Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 10:22 PM GMT

PDB ID	:	1T6U
Title	:	Nickel Superoxide Dismutase (NiSOD) Native 1.30 A Structure
Authors	:	Barondeau, D.P.; Kassmann, C.J.; Bruns, C.K.; Tainer, J.A.; Getzoff, E.D.
Deposited on	:	2004-05-07
Resolution	:	1.30 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:

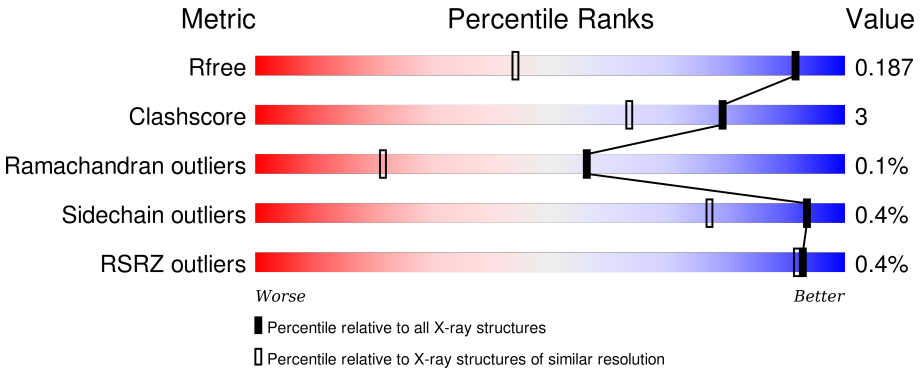
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 1.30 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	1475 (1.34-1.26)
Clashscore	102246	1031 (1.32-1.28)
Ramachandran outliers	100387	1504 (1.34-1.26)
Sidechain outliers	100360	1503 (1.34-1.26)
RSRZ outliers	91569	1476 (1.34-1.26)

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	117	2% 91% 5%
1	B	117	90% 10%
1	C	117	2% 91% 9%
1	D	117	84% 14%
1	E	117	93% 5%

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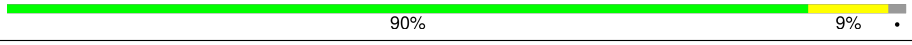
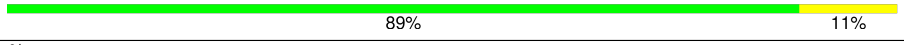
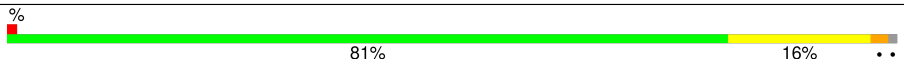
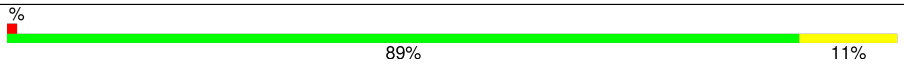
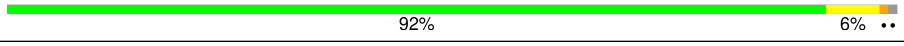
Mol	Chain	Length	Quality of chain
1	F	117	
1	G	117	
1	H	117	
1	I	117	
1	J	117	
1	K	117	
1	L	117	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13192 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 Superoxide dismutase [Ni].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	117	Total	C	N	O	S	0	2	0
			944	600	165	175	4			
1	B	117	Total	C	N	O	S	0	1	0
			940	596	165	175	4			
1	C	117	Total	C	N	O	S	0	2	0
			943	598	165	176	4			
1	D	116	Total	C	N	O	S	0	1	0
			934	593	164	173	4			
1	E	116	Total	C	N	O	S	0	2	0
			938	595	164	175	4			
1	F	115	Total	C	N	O	S	0	1	0
			925	587	162	172	4			
1	G	117	Total	C	N	O	S	0	1	0
			940	596	165	175	4			
1	H	116	Total	C	N	O	S	0	1	0
			934	593	164	173	4			
1	I	117	Total	C	N	O	S	0	2	0
			944	600	165	175	4			
1	J	117	Total	C	N	O	S	0	1	0
			940	596	165	175	4			
1	K	116	Total	C	N	O	S	0	2	0
			938	597	164	173	4			
1	L	116	Total	C	N	O	S	0	1	0
			934	593	164	173	4			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	85	MET	LEU	ENGINEERED	UNP P80735
B	85	MET	LEU	ENGINEERED	UNP P80735
C	85	MET	LEU	ENGINEERED	UNP P80735
D	85	MET	LEU	ENGINEERED	UNP P80735
E	85	MET	LEU	ENGINEERED	UNP P80735

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Chain	Residue	Modelled	Actual	Comment	Reference
F	85	MET	LEU	ENGINEERED	UNP P80735
G	85	MET	LEU	ENGINEERED	UNP P80735
H	85	MET	LEU	ENGINEERED	UNP P80735
I	85	MET	LEU	ENGINEERED	UNP P80735
J	85	MET	LEU	ENGINEERED	UNP P80735
K	85	MET	LEU	ENGINEERED	UNP P80735
L	85	MET	LEU	ENGINEERED	UNP P80735

- Molecule 2 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	155	Total O 155 155	0	0
3	B	169	Total O 169 169	0	0

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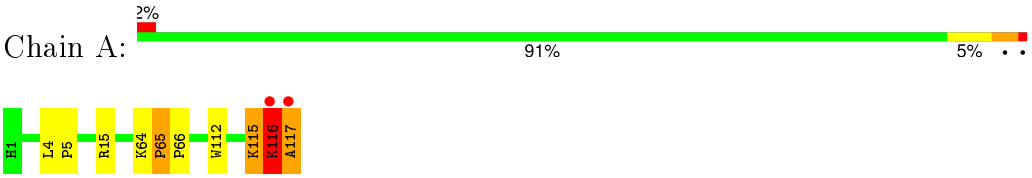
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	C	163	Total O 163 163	0	0
3	D	164	Total O 164 164	0	0
3	E	175	Total O 175 175	0	0
3	F	158	Total O 158 158	0	0
3	G	162	Total O 162 162	0	0
3	H	154	Total O 154 154	0	0
3	I	154	Total O 154 154	0	0
3	J	173	Total O 173 173	0	0
3	K	142	Total O 142 142	0	0
3	L	157	Total O 157 157	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.

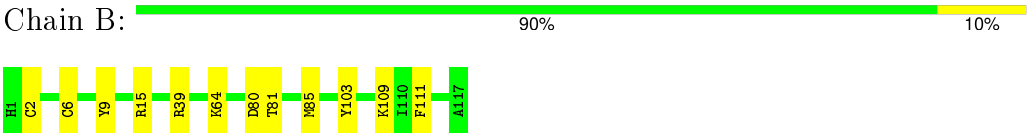
- Molecule 1: Superoxide dismutase [Ni]

Chain A:



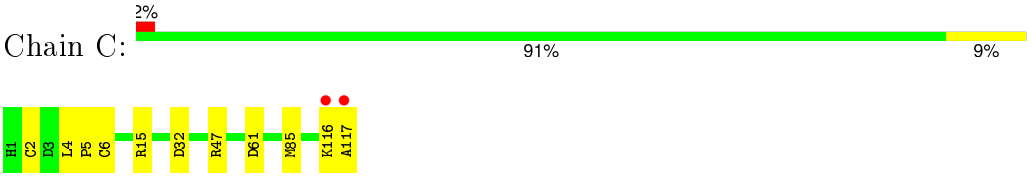
- Molecule 1: Superoxide dismutase [Ni]

Chain B:



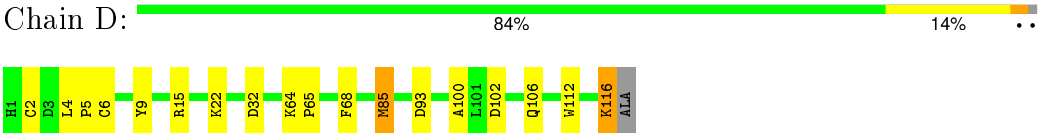
- Molecule 1: Superoxide dismutase [Ni]

Chain C:



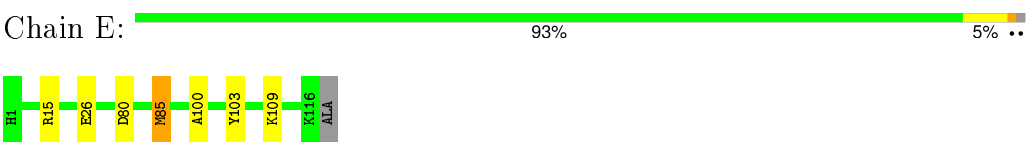
- Molecule 1: Superoxide dismutase [Ni]

Chain D:



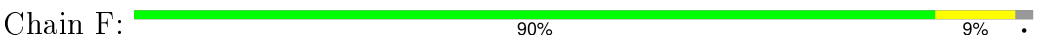
- Molecule 1: Superoxide dismutase [Ni]

Chain E:



- Molecule 1: Superoxide dismutase [Ni]

Chain F:

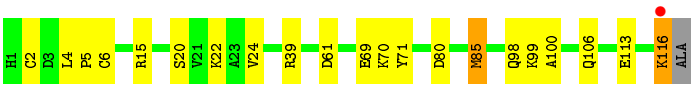


• Molecule 1: Superoxide dismutase [Ni]

Chain G:

• Molecule 1: Superoxide dismutase [Ni]

Chain H:



• Molecule 1: Superoxide dismutase [Ni]

Chain I:

• Molecule 1: Superoxide dismutase [Ni]

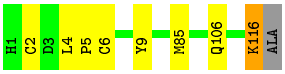
Chain J:

• Molecule 1: Superoxide dismutase [Ni]

Chain K:

• Molecule 1: Superoxide dismutase [Ni]

Chain L:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	64.15Å 111.68Å 106.83Å 90.00° 94.51° 90.00°	Depositor
Resolution (Å)	50.00 – 1.30 49.45 – 1.30	Depositor EDS
% Data completeness (in resolution range)	90.3 (50.00-1.30) 87.3 (49.45-1.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$< I/\sigma(I) >$ [1]	2.17 (at 1.30Å)	Xtriage
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.158 , 0.201 0.152 , 0.187	Depositor DCC
R_{free} test set	15723 reflections (5.14%)	DCC
Wilson B-factor (Å^2)	9.4	Xtriage
Anisotropy	0.152	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 58.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Outliers	4 of 347567 reflections (0.001%)	Xtriage
F_o,F_c correlation	0.98	EDS
Total number of atoms	13192	wwPDB-VP
Average B, all atoms (Å^2)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 45.56 % of the origin peak, indicating pseudo translational symmetry.The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.2982e-04.The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[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: NI

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	0.83	1/975 (0.1%)	1.19	3/1318 (0.2%)
1	B	0.62	0/967	1.11	5/1307 (0.4%)
1	C	0.59	0/974	1.12	6/1318 (0.5%)
1	D	0.58	0/961	1.21	9/1300 (0.7%)
1	E	0.59	0/970	1.16	5/1312 (0.4%)
1	F	0.57	0/952	1.15	2/1289 (0.2%)
1	G	0.58	0/967	1.13	7/1307 (0.5%)
1	H	0.93	1/961 (0.1%)	1.17	6/1300 (0.5%)
1	I	0.58	0/975	1.07	0/1318
1	J	0.64	1/967 (0.1%)	1.12	4/1307 (0.3%)
1	K	0.59	0/969	1.25	9/1311 (0.7%)
1	L	2.19	1/961 (0.1%)	1.07	3/1300 (0.2%)
All	All	0.89	4/11599 (0.0%)	1.15	59/15687 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	116	LYS	C-O	65.24	2.47	1.23
1	H	116	LYS	C-O	22.64	1.66	1.23
1	A	117	ALA	C-OXT	18.44	1.58	1.23
1	J	117	ALA	C-O	7.00	1.36	1.23

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	E	15	ARG	NE-CZ-NH2	-12.58	114.01	120.30
1	K	15	ARG	NE-CZ-NH1	12.42	126.51	120.30
1	D	15	ARG	NE-CZ-NH1	12.09	126.34	120.30
1	B	15	ARG	NE-CZ-NH1	-10.85	114.87	120.30
1	A	116	LYS	C-N-CA	10.72	148.49	121.70
1	A	15	ARG	NE-CZ-NH2	-10.61	115.00	120.30
1	G	15	ARG	NE-CZ-NH1	9.17	124.89	120.30
1	J	15	ARG	NE-CZ-NH1	8.40	124.50	120.30
1	L	85	MET	CG-SD-CE	-8.38	86.79	100.20
1	L	9	TYR	CB-CG-CD2	-8.23	116.06	121.00
1	J	15	ARG	NE-CZ-NH2	-8.11	116.24	120.30
1	C	47	ARG	NE-CZ-NH2	-7.92	116.34	120.30
1	K	15	ARG	NE-CZ-NH2	-7.82	116.39	120.30
1	E	85	MET	CG-SD-CE	-7.50	88.19	100.20
1	G	9	TYR	CB-CG-CD2	-7.42	116.55	121.00
1	D	32	ASP	CB-CG-OD1	7.23	124.80	118.30
1	K	103	TYR	CB-CG-CD1	7.17	125.30	121.00
1	J	39	ARG	NE-CZ-NH2	-7.16	116.72	120.30
1	E	103	TYR	CB-CG-CD2	7.09	125.26	121.00
1	K	39	ARG	NE-CZ-NH1	7.02	123.81	120.30
1	H	80	ASP	CB-CG-OD1	6.77	124.39	118.30
1	G	9	TYR	CB-CG-CD1	6.75	125.05	121.00
1	C	15	ARG	NE-CZ-NH1	6.70	123.65	120.30
1	K	103	TYR	CG-CD2-CE2	6.69	126.65	121.30
1	K	9	TYR	CB-CG-CD2	6.62	124.97	121.00
1	K	99	LYS	CD-CE-NZ	6.60	126.89	111.70
1	A	115	LYS	C-N-CA	6.53	138.02	121.70
1	H	39	ARG	NE-CZ-NH1	6.45	123.52	120.30
1	H	116	LYS	CA-C-O	-6.40	106.67	120.10
1	G	32	ASP	CB-CG-OD2	-6.38	112.55	118.30
1	C	61	ASP	CB-CG-OD2	6.34	124.01	118.30
1	K	15	ARG	CD-NE-CZ	6.27	132.38	123.60
1	D	93	ASP	CB-CG-OD2	-6.27	112.66	118.30
1	K	103	TYR	CD1-CE1-CZ	6.21	125.39	119.80
1	H	85	MET	CG-SD-CE	-6.15	90.37	100.20
1	D	32	ASP	CB-CG-OD2	-6.13	112.79	118.30
1	E	15	ARG	NE-CZ-NH1	6.11	123.35	120.30
1	D	102	ASP	CB-CG-OD2	6.05	123.75	118.30
1	H	15	ARG	NE-CZ-NH2	6.01	123.30	120.30
1	B	39	ARG	NE-CZ-NH2	-5.88	117.36	120.30
1	H	61	ASP	CB-CG-OD1	5.79	123.52	118.30
1	J	10	ASP	CB-CG-OD1	5.67	123.40	118.30
1	B	111	PHE	CB-CG-CD1	-5.55	116.92	120.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	85	MET	CG-SD-CE	-5.53	91.35	100.20
1	D	15	ARG	CD-NE-CZ	5.47	131.25	123.60
1	F	39	ARG	CD-NE-CZ	5.40	131.16	123.60
1	B	9	TYR	CG-CD1-CE1	-5.38	117.00	121.30
1	G	62	TYR	CB-CG-CD2	5.36	124.22	121.00
1	D	9	TYR	CG-CD1-CE1	-5.35	117.02	121.30
1	D	68	PHE	CB-CG-CD1	-5.25	117.13	120.80
1	F	102	ASP	CB-CG-OD1	-5.17	113.64	118.30
1	C	15	ARG	NE-CZ-NH2	-5.15	117.72	120.30
1	G	73	GLU	OE1-CD-OE2	-5.15	117.12	123.30
1	L	116	LYS	CA-C-O	-5.13	109.33	120.10
1	G	102	ASP	CB-CG-OD2	5.09	122.88	118.30
1	E	80	ASP	CB-CG-OD1	5.03	122.83	118.30
1	C	32[A]	ASP	CB-CG-OD1	5.01	122.81	118.30
1	C	32[B]	ASP	CB-CG-OD1	5.01	122.81	118.30
1	B	80	ASP	CB-CG-OD1	5.00	122.80	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	115	LYS	Peptide

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	944	0	927	7	0
1	B	940	0	921	5	0
1	C	943	0	921	4	0
1	D	934	0	916	9	0
1	E	938	0	918	4	0
1	F	925	0	903	5	0
1	G	940	0	921	5	0
1	H	934	0	916	14	0
1	I	944	0	928	8	0
1	J	940	0	921	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	K	938	0	923	7	0
1	L	934	0	916	4	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	155	0	0	2	0
3	B	169	0	0	1	0
3	C	163	0	0	0	0
3	D	164	0	0	3	0
3	E	175	0	0	3	0
3	F	158	0	0	1	0
3	G	162	0	0	2	0
3	H	154	0	0	3	0
3	I	154	0	0	1	0
3	J	173	0	0	3	0
3	K	142	0	0	0	0
3	L	157	0	0	1	0
All	All	13192	0	11031	77	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 3.

All (77) 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:116:LYS:O	1:H:116:LYS:C	1.66	1.32
1:H:113:GLU:HA	1:H:116:LYS:HE2	1.30	1.06
1:K:112:TRP:O	1:K:116:LYS:HD3	1.58	1.02
1:I:20:SER:OG	1:K:1[B]:HIS:CE1	2.38	0.77
1:G:106:GLN:HG2	3:G:2358:HOH:O	1.84	0.76
1:K:112:TRP:O	1:K:116:LYS:CD	2.35	0.72
1:F:70:LYS:HD3	1:F:71:TYR:CE1	2.26	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:64:LYS:HE2	3:J:2910:HOH:O	1.92	0.70
1:L:106:GLN:HG2	3:L:2190:HOH:O	1.92	0.68
1:E:109:LYS:HE2	3:E:1962:HOH:O	1.94	0.66
1:J:64:LYS:CE	3:J:2910:HOH:O	2.43	0.66
1:I:22:LYS:HE2	1:I:26:GLU:OE1	1.97	0.65
1:B:109:LYS:NZ	3:B:2648:HOH:O	2.30	0.64
1:I:23:ALA:O	1:I:27:LYS:HG3	1.99	0.63
1:A:64:LYS:HB3	1:A:65:PRO:CD	2.29	0.63
1:D:22:LYS:NZ	3:D:2799:HOH:O	2.12	0.62
1:J:106:GLN:HG2	3:J:2334:HOH:O	1.99	0.61
1:C:85:MET:HE3	1:C:85:MET:HA	1.82	0.61
1:F:7:GLY:O	3:F:2824:HOH:O	2.16	0.61
1:A:65:PRO:HB2	1:A:66:PRO:HD3	1.81	0.61
1:C:116:LYS:O	1:C:117:ALA:HB2	2.01	0.59
1:A:112:TRP:O	1:A:117:ALA:HB3	2.05	0.57
1:I:85:MET:HE3	1:I:85:MET:HA	1.87	0.56
1:A:64:LYS:HB3	1:A:65:PRO:HD2	1.87	0.56
1:H:99:LYS:HE2	3:H:2468:HOH:O	2.05	0.56
1:E:109:LYS:CE	3:E:1962:HOH:O	2.54	0.55
1:H:113:GLU:HA	1:H:116:LYS:CE	2.21	0.55
1:H:113:GLU:CA	1:H:116:LYS:HE2	2.21	0.54
1:H:85:MET:HE2	1:H:100:ALA:HB1	1.91	0.53
1:L:116:LYS:O	1:L:116:LYS:C	2.47	0.53
1:B:85:MET:HA	1:B:85:MET:HE3	1.92	0.52
1:D:106:GLN:NE2	3:D:1701:HOH:O	2.42	0.51
1:D:112:TRP:O	1:D:116:LYS:CD	2.58	0.51
1:D:112:TRP:O	1:D:116:LYS:HD3	2.10	0.51
1:H:22:LYS:NZ	1:H:98:GLN:OE1	2.24	0.51
1:H:4:LEU:N	1:H:5:PRO:HA	2.28	0.49
1:D:85:MET:HE2	1:D:100:ALA:HB1	1.94	0.48
1:H:116:LYS:O	1:H:116:LYS:CA	2.57	0.48
1:I:29:ALA:HB2	3:I:2049:HOH:O	2.13	0.48
1:H:2:CYS:SG	1:H:6:CYS:HB3	2.52	0.48
1:H:106:GLN:NE2	3:H:2472:HOH:O	2.47	0.48
1:F:2:CYS:SG	1:F:6:CYS:HB3	2.53	0.47
1:H:70:LYS:HE2	1:H:71:TYR:CZ	2.49	0.47
1:E:26[A]:GLU:OE2	3:E:1718:HOH:O	2.20	0.46
1:J:4:LEU:N	1:J:5:PRO:HA	2.30	0.46
1:E:85:MET:HE2	1:E:100:ALA:HB1	1.97	0.46
1:G:85:MET:HE3	1:G:85:MET:HA	1.98	0.46
1:D:4:LEU:HB3	1:D:5:PRO:C	2.36	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:2:CYS:SG	1:D:6:CYS:HB3	2.56	0.44
1:C:4:LEU:HB3	1:C:5:PRO:C	2.37	0.44
1:A:64:LYS:HE2	3:A:2735:HOH:O	2.18	0.44
1:J:20:SER:O	1:J:24:VAL:HG23	2.17	0.44
1:D:22:LYS:HD3	3:D:2799:HOH:O	2.18	0.44
1:B:81:THR:OG1	1:B:103:TYR:HB3	2.18	0.44
1:A:116:LYS:HD2	3:A:2744:HOH:O	2.18	0.43
1:F:4:LEU:HB3	1:F:5:PRO:C	2.39	0.43
1:B:64:LYS:HG3	1:B:64:LYS:HZ3	1.56	0.43
1:J:2:CYS:SG	1:J:6:CYS:HB3	2.57	0.43
1:G:4:LEU:HB3	1:G:5:PRO:C	2.39	0.43
1:G:2:CYS:SG	1:G:6:CYS:HB3	2.59	0.42
1:B:2:CYS:SG	1:B:6:CYS:HB3	2.59	0.42
1:F:4:LEU:N	1:F:5:PRO:HA	2.34	0.42
1:K:2:CYS:SG	1:K:6:CYS:HB3	2.59	0.42
1:H:99:LYS:HD3	3:H:2468:HOH:O	2.18	0.42
1:K:4:LEU:HB3	1:K:5:PRO:C	2.40	0.41
1:K:4:LEU:N	1:K:5:PRO:HA	2.34	0.41
1:C:2:CYS:SG	1:C:6:CYS:HB3	2.59	0.41
1:A:4:LEU:N	1:A:5:PRO:HA	2.35	0.41
1:I:24:VAL:O	1:I:28:MET:HG3	2.20	0.41
1:L:2:CYS:SG	1:L:6:CYS:HB3	2.60	0.41
1:H:20:SER:O	1:H:24:VAL:HG23	2.21	0.41
1:G:116:LYS:HE3	3:G:2037:HOH:O	2.21	0.41
1:K:85:MET:HE2	1:K:100:ALA:HB1	2.02	0.41
1:I:4:LEU:N	1:I:5:PRO:HA	2.36	0.41
1:I:2:CYS:SG	1:I:6:CYS:HB3	2.59	0.41
1:L:4:LEU:HB3	1:L:5:PRO:C	2.42	0.40
1:D:64:LYS:HB3	1:D:65:PRO:HD2	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	116/117 (99%)	114 (98%)	1 (1%)	1 (1%)	21 2
1	B	115/117 (98%)	113 (98%)	2 (2%)	0	100 100
1	C	116/117 (99%)	116 (100%)	0	0	100 100
1	D	114/117 (97%)	111 (97%)	3 (3%)	0	100 100
1	E	115/117 (98%)	114 (99%)	1 (1%)	0	100 100
1	F	113/117 (97%)	113 (100%)	0	0	100 100
1	G	115/117 (98%)	113 (98%)	2 (2%)	0	100 100
1	H	114/117 (97%)	112 (98%)	2 (2%)	0	100 100
1	I	116/117 (99%)	115 (99%)	1 (1%)	0	100 100
1	J	115/117 (98%)	114 (99%)	1 (1%)	0	100 100
1	K	115/117 (98%)	113 (98%)	2 (2%)	0	100 100
1	L	114/117 (97%)	113 (99%)	1 (1%)	0	100 100
All	All	1378/1404 (98%)	1361 (99%)	16 (1%)	1 (0%)	56 21

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	116	LYS

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	100/98 (102%)	99 (99%)	1 (1%)	82 51
1	B	99/98 (101%)	99 (100%)	0	100 100
1	C	100/98 (102%)	100 (100%)	0	100 100
1	D	99/98 (101%)	98 (99%)	1 (1%)	82 51
1	E	100/98 (102%)	100 (100%)	0	100 100
1	F	98/98 (100%)	97 (99%)	1 (1%)	82 51
1	G	99/98 (101%)	99 (100%)	0	100 100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	H	99/98 (101%)	98 (99%)	1 (1%)	82 51
1	I	100/98 (102%)	100 (100%)	0	100 100
1	J	99/98 (101%)	99 (100%)	0	100 100
1	K	100/98 (102%)	99 (99%)	1 (1%)	82 51
1	L	99/98 (101%)	99 (100%)	0	100 100
All	All	1192/1176 (101%)	1187 (100%)	5 (0%)	93 77

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	PRO
1	D	116	LYS
1	F	115	LYS
1	H	69	GLU
1	K	116	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	106	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 12 ligands modelled in this entry, 12 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	117/117 (100%)	-0.58	2 (1%) 73 70	7, 12, 26, 49	0
1	B	117/117 (100%)	-0.63	0 100 100	8, 12, 23, 37	0
1	C	117/117 (100%)	-0.55	2 (1%) 73 70	7, 11, 22, 72	0
1	D	116/117 (99%)	-0.67	0 100 100	8, 12, 22, 62	0
1	E	116/117 (99%)	-0.63	0 100 100	8, 11, 21, 33	0
1	F	115/117 (98%)	-0.57	0 100 100	10, 14, 27, 38	0
1	G	117/117 (100%)	-0.62	0 100 100	8, 12, 26, 38	0
1	H	116/117 (99%)	-0.59	1 (0%) 85 85	8, 14, 27, 61	0
1	I	117/117 (100%)	-0.52	1 (0%) 85 85	8, 12, 28, 42	0
1	J	117/117 (100%)	-0.68	0 100 100	7, 11, 22, 34	0
1	K	116/117 (99%)	-0.61	0 100 100	7, 12, 23, 58	0
1	L	116/117 (99%)	-0.69	0 100 100	7, 11, 21, 35	0
All	All	1397/1404 (99%)	-0.61	6 (0%) 93 92	7, 12, 25, 72	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	117	ALA	8.9
1	A	117	ALA	6.5
1	I	117	ALA	5.4
1	H	116	LYS	2.8
1	C	116	LYS	2.3
1	A	116	LYS	2.3

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 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
2	NI	H	208	1/1	1.00	0.05	-0.58	8,8,8,8	0
2	NI	G	207	1/1	1.00	0.05	-0.66	7,7,7,7	0
2	NI	B	202	1/1	1.00	0.05	-0.73	9,9,9,9	0
2	NI	C	203	1/1	1.00	0.05	-0.79	8,8,8,8	0
2	NI	J	210	1/1	1.00	0.04	-0.82	9,9,9,9	0
2	NI	K	211	1/1	1.00	0.04	-1.00	10,10,10,10	0
2	NI	D	204	1/1	1.00	0.04	-1.24	7,7,7,7	0
2	NI	F	206	1/1	1.00	0.04	-1.33	9,9,9,9	0
2	NI	L	212	1/1	1.00	0.04	-1.35	8,8,8,8	0
2	NI	I	209	1/1	1.00	0.04	-1.80	7,7,7,7	0
2	NI	A	201	1/1	1.00	0.03	-1.91	9,9,9,9	0
2	NI	E	205	1/1	1.00	0.03	-2.60	7,7,7,7	0

6.5 Other polymers (i)

There are no such residues in this entry.