Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 02:29 PM GMT

PDB ID	:	3ZKE
Title	:	Structure of LC8 in complex with Nek9 peptide
Authors	:	Gallego, P.; Velazquez-Campoy, A.; Regue, L.; Roig, J.; Reverter, D.
Deposited on	:	2013-01-22
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:

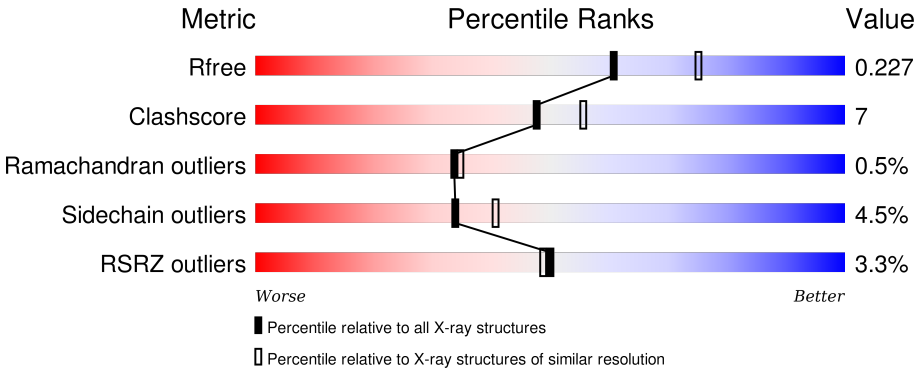
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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))
R_{free}	91344	3774 (2.20-2.20)
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	89	2% 78% 18%
1	C	89	2% 79% 16%
1	E	89	2% 73% 18%
1	G	89	% 72% 18%
1	I	89	2% 78% 16%

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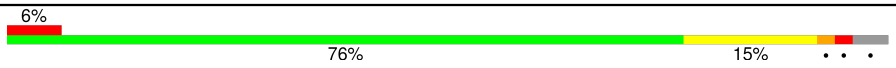
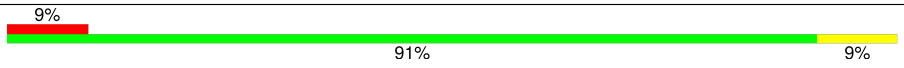
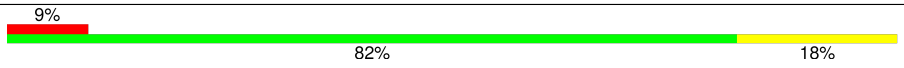
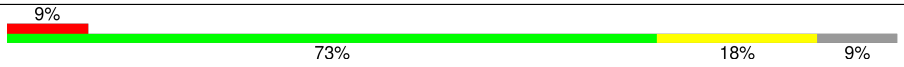
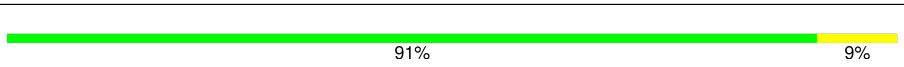
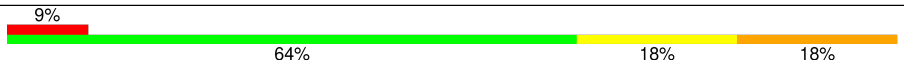
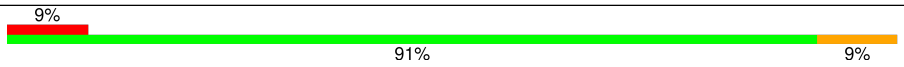
Mol	Chain	Length	Quality of chain
1	K	89	
2	B	11	
2	D	11	
2	F	11	
2	H	11	
2	J	11	
2	L	11	

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 4726 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 DYNEIN LIGHT CHAIN 1, CYTOPLASMIC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	85	Total	C	N	O	S	0	0	0
			695	447	115	129	4			
1	C	85	Total	C	N	O	S	0	0	0
			695	447	115	129	4			
1	E	85	Total	C	N	O	S	0	0	0
			695	447	115	129	4			
1	G	85	Total	C	N	O	S	0	0	0
			695	447	115	129	4			
1	I	85	Total	C	N	O	S	0	0	0
			695	447	115	129	4			
1	K	85	Total	C	N	O	S	0	0	0
			695	447	115	129	4			

- Molecule 2 is a protein called NEK9 PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	11	Total	C	N	O	S	0	0	0
			76	45	15	15	1			
2	D	11	Total	C	N	O	S	0	0	0
			76	45	15	15	1			
2	F	10	Total	C	N	O	S	0	0	0
			69	40	14	14	1			
2	H	11	Total	C	N	O	S	0	0	0
			76	45	15	15	1			
2	J	11	Total	C	N	O	S	0	0	0
			76	45	15	15	1			
2	L	11	Total	C	N	O	S	0	0	0
			76	45	15	15	1			

- Molecule 3 is water.

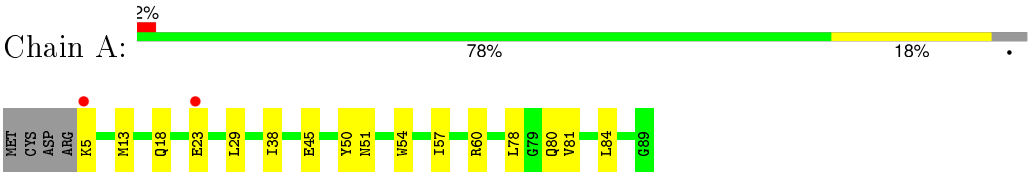
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	18	Total O 18 18	0	0
3	B	3	Total O 3 3	0	0
3	C	13	Total O 13 13	0	0
3	D	3	Total O 3 3	0	0
3	E	18	Total O 18 18	0	0
3	F	3	Total O 3 3	0	0
3	G	18	Total O 18 18	0	0
3	H	3	Total O 3 3	0	0
3	I	12	Total O 12 12	0	0
3	J	6	Total O 6 6	0	0
3	K	9	Total O 9 9	0	0
3	L	1	Total O 1 1	0	0

3 Residue-property plots

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.

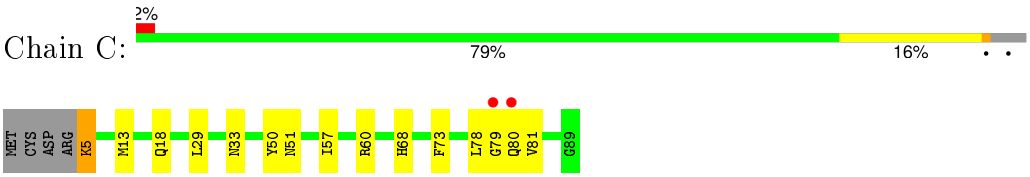
- Molecule 1: DYNEIN LIGHT CHAIN 1, CYTOPLASMIC

Chain A:



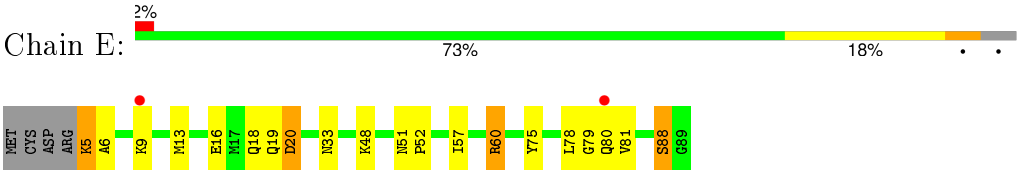
- Molecule 1: DYNEIN LIGHT CHAIN 1, CYTOPLASMIC

Chain C:



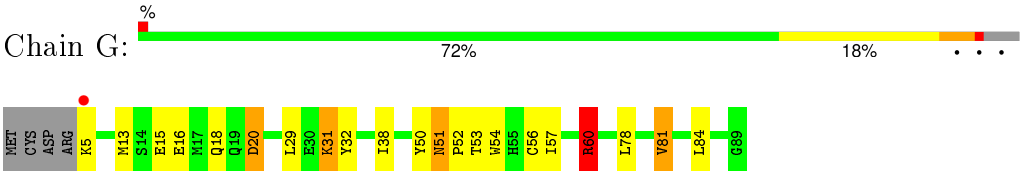
- Molecule 1: DYNEIN LIGHT CHAIN 1, CYTOPLASMIC

Chain E:



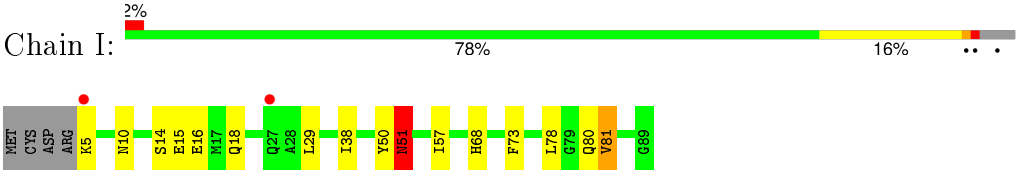
- Molecule 1: DYNEIN LIGHT CHAIN 1, CYTOPLASMIC

Chain G:



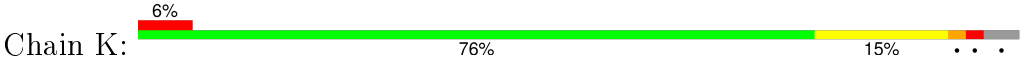
- Molecule 1: DYNEIN LIGHT CHAIN 1, CYTOPLASMIC

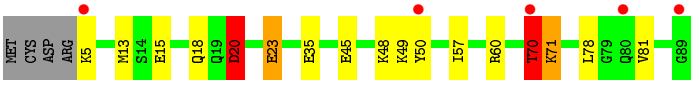
Chain I:



- Molecule 1: DYNEIN LIGHT CHAIN 1, CYTOPLASMIC

Chain K:





• Molecule 2: NEK9 PROTEIN

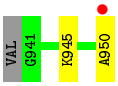
Chain B:

• Molecule 2: NEK9 PROTEIN

Chain D:

• Molecule 2: NEK9 PROTEIN

Chain F:

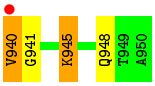


• Molecule 2: NEK9 PROTEIN

Chain H:

• Molecule 2: NEK9 PROTEIN

Chain J:



• Molecule 2: NEK9 PROTEIN

Chain L:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	39.98Å 105.12Å 133.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	82.69 – 2.20 41.34 – 2.20	Depositor EDS
% Data completeness (in resolution range)	96.9 (82.69-2.20) 96.9 (41.34-2.20)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.02 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.196 , 0.228 0.195 , 0.227	Depositor DCC
R_{free} test set	1459 reflections (5.35%)	DCC
Wilson B-factor (Å^2)	35.2	Xtriage
Anisotropy	0.363	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 45.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Outliers	0 of 28736 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	4726	wwPDB-VP
Average B, all atoms (Å^2)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.76% 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

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	1/711 (0.1%)	0.99	3/954 (0.3%)
1	C	0.98	0/711	0.92	1/954 (0.1%)
1	E	1.15	7/711 (1.0%)	1.07	6/954 (0.6%)
1	G	1.09	2/711 (0.3%)	1.06	6/954 (0.6%)
1	I	1.05	2/711 (0.3%)	0.92	1/954 (0.1%)
1	K	1.23	5/711 (0.7%)	1.15	6/954 (0.6%)
2	B	0.91	0/76	0.84	0/100
2	D	0.88	0/76	0.78	0/100
2	F	1.19	0/69	0.79	0/90
2	H	0.88	0/76	0.69	0/100
2	J	1.09	0/76	0.74	0/100
2	L	1.29	2/76 (2.6%)	0.85	0/100
All	All	1.10	19/4715 (0.4%)	1.00	23/6314 (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
2	J	0	1

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	K	45	GLU	CD-OE1	-13.71	1.10	1.25
1	K	45	GLU	CD-OE2	-9.29	1.15	1.25
1	A	54	TRP	CD2-CE2	7.23	1.50	1.41
1	I	51	ASN	CG-ND2	-6.56	1.16	1.32
1	I	51	ASN	CG-OD1	-6.08	1.10	1.24
1	E	75	TYR	CE1-CZ	5.91	1.46	1.38
2	L	943	HIS	CE1-NE2	-5.83	1.19	1.32
1	K	23	GLU	CB-CG	-5.65	1.41	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	943	HIS	CG-ND1	-5.62	1.26	1.38
1	G	15	GLU	CG-CD	-5.62	1.43	1.51
1	K	15	GLU	CG-CD	-5.52	1.43	1.51
1	K	70	THR	CB-CG2	-5.48	1.34	1.52
1	E	20	ASP	CG-OD2	-5.42	1.12	1.25
1	E	88	SER	CB-OG	-5.33	1.35	1.42
1	G	54	TRP	CD2-CE2	5.32	1.47	1.41
1	E	60	ARG	NE-CZ	-5.10	1.26	1.33
1	E	16	GLU	CD-OE2	-5.09	1.20	1.25
1	E	60	ARG	CD-NE	-5.08	1.37	1.46
1	E	60	ARG	CZ-NH1	-5.05	1.26	1.33

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	K	45	GLU	OE1-CD-OE2	-14.90	105.42	123.30
1	E	20	ASP	CB-CG-OD1	8.84	126.25	118.30
1	G	81	VAL	CG1-CB-CG2	-8.25	97.70	110.90
1	K	45	GLU	CG-CD-OE2	7.83	133.95	118.30
1	E	20	ASP	OD1-CG-OD2	-7.31	109.42	123.30
1	A	23	GLU	OE1-CD-OE2	-6.89	115.03	123.30
1	G	5	LYS	CB-CG-CD	-6.70	94.19	111.60
1	E	20	ASP	CB-CG-OD2	6.67	124.31	118.30
1	E	16	GLU	OE1-CD-OE2	-6.60	115.38	123.30
1	G	60	ARG	NE-CZ-NH2	6.25	123.42	120.30
1	G	60	ARG	NE-CZ-NH1	-6.20	117.20	120.30
1	K	20	ASP	CB-CG-OD1	6.05	123.75	118.30
1	K	81	VAL	CG1-CB-CG2	-6.00	101.30	110.90
1	K	5	LYS	CB-CG-CD	-5.87	96.34	111.60
1	A	45	GLU	OE1-CD-OE2	-5.70	116.45	123.30
1	K	23	GLU	CB-CA-C	-5.42	99.57	110.40
1	A	5	LYS	CB-CG-CD	-5.34	97.71	111.60
1	I	5	LYS	CA-CB-CG	-5.33	101.68	113.40
1	G	16	GLU	CG-CD-OE1	5.19	128.69	118.30
1	E	16	GLU	CG-CD-OE1	5.16	128.62	118.30
1	C	60	ARG	CD-NE-CZ	-5.11	116.45	123.60
1	E	9	LYS	CD-CE-NZ	5.09	123.40	111.70
1	G	20	ASP	CB-CG-OD2	5.05	122.84	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	J	940	VAL	Peptide

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	695	0	680	7	0
1	C	695	0	680	15	3
1	E	695	0	680	19	3
1	G	695	0	680	11	2
1	I	695	0	680	12	0
1	K	695	0	680	11	2
2	B	76	0	75	0	0
2	D	76	0	75	3	0
2	F	69	0	66	3	0
2	H	76	0	75	1	0
2	J	76	0	75	4	0
2	L	76	0	75	1	0
3	A	18	0	0	1	0
3	B	3	0	0	0	0
3	C	13	0	0	1	0
3	D	3	0	0	0	0
3	E	18	0	0	2	0
3	F	3	0	0	0	0
3	G	18	0	0	1	0
3	H	3	0	0	0	0
3	I	12	0	0	4	0
3	J	6	0	0	0	0
3	K	9	0	0	1	0
3	L	1	0	0	0	0
All	All	4726	0	4521	68	5

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 7.

All (68) 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:C:80:GLN:HE22	1:E:5:LYS:NZ	1.25	1.31
1:C:79:GLY:O	3:C:2012:HOH:O	1.58	1.22
1:I:78:LEU:O	3:I:2012:HOH:O	1.72	1.06
1:C:80:GLN:NE2	1:E:5:LYS:NZ	2.03	1.06
1:I:16:GLU:OE1	3:I:2004:HOH:O	1.81	0.97
1:C:80:GLN:HE22	1:E:5:LYS:HZ1	1.17	0.93
1:C:80:GLN:HE22	1:E:5:LYS:HZ2	1.10	0.91
1:C:80:GLN:NE2	1:E:5:LYS:HZ2	1.67	0.89
2:D:950:ALA:HB2	2:F:950:ALA:HB2	1.56	0.86
1:G:51:ASN:OD1	3:G:2011:HOH:O	1.96	0.83
2:D:950:ALA:HB2	2:F:950:ALA:CB	2.14	0.77
1:I:29:LEU:HD23	1:I:38:ILE:HD13	1.68	0.73
1:I:81:VAL:HG12	3:I:2012:HOH:O	1.91	0.71
1:C:80:GLN:NE2	1:E:5:LYS:HZ1	1.77	0.70
1:K:13:MET:HG3	1:K:18:GLN:HG3	1.75	0.69
1:K:49:LYS:HG2	1:K:50:TYR:CE2	2.29	0.68
1:K:20:ASP:OD2	1:K:50:TYR:OH	2.12	0.68
1:E:51:ASN:OD1	3:E:2015:HOH:O	2.11	0.68
3:E:2012:HOH:O	2:J:945:LYS:HE3	1.93	0.68
1:E:57:ILE:HG12	1:G:57:ILE:HG12	1.74	0.68
1:I:16:GLU:CD	3:I:2004:HOH:O	2.29	0.65
1:E:80:GLN:OE1	1:E:80:GLN:N	2.24	0.65
1:C:13:MET:HG2	1:C:18:GLN:HG3	1.79	0.63
1:G:29:LEU:HD12	1:G:78:LEU:HD23	1.81	0.62
1:I:50:TYR:O	1:I:51:ASN:HB2	1.99	0.62
1:C:29:LEU:HD12	1:C:78:LEU:HD23	1.83	0.60
1:I:80:GLN:N	1:I:80:GLN:OE1	2.32	0.59
1:G:13:MET:HG3	1:G:18:GLN:HG3	1.86	0.57
2:H:940:VAL:O	2:H:940:VAL:HG22	2.05	0.56
1:A:13:MET:HG3	1:A:18:GLN:HG3	1.90	0.54
1:I:68:HIS:NE2	2:J:945:LYS:HE2	2.22	0.54
1:G:29:LEU:HD23	1:G:38:ILE:HD13	1.90	0.53
1:E:13:MET:HG3	1:E:18:GLN:HG3	1.90	0.53
1:A:60:ARG:HD2	3:A:2012:HOH:O	2.08	0.52
2:J:940:VAL:O	2:J:940:VAL:HG23	2.10	0.52
1:G:57:ILE:HD13	1:G:84:LEU:HD23	1.92	0.51
1:I:57:ILE:HG12	1:K:57:ILE:HG12	1.93	0.50
1:A:57:ILE:HG12	1:C:57:ILE:HG12	1.93	0.49
1:A:50:TYR:O	1:A:51:ASN:HB2	2.13	0.49
1:E:60:ARG:HG2	1:E:80:GLN:O	2.13	0.49
1:C:80:GLN:CD	1:E:5:LYS:HZ1	2.15	0.48
1:G:50:TYR:O	1:G:51:ASN:HB2	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:48:LYS:HG3	3:K:2007:HOH:O	2.12	0.48
1:C:80:GLN:HB3	1:E:79:GLY:HA2	1.95	0.48
1:E:60:ARG:HD3	1:E:80:GLN:HG2	1.95	0.48
1:G:56:CYS:HA	1:G:84:LEU:O	2.14	0.47
1:K:60:ARG:HH11	1:K:60:ARG:HD2	1.50	0.47
1:C:50:TYR:O	1:C:51:ASN:HB2	2.14	0.47
1:K:70:THR:HG23	2:L:943:HIS:CE1	2.50	0.47
1:E:60:ARG:HD2	1:E:60:ARG:HH11	1.43	0.45
1:G:60:ARG:HD2	1:G:60:ARG:HH11	1.56	0.45
1:C:68:HIS:CG	1:C:73:PHE:HB2	2.52	0.45
1:G:50:TYR:O	1:G:51:ASN:CB	2.64	0.44
1:E:48:LYS:O	1:I:10:ASN:HA	2.17	0.44
1:K:78:LEU:HD12	1:K:78:LEU:HA	1.77	0.44
1:A:57:ILE:HD13	1:A:84:LEU:HD23	1.99	0.44
1:K:70:THR:O	1:K:71:LYS:HB2	2.17	0.44
2:D:950:ALA:CB	2:F:950:ALA:HB2	2.39	0.43
1:I:68:HIS:CG	1:I:73:PHE:HB2	2.53	0.43
1:K:49:LYS:HG2	1:K:50:TYR:CD2	2.54	0.42
1:C:5:LYS:HE3	1:E:80:GLN:NE2	2.34	0.42
1:A:29:LEU:HD23	1:A:38:ILE:HD13	2.00	0.42
1:A:29:LEU:HD12	1:A:78:LEU:HD23	2.02	0.41
1:E:6:ALA:HB2	1:E:78:LEU:HD13	2.03	0.41
1:E:51:ASN:HA	1:E:52:PRO:HD3	1.95	0.41
1:G:52:PRO:HA	1:G:53:THR:HA	1.88	0.40
1:I:14:SER:O	1:I:18:GLN:HG3	2.20	0.40
2:J:948:GLN:OE1	1:K:35:GLU:HB2	2.21	0.40

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:31:LYS:NZ	1:K:20:ASP:OD2[3_455]	1.62	0.58
1:G:32:TYR:OH	1:K:23:GLU:OE1[3_455]	1.80	0.40
1:C:33:ASN:OD1	1:E:33:ASN:OD1[4_545]	2.02	0.18
1:C:33:ASN:OD1	1:E:33:ASN:ND2[4_545]	2.09	0.11
1:C:33:ASN:OD1	1:E:33:ASN:CG[4_545]	2.17	0.03

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	83/89 (93%)	79 (95%)	4 (5%)	0	100 100
1	C	83/89 (93%)	80 (96%)	3 (4%)	0	100 100
1	E	83/89 (93%)	80 (96%)	3 (4%)	0	100 100
1	G	83/89 (93%)	80 (96%)	2 (2%)	1 (1%)	16 12
1	I	83/89 (93%)	80 (96%)	2 (2%)	1 (1%)	16 12
1	K	83/89 (93%)	80 (96%)	3 (4%)	0	100 100
2	B	9/11 (82%)	9 (100%)	0	0	100 100
2	D	9/11 (82%)	9 (100%)	0	0	100 100
2	F	8/11 (73%)	8 (100%)	0	0	100 100
2	H	9/11 (82%)	9 (100%)	0	0	100 100
2	J	9/11 (82%)	8 (89%)	0	1 (11%)	0 0
2	L	9/11 (82%)	9 (100%)	0	0	100 100
All	All	551/600 (92%)	531 (96%)	17 (3%)	3 (0%)	34 35

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	J	941	GLY
1	G	51	ASN
1	I	51	ASN

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	74/78 (95%)	72 (97%)	2 (3%)	52 64
1	C	74/78 (95%)	72 (97%)	2 (3%)	52 64
1	E	74/78 (95%)	69 (93%)	5 (7%)	20 21
1	G	74/78 (95%)	70 (95%)	4 (5%)	27 31
1	I	74/78 (95%)	72 (97%)	2 (3%)	52 64
1	K	74/78 (95%)	71 (96%)	3 (4%)	37 45
2	B	8/8 (100%)	7 (88%)	1 (12%)	6 4
2	D	8/8 (100%)	7 (88%)	1 (12%)	6 4
2	F	7/8 (88%)	6 (86%)	1 (14%)	4 3
2	H	8/8 (100%)	8 (100%)	0	100 100
2	J	8/8 (100%)	7 (88%)	1 (12%)	6 4
2	L	8/8 (100%)	8 (100%)	0	100 100
All	All	491/516 (95%)	469 (96%)	22 (4%)	34 41

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	80	GLN
1	A	81	VAL
2	B	940	VAL
1	C	5	LYS
1	C	81	VAL
2	D	940	VAL
1	E	5	LYS
1	E	19	GLN
1	E	20	ASP
1	E	81	VAL
1	E	88	SER
2	F	945	LYS
1	G	20	ASP
1	G	31	LYS
1	G	60	ARG
1	G	81	VAL
1	I	15	GLU
1	I	81	VAL
2	J	945	LYS
1	K	20	ASP
1	K	70	THR
1	K	71	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	C	80	GLN
1	G	51	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

There are no ligands in this entry.

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	85/89 (95%)	-0.23	2 (2%) 62 61	23, 33, 52, 72	0
1	C	85/89 (95%)	-0.19	2 (2%) 62 61	23, 37, 57, 73	0
1	E	85/89 (95%)	-0.22	2 (2%) 62 61	24, 35, 52, 72	0
1	G	85/89 (95%)	-0.26	1 (1%) 81 80	22, 35, 54, 77	0
1	I	85/89 (95%)	-0.07	2 (2%) 62 61	24, 40, 65, 81	0
1	K	85/89 (95%)	-0.04	5 (5%) 26 25	26, 39, 62, 81	0
2	B	11/11 (100%)	-0.11	1 (9%) 11 11	26, 31, 49, 55	0
2	D	11/11 (100%)	-0.21	1 (9%) 11 11	26, 34, 45, 61	0
2	F	10/11 (90%)	-0.13	1 (10%) 9 8	27, 31, 57, 62	0
2	H	11/11 (100%)	-0.29	0 100 100	27, 33, 50, 56	0
2	J	11/11 (100%)	0.32	1 (9%) 11 11	28, 31, 56, 82	0
2	L	11/11 (100%)	0.29	1 (9%) 11 11	38, 42, 68, 69	0
All	All	575/600 (95%)	-0.15	19 (3%) 50 49	22, 37, 59, 82	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	J	940	VAL	6.8
1	I	5	LYS	4.0
1	C	80	GLN	3.5
1	K	5	LYS	3.2
1	E	80	GLN	3.0
1	K	89	GLY	3.0
1	G	5	LYS	2.9
2	F	950	ALA	2.8
1	A	5	LYS	2.7
1	K	50	TYR	2.6
1	K	70	THR	2.5

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Mol	Chain	Res	Type	RSRZ
1	I	27	GLN	2.5
2	B	950	ALA	2.4
1	C	79	GLY	2.4
2	L	950	ALA	2.2
1	E	9	LYS	2.2
1	K	80	GLN	2.2
2	D	940	VAL	2.1
1	A	23	GLU	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)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.