Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:55 AM GMT

PDB ID	:	2JD4
Title	:	MOUSE LAMININ ALPHA1 CHAIN, DOMAINS LG4-5
Authors	:	Harrison, D.; Hussain, S.A.; Combs, A.C.; Ervasti, J.M.; Yurchenco, P.D.; Hohenester, E.
Deposited on	:	2007-01-04
Resolution	:	1.90 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:

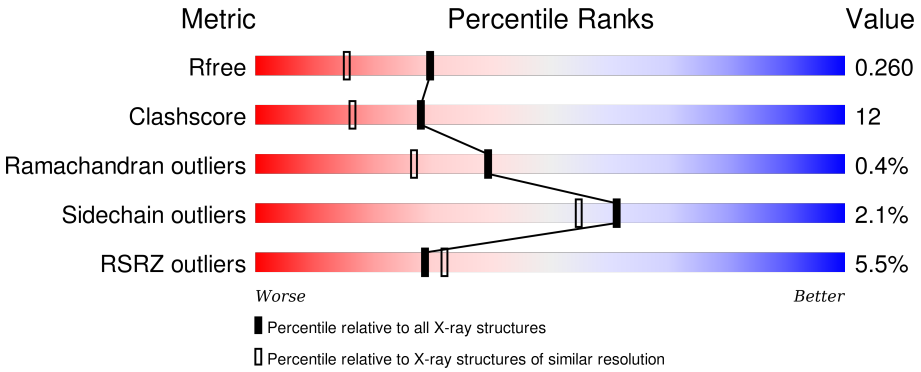
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 1.90 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	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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	383	6% 78% 19%
1	B	383	5% 78% 20%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6115 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 LAMININ SUBUNIT ALPHA-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	376	Total	C	N	O	S	0	0	1
			2890	1814	529	537	10			
1	B	376	Total	C	N	O	S	0	0	1
			2891	1811	533	537	10			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2714	GLN	ASN	ENGINEERED MUTATION	UNP P19137
A	2811	LYS	ASN	ENGINEERED MUTATION	UNP P19137
A	2900	GLN	ASN	ENGINEERED MUTATION	UNP P19137
A	3014	SER	CYS	ENGINEERED MUTATION	UNP P19137
B	2714	GLN	ASN	ENGINEERED MUTATION	UNP P19137
B	2811	LYS	ASN	ENGINEERED MUTATION	UNP P19137
B	2900	GLN	ASN	ENGINEERED MUTATION	UNP P19137
B	3014	SER	CYS	ENGINEERED MUTATION	UNP P19137

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Mg	0	0
			2	2		
2	A	2	Total	Mg	0	0
			2	2		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Cl	0	0
			2	2		
3	A	1	Total	Cl	0	0
			1	1		

- Molecule 4 is water.

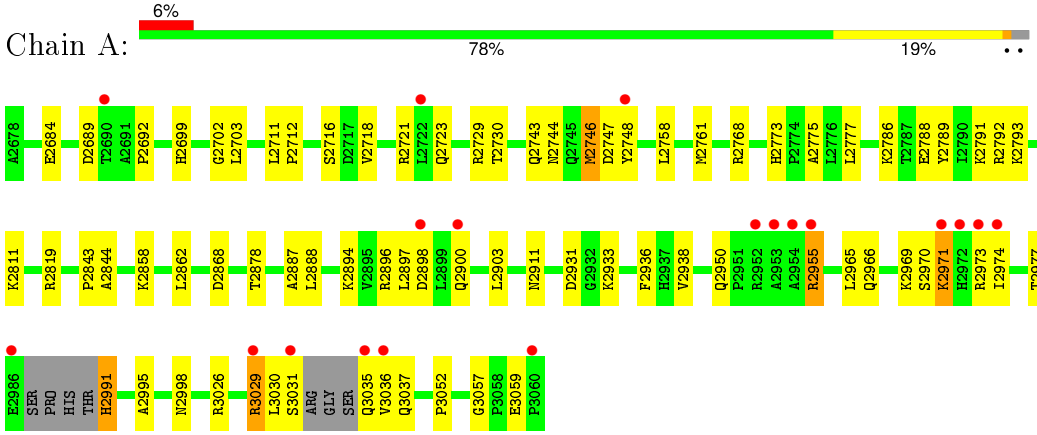
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	161	Total O 161 161	0	0
4	B	166	Total O 166 166	0	0

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.

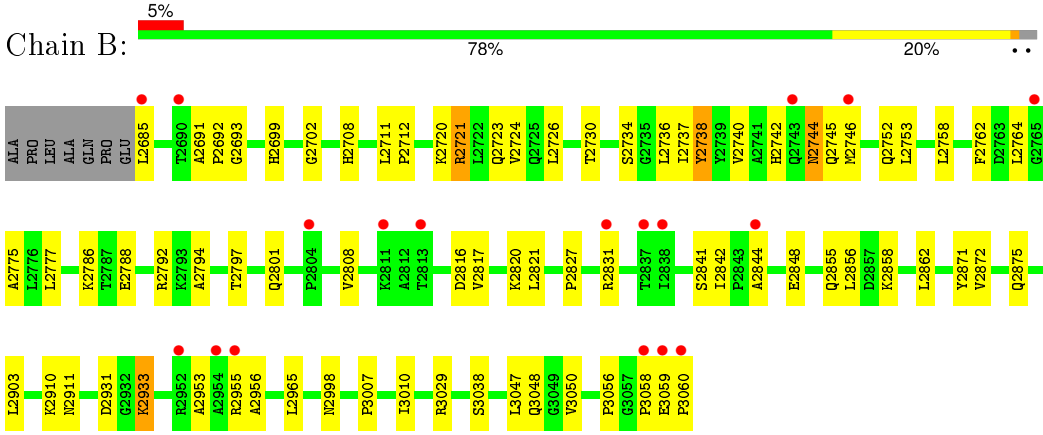
- Molecule 1: LAMININ SUBUNIT ALPHA-1

Chain A:



- Molecule 1: LAMININ SUBUNIT ALPHA-1

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	70.53Å 55.81Å 100.99Å 90.00° 98.48° 90.00°	Depositor
Resolution (Å)	20.00 – 1.90 19.98 – 1.90	Depositor EDS
% Data completeness (in resolution range)	97.0 (20.00-1.90) 97.0 (19.98-1.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.95 (at 1.90Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.230 , 0.262 0.225 , 0.260	Depositor DCC
R_{free} test set	3015 reflections (5.32%)	DCC
Wilson B-factor (Å^2)	24.3	Xtriage
Anisotropy	0.202	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 50.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Outliers	0 of 59669 reflections	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	6115	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.52% 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: MG, CL

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	0.33	0/2946	0.62	0/3987
1	B	0.33	0/2949	0.64	0/3992
All	All	0.33	0/5895	0.63	0/7979

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	2890	0	2887	62	0
1	B	2891	0	2887	73	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	1	0	0	0	0
3	B	2	0	0	0	0
4	A	161	0	0	1	0
4	B	166	0	0	5	0
All	All	6115	0	5774	135	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 12.

All (135) 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:A:2798:VAL:O	1:A:2801:GLN:HG2	1.79	0.83
1:B:2702:GLY:HA2	4:B:2023:HOH:O	1.79	0.81
1:A:3029:ARG:NH2	1:A:3036:VAL:HG11	1.97	0.80
1:A:2791:LYS:HE2	1:A:2791:LYS:HA	1.65	0.79
1:B:2773:HIS:HD2	1:B:2775:ALA:H	1.27	0.78
1:B:2953:ALA:HB3	1:B:2956:ALA:HB2	1.66	0.78
1:B:2742:HIS:HD2	1:B:2744:ASN:H	1.29	0.78
1:A:2729:ARG:HD3	4:A:2002:HOH:O	1.84	0.77
1:B:2786:LYS:HG2	1:B:2797:THR:HB	1.67	0.76
1:A:3029:ARG:HH11	1:A:3029:ARG:HB3	1.54	0.73
1:A:2773:HIS:HD2	1:A:2775:ALA:H	1.38	0.72
1:B:2744:ASN:HD22	1:B:2744:ASN:C	1.92	0.72
1:A:2758:LEU:HD11	1:A:2777:LEU:HD21	1.71	0.72
1:B:2841:SER:HB3	4:B:2023:HOH:O	1.90	0.71
1:A:2712:PRO:HG3	1:A:2862:LEU:HG	1.72	0.70
1:A:2744:ASN:OD1	1:A:2746:MET:HB2	1.92	0.69
1:A:2748:TYR:OH	1:A:2768:ARG:HD3	1.91	0.69
1:B:2699:HIS:ND1	1:B:2858:LYS:HE2	2.09	0.68
1:B:2742:HIS:CD2	1:B:2744:ASN:H	2.12	0.67
1:B:2758:LEU:HG	1:B:2777:LEU:HD13	1.76	0.66
1:B:2786:LYS:CG	1:B:2797:THR:HB	2.26	0.65
1:B:2712:PRO:HG3	1:B:2862:LEU:HD11	1.80	0.64
1:A:2792:ARG:C	1:A:2793:LYS:HD3	2.17	0.63
1:B:2848:GLU:HB2	1:B:2855:GLN:NE2	2.13	0.62
1:B:2712:PRO:HG3	1:B:2862:LEU:CD1	2.30	0.62
1:B:2773:HIS:CD2	1:B:2775:ALA:H	2.16	0.60
1:A:2900:GLN:HB3	1:A:3031:SER:HB2	1.84	0.60
1:A:2703:LEU:HD11	1:A:2868:ASP:HB2	1.84	0.59
1:A:2761:MET:SD	1:A:2768:ARG:HD2	2.42	0.59
1:A:2933:LYS:HD2	1:A:2950:GLN:NE2	2.17	0.59
1:A:2811:LYS:HD3	1:A:2811:LYS:N	2.18	0.58
1:A:2911:ASN:ND2	1:A:2931:ASP:H	2.01	0.58
1:B:2773:HIS:HE1	1:B:2801:GLN:O	1.87	0.57
1:B:2685:LEU:N	1:B:2685:LEU:HD12	2.19	0.57
1:B:2910:LYS:HG3	1:B:2911:ASN:OD1	2.04	0.57
1:A:2703:LEU:CD1	1:A:2868:ASP:HB2	2.36	0.56
1:A:2955:ARG:HG2	1:A:2955:ARG:HH11	1.71	0.55
1:A:2730:THR:HB	1:A:2844:ALA:HB2	1.89	0.55
1:B:2903:LEU:C	1:B:2903:LEU:HD12	2.28	0.54
1:A:2970:SER:OG	1:A:2973:ARG:HB3	2.07	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2744:ASN:ND2	1:B:2746:MET:H	2.05	0.54
1:B:2730:THR:HB	1:B:2844:ALA:HB2	1.90	0.52
1:B:2758:LEU:HG	1:B:2777:LEU:CD1	2.40	0.52
1:B:2898:ASP:HB2	4:B:2096:HOH:O	2.08	0.52
1:A:2773:HIS:CD2	1:A:2775:ALA:H	2.25	0.51
1:A:2699:HIS:HD2	1:A:2858:LYS:NZ	2.08	0.51
1:B:2786:LYS:HG3	4:B:2049:HOH:O	2.10	0.51
1:A:2692:PRO:HG2	1:A:3057:GLY:H	1.75	0.51
1:A:2966:GLN:HB2	1:A:2977:THR:HB	1.91	0.50
1:B:2737:ILE:HD13	1:B:2842:ILE:HG21	1.94	0.50
1:A:2811:LYS:CD	1:A:2811:LYS:N	2.75	0.50
1:A:2938:VAL:HB	1:A:2974:ILE:CD1	2.41	0.49
1:A:2938:VAL:HB	1:A:2974:ILE:HD11	1.94	0.49
1:A:2936:PHE:CE2	1:A:2974:ILE:HG23	2.47	0.49
1:B:2762:PHE:HE1	1:B:2764:LEU:HD21	1.76	0.49
1:B:2762:PHE:CE1	1:B:2764:LEU:HD21	2.47	0.49
1:B:2685:LEU:N	1:B:2955:ARG:HG3	2.28	0.49
1:B:2724:VAL:O	1:B:2786:LYS:HA	2.13	0.49
1:A:3029:ARG:C	1:A:3030:LEU:HD12	2.32	0.49
1:B:2856:LEU:N	1:B:2856:LEU:HD12	2.28	0.49
1:A:2758:LEU:HD11	1:A:2777:LEU:CD2	2.42	0.48
1:B:2730:THR:HB	1:B:2844:ALA:CB	2.44	0.48
1:A:3029:ARG:NH1	1:A:3029:ARG:HB3	2.24	0.48
1:B:2744:ASN:HD21	1:B:2746:MET:HG2	1.78	0.48
1:A:3029:ARG:CZ	1:A:3036:VAL:HG11	2.44	0.47
1:A:2955:ARG:C	1:A:2955:ARG:HD2	2.35	0.47
1:A:2730:THR:HB	1:A:2844:ALA:CB	2.43	0.47
1:A:3035:GLN:OE1	1:A:3035:GLN:HA	2.15	0.47
1:A:3036:VAL:HG12	1:A:3037:GLN:N	2.30	0.47
1:B:3029:ARG:HG2	1:B:3038:SER:HA	1.95	0.47
1:B:2911:ASN:ND2	1:B:2931:ASP:H	2.13	0.47
1:B:2764:LEU:N	1:B:2764:LEU:HD22	2.29	0.47
1:B:3058:PRO:O	1:B:3059:GLU:HB2	2.13	0.47
1:A:2903:LEU:HD12	1:A:2903:LEU:C	2.35	0.47
1:A:2723:GLN:HG3	1:A:2788:GLU:HG2	1.96	0.47
1:B:2744:ASN:C	1:B:2744:ASN:ND2	2.64	0.46
1:B:3007:PRO:HG2	1:B:3010:ILE:CG1	2.45	0.46
1:B:3047:LEU:HD22	1:B:3050:VAL:HB	1.98	0.46
1:A:2878:THR:HG23	1:A:3052:PRO:HA	1.98	0.46
1:A:2811:LYS:H	1:A:2811:LYS:HE2	1.80	0.46
1:A:2743:GLN:HE21	1:A:2743:GLN:HA	1.81	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2716:SER:O	1:A:2819:ARG:NH2	2.44	0.46
1:A:2773:HIS:CE1	1:A:2801:GLN:HG3	2.51	0.45
1:B:2744:ASN:HD22	1:B:2745:GLN:N	2.14	0.45
1:B:2910:LYS:HG2	4:B:2104:HOH:O	2.17	0.45
1:A:2718:VAL:HA	1:A:2721:ARG:O	2.17	0.45
1:A:2894:LYS:HB3	1:A:2896:ARG:HD3	1.99	0.45
1:B:2744:ASN:O	1:B:2745:GLN:HB2	2.17	0.45
1:B:2965:LEU:HD23	1:B:2965:LEU:C	2.37	0.45
1:A:2786:LYS:HD3	1:A:2788:GLU:OE1	2.16	0.44
1:A:2897:LEU:N	1:A:2897:LEU:HD12	2.31	0.44
1:B:2875:GLN:OE1	1:B:3059:GLU:HB3	2.17	0.44
1:B:2699:HIS:CE1	1:B:2858:LYS:HE2	2.53	0.44
1:B:2736:LEU:HD13	1:B:2752:GLN:HG2	1.98	0.44
1:A:2991:HIS:CD2	1:A:2991:HIS:N	2.86	0.44
1:B:2831:ARG:HH11	1:B:2831:ARG:HG3	1.83	0.44
1:B:2711:LEU:HA	1:B:2712:PRO:HD3	1.86	0.44
1:A:2965:LEU:C	1:A:2965:LEU:HD23	2.38	0.44
1:B:2742:HIS:CD2	1:B:2744:ASN:ND2	2.86	0.43
1:B:2821:LEU:C	1:B:2821:LEU:HD23	2.39	0.43
1:B:2734:SER:HA	1:B:2753:LEU:O	2.18	0.43
1:A:2793:LYS:N	1:A:2793:LYS:HD3	2.34	0.43
1:B:2884:GLY:HA3	1:B:3048:GLN:O	2.19	0.43
1:A:2703:LEU:HD11	1:A:2868:ASP:CB	2.49	0.43
1:B:2933:LYS:HE2	1:B:2933:LYS:HB2	1.79	0.43
1:B:2820:LYS:O	1:B:2820:LYS:HG3	2.18	0.43
1:B:2708:HIS:NE2	1:B:2827:PRO:HG3	2.34	0.43
1:B:2871:TYR:CE1	1:B:3056:PRO:HD3	2.54	0.43
1:A:2684:GLU:OE1	1:A:2955:ARG:HG3	2.19	0.42
1:B:2736:LEU:HD12	1:B:2737:ILE:N	2.34	0.42
1:A:2969:LYS:HE3	1:A:2995:ALA:HB2	2.01	0.42
1:A:2970:SER:O	1:A:2971:LYS:C	2.57	0.42
1:B:2720:LYS:O	1:B:2721:ARG:HB2	2.18	0.42
1:A:2791:LYS:HE2	1:A:2791:LYS:CA	2.43	0.42
1:A:2898:ASP:HA	1:A:2969:LYS:O	2.19	0.42
1:B:2744:ASN:HD22	1:B:2746:MET:H	1.66	0.42
1:B:2726:LEU:C	1:B:2726:LEU:HD12	2.40	0.42
1:B:2773:HIS:HD2	1:B:2775:ALA:N	2.07	0.42
1:B:2699:HIS:ND1	1:B:2858:LYS:CE	2.80	0.41
1:B:3059:GLU:HG3	1:B:3060:PRO:N	2.35	0.41
1:B:2792:ARG:C	1:B:2808:VAL:HG23	2.39	0.41
1:B:2693:GLY:C	1:B:2872:VAL:HG23	2.40	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2740:VAL:CG1	1:B:2817:VAL:HG21	2.50	0.41
1:B:2740:VAL:HG12	1:B:2817:VAL:HG21	2.02	0.41
1:A:2703:LEU:O	1:A:3052:PRO:HG2	2.19	0.41
1:A:2702:GLY:HA3	1:A:2843:PRO:HA	2.03	0.41
1:A:2773:HIS:HE1	1:A:2801:GLN:O	2.03	0.41
1:A:2887:ALA:C	1:A:2888:LEU:HD12	2.41	0.41
1:B:2723:GLN:O	1:B:2723:GLN:HG3	2.21	0.41
1:B:2738:TYR:CD1	1:B:2738:TYR:C	2.94	0.41
1:B:2788:GLU:O	1:B:2794:ALA:HA	2.21	0.41
1:B:2742:HIS:HB2	1:B:2816:ASP:O	2.21	0.40
1:B:2691:ALA:HA	1:B:2692:PRO:HD3	1.89	0.40
1:A:2711:LEU:HA	1:A:2712:PRO:HD3	1.86	0.40
1:B:2842:ILE:O	1:B:2842:ILE:HG23	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	370/383 (97%)	352 (95%)	16 (4%)	2 (0%)	34	21
1	B	374/383 (98%)	353 (94%)	20 (5%)	1 (0%)	46	35
All	All	744/766 (97%)	705 (95%)	36 (5%)	3 (0%)	39	27

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2971	LYS
1	A	3059	GLU
1	B	2721	ARG

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	311/318 (98%)	302 (97%)	9 (3%)	50 40
1	B	312/318 (98%)	308 (99%)	4 (1%)	76 73
All	All	623/636 (98%)	610 (98%)	13 (2%)	61 55

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2689	ASP
1	A	2746	MET
1	A	2747	ASP
1	A	2789	TYR
1	A	2955	ARG
1	A	2991	HIS
1	A	2998	ASN
1	A	3026	ARG
1	A	3029	ARG
1	B	2738	TYR
1	B	2744	ASN
1	B	2933	LYS
1	B	2998	ASN

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

Mol	Chain	Res	Type
1	A	2682	GLN
1	A	2699	HIS
1	A	2700	GLN
1	A	2705	GLN
1	A	2743	GLN
1	A	2773	HIS
1	A	2801	GLN
1	A	2855	GLN
1	A	2911	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	2950	GLN
1	A	2991	HIS
1	A	2998	ASN
1	B	2700	GLN
1	B	2705	GLN
1	B	2742	HIS
1	B	2744	ASN
1	B	2754	GLN
1	B	2773	HIS
1	B	2855	GLN
1	B	2911	ASN
1	B	2950	GLN
1	B	2991	HIS
1	B	2998	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 7 ligands modelled in this entry, 7 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	376/383 (98%)	0.37	22 (5%)	26	29	14, 27, 47, 61	0
1	B	376/383 (98%)	0.34	19 (5%)	32	35	14, 25, 41, 62	0
All	All	752/766 (98%)	0.35	41 (5%)	29	32	14, 26, 43, 62	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	3060	PRO	8.5
1	B	2954	ALA	4.6
1	A	2972	HIS	4.6
1	A	3060	PRO	4.6
1	A	2952	ARG	4.4
1	B	3059	GLU	4.3
1	A	2690	THR	4.2
1	A	2973	ARG	4.0
1	A	3035	GLN	3.9
1	B	2685	LEU	3.5
1	A	3029	ARG	3.5
1	B	2955	ARG	3.5
1	A	3036	VAL	3.4
1	B	2837	THR	3.2
1	A	2955	ARG	3.1
1	A	2974	ILE	2.9
1	B	2743	GLN	2.7
1	B	2804	PRO	2.7
1	A	2953	ALA	2.7
1	B	2811	LYS	2.7
1	B	2813	THR	2.7
1	A	2984	ARG	2.6
1	B	2746	MET	2.6
1	A	2985	ALA	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	2971	LYS	2.6
1	A	2954	ALA	2.6
1	A	2986	GLU	2.6
1	A	2898	ASP	2.5
1	A	3031	SER	2.5
1	B	2831	ARG	2.5
1	B	2690	THR	2.5
1	B	2838	ILE	2.4
1	B	2766	LYS	2.4
1	A	2983	VAL	2.2
1	A	2900	GLN	2.2
1	B	2765	GLY	2.1
1	A	2722	LEU	2.1
1	B	2952	ARG	2.0
1	B	2844	ALA	2.0
1	A	2748	TYR	2.0
1	B	3058	PRO	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	CL	A	4063	1/1	0.94	0.12	0.13	47,47,47,47	0
3	CL	B	4063	1/1	0.96	0.09	0.04	21,21,21,21	0
2	MG	B	4062	1/1	0.96	0.06	-1.66	22,22,22,22	0
2	MG	A	4061	1/1	0.93	0.06	-1.76	30,30,30,30	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
2	MG	A	4062	1/1	0.92	0.06	-1.81	26,26,26,26	0
2	MG	B	4061	1/1	0.93	0.06	-2.54	35,35,35,35	0
3	CL	B	4064	1/1	0.98	0.05	-	24,24,24,24	0

6.5 Other polymers (i)

There are no such residues in this entry.