



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:57 PM GMT

PDB ID : 4DX8  
Title : ICAP1 in complex with KRIT1 N-terminus  
Authors : Liu, W.; Draheim, K.; Zhang, R.; Calderwood, D.A.; Boggon, T.J.  
Deposited on : 2012-02-27  
Resolution : 2.54 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) 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

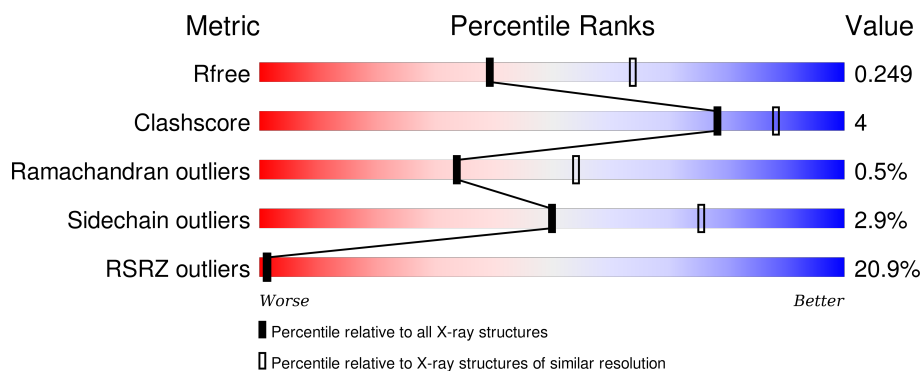
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.54 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	4549 (2.58-2.50)
Clashscore	102246	5292 (2.58-2.50)
Ramachandran outliers	100387	5194 (2.58-2.50)
Sidechain outliers	100360	5196 (2.58-2.50)
RSRZ outliers	91569	4561 (2.58-2.50)

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  $\geq 3$ , 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  $\leq 5\%$ . 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	154	<div> <div>6%</div> <div>82%</div> <div>8%</div> <div>10%</div> </div>
1	B	154	<div> <div>9%</div> <div>81%</div> <div>6%</div> <div>12%</div> </div>
1	D	154	<div> <div>6%</div> <div>76%</div> <div>10%</div> <div>14%</div> </div>
1	E	154	<div> <div>2%</div> <div>79%</div> <div>7%</div> <div>14%</div> </div>
2	H	203	<div> <div>17%</div> <div>69%</div> <div>7%</div> <div>22%</div> </div>

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Mol	Chain	Length	Quality of chain
2	I	203	
2	J	203	
2	K	203	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	BR	A	302	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8128 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 Integrin beta-1-binding protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	139	Total	C	N	O	S	0	0	0
			1080	690	172	212	6			
1	B	136	Total	C	N	O	S	0	0	0
			1061	680	168	207	6			
1	D	133	Total	C	N	O	S	0	0	0
			1038	665	164	203	6			
1	E	133	Total	C	N	O	S	0	0	0
			1038	665	164	203	6			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	47	GLY	-	EXPRESSION TAG	UNP O14713
A	48	SER	-	EXPRESSION TAG	UNP O14713
B	47	GLY	-	EXPRESSION TAG	UNP O14713
B	48	SER	-	EXPRESSION TAG	UNP O14713
D	47	GLY	-	EXPRESSION TAG	UNP O14713
D	48	SER	-	EXPRESSION TAG	UNP O14713
E	47	GLY	-	EXPRESSION TAG	UNP O14713
E	48	SER	-	EXPRESSION TAG	UNP O14713

- Molecule 2 is a protein called Krev interaction trapped protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	159	Total	C	N	O	S	0	0	0
			1278	825	221	225	7			
2	I	111	Total	C	N	O	S	0	0	0
			906	590	159	152	5			
2	J	120	Total	C	N	O	S	0	0	0
			979	634	166	176	3			
2	K	88	Total	C	N	O	S	0	0	0
			718	465	126	125	2			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	-4	GLY	-	EXPRESSION TAG	UNP O00522
H	-3	PRO	-	EXPRESSION TAG	UNP O00522
H	-2	LEU	-	EXPRESSION TAG	UNP O00522
H	-1	GLY	-	EXPRESSION TAG	UNP O00522
H	0	SER	-	EXPRESSION TAG	UNP O00522
I	-4	GLY	-	EXPRESSION TAG	UNP O00522
I	-3	PRO	-	EXPRESSION TAG	UNP O00522
I	-2	LEU	-	EXPRESSION TAG	UNP O00522
I	-1	GLY	-	EXPRESSION TAG	UNP O00522
I	0	SER	-	EXPRESSION TAG	UNP O00522
J	-4	GLY	-	EXPRESSION TAG	UNP O00522
J	-3	PRO	-	EXPRESSION TAG	UNP O00522
J	-2	LEU	-	EXPRESSION TAG	UNP O00522
J	-1	GLY	-	EXPRESSION TAG	UNP O00522
J	0	SER	-	EXPRESSION TAG	UNP O00522
K	-4	GLY	-	EXPRESSION TAG	UNP O00522
K	-3	PRO	-	EXPRESSION TAG	UNP O00522
K	-2	LEU	-	EXPRESSION TAG	UNP O00522
K	-1	GLY	-	EXPRESSION TAG	UNP O00522
K	0	SER	-	EXPRESSION TAG	UNP O00522

- Molecule 3 is BROMIDE ION (three-letter code: BR) (formula: Br).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	H	3	Total Br 3 3	0	0
3	B	3	Total Br 3 3	0	0
3	A	3	Total Br 3 3	0	0
3	D	4	Total Br 4 4	0	0
3	E	3	Total Br 3 3	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	4	Total O 4 4	0	0
4	D	1	Total O 1 1	0	0

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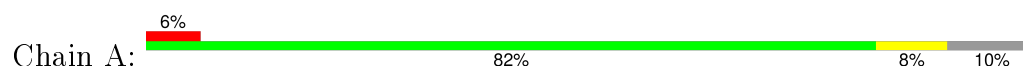
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	7	Total	O	0	0
			7	7		
4	H	1	Total	O	0	0
			1	1		
4	J	1	Total	O	0	0
			1	1		

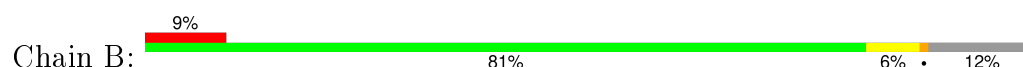
### 3 Residue-property plots [i](#)

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.

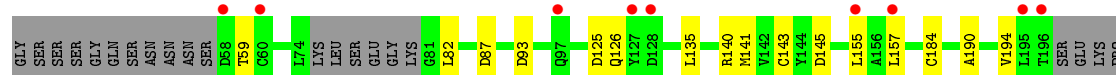
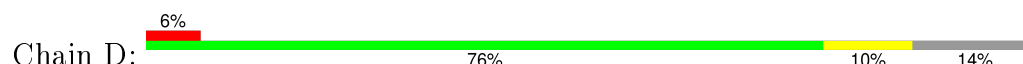
- Molecule 1: Integrin beta-1-binding protein 1



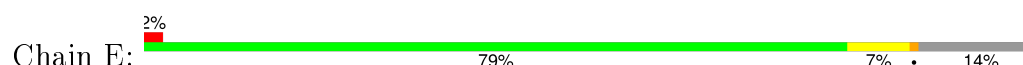
- Molecule 1: Integrin beta-1-binding protein 1



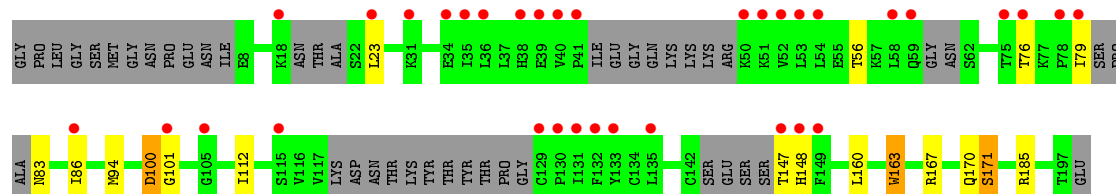
- Molecule 1: Integrin beta-1-binding protein 1



- Molecule 1: Integrin beta-1-binding protein 1



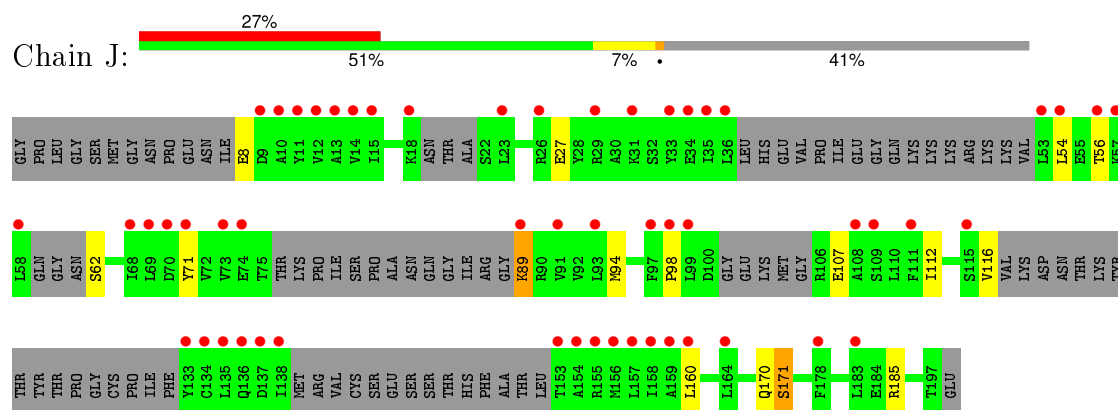
- Molecule 2: Krev interaction trapped protein 1



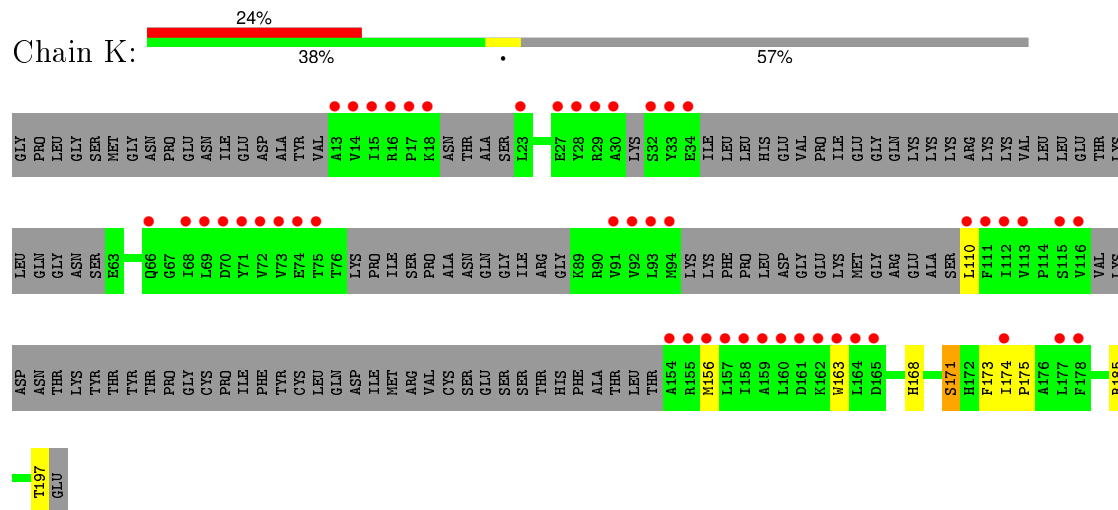
Chain I:



## Chain J:



Chain K:





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	153.90Å 157.98Å 152.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.54 48.79 – 2.54	Depositor EDS
% Data completeness (in resolution range)	100.0 (50.00-2.54) 99.3 (48.79-2.54)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.59 (at 2.54Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, $R_{free}$	0.220 , 0.252 0.218 , 0.249	Depositor DCC
$R_{free}$ test set	3099 reflections (5.35%)	DCC
Wilson B-factor (Å <sup>2</sup> )	83.6	Xtriage
Anisotropy	0.053	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 72.4	EDS
Estimated twinning fraction	0.034 for -k,-h,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 61239 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	8128	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	102.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BR

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	0.43	1/1098 (0.1%)	0.58	1/1486 (0.1%)
1	B	0.41	0/1079	0.54	0/1460
1	D	0.42	0/1056	0.53	0/1430
1	E	0.43	1/1056 (0.1%)	0.55	0/1430
2	H	0.41	1/1299 (0.1%)	0.53	0/1750
2	I	0.47	1/920 (0.1%)	0.49	0/1239
2	J	0.38	0/992	0.50	0/1334
2	K	0.42	1/728 (0.1%)	0.50	0/980
All	All	0.42	5/8228 (0.1%)	0.53	1/11109 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	I	22	SER	CB-OG	7.70	1.52	1.42
1	E	170	TRP	CD2-CE2	5.52	1.48	1.41
1	A	170	TRP	CD2-CE2	5.17	1.47	1.41
2	K	163	TRP	CD2-CE2	5.04	1.47	1.41
2	H	163	TRP	CD2-CE2	5.01	1.47	1.41

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	157	LEU	CA-CB-CG	5.98	129.06	115.30

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	1080	0	1070	5	0
1	B	1061	0	1057	6	0
1	D	1038	0	1028	10	0
1	E	1038	0	1028	8	0
2	H	1278	0	1322	19	0
2	I	906	0	953	5	0
2	J	979	0	1010	10	0
2	K	718	0	740	6	0
3	A	3	0	0	1	0
3	B	3	0	0	1	0
3	D	4	0	0	1	0
3	E	3	0	0	1	0
3	H	3	0	0	0	0
4	A	4	0	0	0	0
4	D	1	0	0	0	0
4	E	7	0	0	0	0
4	H	1	0	0	0	0
4	J	1	0	0	0	0
All	All	8128	0	8208	59	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 (59) 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:H:100:ASP:HB3	2:H:101:GLY:CA	1.77	1.15
2:H:100:ASP:HB3	2:H:101:GLY:HA3	1.30	1.09
2:H:147:THR:HA	2:H:148:HIS:HB2	1.48	0.93
1:D:140:ARG:NH1	3:D:302:BR:BR	2.61	0.88
1:E:140:ARG:NH1	3:E:302:BR:BR	2.62	0.87
1:A:140:ARG:NH1	3:A:302:BR:BR	2.73	0.76
2:H:147:THR:CA	2:H:148:HIS:HB2	2.17	0.73
2:H:100:ASP:HB3	2:H:101:GLY:HA2	1.71	0.73
2:H:100:ASP:CB	2:H:101:GLY:CA	2.61	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:140:ARG:NH1	3:B:302:BR:BR	2.77	0.72
1:D:190:ALA:O	1:D:194:VAL:HG13	1.90	0.71
1:D:143:CYS:HB2	1:D:155:LEU:HD23	1.76	0.67
2:J:94:MET:HE1	2:J:160:LEU:HA	1.78	0.65
2:H:83:ASN:N	2:J:62:SER:HG	1.96	0.64
1:E:93:ASP:OD2	2:K:185:ARG:NH2	2.31	0.64
1:E:160:THR:HG21	1:E:164:ASN:OD1	1.99	0.63
2:H:94:MET:HE1	2:H:160:LEU:HA	1.80	0.63
2:H:94:MET:HE2	2:H:112:ILE:HD11	1.83	0.60
2:I:170:GLN:O	2:I:171:SER:HB3	2.02	0.60
2:H:170:GLN:O	2:H:171:SER:HB3	2.02	0.60
2:H:100:ASP:CB	2:H:101:GLY:HA3	2.18	0.59
1:A:134:ALA:HB1	1:E:134:ALA:HB1	1.84	0.59
1:B:145:ASP:O	2:H:185:ARG:HD2	2.01	0.59
1:D:93:ASP:OD2	2:I:185:ARG:NH2	2.35	0.59
1:B:59:THR:HG21	1:B:116:LYS:HE2	1.83	0.59
1:B:160:THR:HG21	1:B:164:ASN:OD1	2.04	0.57
1:E:145:ASP:O	2:K:185:ARG:HD2	2.08	0.54
1:A:145:ASP:O	2:J:185:ARG:HD2	2.08	0.53
1:D:143:CYS:HB2	1:D:155:LEU:CD2	2.37	0.53
1:D:145:ASP:O	2:I:185:ARG:HD2	2.08	0.53
1:A:93:ASP:OD2	2:J:185:ARG:NH2	2.41	0.52
2:H:147:THR:HA	2:H:148:HIS:CB	2.27	0.52
1:D:135:LEU:CD2	1:D:194:VAL:HG11	2.41	0.51
2:J:170:GLN:O	2:J:171:SER:HB3	2.11	0.50
2:H:163:TRP:O	2:H:167:ARG:HG2	2.12	0.50
1:E:139:ILE:HD11	1:E:160:THR:HG23	1.95	0.48
2:J:89:LYS:HE2	2:J:116:VAL:HG23	1.95	0.48
2:I:65:THR:HG23	2:I:91:VAL:HB	1.95	0.48
1:D:135:LEU:HD23	1:D:194:VAL:HG11	1.95	0.47
2:H:94:MET:HE1	2:H:160:LEU:HD12	1.98	0.46
2:J:94:MET:CE	2:J:112:ILE:HD11	2.46	0.46
1:D:82:LEU:HD12	1:D:87:ASP:HB3	1.99	0.45
2:H:76:THR:HG22	2:H:76:THR:O	2.18	0.44
1:E:93:ASP:CG	2:K:185:ARG:HH22	2.21	0.44
2:H:76:THR:HG21	2:H:86:ILE:HG12	2.00	0.44
1:D:141:MET:CE	1:D:184:CYS:SG	3.06	0.44
2:H:76:THR:HG21	2:H:86:ILE:CG1	2.48	0.44
2:H:170:GLN:O	2:H:171:SER:CB	2.65	0.43
1:B:76:LEU:HD13	1:B:165:GLU:HA	2.01	0.42
2:J:98:PRO:HA	2:J:107:GLU:HG2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:129:CYS:HB3	2:I:130:PRO:HD3	2.00	0.42
2:K:174:ILE:HB	2:K:175:PRO:HD3	2.01	0.42
2:K:171:SER:C	2:K:173:PHE:H	2.23	0.42
2:J:54:LEU:HB3	2:J:71:TYR:HE2	1.85	0.42
1:A:56:ASN:HA	1:A:57:SER:HA	1.91	0.42
2:K:110:LEU:HD22	2:K:156:MET:HB3	2.01	0.42
1:E:160:THR:HG22	1:E:166:GLU:O	2.19	0.41
1:B:111:ILE:HD12	1:B:126:GLN:HG3	2.03	0.41
2:J:94:MET:HE2	2:J:112:ILE:HD11	2.02	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	135/154 (88%)	130 (96%)	5 (4%)	0	100	100
1	B	132/154 (86%)	128 (97%)	4 (3%)	0	100	100
1	D	129/154 (84%)	126 (98%)	2 (2%)	1 (1%)	24	40
1	E	129/154 (84%)	127 (98%)	2 (2%)	0	100	100
2	H	145/203 (71%)	135 (93%)	9 (6%)	1 (1%)	26	44
2	I	97/203 (48%)	94 (97%)	2 (2%)	1 (1%)	19	33
2	J	104/203 (51%)	96 (92%)	7 (7%)	1 (1%)	19	33
2	K	74/203 (36%)	71 (96%)	2 (3%)	1 (1%)	14	23
All	All	945/1428 (66%)	907 (96%)	33 (4%)	5 (0%)	34	54

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	126	GLN

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Mol	Chain	Res	Type
2	K	171	SER
2	H	100	ASP
2	I	171	SER
2	J	171	SER

### 5.3.2 Protein sidechains ⓘ

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	119/132 (90%)	115 (97%)	4 (3%)	44	70
1	B	117/132 (89%)	114 (97%)	3 (3%)	54	79
1	D	114/132 (86%)	111 (97%)	3 (3%)	54	79
1	E	114/132 (86%)	110 (96%)	4 (4%)	43	68
2	H	142/179 (79%)	138 (97%)	4 (3%)	51	76
2	I	102/179 (57%)	100 (98%)	2 (2%)	63	85
2	J	109/179 (61%)	105 (96%)	4 (4%)	41	66
2	K	79/179 (44%)	77 (98%)	2 (2%)	55	80
All	All	896/1244 (72%)	870 (97%)	26 (3%)	50	75

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	76	LEU
1	A	130	LEU
1	A	136	TYR
1	A	175	ASN
1	B	59	THR
1	B	160	THR
1	B	175	ASN
1	D	59	THR
1	D	125	ASP
1	D	157	LEU
1	E	125	ASP

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Mol	Chain	Res	Type
1	E	136	TYR
1	E	160	THR
1	E	175	ASN
2	H	23	LEU
2	H	56	THR
2	H	79	ILE
2	H	171	SER
2	I	76	THR
2	I	197	THR
2	J	8	GLU
2	J	27	GLU
2	J	56	THR
2	J	89	LYS
2	K	168	HIS
2	K	197	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	139/154 (90%)	0.85	10 (7%) 18 21	58, 73, 129, 152	0
1	B	136/154 (88%)	0.85	14 (10%) 9 9	57, 72, 118, 149	0
1	D	133/154 (86%)	0.78	9 (6%) 20 23	59, 73, 113, 158	0
1	E	133/154 (86%)	0.71	3 (2%) 64 69	58, 76, 113, 150	0
2	H	159/203 (78%)	1.16	34 (21%) 1 1	57, 91, 140, 162	0
2	I	111/203 (54%)	2.18	41 (36%) 0 0	57, 143, 208, 254	0
2	J	120/203 (59%)	2.77	54 (45%) 0 0	58, 143, 197, 219	0
2	K	88/203 (43%)	3.18	48 (54%) 0 0	58, 151, 210, 256	0
All	All	1019/1428 (71%)	1.44	213 (20%) 1 1	57, 87, 186, 256	0

All (213) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	K	33	TYR	15.3
2	J	135	LEU	13.6
2	J	54	LEU	10.4
2	J	156	MET	10.3
2	K	156	MET	10.2
2	J	11	TYR	10.1
2	K	158	ILE	10.1
2	K	69	LEU	10.1
2	I	111	PHE	9.9
2	K	157	LEU	9.6
2	J	10	ALA	9.3
2	I	12	VAL	8.9
2	I	110	LEU	8.9
2	J	53	LEU	8.8
2	J	12	VAL	8.8
2	K	14	VAL	8.7

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Mol	Chain	Res	Type	RSRZ
2	K	72	VAL	8.7
2	J	137	ASP	8.6
2	K	74	GLU	8.6
2	K	163	TRP	8.5
2	K	15	ILE	8.3
2	K	160	LEU	8.2
2	K	111	PHE	8.0
2	I	133	TYR	7.9
2	J	56	THR	7.9
2	J	155	ARG	7.8
2	I	156	MET	7.5
2	I	129	CYS	7.4
2	I	135	LEU	7.4
2	J	14	VAL	7.3
2	K	71	TYR	7.0
2	K	154	ALA	7.0
2	J	13	ALA	6.9
2	K	162	LYS	6.9
2	J	15	ILE	6.7
2	J	34	GLU	6.7
2	J	138	ILE	6.6
2	I	157	LEU	6.6
2	J	57	LYS	6.5
2	J	115	SER	6.5
2	J	36	LEU	6.4
1	D	127	TYR	6.2
2	J	74	GLU	6.0
2	I	158	ILE	6.0
2	K	23	LEU	6.0
1	B	127	TYR	5.9
2	J	154	ALA	5.9
2	J	157	LEU	5.8
2	I	71	TYR	5.7
2	J	99	LEU	5.7
2	H	40	VAL	5.7
2	J	33	TYR	5.6
2	I	155	ARG	5.6
2	K	112	ILE	5.6
2	K	91	VAL	5.5
2	H	50	LYS	5.4
2	I	35	ILE	5.4
2	K	159	ALA	5.4

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Mol	Chain	Res	Type	RSRZ
2	I	132	PHE	5.3
2	J	111	PHE	5.3
2	K	68	ILE	5.3
2	K	155	ARG	5.2
2	J	153	THR	5.2
2	I	33	TYR	5.2
1	A	56	ASN	5.1
2	H	149	PHE	5.1
2	K	110	LEU	5.1
2	K	94	MET	5.0
2	J	73	VAL	4.9
2	K	30	ALA	4.9
2	J	69	LEU	4.8
2	I	137	ASP	4.8
2	I	136	GLN	4.7
2	J	108	ALA	4.6
2	K	13	ALA	4.6
1	A	57	SER	4.6
2	K	17	PRO	4.5
2	J	134	CYS	4.4
2	J	133	TYR	4.4
2	J	91	VAL	4.3
2	H	133	TYR	4.2
2	H	132	PHE	4.2
1	B	110	PHE	4.1
2	I	13	ALA	4.1
2	I	76	THR	4.1
2	J	71	TYR	4.1
1	B	81	GLY	4.1
2	J	158	ILE	4.1
2	H	131	ILE	4.1
2	J	35	ILE	4.0
2	J	29	ARG	4.0
2	H	148	HIS	4.0
2	J	97	PHE	4.0
2	J	159	ALA	3.9
2	K	66	GLN	3.9
2	I	15	ILE	3.9
2	I	112	ILE	3.8
2	I	139	MET	3.8
2	H	36	LEU	3.8
2	K	34	GLU	3.7

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Mol	Chain	Res	Type	RSRZ
2	I	78	PRO	3.7
2	H	54	LEU	3.6
1	B	194	VAL	3.6
1	B	77	SER	3.6
2	J	89	LYS	3.6
2	J	9	ASP	3.6
1	A	76	LEU	3.5
2	I	131	ILE	3.5
2	H	51	LYS	3.5
1	A	77	SER	3.5
2	H	53	LEU	3.4
2	I	86	ILE	3.4
2	K	27	GLU	3.4
1	D	60	CYS	3.4
2	K	75	THR	3.3
2	J	136	GLN	3.3
2	H	115	SER	3.3
2	H	38	HIS	3.3
2	J	160	LEU	3.2
1	E	195	LEU	3.2
2	I	160	LEU	3.2
2	H	52	VAL	3.2
2	K	92	VAL	3.2
2	K	18	LYS	3.1
2	J	26	ARG	3.1
2	J	31	LYS	3.1
2	H	41	PRO	3.1
2	H	18	LYS	3.1
2	K	164	LEU	3.0
1	D	196	THR	3.0
2	K	73	VAL	3.0
2	K	113	VAL	3.0
2	H	76	THR	3.0
2	H	35	ILE	3.0
2	I	134	CYS	3.0
2	I	34	GLU	2.9
1	A	97	GLN	2.9
1	D	128	ASP	2.9
2	K	115	SER	2.9
2	J	70	ASP	2.9
2	J	109	SER	2.8
2	I	69	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
2	I	75	THR	2.8
2	H	130	PRO	2.8
2	I	92	VAL	2.8
2	K	174	ILE	2.8
2	I	23	LEU	2.7
2	K	165	ASP	2.7
2	H	34	GLU	2.7
2	I	31	LYS	2.7
2	J	23	LEU	2.7
1	B	76	LEU	2.6
2	K	32	SER	2.6
2	H	129	CYS	2.6
2	H	135	LEU	2.6
2	I	93	LEU	2.6
2	J	98	PRO	2.6
1	B	112	MET	2.6
2	K	161	ASP	2.6
2	H	58	LEU	2.6
2	H	75	THR	2.6
2	I	140	ARG	2.5
2	J	93	LEU	2.5
2	H	59	GLN	2.5
2	K	16	ARG	2.5
2	K	116	VAL	2.5
2	J	68	ILE	2.5
2	H	23	LEU	2.5
2	I	36	LEU	2.5
1	A	157	LEU	2.4
1	D	58	ASP	2.4
2	K	178	PHE	2.4
1	A	55	ASN	2.4
2	I	164	LEU	2.4
2	K	177	LEU	2.4
2	K	28	TYR	2.4
1	B	58	ASP	2.4
2	H	31	LYS	2.4
2	I	79	ILE	2.4
2	I	94	MET	2.4
2	I	138	ILE	2.4
2	H	86	ILE	2.3
2	J	18	LYS	2.3
1	D	157	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
2	H	147	THR	2.3
2	H	39	GLU	2.3
2	H	101	GLY	2.3
1	B	155	LEU	2.3
2	I	74	GLU	2.3
1	E	59	THR	2.3
1	B	64	ARG	2.3
1	B	107	GLU	2.2
2	H	105	GLY	2.2
1	A	127	TYR	2.2
2	H	79	ILE	2.2
2	K	29	ARG	2.2
1	D	97	GLN	2.2
1	A	194	VAL	2.2
1	B	124	SER	2.2
1	D	155	LEU	2.2
2	K	93	LEU	2.2
1	E	60	CYS	2.2
2	K	70	ASP	2.2
1	B	130	LEU	2.2
1	A	124	SER	2.1
1	B	111	ILE	2.1
2	J	164	LEU	2.1
1	D	195	LEU	2.1
2	J	183	LEU	2.1
2	J	58	LEU	2.0
2	I	77	LYS	2.0
2	H	78	PRO	2.0
2	J	178	PHE	2.0

## 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 [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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	BR	A	302	1/1	0.99	0.25	3.01	97,97,97,97	0
3	BR	B	302	1/1	0.97	0.21	0.64	105,105,105,105	0
3	BR	E	302	1/1	0.96	0.23	0.55	92,92,92,92	0
3	BR	D	302	1/1	0.98	0.19	0.14	95,95,95,95	0
3	BR	D	304	1/1	0.50	0.18	-0.20	166,166,166,166	0
3	BR	A	303	1/1	0.87	0.17	-0.45	162,162,162,162	0
3	BR	H	302	1/1	0.76	0.09	-	146,146,146,146	0
3	BR	D	303	1/1	0.96	0.11	-	133,133,133,133	0
3	BR	A	301	1/1	0.94	0.10	-	119,119,119,119	0
3	BR	H	303	1/1	0.78	0.07	-	161,161,161,161	0
3	BR	E	301	1/1	0.88	0.11	-	127,127,127,127	0
3	BR	H	301	1/1	0.94	0.17	-	135,135,135,135	0
3	BR	D	301	1/1	0.95	0.11	-	120,120,120,120	0
3	BR	E	303	1/1	0.95	0.34	-	139,139,139,139	1
3	BR	B	301	1/1	0.95	0.06	-	121,121,121,121	0
3	BR	B	303	1/1	0.76	0.12	-	146,146,146,146	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.