Full wwPDB X-ray Structure Validation Report ⓘ

Oct 11, 2023 – 02:43 AM EDT

PDB ID	:	7K26
Title	:	Crystal structure of Human H-chain Ferritin variant infused with Sodium Acrylate
Authors	:	Bailey, J.B.; Zhang, L.
Deposited on	:	2020-09-08
Resolution	:	2.70 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:

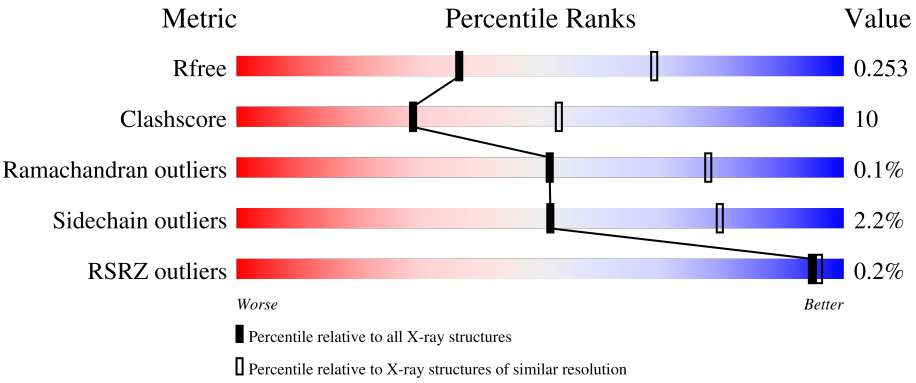
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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 <=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	182	70% 25%
1	B	182	79% 15% 5%
1	C	182	75% 19%
1	D	182	% 85% 10%
1	E	182	73% 22% 5%

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Mol	Chain	Length	Quality of chain
1	F	182	73% 20% • 5%
1	G	182	77% 19% •
1	H	182	% 74% 20% • 5%
1	I	182	77% 17% • 5%
1	J	182	% 75% 20% • 5%
1	K	182	76% 20% •
1	L	182	78% 16% • 5%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 17398 atoms, of which 98 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 Ferritin heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	174	Total	C	N	O	S	0	0	0
			1406	880	244	278	4			
1	B	173	Total	C	N	O	S	0	2	0
			1425	894	250	277	4			
1	C	174	Total	C	N	O	S	0	1	0
			1415	886	247	278	4			
1	D	176	Total	C	N	O	S	0	0	0
			1420	889	245	282	4			
1	E	173	Total	C	N	O	S	0	0	0
			1396	877	241	274	4			
1	F	173	Total	C	N	O	S	0	0	0
			1403	881	245	273	4			
1	G	174	Total	C	N	O	S	0	0	0
			1406	881	243	278	4			
1	H	173	Total	C	N	O	S	0	0	0
			1394	875	241	274	4			
1	I	173	Total	C	N	O	S	0	0	0
			1389	872	242	271	4			
1	J	173	Total	C	N	O	S	0	1	0
			1392	873	239	276	4			
1	K	175	Total	C	N	O	S	0	1	0
			1412	884	243	281	4			
1	L	173	Total	C	N	O	S	0	0	0
			1401	879	243	275	4			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	86	GLN	LYS	conflict	UNP P02794
A	90	GLU	CYS	conflict	UNP P02794
A	102	ALA	CYS	conflict	UNP P02794
A	130	ALA	CYS	conflict	UNP P02794
B	86	GLN	LYS	conflict	UNP P02794

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Chain	Residue	Modelled	Actual	Comment	Reference
B	90	GLU	CYS	conflict	UNP P02794
B	102	ALA	CYS	conflict	UNP P02794
B	130	ALA	CYS	conflict	UNP P02794
C	86	GLN	LYS	conflict	UNP P02794
C	90	GLU	CYS	conflict	UNP P02794
C	102	ALA	CYS	conflict	UNP P02794
C	130	ALA	CYS	conflict	UNP P02794
D	86	GLN	LYS	conflict	UNP P02794
D	90	GLU	CYS	conflict	UNP P02794
D	102	ALA	CYS	conflict	UNP P02794
D	130	ALA	CYS	conflict	UNP P02794
E	86	GLN	LYS	conflict	UNP P02794
E	90	GLU	CYS	conflict	UNP P02794
E	102	ALA	CYS	conflict	UNP P02794
E	130	ALA	CYS	conflict	UNP P02794
F	86	GLN	LYS	conflict	UNP P02794
F	90	GLU	CYS	conflict	UNP P02794
F	102	ALA	CYS	conflict	UNP P02794
F	130	ALA	CYS	conflict	UNP P02794
G	86	GLN	LYS	conflict	UNP P02794
G	90	GLU	CYS	conflict	UNP P02794
G	102	ALA	CYS	conflict	UNP P02794
G	130	ALA	CYS	conflict	UNP P02794
H	86	GLN	LYS	conflict	UNP P02794
H	90	GLU	CYS	conflict	UNP P02794
H	102	ALA	CYS	conflict	UNP P02794
H	130	ALA	CYS	conflict	UNP P02794
I	86	GLN	LYS	conflict	UNP P02794
I	90	GLU	CYS	conflict	UNP P02794
I	102	ALA	CYS	conflict	UNP P02794
I	130	ALA	CYS	conflict	UNP P02794
J	86	GLN	LYS	conflict	UNP P02794
J	90	GLU	CYS	conflict	UNP P02794
J	102	ALA	CYS	conflict	UNP P02794
J	130	ALA	CYS	conflict	UNP P02794
K	86	GLN	LYS	conflict	UNP P02794
K	90	GLU	CYS	conflict	UNP P02794
K	102	ALA	CYS	conflict	UNP P02794
K	130	ALA	CYS	conflict	UNP P02794
L	86	GLN	LYS	conflict	UNP P02794
L	90	GLU	CYS	conflict	UNP P02794
L	102	ALA	CYS	conflict	UNP P02794

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Chain	Residue	Modelled	Actual	Comment	Reference
L	130	ALA	CYS	conflict	UNP P02794

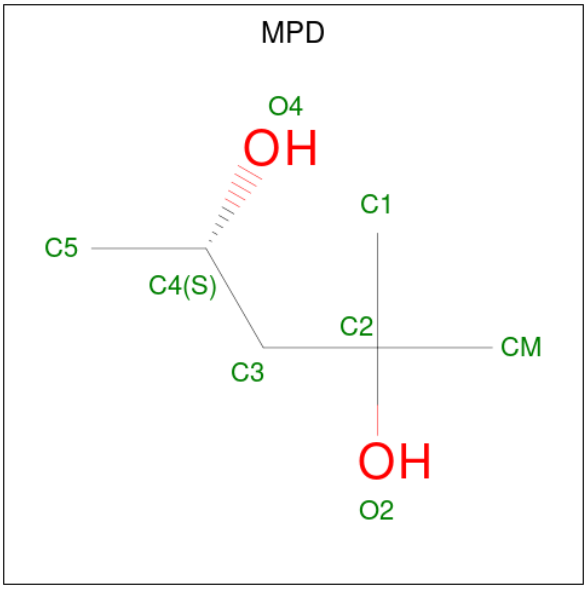
- Molecule 2 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Fe 1 1	0	0
2	B	1	Total Fe 1 1	0	0
2	C	1	Total Fe 1 1	0	0
2	D	1	Total Fe 1 1	0	0
2	E	1	Total Fe 1 1	0	0
2	F	1	Total Fe 1 1	0	0
2	G	1	Total Fe 1 1	0	0
2	H	1	Total Fe 1 1	0	0
2	I	1	Total Fe 1 1	0	0
2	J	1	Total Fe 1 1	0	0
2	K	1	Total Fe 1 1	0	0
2	L	1	Total Fe 1 1	0	0

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total Na 2 2	0	0
3	B	2	Total Na 2 2	0	0
3	C	1	Total Na 1 1	0	0
3	H	1	Total Na 1 1	0	0
3	L	1	Total Na 1 1	0	0

- Molecule 4 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	H	O	0	0
			22	6	14	2		
4	D	1	Total	C	H	O	0	0
			22	6	14	2		
4	E	1	Total	C	H	O	0	0
			22	6	14	2		
4	F	1	Total	C	H	O	0	0
			22	6	14	2		
4	G	1	Total	C	H	O	0	0
			22	6	14	2		
4	K	1	Total	C	H	O	0	0
			22	6	14	2		
4	L	1	Total	C	H	O	0	0
			22	6	14	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	36	Total	O	0	0
			36	36		
5	B	43	Total	O	0	0
			43	43		
5	C	31	Total	O	0	0
			31	31		

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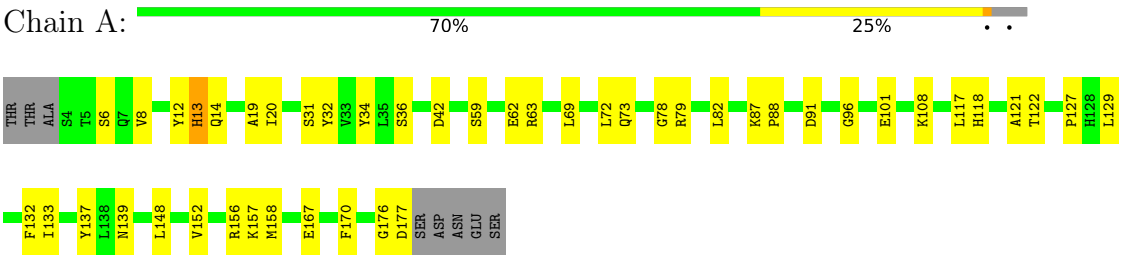
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	D	55	Total O 55 55	0	0
5	E	27	Total O 27 27	0	0
5	F	29	Total O 29 29	0	0
5	G	24	Total O 24 24	0	0
5	H	24	Total O 24 24	0	0
5	I	22	Total O 22 22	0	0
5	J	21	Total O 21 21	0	0
5	K	31	Total O 31 31	0	0
5	L	23	Total O 23 23	0	0

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.

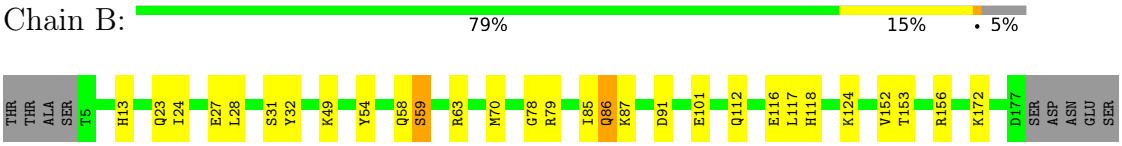
• Molecule 1: Ferritin heavy chain

Chain A:



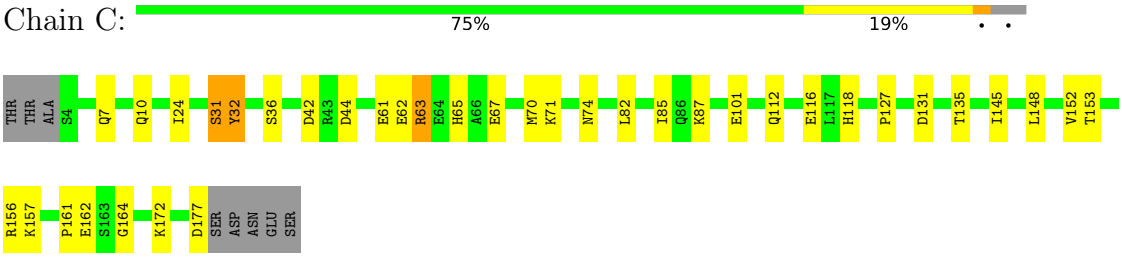
• Molecule 1: Ferritin heavy chain

Chain B:



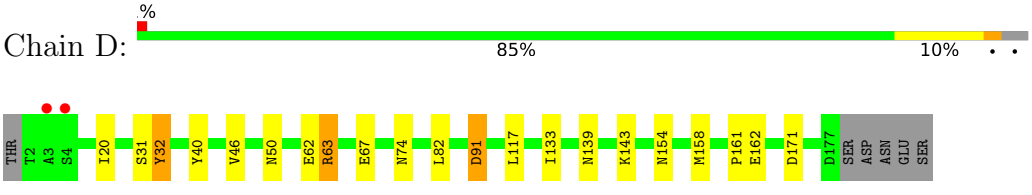
• Molecule 1: Ferritin heavy chain

Chain C:



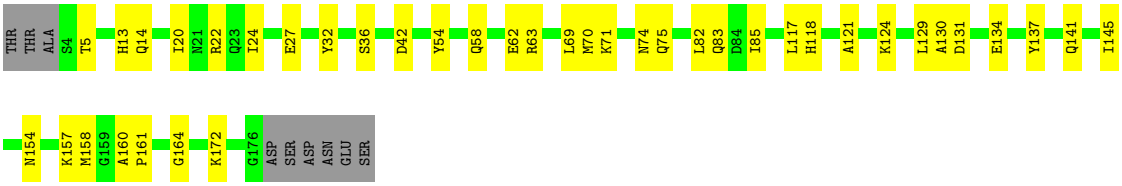
• Molecule 1: Ferritin heavy chain

Chain D:



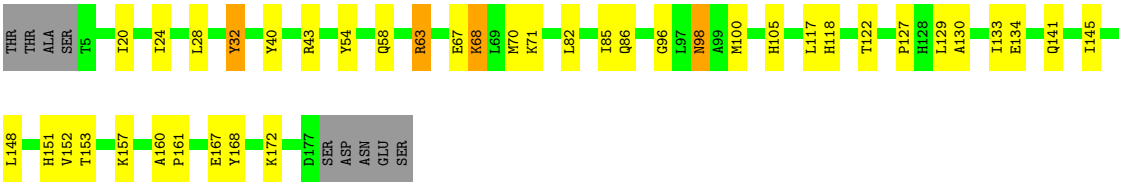
• Molecule 1: Ferritin heavy chain

Chain E:



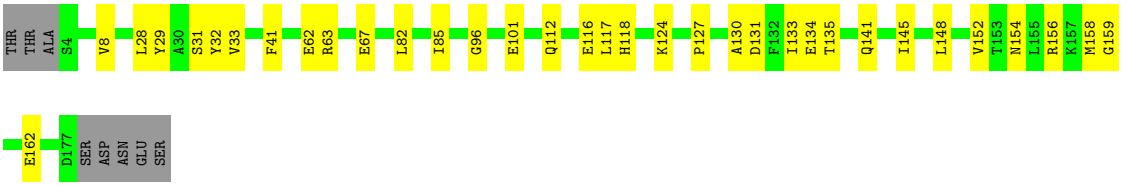
• Molecule 1: Ferritin heavy chain

Chain F:



• Molecule 1: Ferritin heavy chain

Chain G:

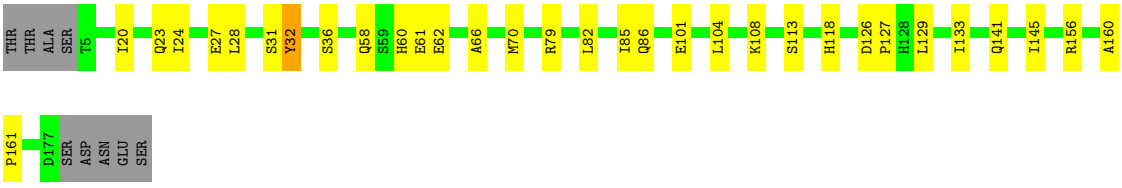


• Molecule 1: Ferritin heavy chain

Chain H:

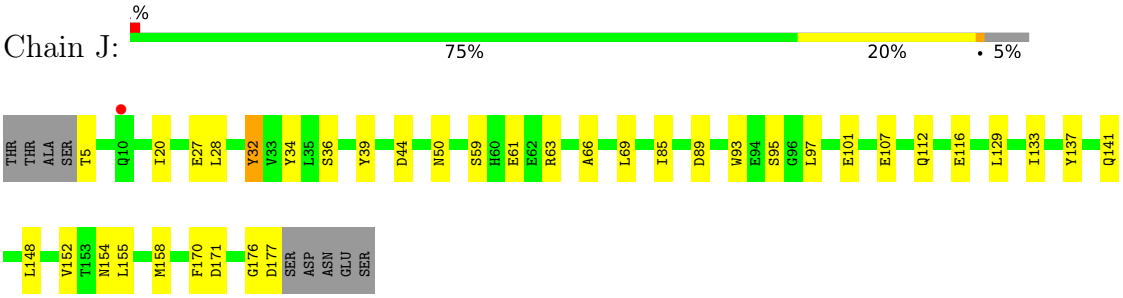
• Molecule 1: Ferritin heavy chain

Chain I:



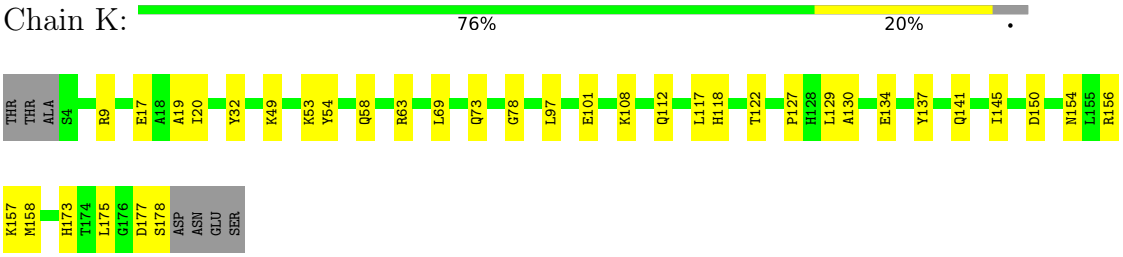
• Molecule 1: Ferritin heavy chain

Chain J:



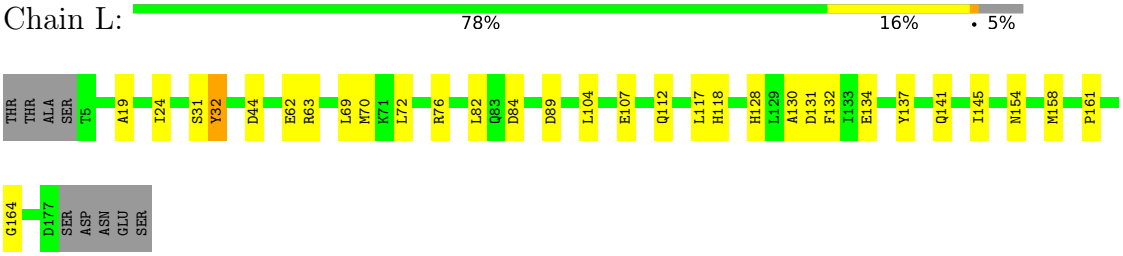
• Molecule 1: Ferritin heavy chain

Chain K:



• Molecule 1: Ferritin heavy chain

Chain L:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	131.57Å 131.57Å 306.56Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	53.99 – 2.70 91.44 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.1 (53.99-2.70) 99.2 (91.44-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.96 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.198 , 0.254 0.198 , 0.253	Depositor DCC
R_{free} test set	4417 reflections (5.19%)	wwPDB-VP
Wilson B-factor (Å^2)	49.1	Xtriage
Anisotropy	0.478	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 39.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	0.008 for -h,-k,l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	17398	wwPDB-VP
Average B, all atoms (Å^2)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.96% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: FE, MPD, NA

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	0.35	0/1435	0.52	0/1941
1	B	0.30	0/1460	0.47	0/1969
1	C	0.33	0/1447	0.49	0/1955
1	D	0.32	0/1449	0.46	0/1960
1	E	0.30	0/1425	0.46	0/1926
1	F	0.32	0/1432	0.46	0/1934
1	G	0.30	0/1435	0.45	0/1940
1	H	0.34	0/1423	0.54	2/1925 (0.1%)
1	I	0.30	0/1418	0.48	0/1918
1	J	0.35	0/1424	0.48	0/1929
1	K	0.32	0/1444	0.45	0/1954
1	L	0.29	0/1430	0.44	0/1933
All	All	0.32	0/17222	0.47	2/23284 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	H	108	LYS	CD-CE-NZ	-7.67	94.07	111.70
1	H	86	GLN	CA-CB-CG	-5.45	101.42	113.40

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	1406	0	1304	40	0
1	B	1425	0	1357	18	0
1	C	1415	0	1325	37	0
1	D	1420	0	1318	19	0
1	E	1396	0	1305	36	0
1	F	1403	0	1320	37	0
1	G	1406	0	1307	27	0
1	H	1394	0	1291	43	0
1	I	1389	0	1297	32	0
1	J	1392	0	1275	28	0
1	K	1412	0	1302	28	0
1	L	1401	0	1306	27	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
2	K	1	0	0	0	0
2	L	1	0	0	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	1	0	0	0	0
3	H	1	0	0	0	0
3	L	1	0	0	0	0
4	B	8	14	14	0	0
4	D	8	14	14	1	0
4	E	8	14	14	4	0
4	F	8	14	14	2	0
4	G	8	14	14	0	0
4	K	8	14	14	1	0
4	L	8	14	14	1	0
5	A	36	0	0	1	0
5	B	43	0	0	2	0
5	C	31	0	0	2	0
5	D	55	0	0	0	0
5	E	27	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	F	29	0	0	3	0
5	G	24	0	0	0	0
5	H	24	0	0	1	0
5	I	22	0	0	1	0
5	J	21	0	0	2	0
5	K	31	0	0	2	1
5	L	23	0	0	1	0
All	All	17300	98	15805	315	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 10.

All (315) 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:H:108:LYS:HE2	1:H:145:ILE:HD13	1.37	1.06
1:I:32:TYR:OH	1:L:82:LEU:HB3	1.55	1.03
1:B:91:ASP:OD2	5:B:301:HOH:O	1.80	0.99
1:F:168:TYR:CE1	1:F:172:LYS:HE3	2.00	0.96
1:A:87:LYS:CE	1:D:82:LEU:O	2.15	0.95
1:C:172:LYS:NZ	5:C:301:HOH:O	1.98	0.94
1:F:168:TYR:CZ	1:F:172:LYS:HE3	2.02	0.94
1:H:117:LEU:HD22	1:H:133:ILE:CD1	1.97	0.94
1:H:117:LEU:HD22	1:H:133:ILE:HD11	1.48	0.94
1:H:108:LYS:CE	1:H:145:ILE:HD13	1.97	0.94
1:A:101:GLU:OE2	1:A:156:ARG:NH2	2.03	0.92
1:I:104:LEU:HD11	1:I:145:ILE:HG23	1.50	0.90
1:A:87:LYS:HE2	1:D:82:LEU:O	1.69	0.90
1:H:12:TYR:CD1	1:H:76:ARG:HD3	2.13	0.83
1:G:28:LEU:HB3	1:G:85:ILE:HD13	1.60	0.83
1:I:104:LEU:CD1	1:I:145:ILE:HG23	2.08	0.82
1:I:108:LYS:HG2	1:I:145:ILE:HG12	1.63	0.81
1:L:31:SER:HB2	1:L:62:GLU:HB2	1.61	0.81
1:C:63[B]:ARG:HH21	1:E:63:ARG:HG2	1.47	0.79
1:H:28:LEU:HB3	1:H:85:ILE:CD1	2.13	0.78
1:G:63:ARG:HD3	1:J:63:ARG:HH11	1.47	0.78
1:H:28:LEU:HB3	1:H:85:ILE:HD13	1.64	0.77
1:F:28:LEU:HB3	1:F:85:ILE:HD13	1.63	0.77
1:E:154:ASN:O	1:E:158:MET:HG3	1.86	0.76
1:K:118:HIS:O	1:K:122:THR:HG23	1.86	0.75
1:B:85:ILE:O	1:B:86:GLN:HB2	1.86	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:117:LEU:HD22	1:G:133:ILE:HD11	1.69	0.74
1:L:141:GLN:O	1:L:145:ILE:HG13	1.87	0.74
1:A:87:LYS:NZ	1:A:91:ASP:OD2	2.17	0.73
1:H:117:LEU:CD2	1:H:133:ILE:HD11	2.17	0.73
1:J:176:GLY:O	1:J:177:ASP:HB2	1.88	0.72
1:J:61[A]:GLU:OE2	5:J:301:HOH:O	2.07	0.72
1:A:78:GLY:O	1:A:79:ARG:NE	2.18	0.71
1:E:172:LYS:HG2	5:E:304:HOH:O	1.91	0.70
1:I:36:SER:HB2	1:L:82:LEU:CD1	2.22	0.70
1:K:130:ALA:O	1:K:134:GLU:HG3	1.92	0.70
1:F:82:LEU:O	1:H:87:LYS:HE3	1.92	0.69
1:H:131:ASP:O	1:H:135:THR:HG23	1.92	0.69
1:A:176:GLY:O	1:A:177:ASP:HB2	1.93	0.69
1:B:172:LYS:HE2	5:B:320:HOH:O	1.93	0.69
1:A:87:LYS:HE3	1:D:82:LEU:O	1.91	0.69
1:C:63[A]:ARG:NH2	4:E:202:MPD:H12	2.09	0.68
1:J:107:GLU:OE2	1:J:141:GLN:NE2	2.26	0.68
1:A:63:ARG:HH21	1:D:63:ARG:HG2	1.57	0.68
1:H:28:LEU:HD13	1:H:85:ILE:HD11	1.75	0.68
1:H:76:ARG:HH12	1:H:128:HIS:HB3	1.59	0.67
1:J:112:GLN:O	1:J:116:GLU:HG3	1.94	0.67
1:H:63:ARG:HD3	1:H:67:GLU:OE2	1.94	0.67
1:L:107:GLU:OE2	1:L:141:GLN:NE2	2.20	0.66
1:L:130:ALA:O	1:L:134:GLU:HG3	1.95	0.66
1:G:31:SER:HB2	1:G:62:GLU:HB2	1.77	0.66
1:L:112:GLN:O	1:L:112:GLN:HG3	1.94	0.66
1:B:28:LEU:HB3	1:B:85:ILE:HD12	1.78	0.65
1:H:12:TYR:HD1	1:H:76:ARG:HD3	1.59	0.65
1:G:63:ARG:HD3	1:J:63:ARG:NH1	2.12	0.65
1:I:126:ASP:OD1	5:I:301:HOH:O	2.14	0.65
1:H:161:PRO:HD2	1:H:162:GLU:OE1	1.98	0.64
1:K:19:ALA:HB1	1:K:117:LEU:HD13	1.80	0.64
1:E:14:GLN:NE2	5:E:303:HOH:O	2.27	0.64
1:I:36:SER:HB2	1:L:82:LEU:HD11	1.80	0.63
1:E:130:ALA:O	1:E:134:GLU:HG2	1.98	0.63
1:I:32:TYR:CE1	1:L:82:LEU:HD13	2.33	0.63
1:K:20:ILE:HD13	1:K:117:LEU:HD21	1.80	0.62
1:E:121:ALA:HB2	1:E:129:LEU:HD23	1.80	0.62
1:H:108:LYS:HE2	1:H:145:ILE:HG21	1.81	0.62
1:K:73:GLN:OE1	1:K:78:GLY:HA3	2.00	0.62
1:H:117:LEU:CD2	1:H:133:ILE:CD1	2.75	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:31:SER:HB2	1:I:62:GLU:HB2	1.81	0.61
1:F:100:MET:HE2	1:F:151:HIS:CB	2.30	0.61
1:K:54:TYR:O	1:K:58:GLN:HG2	2.00	0.61
1:A:20:ILE:HD13	1:A:117:LEU:HD21	1.81	0.61
1:E:141:GLN:O	1:E:145:ILE:HG12	1.99	0.61
1:C:63[B]:ARG:NH2	1:E:63:ARG:HG2	2.16	0.60
1:D:154:ASN:O	1:D:158:MET:HG3	2.01	0.60
4:D:202:MPD:O4	4:D:202:MPD:HM1	2.00	0.60
1:H:76:ARG:NH1	1:H:128:HIS:HB3	2.15	0.60
1:A:13:HIS:ND1	1:A:14:GLN:N	2.49	0.60
1:D:117:LEU:HD22	1:D:133:ILE:HD11	1.82	0.60
1:C:172:LYS:CD	1:K:173:HIS:HB3	2.32	0.60
1:K:154:ASN:O	1:K:158:MET:HG3	2.00	0.60
1:A:78:GLY:C	1:A:79:ARG:HE	2.05	0.60
1:I:82:LEU:HB3	1:L:32:TYR:OH	2.01	0.59
1:I:23:GLN:O	1:I:27:GLU:HG2	2.02	0.59
1:L:19:ALA:HB1	1:L:117:LEU:CD1	2.31	0.59
1:B:13:HIS:CE1	1:B:124:LYS:HE3	2.37	0.59
1:G:131:ASP:O	1:G:135:THR:HG23	2.03	0.59
1:H:78:GLY:O	1:H:79:ARG:NH1	2.36	0.59
1:H:63:ARG:NH2	5:H:302:HOH:O	2.35	0.59
1:L:72:LEU:HD22	1:L:132:PHE:CD1	2.38	0.59
1:B:152:VAL:O	1:B:156:ARG:HG3	2.02	0.58
1:K:101:GLU:OE2	1:K:156:ARG:NH2	2.33	0.58
1:A:118:HIS:O	1:A:122:THR:HG23	2.04	0.58
1:A:121:ALA:HB2	1:A:129:LEU:HD23	1.86	0.58
1:J:28:LEU:HB3	1:J:85:ILE:HD12	1.86	0.57
1:E:121:ALA:CB	1:E:129:LEU:HD23	2.35	0.57
1:F:63:ARG:NH1	1:F:67:GLU:OE2	2.37	0.57
1:K:63:ARG:HD3	5:K:318:HOH:O	2.05	0.57
1:G:82:LEU:HB3	1:J:32:TYR:OH	2.04	0.57
1:C:131:ASP:O	1:C:135:THR:HG23	2.04	0.57
1:A:36:SER:HB2	1:D:82:LEU:CD1	2.35	0.57
1:B:31:SER:O	1:B:59:SER:HB2	2.05	0.57
1:C:172:LYS:HD2	1:K:173:HIS:HB3	1.86	0.57
1:D:20:ILE:HD13	1:D:117:LEU:HD11	1.87	0.57
1:C:161:PRO:HD2	1:C:162:GLU:OE1	2.05	0.56
1:G:112:GLN:O	1:G:116:GLU:HG2	2.06	0.56
1:G:148:LEU:O	1:G:152:VAL:HG23	2.06	0.56
1:A:158:MET:HE3	1:A:170:PHE:HB2	1.88	0.55
1:C:82:LEU:CD1	1:E:36:SER:HB2	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:69:LEU:HG	1:J:137:TYR:OH	2.06	0.55
1:E:71:LYS:HE2	1:E:75:GLN:NE2	2.22	0.55
1:F:20:ILE:HD13	1:F:117:LEU:HD11	1.89	0.55
1:A:19:ALA:HB1	1:A:117:LEU:HD13	1.87	0.55
1:L:31:SER:HB2	1:L:62:GLU:CB	2.35	0.55
1:F:82:LEU:CD1	1:H:36:SER:HB2	2.37	0.55
1:L:24:ILE:HD13	1:L:70:MET:HG2	1.88	0.55
1:E:131:ASP:HA	1:E:134:GLU:CG	2.36	0.55
1:K:63:ARG:NH1	4:K:202:MPD:O2	2.40	0.55
1:G:82:LEU:CD1	1:J:36:SER:HB2	2.37	0.54
1:A:118:HIS:CE1	1:C:127:PRO:HB3	2.43	0.54
1:I:24:ILE:HD13	1:I:70:MET:HG2	1.89	0.54
1:I:28:LEU:HD23	1:I:85:ILE:CD1	2.37	0.54
1:A:96:GLY:N	1:A:167:GLU:OE2	2.40	0.54
1:E:20:ILE:HD13	1:E:117:LEU:HD11	1.90	0.54
1:F:118:HIS:O	1:F:122:THR:HG23	2.07	0.54
1:C:112:GLN:O	1:C:116:GLU:HG2	2.08	0.54
1:I:141:GLN:O	1:I:145:ILE:HD12	2.07	0.54
1:L:72:LEU:HD22	1:L:132:PHE:CE1	2.43	0.54
1:C:118:HIS:CE1	1:G:127:PRO:HB3	2.43	0.54
1:E:54:TYR:O	1:E:58:GLN:HG2	2.08	0.54
1:G:117:LEU:HD22	1:G:133:ILE:CD1	2.36	0.54
1:I:108:LYS:CG	1:I:145:ILE:HG12	2.36	0.54
1:J:28:LEU:HD13	1:J:85:ILE:HD11	1.89	0.54
1:H:28:LEU:HB3	1:H:85:ILE:HD11	1.87	0.53
1:I:28:LEU:HD12	1:I:66:ALA:CB	2.38	0.53
1:C:63[A]:ARG:HD2	1:C:67:GLU:OE1	2.08	0.53
1:C:87:LYS:HD2	1:E:83:GLN:HA	1.91	0.53
1:F:40:TYR:O	1:F:43:ARG:HG3	2.08	0.53
1:J:154:ASN:O	1:J:158:MET:HG3	2.08	0.53
1:C:164:GLY:HA3	1:K:157:LYS:O	2.08	0.53
1:F:100:MET:CE	1:F:148:LEU:HD23	2.38	0.53
1:G:29:TYR:O	1:G:33:VAL:HG23	2.08	0.53
1:G:159:GLY:O	1:G:162:GLU:HB2	2.09	0.53
1:F:24:ILE:HD13	1:F:70:MET:HG2	1.90	0.53
1:G:82:LEU:HD12	1:J:36:SER:HB2	1.91	0.53
1:C:145:ILE:HG22	1:G:8:VAL:HB	1.90	0.52
1:C:63[A]:ARG:HH21	4:E:202:MPD:H12	1.75	0.52
1:A:108:LYS:HG2	1:C:10:GLN:NE2	2.23	0.52
1:A:108:LYS:HE3	1:C:7:GLN:O	2.08	0.52
1:G:130:ALA:O	1:G:134:GLU:HG3	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:20:ILE:O	1:I:24:ILE:HG13	2.10	0.52
1:F:130:ALA:O	1:F:134:GLU:HG3	2.10	0.52
1:C:63[A]:ARG:NH1	1:C:67:GLU:OE1	2.43	0.51
1:B:23:GLN:O	1:B:27:GLU:HG2	2.10	0.51
1:K:20:ILE:CD1	1:K:117:LEU:HD21	2.41	0.51
1:C:148:LEU:O	1:C:152:VAL:HG22	2.10	0.51
1:G:154:ASN:O	1:G:158:MET:HG3	2.10	0.51
1:I:129:LEU:O	1:I:133:ILE:HG12	2.11	0.51
1:F:86:GLN:N	5:F:301:HOH:O	2.23	0.51
1:C:82:LEU:HD12	1:E:36:SER:HB2	1.92	0.50
1:C:42:ASP:HB3	1:E:74:ASN:HD22	1.76	0.50
1:J:148:LEU:O	1:J:152:VAL:HG22	2.11	0.50
1:A:176:GLY:O	1:A:177:ASP:CB	2.58	0.50
1:A:72:LEU:HD21	1:A:129:LEU:CD1	2.41	0.50
1:F:96:GLY:N	1:F:167:GLU:OE1	2.44	0.50
1:F:100:MET:HE2	1:F:151:HIS:HB3	1.93	0.50
1:K:49:LYS:O	1:K:53:LYS:HG3	2.11	0.50
1:A:6:SER:OG	1:A:8:VAL:HG22	2.12	0.50
1:F:153:THR:O	1:F:157:LYS:HG3	2.12	0.50
1:G:141:GLN:O	1:G:145:ILE:HG13	2.11	0.50
1:F:63:ARG:NH2	1:H:59:SER:OG	2.39	0.49
1:A:129:LEU:O	1:A:133:ILE:HG12	2.12	0.49
1:F:82:LEU:HD12	1:H:36:SER:HB2	1.93	0.49
1:E:13:HIS:ND1	1:E:124:LYS:HD2	2.26	0.49
1:C:36:SER:HB2	1:E:82:LEU:CD1	2.42	0.49
1:B:24:ILE:HD13	1:B:70:MET:HG2	1.93	0.49
1:A:69:LEU:HG	1:A:137:TYR:OH	2.12	0.49
1:C:82:LEU:HA	5:C:320:HOH:O	2.12	0.49
1:E:24:ILE:HD13	1:E:70:MET:HG2	1.94	0.49
1:A:127:PRO:HB3	1:G:118:HIS:CE1	2.48	0.49
1:E:117:LEU:O	1:E:117:LEU:HD23	2.13	0.49
1:D:139:ASN:O	1:D:143:LYS:HG3	2.13	0.48
1:H:41:PHE:HA	1:H:46:VAL:HG11	1.95	0.48
1:J:129:LEU:O	1:J:133:ILE:HG12	2.13	0.48
1:B:101:GLU:OE2	1:B:156:ARG:NH1	2.46	0.48
1:C:85:ILE:HB	1:E:85:ILE:HB	1.94	0.48
1:D:50:ASN:HB2	1:D:171:ASP:OD2	2.13	0.48
1:L:19:ALA:HB1	1:L:117:LEU:HD13	1.94	0.48
1:C:74:ASN:HD22	1:E:42:ASP:HB3	1.77	0.48
1:E:131:ASP:HB2	1:K:134:GLU:OE2	2.13	0.48
1:I:28:LEU:HD12	1:I:66:ALA:HB2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:108:LYS:NZ	1:H:145:ILE:HD13	2.27	0.48
1:F:82:LEU:HA	5:F:309:HOH:O	2.14	0.48
1:J:28:LEU:CB	1:J:85:ILE:HD12	2.43	0.48
1:K:97:LEU:HD11	1:K:156:ARG:CG	2.43	0.48
1:F:100:MET:HE1	1:F:148:LEU:HD23	1.94	0.48
1:G:28:LEU:HD13	1:G:85:ILE:HD11	1.95	0.48
1:C:44:ASP:HA	1:K:150:ASP:OD1	2.14	0.48
1:A:34:TYR:CB	1:A:59:SER:HB2	2.44	0.48
1:D:63:ARG:NH1	1:D:67:GLU:OE2	2.43	0.48
1:E:13:HIS:CG	1:E:124:LYS:HD2	2.49	0.48
1:K:129:LEU:HD12	1:K:129:LEU:O	2.13	0.48
1:A:72:LEU:HD21	1:A:129:LEU:HD12	1.95	0.48
1:K:178:SER:O	1:K:178:SER:OG	2.26	0.47
1:A:157:LYS:O	1:E:164:GLY:HA3	2.14	0.47
1:G:28:LEU:HB3	1:G:85:ILE:CD1	2.40	0.47
1:L:141:GLN:HE21	1:L:141:GLN:HA	1.79	0.47
1:I:28:LEU:CD1	1:I:66:ALA:CB	2.92	0.47
1:A:158:MET:HE1	1:A:170:PHE:HD1	1.80	0.47
1:B:118:HIS:CE1	1:F:127:PRO:HB3	2.49	0.47
1:I:61:GLU:OE1	1:I:61:GLU:HA	2.14	0.47
1:E:22:ARG:NH1	5:E:301:HOH:O	1.99	0.47
1:I:28:LEU:HD23	1:I:85:ILE:HD11	1.96	0.47
1:I:28:LEU:CD1	1:I:66:ALA:HB1	2.45	0.47
1:A:82:LEU:HD13	1:D:32:TYR:CE1	2.50	0.47
1:H:108:LYS:NZ	1:H:145:ILE:CD1	2.79	0.46
1:C:172:LYS:HD3	1:K:173:HIS:HB3	1.98	0.46
1:F:168:TYR:CD1	1:F:172:LYS:HE3	2.48	0.46
1:C:153:THR:HG22	1:C:157:LYS:HD2	1.99	0.45
1:I:101:GLU:OE2	1:I:156:ARG:NH2	2.36	0.45
1:A:59:SER:OG	1:D:63:ARG:NH2	2.48	0.45
1:B:49:LYS:O	1:B:49:LYS:HD3	2.16	0.45
1:J:5:THR:N	5:J:303:HOH:O	2.49	0.45
1:K:141:GLN:O	1:K:145:ILE:HG13	2.17	0.45
1:K:108:LYS:NZ	5:K:303:HOH:O	2.47	0.45
4:L:203:MPD:O4	4:L:203:MPD:O2	2.24	0.45
1:A:31:SER:HB2	1:A:62:GLU:HB2	1.98	0.45
1:J:97:LEU:HD21	1:J:101:GLU:OE2	2.17	0.45
1:C:32:TYR:OH	1:E:82:LEU:HB3	2.17	0.45
1:C:42:ASP:HB3	1:E:74:ASN:ND2	2.32	0.45
1:H:162:GLU:OE1	1:H:162:GLU:N	2.49	0.45
1:I:60:HIS:CE1	1:L:63:ARG:HH11	2.35	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:ASN:N	5:A:303:HOH:O	2.50	0.44
1:F:86:GLN:HG2	5:F:301:HOH:O	2.17	0.44
1:I:86:GLN:HA	1:L:84:ASP:OD2	2.17	0.44
1:K:54:TYR:HB2	1:K:175:LEU:HD22	1.99	0.44
1:C:31:SER:HB2	1:C:62:GLU:HB2	1.99	0.44
1:H:148:LEU:O	1:H:152:VAL:HG13	2.17	0.44
1:B:78:GLY:O	1:B:79:ARG:NH1	2.50	0.44
1:K:69:LEU:HG	1:K:137:TYR:OH	2.18	0.44
1:H:60:HIS:O	1:H:64:GLU:HG3	2.16	0.44
1:B:54:TYR:O	1:B:58:GLN:HG2	2.18	0.44
1:D:117:LEU:HD22	1:D:133:ILE:CD1	2.48	0.44
1:F:98:ASN:C	1:F:98:ASN:HD22	2.17	0.44
1:H:155:LEU:HD21	1:H:170:PHE:CD2	2.52	0.44
1:K:97:LEU:HD11	1:K:156:ARG:HG2	1.98	0.44
1:A:12:TYR:OH	1:A:73:GLN:OE1	2.32	0.44
1:L:69:LEU:HG	1:L:137:TYR:OH	2.17	0.44
1:A:87:LYS:HG3	1:A:88:PRO:HD2	1.99	0.44
1:J:93:TRP:O	1:J:95:SER:N	2.47	0.44
1:H:69:LEU:HG	1:H:137:TYR:OH	2.18	0.43
1:C:87:LYS:HB2	1:C:87:LYS:HE2	1.60	0.43
1:E:157:LYS:O	1:L:164:GLY:HA3	2.18	0.43
1:H:31:SER:HB2	1:H:62:GLU:CB	2.47	0.43
1:E:69:LEU:HG	1:E:137:TYR:OH	2.17	0.43
1:L:154:ASN:O	1:L:158:MET:HG3	2.19	0.43
1:D:91:ASP:OD2	1:D:91:ASP:N	2.50	0.43
1:A:148:LEU:O	1:A:152:VAL:HG23	2.18	0.43
1:D:40:TYR:CZ	1:D:46:VAL:HG21	2.53	0.43
1:F:54:TYR:O	1:F:58:GLN:HG2	2.19	0.43
1:F:148:LEU:O	1:F:152:VAL:HG23	2.19	0.42
1:G:124:LYS:HA	1:G:124:LYS:HD3	1.63	0.42
1:H:31:SER:HB2	1:H:62:GLU:HB2	2.01	0.42
4:F:202:MPD:C1	1:H:31:SER:OG	2.67	0.42
1:J:34:TYR:CB	1:J:59:SER:HB2	2.49	0.42
1:J:50:ASN:HB2	1:J:171:ASP:OD2	2.19	0.42
1:F:129:LEU:O	1:F:133:ILE:HG12	2.20	0.42
1:B:85:ILE:O	1:B:86:GLN:CB	2.60	0.42
1:F:63:ARG:NH1	1:H:35:LEU:HD13	2.34	0.42
1:I:28:LEU:HD23	1:I:85:ILE:HD13	2.02	0.42
1:J:176:GLY:O	1:J:177:ASP:CB	2.63	0.42
1:F:168:TYR:CE2	1:F:172:LYS:HE3	2.52	0.42
1:I:118:HIS:CE1	1:K:127:PRO:HB3	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:104:LEU:CD1	1:L:145:ILE:HG23	2.49	0.41
1:B:112:GLN:O	1:B:116:GLU:HG3	2.20	0.41
1:B:153:THR:HG21	1:H:45:ASP:C	2.41	0.41
1:L:76:ARG:HD3	1:L:128:HIS:CE1	2.56	0.41
1:L:161:PRO:HG2	5:L:317:HOH:O	2.20	0.41
1:D:161:PRO:HD2	1:D:162:GLU:OE1	2.19	0.41
1:F:32:TYR:CE1	1:H:82:LEU:HD13	2.55	0.41
1:F:63:ARG:HH12	1:H:35:LEU:HD13	1.85	0.41
1:F:160:ALA:HB1	1:F:161:PRO:HA	2.01	0.41
1:G:41:PHE:CZ	1:G:96:GLY:HA2	2.56	0.41
1:I:160:ALA:HB1	1:I:161:PRO:HA	2.03	0.41
1:A:158:MET:HE1	1:A:170:PHE:CD1	2.55	0.41
1:E:118:HIS:CE1	1:I:127:PRO:HB3	2.55	0.41
1:H:127:PRO:HB3	1:L:118:HIS:CE1	2.54	0.41
1:H:108:LYS:HE2	1:H:145:ILE:CG2	2.49	0.41
1:G:82:LEU:HD13	1:J:32:TYR:CE1	2.56	0.41
1:G:101:GLU:OE2	1:G:156:ARG:NH2	2.53	0.41
1:H:79:ARG:HA	1:H:79:ARG:HD3	1.84	0.41
1:I:58:GLN:O	1:I:61:GLU:HB2	2.21	0.41
1:A:132:PHE:CE1	1:A:137:TYR:HE1	2.39	0.41
1:C:63[A]:ARG:CZ	4:E:202:MPD:H12	2.51	0.41
1:F:20:ILE:HD13	1:F:20:ILE:HA	1.94	0.41
1:A:42:ASP:HB3	1:D:74:ASN:ND2	2.36	0.41
1:D:31:SER:HB2	1:D:62:GLU:HB2	2.02	0.41
1:F:32:TYR:OH	1:H:82:LEU:HB3	2.21	0.41
1:J:20:ILE:HD13	1:J:20:ILE:HA	1.90	0.41
1:J:97:LEU:CD2	1:J:101:GLU:OE2	2.69	0.41
1:F:68:LYS:HB2	1:F:68:LYS:HE3	1.79	0.40
1:G:67:GLU:HB3	1:J:39:TYR:OH	2.21	0.40
1:B:28:LEU:CB	1:B:85:ILE:HD12	2.47	0.40
1:C:24:ILE:HD13	1:C:70:MET:HG2	2.04	0.40
1:F:141:GLN:O	1:F:145:ILE:HG13	2.22	0.40
1:L:131:ASP:HA	1:L:134:GLU:OE1	2.21	0.40
4:F:202:MPD:HM1	4:F:202:MPD:H4	1.92	0.40
1:J:155:LEU:HD21	1:J:170:PHE:CD2	2.57	0.40
1:C:101:GLU:OE2	1:C:156:ARG:NH1	2.54	0.40
1:E:63:ARG:NH1	4:E:202:MPD:H53	2.37	0.40
1:E:131:ASP:HA	1:E:134:GLU:HG3	2.04	0.40
1:E:160:ALA:HB1	1:E:161:PRO:HA	2.03	0.40
1:I:79:ARG:HA	1:I:79:ARG:HD3	1.89	0.40
1:E:27:GLU:OE1	1:E:62:GLU:OE2	2.39	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:27:GLU:HB2	1:J:66:ALA:HB2	2.02	0.40
1:K:9:ARG:NH2	1:K:17:GLU:OE1	2.42	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)
5:K:318:HOH:O	5:K:318:HOH:O[4_555]	2.02	0.18

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	172/182 (94%)	167 (97%)	5 (3%)	0	100	100
1	B	173/182 (95%)	169 (98%)	3 (2%)	1 (1%)	25	50
1	C	173/182 (95%)	169 (98%)	4 (2%)	0	100	100
1	D	174/182 (96%)	171 (98%)	3 (2%)	0	100	100
1	E	171/182 (94%)	165 (96%)	5 (3%)	1 (1%)	25	50
1	F	171/182 (94%)	165 (96%)	6 (4%)	0	100	100
1	G	172/182 (94%)	165 (96%)	7 (4%)	0	100	100
1	H	171/182 (94%)	168 (98%)	3 (2%)	0	100	100
1	I	171/182 (94%)	165 (96%)	6 (4%)	0	100	100
1	J	172/182 (94%)	168 (98%)	4 (2%)	0	100	100
1	K	174/182 (96%)	167 (96%)	6 (3%)	1 (1%)	25	50
1	L	171/182 (94%)	167 (98%)	4 (2%)	0	100	100
All	All	2065/2184 (95%)	2006 (97%)	56 (3%)	3 (0%)	51	78

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	5	THR
1	B	86	GLN
1	K	177	ASP

5.3.2 Protein sidechains (i)

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	146/160 (91%)	144 (99%)	2 (1%)	67	86
1	B	151/160 (94%)	145 (96%)	6 (4%)	31	60
1	C	148/160 (92%)	140 (95%)	8 (5%)	22	47
1	D	147/160 (92%)	144 (98%)	3 (2%)	55	81
1	E	145/160 (91%)	144 (99%)	1 (1%)	84	94
1	F	146/160 (91%)	140 (96%)	6 (4%)	30	59
1	G	146/160 (91%)	145 (99%)	1 (1%)	84	94
1	H	143/160 (89%)	140 (98%)	3 (2%)	53	80
1	I	144/160 (90%)	142 (99%)	2 (1%)	67	86
1	J	142/160 (89%)	139 (98%)	3 (2%)	53	80
1	K	146/160 (91%)	144 (99%)	2 (1%)	67	86
1	L	145/160 (91%)	142 (98%)	3 (2%)	53	80
All	All	1749/1920 (91%)	1709 (98%)	40 (2%)	52	78

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	HIS
1	A	32	TYR
1	B	32	TYR
1	B	59	SER
1	B	63[A]	ARG
1	B	63[B]	ARG
1	B	87	LYS
1	B	117	LEU

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Mol	Chain	Res	Type
1	C	31	SER
1	C	32	TYR
1	C	61	GLU
1	C	63[A]	ARG
1	C	63[B]	ARG
1	C	65	HIS
1	C	71	LYS
1	C	177	ASP
1	D	32	TYR
1	D	63	ARG
1	D	91	ASP
1	E	32	TYR
1	F	32	TYR
1	F	63	ARG
1	F	68	LYS
1	F	71	LYS
1	F	98	ASN
1	F	105	HIS
1	G	32	TYR
1	H	32	TYR
1	H	36	SER
1	H	89	ASP
1	I	32	TYR
1	I	113	SER
1	J	32	TYR
1	J	44	ASP
1	J	89	ASP
1	K	32	TYR
1	K	112	GLN
1	L	32	TYR
1	L	44	ASP
1	L	89	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	105	HIS
1	C	10	GLN
1	E	75	GLN
1	F	109	ASN
1	H	111	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 26 ligands modelled in this entry, 19 are monoatomic - leaving 7 for Mogul analysis.

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	$\#\|Z\| > 2$
4	MPD	K	202	-	7,7,7	0.26	0	9,10,10	0.18	0
4	MPD	G	202	-	7,7,7	0.28	0	9,10,10	0.37	0
4	MPD	F	202	-	7,7,7	0.23	0	9,10,10	0.84	1 (11%)
4	MPD	L	203	-	7,7,7	0.31	0	9,10,10	0.71	0
4	MPD	E	202	-	7,7,7	0.24	0	9,10,10	0.52	0
4	MPD	D	202	-	7,7,7	0.25	0	9,10,10	0.69	0
4	MPD	B	204	-	7,7,7	0.24	0	9,10,10	0.72	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
4	MPD	K	202	-	-	0/5/5/5	-
4	MPD	G	202	-	-	0/5/5/5	-
4	MPD	F	202	-	-	0/5/5/5	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	MPD	L	203	-	-	2/5/5/5	-
4	MPD	E	202	-	-	0/5/5/5	-
4	MPD	D	202	-	-	1/5/5/5	-
4	MPD	B	204	-	-	0/5/5/5	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	F	202	MPD	CM-C2-C1	-2.23	105.92	110.57

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	L	203	MPD	C2-C3-C4-O4
4	L	203	MPD	C2-C3-C4-C5
4	D	202	MPD	CM-C2-C3-C4

There are no ring outliers.

5 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	K	202	MPD	1	0
4	F	202	MPD	2	0
4	L	203	MPD	1	0
4	E	202	MPD	4	0
4	D	202	MPD	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, 95^{th} 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	174/182 (95%)	-0.12	0 100 100	44, 54, 74, 94	0
1	B	173/182 (95%)	-0.19	0 100 100	39, 50, 64, 81	0
1	C	174/182 (95%)	-0.16	0 100 100	49, 59, 78, 95	0
1	D	176/182 (96%)	-0.16	2 (1%) 80 82	42, 51, 70, 110	0
1	E	173/182 (95%)	-0.09	0 100 100	48, 59, 74, 100	0
1	F	173/182 (95%)	-0.22	0 100 100	41, 53, 71, 95	0
1	G	174/182 (95%)	-0.12	0 100 100	47, 59, 81, 114	0
1	H	173/182 (95%)	-0.16	2 (1%) 79 80	39, 60, 80, 96	0
1	I	173/182 (95%)	-0.03	0 100 100	51, 64, 82, 97	0
1	J	173/182 (95%)	-0.22	1 (0%) 89 91	49, 65, 86, 94	0
1	K	175/182 (96%)	-0.23	0 100 100	51, 59, 76, 124	0
1	L	173/182 (95%)	-0.17	0 100 100	43, 64, 84, 110	0
All	All	2084/2184 (95%)	-0.16	5 (0%) 95 96	39, 58, 80, 124	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	108	LYS	2.8
1	H	112	GLN	2.2
1	J	10	GLN	2.2
1	D	3	ALA	2.0
1	D	4	SER	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates (i)

There are no monosaccharides 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, 95^{th} 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
2	FE	I	201	1/1	0.84	0.10	103,103,103,103	1
3	NA	B	202	1/1	0.84	0.22	54,54,54,54	0
2	FE	C	201	1/1	0.86	0.06	107,107,107,107	1
3	NA	L	202	1/1	0.87	0.14	102,102,102,102	0
2	FE	D	201	1/1	0.90	0.12	97,97,97,97	1
2	FE	G	201	1/1	0.91	0.08	104,104,104,104	1
2	FE	K	201	1/1	0.91	0.10	105,105,105,105	1
2	FE	F	201	1/1	0.92	0.18	105,105,105,105	1
2	FE	B	201	1/1	0.92	0.09	102,102,102,102	1
2	FE	E	201	1/1	0.93	0.09	93,93,93,93	1
2	FE	L	201	1/1	0.93	0.17	108,108,108,108	1
3	NA	C	202	1/1	0.94	0.20	48,48,48,48	0
3	NA	H	202	1/1	0.95	0.14	74,74,74,74	0
2	FE	H	201	1/1	0.95	0.13	127,127,127,127	1
4	MPD	E	202	8/8	0.95	0.38	45,64,74,74	0
4	MPD	G	202	8/8	0.95	0.47	67,84,97,97	0
4	MPD	L	203	8/8	0.95	0.38	58,72,87,87	0
4	MPD	D	202	8/8	0.96	0.42	41,55,63,63	0
3	NA	B	203	1/1	0.96	0.13	35,35,35,35	0
4	MPD	F	202	8/8	0.96	0.44	73,89,91,93	0
2	FE	J	201	1/1	0.96	0.12	118,118,118,118	1
4	MPD	B	204	8/8	0.96	0.28	48,59,64,64	22
2	FE	A	201	1/1	0.97	0.15	83,83,83,83	1
3	NA	A	203	1/1	0.98	0.13	32,32,32,32	0
4	MPD	K	202	8/8	0.98	0.34	36,45,48,48	22
3	NA	A	202	1/1	0.98	0.14	81,81,81,81	0

6.5 Other polymers (i)

There are no such residues in this entry.