Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 06:32 PM GMT

PDB ID	:	1BAO
Title	:	THE CONTRIBUTION OF BURIED HYDROGEN BONDS TO PROTEIN STABILITY: THE CRYSTAL STRUCTURES OF TWO BARNASE MUTANTS
Authors	:	Chen, Y.W.; Fersht, A.R.; Henrick, K.
Deposited on	:	1993-05-19
Resolution	:	2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

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

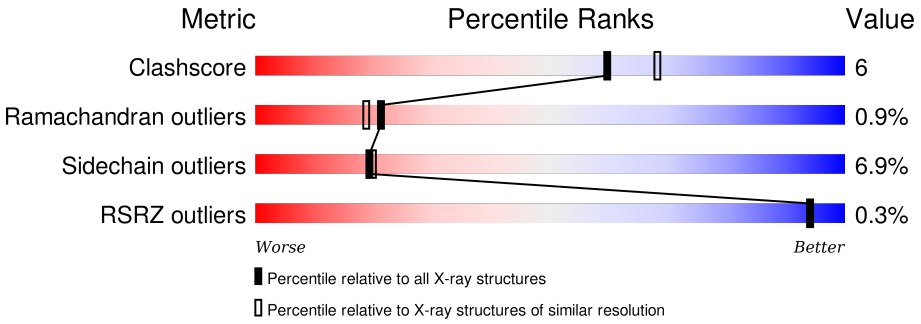
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 2.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	110	55% 33% 11%
1	B	110	% 59% 35%
1	C	110	66% 26% 5%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2752 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 BARNASE.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	108	Total	C	N	O	0	0	0
			853	541	148	164			
1	B	108	Total	C	N	O	0	0	1
			842	535	144	163			
1	C	107	Total	C	N	O	0	0	0
			843	535	145	163			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	78	PHE	TYR	CONFLICT	UNP P00648
B	78	PHE	TYR	CONFLICT	UNP P00648
C	78	PHE	TYR	CONFLICT	UNP P00648

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	1	Total	Zn	0	0
			1	1		

- Molecule 3 is water.

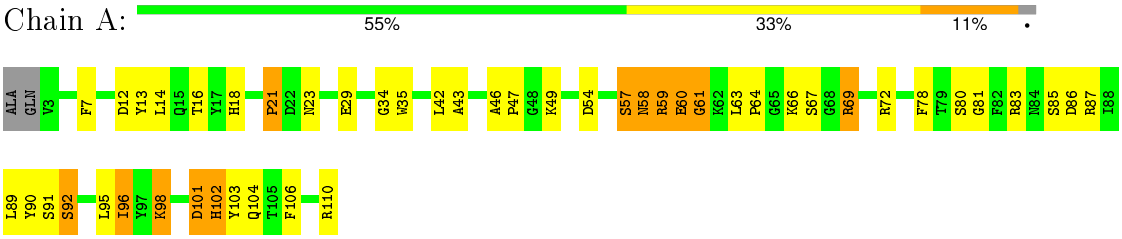
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	72	Total	O	0	0
			72	72		
3	B	76	Total	O	0	0
			76	76		
3	C	65	Total	O	0	0
			65	65		

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.

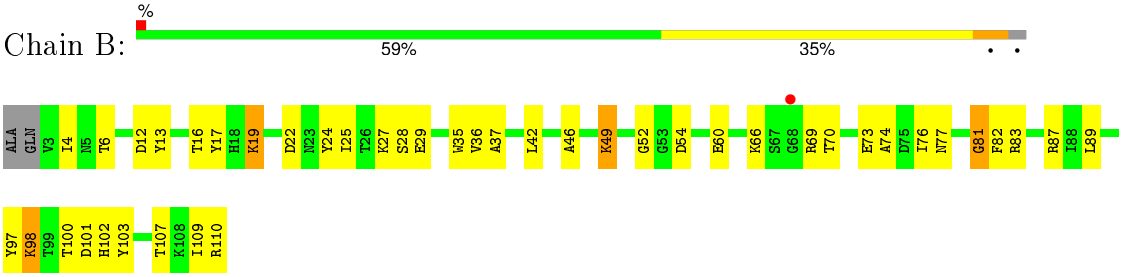
- Molecule 1: BARNASE

Chain A:



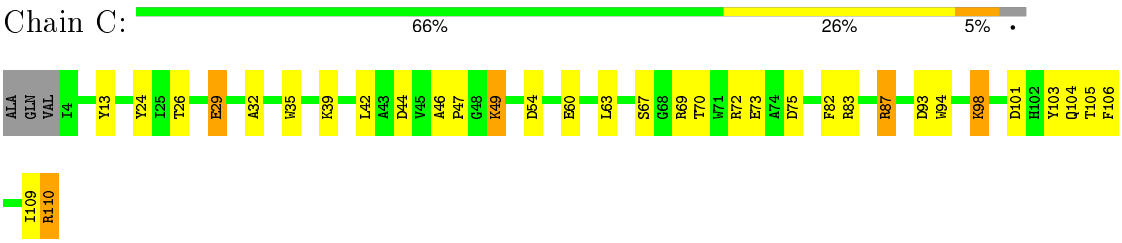
- Molecule 1: BARNASE

Chain B:



- Molecule 1: BARNASE

Chain C:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α, β, γ	58.81Å 58.81Å 81.90Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	6.00 – 2.20 13.92 – 2.20	Depositor EDS
% Data completeness (in resolution range)	(Not available) (6.00-2.20) 93.0 (13.92-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.35 (at 2.20Å)	Xtriage
Refinement program	PROLSQ, X-PLOR	Depositor
R, R_{free}	0.152 , (Not available) 0.145 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å^2)	16.9	Xtriage
Anisotropy	0.197	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 85.0	EDS
Estimated twinning fraction	0.006 for -h,-k,l 0.057 for h,-h-k,-l 0.028 for -k,-h,-l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Outliers	0 of 14908 reflections	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	2752	wwPDB-VP
Average B, all atoms (Å^2)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.03% 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: ZN

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	1.09	1/874 (0.1%)	2.45	46/1182 (3.9%)
1	B	1.08	0/863	2.20	38/1168 (3.3%)
1	C	1.08	0/864	2.48	37/1170 (3.2%)
All	All	1.08	1/2601 (0.0%)	2.38	121/3520 (3.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	29	GLU	CD-OE1	-5.16	1.20	1.25

All (121) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	110	ARG	NE-CZ-NH2	29.07	134.83	120.30
1	A	87	ARG	NE-CZ-NH2	19.52	130.06	120.30
1	A	72	ARG	NE-CZ-NH1	19.14	129.87	120.30
1	C	87	ARG	NE-CZ-NH2	17.83	129.22	120.30
1	C	72	ARG	NE-CZ-NH1	15.25	127.92	120.30
1	C	72	ARG	NE-CZ-NH2	-14.10	113.25	120.30
1	C	24	TYR	CB-CG-CD2	-13.30	113.02	121.00
1	A	72	ARG	NE-CZ-NH2	-12.79	113.90	120.30
1	B	87	ARG	NE-CZ-NH1	12.51	126.56	120.30
1	A	54	ASP	CB-CG-OD2	11.94	129.04	118.30
1	C	83	ARG	NE-CZ-NH1	11.81	126.21	120.30
1	B	83	ARG	NE-CZ-NH2	10.68	125.64	120.30
1	A	110	ARG	CD-NE-CZ	10.50	138.30	123.60
1	B	69	ARG	NE-CZ-NH1	10.46	125.53	120.30
1	B	69	ARG	NE-CZ-NH2	-10.01	115.29	120.30
1	C	44	ASP	CB-CG-OD2	10.01	127.31	118.30
1	C	110	ARG	NH1-CZ-NH2	-9.58	108.86	119.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	12	ASP	CB-CG-OD2	-9.53	109.72	118.30
1	A	60	GLU	CG-CD-OE2	9.44	137.17	118.30
1	C	83	ARG	NH1-CZ-NH2	-9.42	109.03	119.40
1	A	59	ARG	NE-CZ-NH1	9.32	124.96	120.30
1	C	13	TYR	CB-CG-CD1	9.30	126.58	121.00
1	A	87	ARG	NE-CZ-NH1	-9.14	115.73	120.30
1	C	83	ARG	NE-CZ-NH2	8.87	124.73	120.30
1	C	24	TYR	CB-CG-CD1	8.76	126.25	121.00
1	A	90	TYR	CB-CG-CD1	8.60	126.16	121.00
1	B	29	GLU	OE1-CD-OE2	-8.59	112.99	123.30
1	A	59	ARG	CD-NE-CZ	8.07	134.90	123.60
1	B	37	ALA	N-CA-CB	8.05	121.37	110.10
1	A	86	ASP	CB-CG-OD1	8.01	125.51	118.30
1	C	110	ARG	NE-CZ-NH1	-7.98	116.31	120.30
1	B	24	TYR	CB-CG-CD2	7.73	125.64	121.00
1	A	103	TYR	CB-CG-CD2	7.70	125.62	121.00
1	A	13	TYR	CB-CG-CD2	-7.68	116.39	121.00
1	A	69	ARG	NE-CZ-NH1	7.60	124.10	120.30
1	A	110	ARG	NE-CZ-NH2	-7.55	116.53	120.30
1	A	59	ARG	NE-CZ-NH2	-7.45	116.57	120.30
1	B	110	ARG	NE-CZ-NH2	-7.45	116.58	120.30
1	A	60	GLU	OE1-CD-OE2	-7.44	114.37	123.30
1	C	101	ASP	CB-CG-OD2	7.26	124.83	118.30
1	B	74	ALA	CB-CA-C	7.25	120.98	110.10
1	C	13	TYR	CB-CG-CD2	-7.11	116.73	121.00
1	A	103	TYR	CB-CG-CD1	-7.05	116.77	121.00
1	C	93	ASP	CB-CG-OD1	6.98	124.58	118.30
1	A	90	TYR	CB-CG-CD2	-6.79	116.93	121.00
1	C	32	ALA	N-CA-CB	6.77	119.58	110.10
1	B	82	PHE	CB-CG-CD2	6.74	125.52	120.80
1	A	60	GLU	CA-CB-CG	6.74	128.22	113.40
1	B	83	ARG	NH1-CZ-NH2	-6.71	112.01	119.40
1	B	97	TYR	CB-CG-CD1	6.58	124.95	121.00
1	A	106	PHE	CB-CG-CD1	6.57	125.40	120.80
1	C	63	LEU	CA-CB-CG	6.54	130.34	115.30
1	A	61	GLY	C-N-CA	6.46	137.86	121.70
1	A	66	LYS	N-CA-CB	6.36	122.05	110.60
1	B	100	THR	CA-CB-CG2	6.31	121.23	112.40
1	C	82	PHE	CB-CG-CD2	-6.30	116.39	120.80
1	C	106	PHE	CB-CG-CD1	-6.27	116.41	120.80
1	C	60	GLU	OE1-CD-OE2	-6.26	115.78	123.30
1	B	54	ASP	CB-CG-OD2	6.24	123.91	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	81	GLY	CA-C-O	6.14	131.66	120.60
1	A	54	ASP	OD1-CG-OD2	-6.08	111.74	123.30
1	B	17	TYR	CB-CG-CD2	-6.03	117.38	121.00
1	C	75	ASP	CB-CG-OD1	6.03	123.72	118.30
1	A	96	ILE	CB-CG1-CD1	5.97	130.62	113.90
1	B	6	THR	O-C-N	5.96	132.23	122.70
1	C	87	ARG	NH1-CZ-NH2	-5.91	112.90	119.40
1	A	106	PHE	CB-CG-CD2	-5.89	116.68	120.80
1	A	81	GLY	O-C-N	-5.88	113.28	122.70
1	A	57	SER	CA-C-N	5.86	130.09	117.20
1	B	12	ASP	CB-CG-OD1	5.85	123.57	118.30
1	B	69	ARG	N-CA-CB	5.84	121.11	110.60
1	B	52	GLY	C-N-CA	5.82	134.51	122.30
1	C	54	ASP	CB-CG-OD1	-5.80	113.08	118.30
1	A	13	TYR	O-C-N	-5.80	113.42	122.70
1	A	34	GLY	CA-C-O	-5.80	110.16	120.60
1	A	87	ARG	CD-NE-CZ	5.77	131.68	123.60
1	A	101	ASP	CB-CG-OD1	5.76	123.48	118.30
1	B	35	TRP	CG-CD2-CE3	-5.71	128.76	133.90
1	B	70	THR	CB-CA-C	5.70	126.98	111.60
1	B	22	ASP	CB-CG-OD1	5.63	123.37	118.30
1	A	80	SER	CA-CB-OG	-5.60	96.08	111.20
1	C	94	TRP	N-CA-C	5.60	126.11	111.00
1	C	39	LYS	CG-CD-CE	5.57	128.62	111.90
1	B	66	LYS	N-CA-CB	5.55	120.59	110.60
1	A	95	LEU	CA-CB-CG	5.54	128.04	115.30
1	B	87	ARG	NH1-CZ-NH2	-5.53	113.32	119.40
1	C	29	GLU	OE1-CD-OE2	-5.53	116.67	123.30
1	A	83	ARG	NE-CZ-NH2	-5.51	117.55	120.30
1	A	85	SER	N-CA-CB	-5.46	102.32	110.50
1	C	69	ARG	NE-CZ-NH1	5.44	123.02	120.30
1	B	36	VAL	CA-CB-CG2	5.42	119.03	110.90
1	B	107	THR	CA-CB-CG2	5.40	119.96	112.40
1	A	29	GLU	CG-CD-OE1	5.38	129.06	118.30
1	A	102	HIS	O-C-N	-5.33	114.18	122.70
1	B	81	GLY	O-C-N	5.29	131.16	122.70
1	C	49	LYS	CA-CB-CG	5.27	125.00	113.40
1	C	110	ARG	N-CA-CB	5.26	120.06	110.60
1	A	12	ASP	CB-CG-OD2	-5.25	113.57	118.30
1	C	44	ASP	CB-CG-OD1	-5.25	113.58	118.30
1	C	98	LYS	CA-C-O	-5.24	109.09	120.10
1	A	69	ARG	NE-CZ-NH2	-5.23	117.68	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	42	LEU	CB-CG-CD2	-5.23	102.11	111.00
1	B	97	TYR	CB-CG-CD2	-5.22	117.87	121.00
1	A	35	TRP	CE3-CZ3-CH2	-5.19	115.49	121.20
1	B	17	TYR	CA-CB-CG	-5.19	103.53	113.40
1	B	16	THR	O-C-N	5.16	130.96	122.70
1	A	13	TYR	CD1-CE1-CZ	-5.16	115.16	119.80
1	A	21	PRO	O-C-N	-5.16	114.44	122.70
1	B	4	ILE	CA-C-N	5.16	128.55	117.20
1	A	16	THR	CA-CB-OG1	-5.16	98.17	109.00
1	B	28	SER	N-CA-CB	-5.15	102.78	110.50
1	C	105	THR	C-N-CA	5.13	134.53	121.70
1	B	24	TYR	CG-CD1-CE1	5.11	125.39	121.30
1	A	60	GLU	CB-CG-CD	5.10	127.98	114.20
1	B	66	LYS	CB-CG-CD	5.10	124.85	111.60
1	C	29	GLU	CB-CG-CD	5.08	127.92	114.20
1	B	28	SER	CA-C-O	5.07	130.75	120.10
1	C	104	GLN	C-N-CA	5.06	134.35	121.70
1	C	72	ARG	CA-CB-CG	5.04	124.50	113.40
1	C	32	ALA	CB-CA-C	-5.04	102.54	110.10
1	B	110	ARG	N-CA-CB	5.03	119.65	110.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	853	0	809	14	0
1	B	842	0	792	9	0
1	C	843	0	800	6	0
2	C	1	0	0	0	0
3	A	72	0	0	1	0
3	B	76	0	0	0	0
3	C	65	0	0	0	0
All	All	2752	0	2401	28	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 6.

All (28) 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:63:LEU:HD11	1:A:89:LEU:HG	1.75	0.68
1:B:76:ILE:O	1:B:77:ASN:HB2	2.05	0.57
1:B:46:ALA:O	1:B:49:LYS:HB2	2.09	0.52
1:C:26:THR:OG1	1:C:29:GLU:HG3	2.11	0.51
1:A:7:PHE:CE1	1:A:98:LYS:HB2	2.47	0.50
1:A:42:LEU:HD23	1:A:78:PHE:CE2	2.47	0.50
1:B:101:ASP:O	1:B:102:HIS:HB2	2.10	0.49
1:B:73:GLU:HB3	1:B:89:LEU:HD13	1.94	0.49
1:A:46:ALA:N	1:A:47:PRO:HD3	2.30	0.46
1:A:57:SER:O	1:A:58:ASN:CB	2.65	0.45
1:C:87:ARG:HG3	1:C:103:TYR:CE1	2.51	0.44
1:B:25:ILE:HG12	1:B:49:LYS:HG3	1.99	0.44
1:A:46:ALA:N	1:A:47:PRO:CD	2.80	0.44
1:C:46:ALA:N	1:C:47:PRO:HD3	2.33	0.44
1:B:81:GLY:HA2	1:C:47:PRO:HB2	1.99	0.43
1:B:13:TYR:HE1	1:B:19:LYS:HD2	1.82	0.43
1:B:98:LYS:HG2	1:B:109:ILE:CG2	2.48	0.43
1:A:63:LEU:HA	1:A:64:PRO:HD3	1.84	0.43
1:A:43:ALA:HA	3:A:142:HOH:O	2.19	0.43
1:A:69:ARG:HG3	1:A:91:SER:HB2	2.02	0.42
1:C:109:ILE:O	1:C:110:ARG:HD3	2.20	0.42
1:A:57:SER:O	1:A:58:ASN:HB3	2.20	0.41
1:A:21:PRO:HB2	1:A:23:ASN:OD1	2.21	0.41
1:B:98:LYS:HG2	1:B:109:ILE:HG21	2.02	0.41
1:A:101:ASP:O	1:A:104:GLN:HG3	2.20	0.41
1:C:35:TRP:CD1	1:C:42:LEU:HB2	2.56	0.41
1:A:69:ARG:HA	1:A:92:SER:OG	2.21	0.40
1:A:14:LEU:O	1:A:18:HIS:HA	2.21	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	106/110 (96%)	98 (92%)	6 (6%)	2 (2%)	10 6
1	B	106/110 (96%)	98 (92%)	7 (7%)	1 (1%)	21 19
1	C	105/110 (96%)	101 (96%)	4 (4%)	0	100 100
All	All	317/330 (96%)	297 (94%)	17 (5%)	3 (1%)	21 19

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	58	ASN
1	B	60	GLU
1	A	61	GLY

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	87/92 (95%)	79 (91%)	8 (9%)	11 11
1	B	85/92 (92%)	80 (94%)	5 (6%)	24 27
1	C	87/92 (95%)	82 (94%)	5 (6%)	25 29
All	All	259/276 (94%)	241 (93%)	18 (7%)	19 20

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	LYS
1	A	59	ARG
1	A	60	GLU
1	A	67	SER
1	A	92	SER
1	A	96	ILE

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Mol	Chain	Res	Type
1	A	98	LYS
1	A	102	HIS
1	B	19	LYS
1	B	27	LYS
1	B	49	LYS
1	B	98	LYS
1	B	103	TYR
1	C	49	LYS
1	C	67	SER
1	C	70	THR
1	C	73	GLU
1	C	98	LYS

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

Mol	Chain	Res	Type
1	A	18	HIS
1	B	102	HIS

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 1 ligands modelled in this entry, 1 is 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	108/110 (98%)	-0.67	0 100 100	2, 14, 28, 41	0
1	B	108/110 (98%)	-0.75	1 (0%) 85 85	3, 13, 27, 35	0
1	C	107/110 (97%)	-0.77	0 100 100	5, 13, 23, 30	0
All	All	323/330 (97%)	-0.73	1 (0%) 94 94	2, 13, 27, 41	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	68	GLY	2.1

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

There are no carbohydrates in this entry.

6.4 Ligands

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
2	ZN	C	111	1/1	1.00	0.03	-5.74	18,18,18,18	0

6.5 Other polymers (i)

There are no such residues in this entry.