98%)	182 (90%)	19 (10%)	8	8
2	C	124/125 (99%)	116 (94%)	8 (6%)	15	18
2	D	123/125 (98%)	118 (96%)	5 (4%)	27	37
All	All	652/658 (99%)	602 (92%)	50 (8%)	13	13

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

Mol	Chain	Res	Type
1	A	17	VAL
1	A	20	THR
1	A	23[A]	VAL
1	A	23[B]	VAL
1	A	24	ARG
1	A	33	LEU
1	A	37	SER
1	A	37(A)	PRO
1	A	37(B)	THR
1	A	37(C)	GLN
1	A	37(D)	ARG
1	A	93	GLN
1	A	96	MET
1	A	116	ASP
1	A	156	GLN
1	A	162	LEU
1	A	171	ARG
1	A	186	GLU

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Mol	Chain	Res	Type
1	B	17	VAL
1	B	23	VAL
1	B	32	THR
1	B	37	SER
1	B	37(A)	PRO
1	B	37(B)	THR
1	B	37(C)	GLN
1	B	37(D)	ARG
1	B	63	LYS
1	B	66	ARG
1	B	93	GLN
1	B	96	MET
1	B	148	ARG
1	B	151	ILE
1	B	156	GLN
1	B	166	GLU
1	B	173	ARG
1	B	186[A]	GLU
1	B	186[B]	GLU
2	C	23	GLN
2	C	28	THR
2	C	31	GLU
2	C	61	ASN
2	C	95	LYS
2	C	107	LEU
2	C	121	ASP
2	C	122	ASN
2	D	47	GLU
2	D	78	SER
2	D	89	ASP
2	D	91	LYS
2	D	107	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	GLN
1	A	73	ASN
1	A	91	HIS
1	A	118	GLN
1	A	156	GLN
1	A	243	GLN

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Mol	Chain	Res	Type
1	B	37(C)	GLN
1	B	48	GLN
1	B	86	GLN
1	B	93	GLN
1	B	113	ASN
1	B	156	GLN
1	B	169	GLN
1	B	178	HIS
1	B	202(B)	ASN
2	C	5	GLN
2	C	51	ASN
2	C	61	ASN
2	C	122	ASN
2	D	15	GLN
2	D	44	GLN
2	D	53	HIS
2	D	61	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](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	A	1	1	14,14,15	0.85	0	17,19,21	0.81	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	A	1	1	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1	NAG	C3-C2-N2-C7

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1	NAG	1	0

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	237/238 (99%)	0.51	12 (5%) 33 30	23, 35, 55, 67	1 (0%)
1	B	233/238 (97%)	0.52	12 (5%) 33 29	24, 36, 60, 76	2 (0%)
2	C	140/142 (98%)	0.51	5 (3%) 46 43	26, 38, 61, 71	0
2	D	139/142 (97%)	0.53	4 (2%) 53 50	24, 39, 59, 73	0
All	All	749/760 (98%)	0.52	33 (4%) 39 36	23, 37, 59, 76	3 (0%)

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	4	VAL	5.6
1	A	37(A)	PRO	4.2
1	B	178	HIS	3.9
2	C	3	SER	3.8
1	B	146	LEU	3.7
1	B	151	ILE	3.5
2	D	90	GLY	3.3
1	B	148	ARG	3.3
2	D	91	LYS	3.2
1	A	33	LEU	3.2
1	B	37(A)	PRO	3.1
1	A	59(A)	TYR	3.0
1	B	128	GLY	3.0
2	D	89	ASP	2.6
1	A	244	ALA	2.5
1	A	80	ASP	2.5
1	A	59(B)	GLY	2.4
1	A	146	LEU	2.4
1	B	35	THR	2.4
1	B	143	TYR	2.3
1	B	145	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
2	C	89	ASP	2.2
2	C	121	ASP	2.2
1	B	59(A)	TYR	2.2
1	B	222	ARG	2.2
2	C	95	LYS	2.2
1	A	24	ARG	2.1
2	C	9	LYS	2.1
1	A	37(B)	THR	2.1
1	A	47	ASN	2.1
1	B	133	TYR	2.1
1	A	37	SER	2.0
1	A	151	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](#)

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($\text{\AA}^2$)	Q<0.9
3	NAG	A	1	14/15	0.58	0.16	49,52,54,54	0

6.5 Other polymers [i](#)

There are no such residues in this entry.