Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:59 AM GMT

PDB ID	:	2EX3
Title	:	Bacteriophage phi29 DNA polymerase bound to terminal protein
Authors	:	Kamtekar, S.; Berman, A.J.; Wang, J.; de Vega, M.; Blanco, L.; Salas, M.; Steitz, T.A.
Deposited on	:	2005-11-07
Resolution	:	3.00 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see references ①) were used in the production of this report:

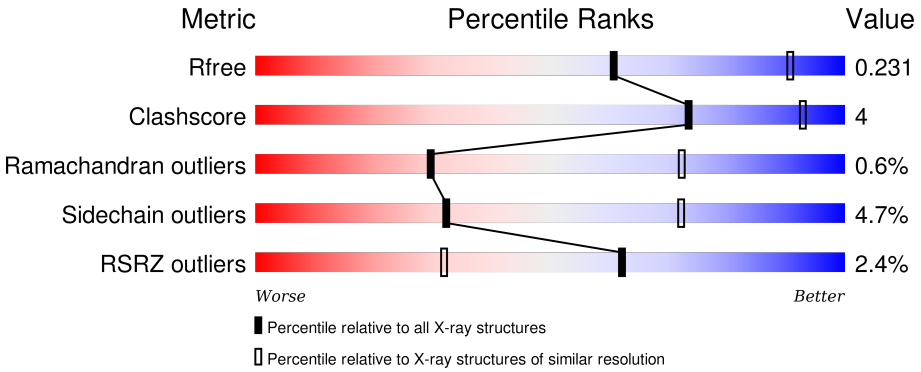
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

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 3.00 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}	91344	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

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. 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	575	2% 81% 18%
1	C	575	2% 82% 16%
1	E	575	2% 86% 12%
1	G	575	% 86% 11%
1	I	575	2% 87% 11%

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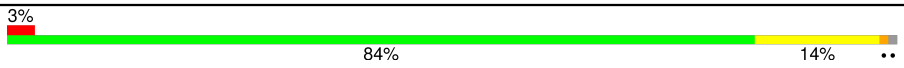
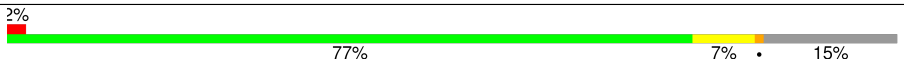
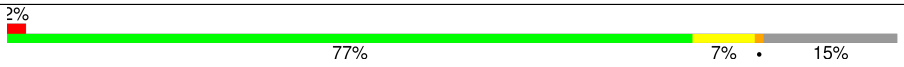
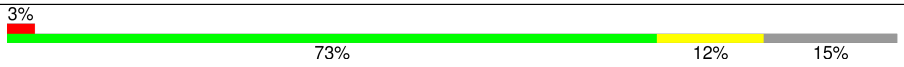
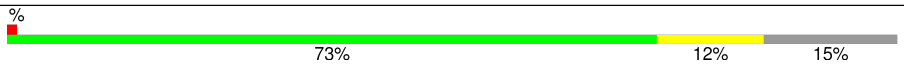
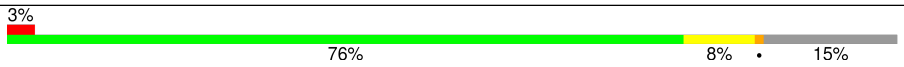
Mol	Chain	Length	Quality of chain
1	K	575	
2	B	230	
2	D	230	
2	F	230	
2	H	230	
2	J	230	
2	L	230	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 37108 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 DNA polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	570	Total	C	N	O	S	0	0	0
			4661	3036	753	851	21			
1	C	570	Total	C	N	O	S	0	0	0
			4661	3036	753	851	21			
1	E	570	Total	C	N	O	S	0	0	0
			4661	3036	753	851	21			
1	G	570	Total	C	N	O	S	0	0	0
			4661	3036	753	851	21			
1	I	570	Total	C	N	O	S	0	0	0
			4661	3036	753	851	21			
1	K	570	Total	C	N	O	S	0	0	0
			4661	3036	753	851	21			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	12	ALA	ASP	engineered	UNP P03680
A	66	ALA	ASP	engineered	UNP P03680
C	12	ALA	ASP	engineered	UNP P03680
C	66	ALA	ASP	engineered	UNP P03680
E	12	ALA	ASP	engineered	UNP P03680
E	66	ALA	ASP	engineered	UNP P03680
G	12	ALA	ASP	engineered	UNP P03680
G	66	ALA	ASP	engineered	UNP P03680
I	12	ALA	ASP	engineered	UNP P03680
I	66	ALA	ASP	engineered	UNP P03680
K	12	ALA	ASP	engineered	UNP P03680
K	66	ALA	ASP	engineered	UNP P03680

- Molecule 2 is a protein called DNA terminal protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	196	Total	C	N	O	S	0	0	0
			1521	959	261	296	5			
2	D	196	Total	C	N	O	S	0	0	0
			1521	959	261	296	5			
2	F	196	Total	C	N	O	S	0	0	0
			1521	959	261	296	5			
2	H	196	Total	C	N	O	S	0	0	0
			1521	959	261	296	5			
2	J	196	Total	C	N	O	S	0	0	0
			1521	959	261	296	5			
2	L	196	Total	C	N	O	S	0	0	0
			1521	959	261	296	5			

- Molecule 3 is LEAD (II) ION (three-letter code: PB) (formula: Pb).

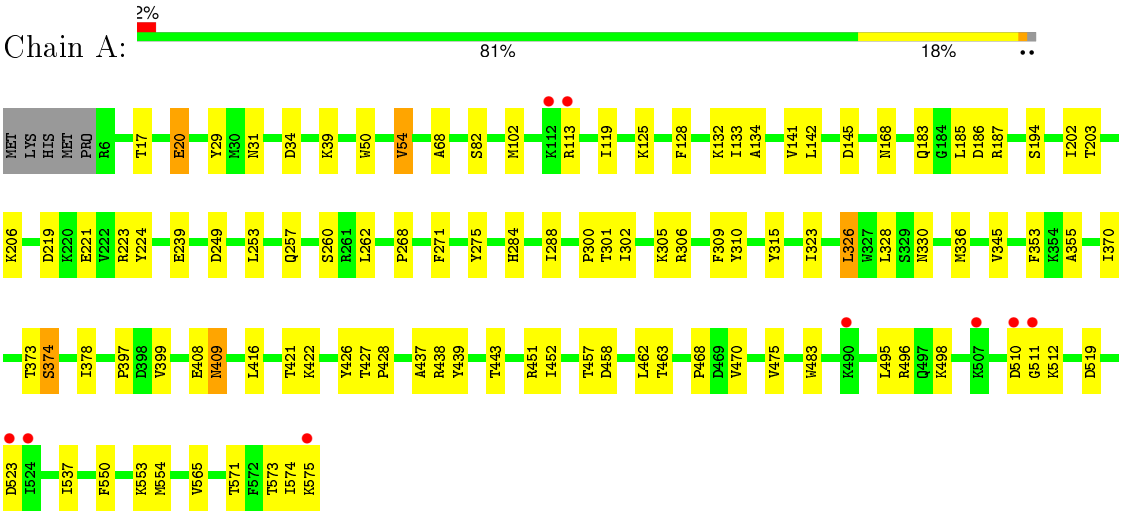
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	3	Total	Pb	0	0
			3	3		
3	K	2	Total	Pb	0	0
			2	2		
3	E	3	Total	Pb	0	0
			3	3		
3	I	2	Total	Pb	0	0
			2	2		
3	C	3	Total	Pb	0	0
			3	3		
3	A	3	Total	Pb	0	0
			3	3		

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

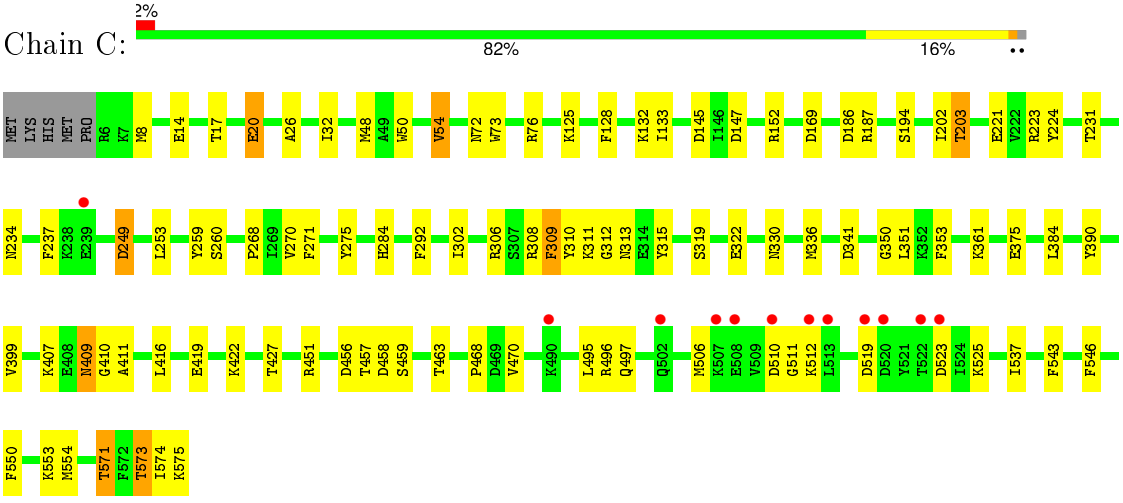
- Molecule 1: DNA polymerase

Chain A:



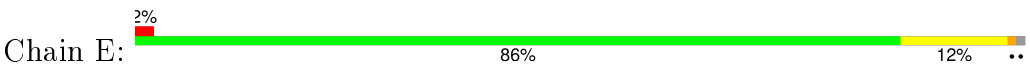
- Molecule 1: DNA polymerase

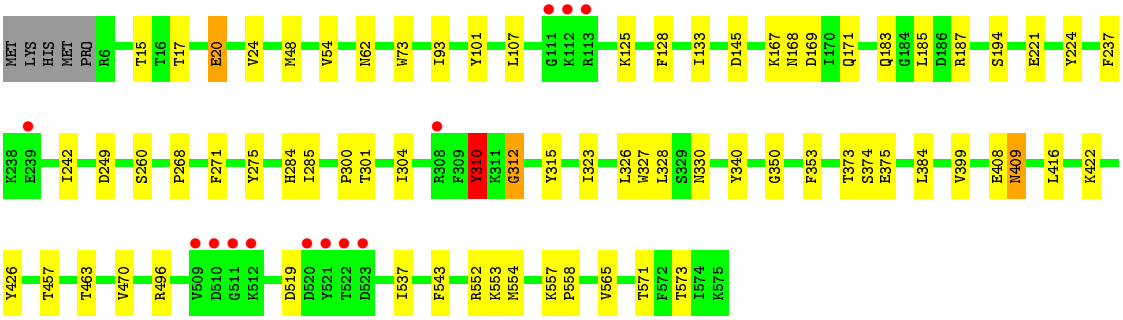
Chain C:



- Molecule 1: DNA polymerase

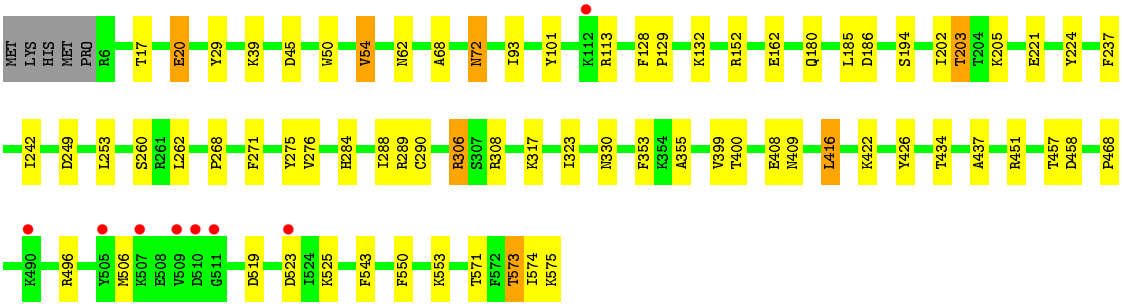
Chain E:





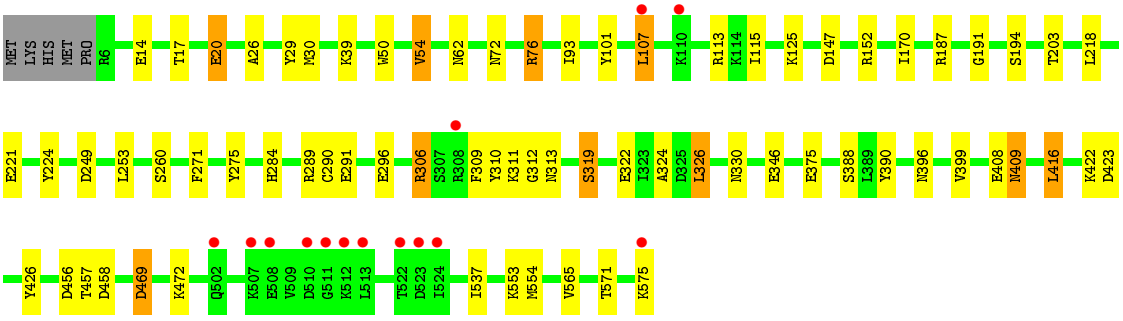
• Molecule 1: DNA polymerase

Chain G:



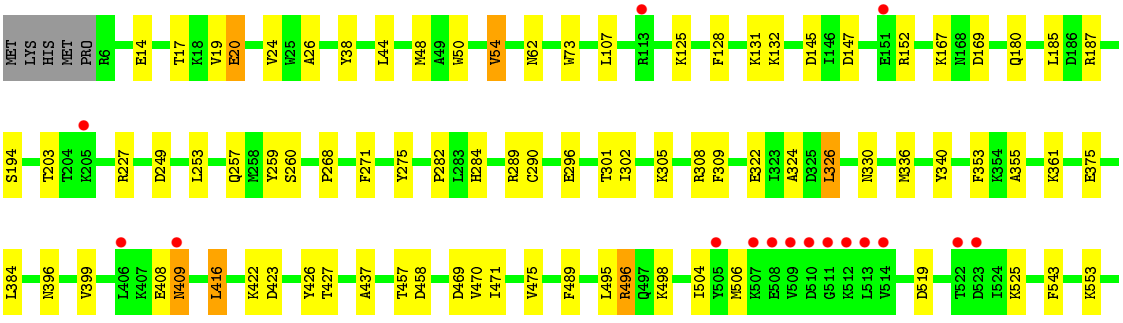
• Molecule 1: DNA polymerase

Chain I:



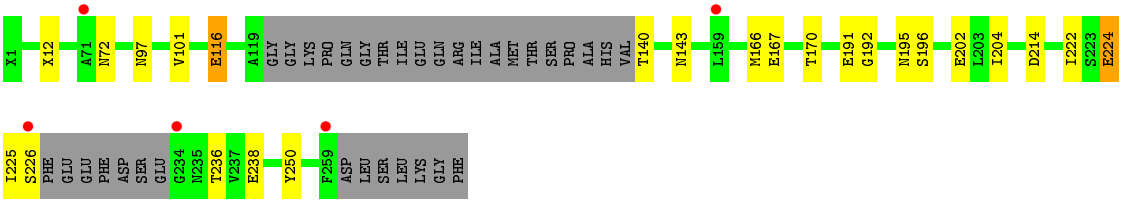
• Molecule 1: DNA polymerase

Chain K:



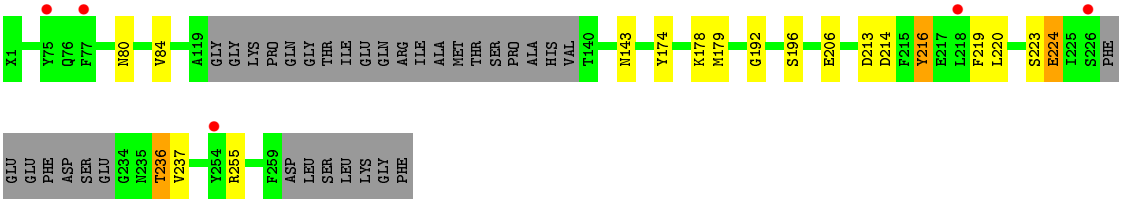
• Molecule 2: DNA terminal protein

Chain B:



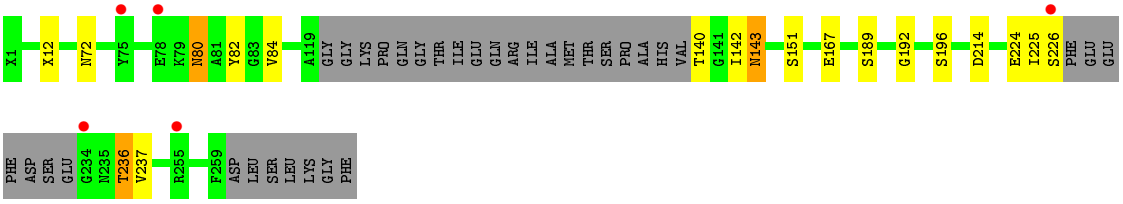
• Molecule 2: DNA terminal protein

Chain D:



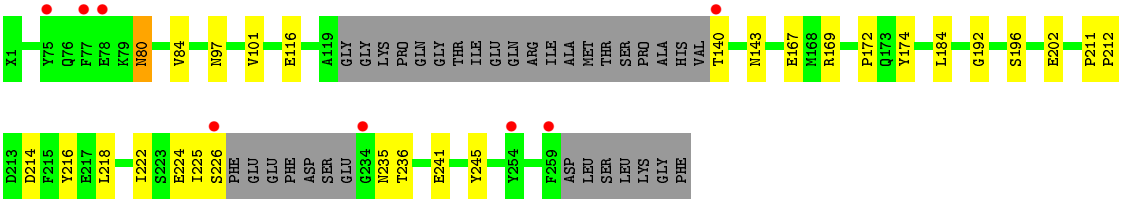
• Molecule 2: DNA terminal protein

Chain F:



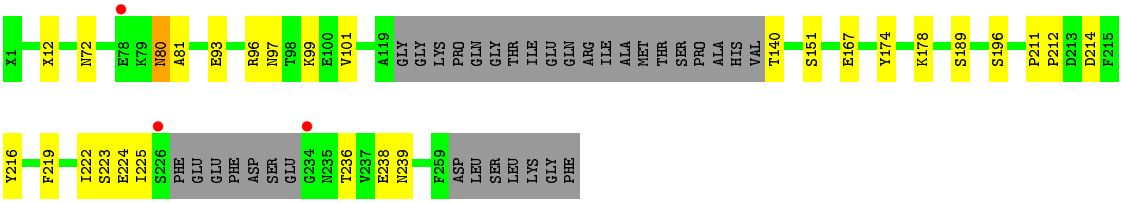
• Molecule 2: DNA terminal protein

Chain H:



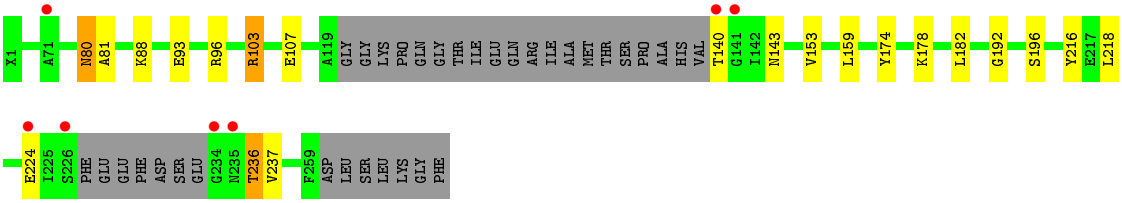
• Molecule 2: DNA terminal protein

Chain J:



- Molecule 2: DNA terminal protein

Chain L:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	304.93Å 220.28Å 217.16Å 90.00° 45.40° 90.00°	Depositor
Resolution (Å)	48.85 – 3.00 48.83 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.7 (48.85-3.00) 99.7 (48.83-3.00)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.43 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.2.0011	Depositor
R, R_{free}	0.200 , 0.229 0.204 , 0.231	Depositor DCC
R_{free} test set	20161 reflections (11.07%)	DCC
Wilson B-factor (Å^2)	91.1	Xtriage
Anisotropy	0.010	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 97.4	EDS
Estimated twinning fraction	0.006 for -h+2*l,k,-h+l 0.011 for -h-k+l,-l,-k 0.006 for -h+k+l,l,k 0.008 for -k-l,-h+l,-l 0.008 for k-l,h-l,-l 0.039 for k+l,h-l,k 0.044 for h+k-l,l,h-l 0.038 for -k+l,-h+l,-k 0.036 for h-k-l,-l,h-l 0.010 for h,-k,h-l 0.023 for h-2*l,-k,-l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.31$	Xtriage
Outliers	0 of 202286 reflections	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	37108	wwPDB-VP
Average B, all atoms (Å^2)	99.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.97% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: PB

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	0.38	0/4780	0.53	0/6448
1	C	0.37	0/4780	0.52	0/6448
1	E	0.37	0/4780	0.52	0/6448
1	G	0.37	0/4780	0.53	0/6448
1	I	0.37	0/4780	0.51	0/6448
1	K	0.37	0/4780	0.51	0/6448
2	B	0.36	0/1373	0.50	0/1837
2	D	0.38	0/1373	0.50	0/1837
2	F	0.36	0/1373	0.50	0/1837
2	H	0.35	0/1373	0.50	0/1837
2	J	0.38	0/1373	0.49	0/1837
2	L	0.37	0/1373	0.48	0/1837
All	All	0.37	0/36918	0.51	0/49710

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

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	4661	0	4668	53	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	4661	0	4668	51	0
1	E	4661	0	4668	35	0
1	G	4661	0	4668	35	0
1	I	4661	0	4668	36	0
1	K	4661	0	4668	43	0
2	B	1521	0	1366	13	0
2	D	1521	0	1365	13	0
2	F	1521	0	1366	11	0
2	H	1521	0	1365	15	0
2	J	1521	0	1366	14	0
2	L	1521	0	1366	14	0
3	A	3	0	0	0	0
3	C	3	0	0	0	0
3	E	3	0	0	0	0
3	G	3	0	0	0	0
3	I	2	0	0	0	0
3	K	2	0	0	0	0
All	All	37108	0	36202	303	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 4.

All (303) 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:L:80:ASN:HD22	2:L:81:ALA:H	1.28	0.81
1:E:310:TYR:CE1	1:E:312:GLY:HA2	2.20	0.76
1:E:399:VAL:HG21	1:E:422:LYS:HB2	1.69	0.75
1:I:309:PHE:CE1	1:I:322:GLU:HG2	2.24	0.72
1:A:370:ILE:O	1:A:374:SER:HB2	1.89	0.72
1:I:284:HIS:H	1:I:330:ASN:ND2	1.89	0.71
1:G:284:HIS:H	1:G:330:ASN:HD22	1.39	0.70
1:C:399:VAL:O	1:C:399:VAL:HG12	1.91	0.69
1:K:308:ARG:HB3	2:L:216:TYR:HE1	1.58	0.69
1:G:306:ARG:HG2	2:H:224:GLU:HG2	1.74	0.68
1:K:308:ARG:HB3	2:L:216:TYR:CE1	2.29	0.68
1:E:310:TYR:HE1	1:E:312:GLY:HA2	1.61	0.66
1:C:409:ASN:H	1:C:409:ASN:HD22	1.44	0.66
1:C:309:PHE:CE1	1:C:322:GLU:HG2	2.31	0.65
1:C:284:HIS:H	1:C:330:ASN:HD22	1.45	0.63
2:D:174:TYR:CE1	2:D:178:LYS:HE3	2.34	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:409:ASN:HD22	1:I:409:ASN:H	1.48	0.62
1:A:284:HIS:H	1:A:330:ASN:HD22	1.48	0.61
1:C:253:LEU:HD22	1:C:458:ASP:HB3	1.81	0.61
1:I:399:VAL:HG12	1:I:399:VAL:O	2.01	0.61
1:E:496:ARG:NH1	2:F:192:GLY:O	2.34	0.60
1:K:399:VAL:HG21	1:K:422:LYS:HB2	1.83	0.60
2:J:97:ASN:O	2:J:101:VAL:HG23	2.01	0.60
1:E:221:GLU:HG2	1:E:327:TRP:CH2	2.36	0.60
1:K:308:ARG:HA	2:L:182:LEU:HD13	1.84	0.59
2:J:80:ASN:HD22	2:J:81:ALA:H	1.48	0.59
1:C:50:TRP:O	1:C:54:VAL:HG22	2.03	0.59
1:C:259:TYR:O	1:C:361:LYS:HE2	2.03	0.59
1:A:537:ILE:HG12	1:A:554:MET:HG3	1.85	0.59
2:F:225:ILE:HD12	2:F:226:SER:H	1.68	0.58
1:A:20:GLU:CD	1:A:20:GLU:H	2.07	0.58
1:G:253:LEU:HD22	1:G:458:ASP:HB3	1.84	0.57
1:K:284:HIS:H	1:K:330:ASN:ND2	2.01	0.57
2:D:80:ASN:HD22	2:D:84:VAL:HB	1.69	0.57
1:A:68:ALA:CB	1:A:565:VAL:HG23	2.35	0.56
1:G:50:TRP:O	1:G:54:VAL:HG22	2.05	0.56
1:E:301:THR:HB	1:E:340:TYR:HE1	1.69	0.56
2:J:80:ASN:HD21	2:J:151:SER:HA	1.70	0.56
1:C:14:GLU:HB2	1:C:26:ALA:HB3	1.88	0.56
1:K:409:ASN:H	1:K:409:ASN:HD22	1.54	0.55
1:G:284:HIS:H	1:G:330:ASN:ND2	2.04	0.55
1:G:323:ILE:HD12	2:H:143:ASN:HB3	1.87	0.55
1:I:253:LEU:HD22	1:I:458:ASP:HB3	1.89	0.55
1:I:20:GLU:H	1:I:20:GLU:CD	2.08	0.55
1:E:408:GLU:CD	1:E:408:GLU:H	2.10	0.55
1:E:268:PRO:HB3	1:E:353:PHE:CE2	2.40	0.55
1:G:496:ARG:NH1	2:H:192:GLY:O	2.40	0.55
2:F:80:ASN:HB3	2:F:82:TYR:H	1.71	0.54
2:J:12:UNK:C	2:J:72:ASN:H	2.21	0.54
1:C:223:ARG:NH2	1:C:427:THR:HG21	2.23	0.54
1:A:399:VAL:HG12	1:A:399:VAL:O	2.08	0.54
1:G:268:PRO:HB3	1:G:353:PHE:CE2	2.42	0.54
2:L:174:TYR:CE1	2:L:178:LYS:HE2	2.42	0.54
1:E:552:ARG:HB3	1:E:554:MET:HG2	1.90	0.54
1:I:284:HIS:H	1:I:330:ASN:HD22	1.56	0.54
1:I:306:ARG:O	1:I:306:ARG:HG3	2.08	0.53
1:A:323:ILE:HD12	2:B:143:ASN:HB3	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:309:PHE:CD1	1:K:322:GLU:HG2	2.44	0.53
2:F:225:ILE:HD12	2:F:226:SER:N	2.24	0.53
1:K:284:HIS:CE1	1:K:330:ASN:HB3	2.44	0.53
1:E:553:LYS:HA	1:E:571:THR:HA	1.90	0.53
1:A:253:LEU:HD22	1:A:458:ASP:HB3	1.90	0.52
1:K:50:TRP:O	1:K:54:VAL:HG22	2.09	0.52
1:C:221:GLU:HA	1:C:224:TYR:CD1	2.44	0.52
1:C:399:VAL:O	1:C:399:VAL:CG1	2.57	0.52
1:C:231:THR:O	1:C:315:TYR:HE1	1.92	0.52
2:H:241:GLU:O	2:H:245:TYR:HD1	1.93	0.52
1:I:93:ILE:HB	1:I:101:TYR:HB2	1.90	0.52
1:A:510:ASP:O	1:A:512:LYS:N	2.41	0.51
2:F:80:ASN:HB2	2:F:84:VAL:H	1.75	0.51
1:I:14:GLU:HB2	1:I:26:ALA:HB3	1.91	0.51
1:C:268:PRO:HB3	1:C:353:PHE:CE2	2.45	0.51
1:A:202:ILE:HG23	1:A:206:LYS:HD3	1.92	0.51
2:F:140:THR:HG22	2:F:142:ILE:HG13	1.93	0.51
1:A:125:LYS:HD3	1:A:187:ARG:O	2.11	0.51
1:E:537:ILE:HG12	1:E:554:MET:HG3	1.93	0.51
1:I:537:ILE:HG12	1:I:554:MET:HG3	1.91	0.51
1:A:68:ALA:HB3	1:A:565:VAL:HG23	1.93	0.51
1:I:147:ASP:O	1:I:152:ARG:NH2	2.43	0.51
1:I:324:ALA:HB1	1:I:326:LEU:HD21	1.92	0.51
1:C:575:LYS:HG2	2:D:196:SER:HB2	1.92	0.51
1:C:506:MET:HG3	1:C:525:LYS:HB2	1.93	0.50
2:H:225:ILE:HD12	2:H:226:SER:N	2.26	0.50
1:C:495:LEU:HG	1:C:546:PHE:CE2	2.46	0.50
1:G:451:ARG:NH1	1:G:468:PRO:HG3	2.26	0.50
1:A:309:PHE:HD2	1:A:310:TYR:CD2	2.29	0.50
1:E:300:PRO:HB2	1:E:315:TYR:HB3	1.92	0.50
1:C:311:LYS:O	1:C:313:ASN:N	2.44	0.50
1:A:29:TYR:CZ	1:A:39:LYS:HB3	2.47	0.50
1:K:301:THR:HB	1:K:340:TYR:HE1	1.75	0.50
1:K:324:ALA:HB1	1:K:326:LEU:HD21	1.93	0.49
1:K:253:LEU:O	1:K:257:GLN:HG2	2.12	0.49
1:I:221:GLU:HA	1:I:224:TYR:CD1	2.47	0.49
1:K:125:LYS:HD3	1:K:187:ARG:O	2.12	0.49
1:A:50:TRP:O	1:A:54:VAL:HG22	2.13	0.49
2:L:80:ASN:HD22	2:L:81:ALA:N	2.04	0.49
2:L:103:ARG:O	2:L:107:GLU:HG2	2.11	0.49
1:K:147:ASP:O	1:K:152:ARG:NH2	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:253:LEU:HD21	1:G:437:ALA:HB1	1.96	0.48
1:E:284:HIS:H	1:E:330:ASN:HD22	1.60	0.48
1:G:271:PHE:CZ	1:G:275:TYR:HB2	2.48	0.48
1:I:125:LYS:HD3	1:I:187:ARG:O	2.14	0.48
1:I:408:GLU:H	1:I:408:GLU:CD	2.17	0.48
1:G:399:VAL:HG21	1:G:422:LYS:HB2	1.95	0.48
1:K:489:PHE:HB3	1:K:504:ILE:HD13	1.95	0.48
1:K:408:GLU:H	1:K:408:GLU:CD	2.17	0.48
1:K:496:ARG:NH1	2:L:192:GLY:O	2.45	0.48
2:L:80:ASN:ND2	2:L:81:ALA:H	2.06	0.48
1:I:29:TYR:CZ	1:I:39:LYS:HB3	2.49	0.48
1:I:289:ARG:O	1:I:290:CYS:HB3	2.13	0.48
1:I:399:VAL:HG21	1:I:422:LYS:HB2	1.96	0.47
1:E:20:GLU:H	1:E:20:GLU:CD	2.17	0.47
1:G:129:PRO:HD2	1:G:132:LYS:HB3	1.96	0.47
1:E:573:THR:HB	2:F:196:SER:HB3	1.96	0.47
1:E:409:ASN:H	1:E:409:ASN:HD22	1.62	0.47
1:A:496:ARG:NH1	2:B:192:GLY:O	2.45	0.47
1:E:125:LYS:HD3	1:E:187:ARG:O	2.14	0.47
1:A:50:TRP:CZ2	1:A:54:VAL:HG11	2.49	0.47
1:C:451:ARG:NH1	1:C:468:PRO:HG3	2.29	0.47
1:A:452:ILE:HA	1:A:462:LEU:HD23	1.97	0.47
1:C:537:ILE:HG12	1:C:554:MET:HG3	1.97	0.47
2:F:80:ASN:HD21	2:F:151:SER:HA	1.80	0.47
1:A:408:GLU:H	1:A:408:GLU:CD	2.18	0.47
1:I:575:LYS:HD2	1:I:575:LYS:HA	1.67	0.47
1:I:284:HIS:CE1	1:I:330:ASN:HB3	2.49	0.47
1:C:399:VAL:HG21	1:C:422:LYS:HB2	1.97	0.47
1:E:284:HIS:H	1:E:330:ASN:ND2	2.13	0.47
1:G:237:PHE:HD2	1:G:242:ILE:HG21	1.80	0.47
1:A:575:LYS:HB2	2:B:195:ASN:HA	1.98	0.46
2:B:166:MET:O	2:B:170:THR:HG23	2.15	0.46
1:K:128:PHE:HB3	1:K:132:LYS:HG2	1.96	0.46
1:I:575:LYS:HG2	2:J:196:SER:HB2	1.97	0.46
1:E:48:MET:HG3	1:E:73:TRP:CD2	2.50	0.46
1:C:202:ILE:O	1:C:203:THR:HB	2.16	0.46
1:I:50:TRP:O	1:I:54:VAL:HG22	2.16	0.46
1:C:76:ARG:NH2	1:C:410:GLY:O	2.49	0.46
1:K:20:GLU:CD	1:K:20:GLU:H	2.20	0.46
1:C:302:ILE:HD11	1:C:336:MET:HG3	1.98	0.46
1:G:553:LYS:HA	1:G:571:THR:HA	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:180:GLN:HG3	1:G:185:LEU:HD12	1.98	0.45
1:A:451:ARG:NH1	1:A:468:PRO:HG3	2.31	0.45
1:C:553:LYS:HA	1:C:571:THR:HA	1.98	0.45
2:B:116:GLU:HG3	2:B:116:GLU:H	1.46	0.45
2:L:153:VAL:HG11	2:L:159:LEU:HB2	1.98	0.45
1:A:288:ILE:HD11	1:A:345:VAL:HG13	1.99	0.45
1:E:237:PHE:HD2	1:E:242:ILE:HG21	1.81	0.45
1:C:8:MET:HB3	1:C:32:ILE:HD12	1.97	0.45
1:E:271:PHE:CZ	1:E:350:GLY:HA3	2.51	0.45
1:K:48:MET:HG3	1:K:73:TRP:CD2	2.51	0.45
2:J:238:GLU:HG3	2:J:239:ASN:N	2.32	0.45
1:G:575:LYS:HA	1:G:575:LYS:HD2	1.76	0.45
1:G:400:THR:HG22	1:G:416:LEU:HD22	1.97	0.45
2:H:97:ASN:O	2:H:101:VAL:HG23	2.17	0.45
2:L:236:THR:HG22	2:L:237:VAL:H	1.82	0.45
1:G:128:PHE:HB3	1:G:132:LYS:HG2	1.98	0.45
1:G:575:LYS:HG2	2:H:196:SER:HB2	1.99	0.45
1:K:14:GLU:HB2	1:K:26:ALA:HB3	1.97	0.45
1:I:72:ASN:OD1	1:I:76:ARG:NH2	2.49	0.45
1:A:409:ASN:H	1:A:409:ASN:HD22	1.65	0.44
1:A:268:PRO:HB3	1:A:353:PHE:CE2	2.52	0.44
2:J:93:GLU:HA	2:J:96:ARG:NH1	2.32	0.44
1:K:38:TYR:CE1	1:K:167:LYS:HD2	2.51	0.44
1:G:68:ALA:O	1:G:72:ASN:HB2	2.17	0.44
1:G:29:TYR:CZ	1:G:39:LYS:HB3	2.53	0.44
1:G:550:PHE:HB3	1:G:574:ILE:HD12	2.00	0.44
1:G:20:GLU:CD	1:G:20:GLU:H	2.19	0.44
1:K:259:TYR:CZ	1:K:361:LYS:HD3	2.52	0.44
1:C:456:ASP:O	1:C:458:ASP:N	2.49	0.44
1:K:284:HIS:H	1:K:330:ASN:HD22	1.64	0.44
1:I:469:ASP:HA	1:I:472:LYS:HG3	2.00	0.44
2:J:174:TYR:CE1	2:J:178:LYS:HE2	2.53	0.44
2:D:179:MET:HG2	2:D:216:TYR:CE1	2.53	0.44
2:B:222:ILE:HD11	2:B:250:TYR:HB3	1.99	0.44
1:K:506:MET:HG3	1:K:525:LYS:HB2	2.00	0.44
1:A:397:PRO:HB2	1:A:421:THR:HG23	2.00	0.44
2:B:191:GLU:HG2	2:B:204:ILE:HD13	2.00	0.44
1:K:575:LYS:HG2	2:L:196:SER:HB2	1.99	0.43
2:B:97:ASN:O	2:B:101:VAL:HG23	2.18	0.43
2:H:169:ARG:HG2	2:H:174:TYR:CZ	2.53	0.43
1:G:202:ILE:O	1:G:203:THR:CB	2.66	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:20:GLU:H	1:C:20:GLU:CD	2.20	0.43
1:C:550:PHE:HB3	1:C:574:ILE:HD12	1.99	0.43
1:C:125:LYS:HD3	1:C:187:ARG:O	2.17	0.43
2:J:80:ASN:HD22	2:J:81:ALA:N	2.16	0.43
2:J:219:PHE:CE1	2:J:223:SER:HB2	2.52	0.43
2:D:236:THR:HG22	2:D:237:VAL:H	1.84	0.43
2:H:80:ASN:HB3	2:H:84:VAL:H	1.84	0.43
1:E:183:GLN:HB2	1:E:185:LEU:HG	2.01	0.43
1:A:253:LEU:O	1:A:257:GLN:HG2	2.18	0.43
1:I:291:GLU:HA	1:I:319:SER:OG	2.18	0.43
1:I:456:ASP:O	1:I:458:ASP:N	2.52	0.43
1:A:399:VAL:HG21	1:A:422:LYS:HB2	2.01	0.43
1:K:19:VAL:HG13	1:K:561:VAL:HG11	2.00	0.43
1:K:253:LEU:HD21	1:K:437:ALA:HB1	2.00	0.43
1:K:575:LYS:HA	1:K:575:LYS:HD2	1.85	0.43
1:A:128:PHE:HB3	1:A:132:LYS:HG2	1.99	0.43
1:A:427:THR:N	1:A:428:PRO:CD	2.82	0.43
1:C:308:ARG:HG2	2:D:216:TYR:OH	2.19	0.43
1:K:573:THR:HB	2:L:196:SER:HB3	2.01	0.43
1:I:553:LYS:HA	1:I:571:THR:HA	2.00	0.43
1:C:202:ILE:O	1:C:203:THR:CB	2.67	0.42
1:C:271:PHE:CZ	1:C:275:TYR:HB2	2.54	0.42
1:E:399:VAL:CG2	1:E:422:LYS:HB2	2.45	0.42
1:G:573:THR:HB	2:H:196:SER:HB3	2.00	0.42
1:I:311:LYS:O	1:I:313:ASN:N	2.47	0.42
1:A:102:MET:HA	1:A:119:ILE:O	2.19	0.42
1:C:302:ILE:HD11	1:C:336:MET:SD	2.59	0.42
2:B:225:ILE:HD12	2:B:226:SER:N	2.35	0.42
1:C:48:MET:HG3	1:C:73:TRP:CD2	2.54	0.42
1:E:284:HIS:CE1	1:E:330:ASN:HB3	2.54	0.42
1:K:495:LEU:O	1:K:496:ARG:HB3	2.19	0.42
1:K:553:LYS:HA	1:K:571:THR:HA	2.01	0.42
1:A:439:TYR:O	1:A:443:THR:HB	2.19	0.42
1:A:31:ASN:HB3	1:A:34:ASP:O	2.19	0.42
1:C:510:ASP:O	1:C:512:LYS:N	2.46	0.42
1:A:475:VAL:HA	1:A:483:TRP:O	2.20	0.42
1:I:30:MET:HB2	1:I:170:ILE:HD12	2.01	0.42
1:C:306:ARG:NH2	2:D:220:LEU:O	2.52	0.42
1:C:573:THR:HB	2:D:196:SER:HB3	2.01	0.42
1:K:471:ILE:O	1:K:475:VAL:HG23	2.20	0.42
1:A:302:ILE:HD11	1:A:336:MET:HG3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:191:GLY:HA2	1:I:388:SER:HB3	2.01	0.42
1:G:289:ARG:O	1:G:290:CYS:HB3	2.19	0.42
1:K:289:ARG:O	1:K:290:CYS:HB3	2.19	0.42
1:A:553:LYS:HA	1:A:571:THR:HA	2.01	0.42
1:A:573:THR:HB	2:B:196:SER:HB3	2.01	0.42
2:H:218:LEU:HD13	2:H:222:ILE:HD13	2.00	0.42
1:A:378:ILE:H	1:A:378:ILE:HG12	1.69	0.42
1:K:271:PHE:CZ	1:K:275:TYR:HB2	2.55	0.42
1:A:142:LEU:HD21	2:H:172:PRO:HB2	2.01	0.42
1:C:234:ASN:HB3	1:C:237:PHE:HD1	1.84	0.42
1:E:304:ILE:HG12	1:E:326:LEU:HD21	2.00	0.42
1:A:326:LEU:HD12	1:A:328:LEU:HD11	2.02	0.42
1:K:180:GLN:HG2	1:K:185:LEU:HB2	2.01	0.42
1:I:306:ARG:HH21	2:J:224:GLU:HG2	1.85	0.41
1:C:292:PHE:N	1:C:319:SER:OG	2.49	0.41
1:A:134:ALA:HB1	1:A:141:VAL:HB	2.02	0.41
1:C:128:PHE:HB2	1:C:133:ILE:HG13	2.02	0.41
1:A:306:ARG:HD3	2:B:224:GLU:HB2	2.01	0.41
2:F:236:THR:HG22	2:F:237:VAL:H	1.85	0.41
1:K:24:VAL:HG21	1:K:44:LEU:HD13	2.02	0.41
1:C:284:HIS:H	1:C:330:ASN:ND2	2.13	0.41
1:A:183:GLN:HB2	1:A:185:LEU:HG	2.01	0.41
1:K:227:ARG:HG3	1:K:305:LYS:NZ	2.35	0.41
1:E:128:PHE:HB2	1:E:133:ILE:HG13	2.02	0.41
1:G:93:ILE:HB	1:G:101:TYR:HB2	2.02	0.41
1:I:271:PHE:CZ	1:I:275:TYR:HB2	2.55	0.41
1:A:239:GLU:HG2	1:A:495:LEU:HA	2.01	0.41
1:G:408:GLU:CD	1:G:408:GLU:H	2.23	0.41
1:A:438:ARG:HH11	1:A:438:ARG:HG3	1.85	0.41
2:H:80:ASN:CB	2:H:84:VAL:H	2.33	0.41
1:E:167:LYS:NZ	1:E:171:GLN:HE22	2.19	0.41
1:A:271:PHE:CZ	1:A:275:TYR:HB2	2.55	0.41
2:D:206:GLU:CG	2:D:255:ARG:HH21	2.34	0.41
2:L:93:GLU:HA	2:L:96:ARG:NH1	2.35	0.41
1:G:221:GLU:HA	1:G:224:TYR:CD1	2.54	0.41
1:K:282:PRO:HG2	1:K:355:ALA:HB3	2.01	0.41
1:E:323:ILE:HD12	2:F:143:ASN:HB3	2.02	0.41
2:F:12:UNK:C	2:F:72:ASN:H	2.34	0.41
2:B:191:GLU:HG2	2:B:204:ILE:CD1	2.50	0.41
1:E:93:ILE:HB	1:E:101:TYR:HB2	2.03	0.41
1:C:270:VAL:HG22	1:C:351:LEU:CD2	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:496:ARG:NH1	2:D:192:GLY:O	2.49	0.41
1:E:221:GLU:HG2	1:E:327:TRP:CZ3	2.55	0.41
1:A:128:PHE:HB2	1:A:133:ILE:HG13	2.02	0.41
2:D:213:ASP:OD1	2:D:214:ASP:N	2.53	0.41
1:I:416:LEU:HD23	1:I:416:LEU:HA	1.91	0.41
1:E:15:THR:HG22	1:E:24:VAL:HA	2.03	0.41
1:I:306:ARG:HG2	2:J:224:GLU:OE2	2.21	0.41
1:A:253:LEU:HD21	1:A:437:ALA:HB1	2.03	0.41
1:C:306:ARG:HE	2:D:224:GLU:HB2	1.86	0.41
1:C:292:PHE:HA	1:C:341:ASP:O	2.21	0.41
2:D:219:PHE:CE1	2:D:223:SER:HB2	2.55	0.41
2:J:211:PRO:HA	2:J:212:PRO:HD3	1.91	0.41
2:J:222:ILE:O	2:J:225:ILE:HG13	2.20	0.41
1:A:550:PHE:HB3	1:A:574:ILE:HD12	2.03	0.41
1:E:557:LYS:HA	1:E:558:PRO:HD3	1.93	0.41
1:C:271:PHE:CZ	1:C:350:GLY:HA3	2.56	0.41
2:B:12:UNK:C	2:B:72:ASN:H	2.33	0.41
1:C:249:ASP:OD2	1:C:459:SER:HB3	2.21	0.41
1:G:308:ARG:HG2	2:H:216:TYR:CE1	2.56	0.40
1:E:221:GLU:HA	1:E:224:TYR:CD1	2.55	0.40
1:A:575:LYS:HD2	1:A:575:LYS:HA	1.89	0.40
1:C:128:PHE:HB3	1:C:132:LYS:HG2	2.04	0.40
1:A:262:LEU:HG	1:A:355:ALA:HB1	2.04	0.40
1:C:147:ASP:O	1:C:152:ARG:NH2	2.54	0.40
1:K:302:ILE:HD11	1:K:336:MET:HG3	2.03	0.40
1:A:219:ASP:OD1	1:A:223:ARG:NH2	2.53	0.40
1:K:268:PRO:HB3	1:K:353:PHE:CE2	2.56	0.40
1:G:152:ARG:NH1	1:G:162:GLU:OE2	2.53	0.40
1:E:271:PHE:CZ	1:E:275:TYR:HB2	2.57	0.40
1:C:496:ARG:O	1:C:497:GLN:C	2.60	0.40
1:E:285:ILE:HA	1:E:328:LEU:O	2.21	0.40
1:K:416:LEU:HD23	1:K:416:LEU:HA	1.97	0.40
1:A:221:GLU:HA	1:A:224:TYR:CD1	2.57	0.40
1:G:506:MET:HG3	1:G:525:LYS:HB2	2.03	0.40
1:I:107:LEU:HB2	1:I:115:ILE:HG22	2.03	0.40
1:K:253:LEU:HD22	1:K:458:ASP:HB3	2.03	0.40
1:G:203:THR:HG22	1:G:205:LYS:H	1.86	0.40
1:C:407:LYS:HB2	1:C:411:ALA:O	2.21	0.40
1:G:262:LEU:HG	1:G:355:ALA:HB1	2.02	0.40
2:H:211:PRO:HA	2:H:212:PRO:HD3	1.94	0.40
1:A:300:PRO:HB2	1:A:315:TYR:HB3	2.03	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	568/575 (99%)	547 (96%)	17 (3%)	4 (1%)	26	70
1	C	568/575 (99%)	543 (96%)	21 (4%)	4 (1%)	26	70
1	E	568/575 (99%)	540 (95%)	23 (4%)	5 (1%)	21	64
1	G	568/575 (99%)	549 (97%)	15 (3%)	4 (1%)	26	70
1	I	568/575 (99%)	543 (96%)	20 (4%)	5 (1%)	21	64
1	K	568/575 (99%)	543 (96%)	19 (3%)	6 (1%)	17	58
2	B	156/230 (68%)	149 (96%)	7 (4%)	0	100	100
2	D	156/230 (68%)	148 (95%)	8 (5%)	0	100	100
2	F	156/230 (68%)	149 (96%)	7 (4%)	0	100	100
2	H	156/230 (68%)	149 (96%)	7 (4%)	0	100	100
2	J	156/230 (68%)	148 (95%)	8 (5%)	0	100	100
2	L	156/230 (68%)	148 (95%)	8 (5%)	0	100	100
All	All	4344/4830 (90%)	4156 (96%)	160 (4%)	28 (1%)	30	72

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	310	TYR
1	A	426	TYR
1	C	312	GLY
1	C	511	GLY
1	E	312	GLY
1	E	457	THR
1	G	203	THR
1	I	62	ASN
1	I	312	GLY

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Mol	Chain	Res	Type
1	I	457	THR
1	A	457	THR
1	A	511	GLY
1	C	203	THR
1	E	426	TYR
1	K	62	ASN
1	K	203	THR
1	K	457	THR
1	A	203	THR
1	C	457	THR
1	G	457	THR
1	I	203	THR
1	I	426	TYR
1	K	426	TYR
1	K	496	ARG
1	G	426	TYR
1	E	62	ASN
1	G	62	ASN
1	K	562	PRO

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	501/506 (99%)	478 (95%)	23 (5%)	33	73
1	C	501/506 (99%)	476 (95%)	25 (5%)	30	70
1	E	501/506 (99%)	479 (96%)	22 (4%)	35	74
1	G	501/506 (99%)	480 (96%)	21 (4%)	36	76
1	I	501/506 (99%)	477 (95%)	24 (5%)	31	71
1	K	501/506 (99%)	477 (95%)	24 (5%)	31	71
2	B	147/175 (84%)	139 (95%)	8 (5%)	27	66
2	D	147/175 (84%)	143 (97%)	4 (3%)	52	85
2	F	147/175 (84%)	140 (95%)	7 (5%)	31	71

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	H	147/175 (84%)	138 (94%)	9 (6%)	23 61
2	J	147/175 (84%)	139 (95%)	8 (5%)	27 66
2	L	147/175 (84%)	139 (95%)	8 (5%)	27 66
All	All	3888/4086 (95%)	3705 (95%)	183 (5%)	32 72

All (183) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	THR
1	A	20	GLU
1	A	54	VAL
1	A	82	SER
1	A	113	ARG
1	A	145	ASP
1	A	168	ASN
1	A	186	ASP
1	A	194	SER
1	A	249	ASP
1	A	260	SER
1	A	301	THR
1	A	305	LYS
1	A	326	LEU
1	A	373	THR
1	A	374	SER
1	A	409	ASN
1	A	416	LEU
1	A	463	THR
1	A	470	VAL
1	A	498	LYS
1	A	519	ASP
1	A	523	ASP
2	B	116	GLU
2	B	140	THR
2	B	167	GLU
2	B	202	GLU
2	B	214	ASP
2	B	224	GLU
2	B	236	THR
2	B	238	GLU
1	C	17	THR
1	C	20	GLU

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Mol	Chain	Res	Type
1	C	54	VAL
1	C	72	ASN
1	C	145	ASP
1	C	169	ASP
1	C	186	ASP
1	C	194	SER
1	C	249	ASP
1	C	260	SER
1	C	309	PHE
1	C	310	TYR
1	C	375	GLU
1	C	384	LEU
1	C	390	TYR
1	C	409	ASN
1	C	416	LEU
1	C	419	GLU
1	C	463	THR
1	C	470	VAL
1	C	519	ASP
1	C	523	ASP
1	C	543	PHE
1	C	571	THR
1	C	573	THR
2	D	143	ASN
2	D	216	TYR
2	D	224	GLU
2	D	236	THR
1	E	17	THR
1	E	20	GLU
1	E	54	VAL
1	E	107	LEU
1	E	145	ASP
1	E	168	ASN
1	E	169	ASP
1	E	194	SER
1	E	249	ASP
1	E	260	SER
1	E	310	TYR
1	E	373	THR
1	E	374	SER
1	E	375	GLU
1	E	384	LEU

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Mol	Chain	Res	Type
1	E	409	ASN
1	E	416	LEU
1	E	463	THR
1	E	470	VAL
1	E	519	ASP
1	E	543	PHE
1	E	565	VAL
2	F	80	ASN
2	F	143	ASN
2	F	167	GLU
2	F	189	SER
2	F	214	ASP
2	F	224	GLU
2	F	236	THR
1	G	17	THR
1	G	20	GLU
1	G	45	ASP
1	G	54	VAL
1	G	72	ASN
1	G	113	ARG
1	G	186	ASP
1	G	194	SER
1	G	249	ASP
1	G	260	SER
1	G	276	VAL
1	G	288	ILE
1	G	306	ARG
1	G	317	LYS
1	G	409	ASN
1	G	416	LEU
1	G	434	THR
1	G	519	ASP
1	G	523	ASP
1	G	543	PHE
1	G	573	THR
2	H	80	ASN
2	H	116	GLU
2	H	140	THR
2	H	167	GLU
2	H	184	LEU
2	H	202	GLU
2	H	214	ASP

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Mol	Chain	Res	Type
2	H	235	ASN
2	H	236	THR
1	I	17	THR
1	I	20	GLU
1	I	54	VAL
1	I	76	ARG
1	I	107	LEU
1	I	113	ARG
1	I	194	SER
1	I	218	LEU
1	I	249	ASP
1	I	260	SER
1	I	296	GLU
1	I	306	ARG
1	I	310	TYR
1	I	319	SER
1	I	326	LEU
1	I	346	GLU
1	I	375	GLU
1	I	390	TYR
1	I	396	ASN
1	I	409	ASN
1	I	416	LEU
1	I	423	ASP
1	I	469	ASP
1	I	565	VAL
2	J	80	ASN
2	J	99	LYS
2	J	140	THR
2	J	167	GLU
2	J	189	SER
2	J	214	ASP
2	J	216	TYR
2	J	236	THR
1	K	17	THR
1	K	20	GLU
1	K	54	VAL
1	K	107	LEU
1	K	131	LYS
1	K	145	ASP
1	K	169	ASP
1	K	194	SER

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Mol	Chain	Res	Type
1	K	249	ASP
1	K	260	SER
1	K	296	GLU
1	K	326	LEU
1	K	375	GLU
1	K	384	LEU
1	K	396	ASN
1	K	409	ASN
1	K	416	LEU
1	K	423	ASP
1	K	427	THR
1	K	469	ASP
1	K	470	VAL
1	K	498	LYS
1	K	519	ASP
1	K	543	PHE
2	L	80	ASN
2	L	88	LYS
2	L	103	ARG
2	L	140	THR
2	L	143	ASN
2	L	218	LEU
2	L	224	GLU
2	L	236	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (42) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	171	GLN
1	A	330	ASN
1	A	409	ASN
1	A	485	HIS
2	B	235	ASN
1	C	171	GLN
1	C	330	ASN
1	C	380	GLN
1	C	409	ASN
1	C	485	HIS
2	D	235	ASN
1	E	171	GLN
1	E	303	GLN
1	E	330	ASN

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Mol	Chain	Res	Type
1	E	344	ASN
1	E	409	ASN
1	E	485	HIS
2	F	235	ASN
1	G	171	GLN
1	G	313	ASN
1	G	330	ASN
1	G	380	GLN
1	G	409	ASN
1	G	485	HIS
1	G	497	GLN
2	H	80	ASN
2	H	235	ASN
1	I	171	GLN
1	I	284	HIS
1	I	330	ASN
1	I	409	ASN
1	I	485	HIS
2	J	80	ASN
2	J	183	GLN
2	J	235	ASN
1	K	171	GLN
1	K	284	HIS
1	K	330	ASN
1	K	409	ASN
2	L	80	ASN
2	L	183	GLN
2	L	235	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 16 ligands modelled in this entry, 16 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

There are no such residues in this entry.

5.8 Polymer linkage issues

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	570/575 (99%)	0.07	9 (1%) 74 47	86, 99, 112, 132	0
1	C	570/575 (99%)	0.17	12 (2%) 67 36	84, 99, 112, 158	0
1	E	570/575 (99%)	0.08	13 (2%) 64 33	85, 99, 111, 143	0
1	G	570/575 (99%)	0.09	8 (1%) 78 51	85, 98, 111, 140	0
1	I	570/575 (99%)	0.07	14 (2%) 61 30	86, 99, 111, 148	0
1	K	570/575 (99%)	0.12	16 (2%) 56 27	86, 99, 112, 145	0
2	B	162/230 (70%)	0.21	5 (3%) 52 24	76, 99, 113, 124	0
2	D	162/230 (70%)	0.27	5 (3%) 52 24	83, 98, 114, 125	0
2	F	162/230 (70%)	0.25	5 (3%) 52 24	78, 99, 114, 125	0
2	H	162/230 (70%)	0.18	8 (4%) 33 13	80, 99, 114, 125	0
2	J	162/230 (70%)	0.20	3 (1%) 70 41	82, 98, 114, 125	0
2	L	162/230 (70%)	0.27	7 (4%) 39 16	82, 98, 115, 124	0
All	All	4392/4830 (90%)	0.13	105 (2%) 62 32	76, 99, 112, 158	0

All (105) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	226	SER	5.2
2	J	234	GLY	5.1
2	F	226	SER	4.6
2	B	226	SER	4.5
2	L	71	ALA	4.5
1	I	523	ASP	4.2
1	K	113	ARG	4.2
2	F	234	GLY	4.1
1	A	510	ASP	4.0
1	E	510	ASP	4.0
1	A	507	LYS	3.9

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Mol	Chain	Res	Type	RSRZ
2	D	226	SER	3.8
1	I	507	LYS	3.7
1	I	513	LEU	3.7
1	C	520	ASP	3.7
1	E	112	LYS	3.7
1	C	513	LEU	3.6
1	C	522	THR	3.6
1	G	507	LYS	3.5
1	I	511	GLY	3.5
1	K	511	GLY	3.5
1	A	490	LYS	3.4
1	I	575	LYS	3.4
1	K	510	ASP	3.3
1	K	509	VAL	3.3
1	E	509	VAL	3.3
1	G	510	ASP	3.3
2	L	140	THR	3.2
1	A	112	LYS	3.1
1	A	575	LYS	3.0
2	L	226	SER	3.0
1	I	510	ASP	2.9
1	A	511	GLY	2.9
1	K	507	LYS	2.9
1	K	512	LYS	2.9
1	C	510	ASP	2.9
1	K	508	GLU	2.9
1	A	523	ASP	2.9
1	I	524	ILE	2.8
2	H	259	PHE	2.8
1	C	523	ASP	2.8
1	C	508	GLU	2.8
1	K	523	ASP	2.8
2	D	75	TYR	2.7
1	I	308	ARG	2.7
1	G	511	GLY	2.7
1	C	490	LYS	2.7
1	G	112	LYS	2.7
1	E	520	ASP	2.6
1	K	522	THR	2.6
2	H	78	GLU	2.6
1	G	509	VAL	2.6
1	E	111	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
1	E	523	ASP	2.6
2	F	75	TYR	2.6
2	J	78	GLU	2.6
1	I	522	THR	2.5
2	L	235	ASN	2.5
1	C	502	GLN	2.5
1	E	512	LYS	2.5
2	F	78	GLU	2.5
1	K	205	LYS	2.5
1	A	113	ARG	2.5
1	E	308	ARG	2.5
1	I	512	LYS	2.5
2	H	254	TYR	2.5
2	H	77	PHE	2.5
1	G	490	LYS	2.5
1	E	511	GLY	2.4
1	E	521	TYR	2.4
2	L	224	GLU	2.3
2	B	71	ALA	2.3
1	K	514	VAL	2.3
2	J	226	SER	2.3
1	K	505	TYR	2.3
2	D	218	LEU	2.3
2	L	141	GLY	2.3
2	B	259	PHE	2.3
1	C	519	ASP	2.2
1	E	113	ARG	2.2
1	C	239	GLU	2.2
1	K	406	LEU	2.2
1	E	522	THR	2.2
1	C	512	LYS	2.2
2	H	234	GLY	2.2
2	B	234	GLY	2.2
1	E	239	GLU	2.2
1	I	508	GLU	2.2
2	D	254	TYR	2.2
1	C	507	LYS	2.2
1	A	524	ILE	2.1
2	B	159	LEU	2.1
2	L	234	GLY	2.1
1	I	502	GLN	2.1
2	H	75	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
1	I	110	LYS	2.1
2	F	255	ARG	2.0
1	K	151	GLU	2.0
1	K	409	ASN	2.0
2	H	140	THR	2.0
2	D	77	PHE	2.0
1	G	505	TYR	2.0
1	K	513	LEU	2.0
1	G	523	ASP	2.0
1	I	107	LEU	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å^2)	Q<0.9
3	PB	C	705	1/1	0.77	0.17	-0.81	162,162,162,162	1
3	PB	A	702	1/1	0.66	0.16	-1.68	165,165,165,165	1
3	PB	I	714	1/1	0.81	0.10	-1.79	180,180,180,180	1
3	PB	E	709	1/1	0.80	0.15	-1.98	182,182,182,182	1
3	PB	G	711	1/1	0.94	0.08	-2.12	156,156,156,156	1
3	PB	C	704	1/1	0.91	0.10	-	146,146,146,146	1
3	PB	A	701	1/1	0.94	0.09	-	157,157,157,157	1
3	PB	I	713	1/1	0.88	0.05	-	155,155,155,155	1
3	PB	E	708	1/1	0.95	0.09	-	160,160,160,160	1
3	PB	C	706	1/1	0.82	0.42	-	191,191,191,191	1
3	PB	K	716	1/1	0.74	0.07	-	160,160,160,160	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
3	PB	E	707	1/1	0.73	0.18	-	174,174,174,174	1
3	PB	G	712	1/1	0.97	0.05	-	106,106,106,106	1
3	PB	A	703	1/1	0.98	0.09	-	103,103,103,103	1
3	PB	G	710	1/1	0.92	0.05	-	158,158,158,158	1
3	PB	K	715	1/1	0.78	0.09	-	175,175,175,175	1

6.5 Other polymers

There are no such residues in this entry.