



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 01:03 PM GMT

PDB ID : 1A38  
Title : 14-3-3 PROTEIN ZETA BOUND TO R18 PEPTIDE  
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Deposited on : 1998-01-28  
Resolution : 3.35 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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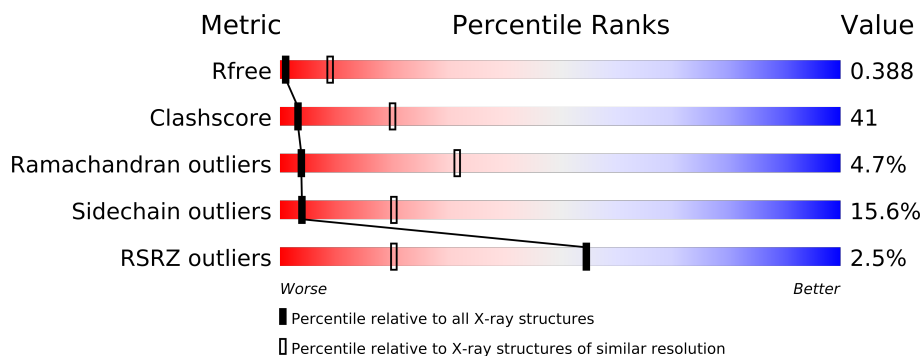
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 3.35 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))
$R_{free}$	66092	1141 (3.50-3.22)
Clashscore	79885	1030 (3.48-3.24)
Ramachandran outliers	78287	1008 (3.48-3.24)
Sidechain outliers	78261	1007 (3.48-3.24)
RSRZ outliers	66119	1141 (3.50-3.22)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	245	
1	B	245	
2	P	20	
2	Q	20	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3554 atoms, of which 0 are hydrogen and 0 are deuterium.

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 14-3-3 PROTEIN ZETA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	216	Total	C	N	O	S	0	0	0
			1730	1085	293	342	10			
1	B	216	Total	C	N	O	S	0	0	0
			1730	1085	293	342	10			

- Molecule 2 is a protein called R18 PEPTIDE (PHCVPRDLSWLDLEANMCLP).

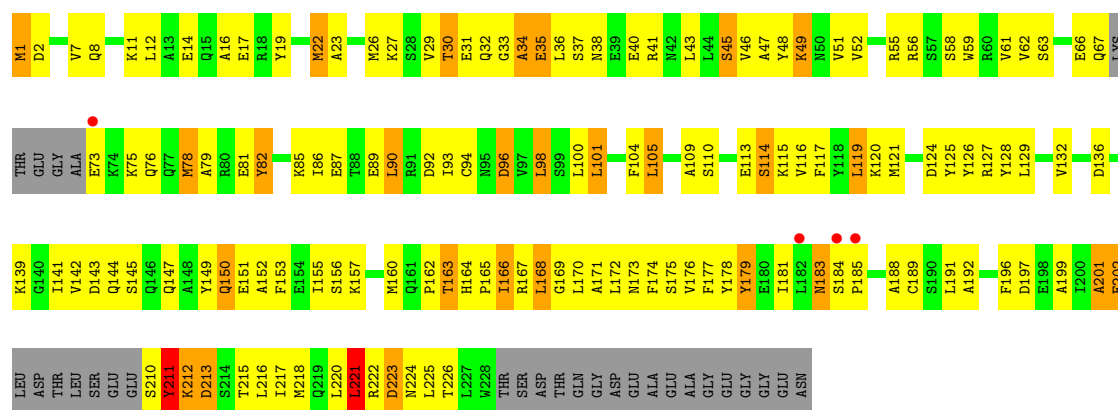
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	P	5	Total	C	N	O	0	0	0
			47	32	6	9			
2	Q	5	Total	C	N	O	0	0	0
			47	32	6	9			

### 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.

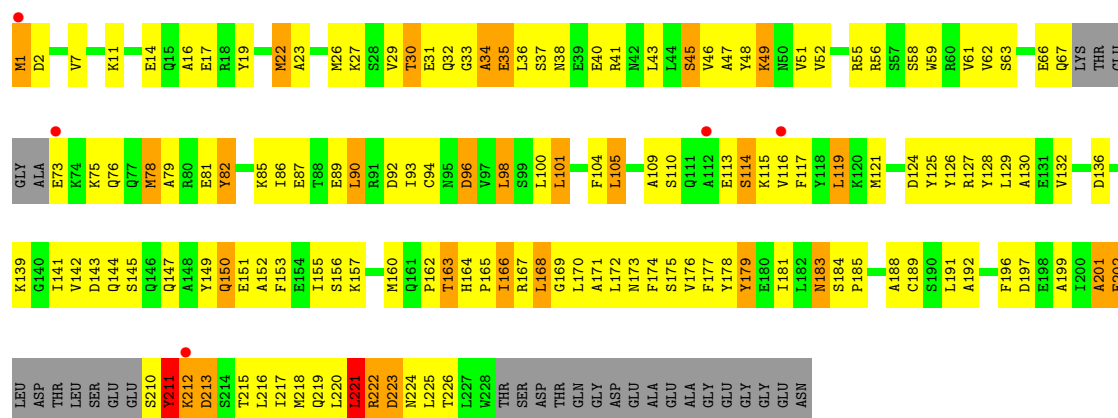
#### • Molecule 1: 14-3-3 PROTEIN ZETA

Chain A: 



#### • Molecule 1: 14-3-3 PROTEIN ZETA

Chain B: 



#### • Molecule 2: R18 PEPTIDE (PHCVPRDLSWLDLEANMCLP)

Chain P: 



#### • Molecule 2: R18 PEPTIDE (PHCVPRDLSWLDLEANMCLP)

Chain Q: 

PHE	HIS	CYS	VAL	PRO	ARG	ASP	LEU	SER	W1	L2	D3	L4	E5	ALA	ASN	MET	CYS	LEU	PRO
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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	95.52Å 95.52Å 235.36Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 3.35 29.04 – 3.10	Depositor EDS
% Data completeness (in resolution range)	96.8 (20.00-3.35) 95.4 (29.04-3.10)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.52 (at 3.11Å)	Xtriage
Refinement program	X-PLOR 3.8	Depositor
R, $R_{free}$	0.330 , 0.390 0.370 , 0.388	Depositor DCC
$R_{free}$ test set	2087 reflections (11.05%)	DCC
Wilson B-factor (Å <sup>2</sup> )	64.4	Xtriage
Anisotropy	0.511	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 27.4	EDS
Estimated twinning fraction	0.449 for h,-h-k,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.42$ , $\langle L^2 \rangle = 0.25$	Xtriage
Outliers	0 of 20975 reflections	Xtriage
$F_o, F_c$ correlation	0.68	EDS
Total number of atoms	3554	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.09% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 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	0.55	0/1751	0.92	3/2348 (0.1%)
1	B	0.55	0/1751	0.92	3/2348 (0.1%)
2	P	0.71	0/48	0.88	0/65
2	Q	0.72	0/48	0.88	0/65
All	All	0.55	0/3598	0.92	6/4826 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	221	LEU	CA-CB-CG	7.83	133.32	115.30
1	A	221	LEU	CA-CB-CG	7.83	133.31	115.30
1	B	202	GLU	N-CA-C	5.43	125.67	111.00
1	A	202	GLU	N-CA-C	5.42	125.64	111.00
1	B	19	TYR	N-CA-C	5.31	125.34	111.00
1	A	19	TYR	N-CA-C	5.30	125.32	111.00

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1730	0	1726	144	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1730	0	1726	145	0
2	P	47	0	44	7	0
2	Q	47	0	44	7	0
All	All	3554	0	3540	290	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 41.

All (290) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:220:LEU:HD23	2:P:4:LEU:HD11	1.49	0.94
1:B:220:LEU:HD23	2:Q:4:LEU:HD11	1.49	0.92
1:A:89:GLU:O	1:A:93:ILE:HG12	1.77	0.85
1:B:89:GLU:O	1:B:93:ILE:HG12	1.77	0.84
1:B:178:TYR:HB2	1:B:188:ALA:HB2	1.59	0.84
1:A:178:TYR:HB2	1:A:188:ALA:HB2	1.59	0.83
1:B:30:THR:OG1	1:B:100:LEU:HD23	1.78	0.83
1:A:30:THR:OG1	1:A:100:LEU:HD23	1.78	0.82
1:A:52:VAL:O	1:A:56:ARG:HG3	1.80	0.82
1:B:52:VAL:O	1:B:56:ARG:HG3	1.80	0.81
1:A:55:ARG:HB3	1:A:90:LEU:HD22	1.64	0.80
1:A:26:MET:O	1:A:30:THR:HG23	1.83	0.79
1:B:101:LEU:HD12	1:B:105:LEU:HD23	1.63	0.79
1:A:101:LEU:HD12	1:A:105:LEU:HD23	1.63	0.79
1:B:55:ARG:HB3	1:B:90:LEU:HD22	1.64	0.79
1:A:48:TYR:O	1:A:52:VAL:HG12	1.83	0.79
1:B:26:MET:O	1:B:30:THR:HG23	1.83	0.78
1:B:48:TYR:O	1:B:52:VAL:HG12	1.83	0.78
1:B:177:PHE:O	1:B:181:ILE:HB	1.83	0.78
1:B:75:LYS:O	1:B:78:MET:HG3	1.84	0.78
1:A:75:LYS:O	1:A:78:MET:HG3	1.84	0.78
1:A:177:PHE:O	1:A:181:ILE:HB	1.83	0.76
1:B:52:VAL:HG11	1:B:125:TYR:CE1	2.23	0.73
1:A:217:ILE:O	1:A:221:LEU:HD23	1.89	0.73
1:B:7:VAL:O	1:B:11:LYS:HG2	1.88	0.72
1:A:7:VAL:O	1:A:11:LYS:HG2	1.88	0.72
1:B:151:GLU:O	1:B:155:ILE:HG13	1.89	0.72
1:A:151:GLU:O	1:A:155:ILE:HG13	1.89	0.72
1:A:52:VAL:HG11	1:A:125:TYR:CE1	2.23	0.72
1:B:217:ILE:O	1:B:221:LEU:HD23	1.89	0.72
1:B:100:LEU:HD12	1:B:100:LEU:H	1.57	0.70
1:A:52:VAL:HG11	1:A:125:TYR:HE1	1.58	0.69

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:100:LEU:H	1:A:100:LEU:HD12	1.57	0.69
1:B:220:LEU:CD2	2:Q:4:LEU:HD11	2.22	0.69
1:B:116:VAL:HG23	1:B:152:ALA:HB1	1.75	0.69
1:A:220:LEU:CD2	2:P:4:LEU:HD11	2.22	0.68
1:A:116:VAL:HG23	1:A:152:ALA:HB1	1.75	0.68
1:B:52:VAL:HG11	1:B:125:TYR:HE1	1.58	0.68
1:B:115:LYS:O	1:B:119:LEU:HD12	1.95	0.67
1:A:115:LYS:O	1:A:119:LEU:HD12	1.95	0.67
1:A:58:SER:O	1:A:62:VAL:HG12	1.95	0.66
1:B:58:SER:O	1:B:62:VAL:HG12	1.95	0.66
1:B:165:PRO:HG3	1:B:213:ASP:HB2	1.77	0.66
1:B:173:ASN:O	1:B:176:VAL:HG12	1.96	0.66
1:A:173:ASN:O	1:A:176:VAL:HG12	1.96	0.66
1:A:101:LEU:CD1	1:A:105:LEU:HD23	2.25	0.65
1:B:101:LEU:CD1	1:B:105:LEU:HD23	2.25	0.65
1:A:165:PRO:HG3	1:A:213:ASP:HB2	1.77	0.65
1:B:43:LEU:O	1:B:46:VAL:HG12	1.98	0.64
1:B:22:MET:SD	1:B:47:ALA:HB2	2.38	0.64
1:B:45:SER:HG	1:B:117:PHE:HZ	1.46	0.63
1:A:101:LEU:HA	1:A:105:LEU:HB2	1.81	0.63
1:A:35:GLU:OE1	1:A:35:GLU:HA	1.99	0.63
2:P:1:TRP:HD1	2:P:2:LEU:H	1.46	0.63
1:A:22:MET:SD	1:A:47:ALA:HB2	2.38	0.63
1:B:101:LEU:HA	1:B:105:LEU:HB2	1.81	0.63
1:A:185:PRO:O	1:A:189:CYS:SG	2.55	0.62
1:B:101:LEU:HD21	1:B:121:MET:HG2	1.80	0.62
1:A:143:ASP:O	1:A:147:GLN:HG3	2.00	0.62
1:A:43:LEU:O	1:A:46:VAL:HG12	1.98	0.62
1:A:192:ALA:HB1	1:A:225:LEU:HD11	1.82	0.62
1:B:143:ASP:O	1:B:147:GLN:HG3	2.00	0.62
1:B:59:TRP:HE1	1:B:87:GLU:HG3	1.65	0.62
1:B:192:ALA:HB1	1:B:225:LEU:HD11	1.82	0.61
1:A:101:LEU:HD21	1:A:121:MET:HG2	1.80	0.61
1:A:59:TRP:HE1	1:A:87:GLU:HG3	1.65	0.61
1:A:94:CYS:O	1:A:98:LEU:HB2	2.00	0.61
1:B:94:CYS:O	1:B:98:LEU:HB2	2.00	0.61
2:Q:1:TRP:HD1	2:Q:2:LEU:H	1.47	0.61
1:B:35:GLU:HA	1:B:35:GLU:OE1	1.99	0.61
1:B:66:GLU:HB2	1:B:79:ALA:HB1	1.81	0.61
1:B:185:PRO:O	1:B:189:CYS:SG	2.55	0.61
1:A:201:ALA:O	1:A:202:GLU:HB2	2.01	0.61
1:A:116:VAL:HA	1:A:152:ALA:HB1	1.83	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:66:GLU:HB2	1:A:79:ALA:HB1	1.81	0.61
1:B:116:VAL:HA	1:B:152:ALA:HB1	1.83	0.61
1:A:51:VAL:HG13	1:A:55:ARG:NH1	2.17	0.59
1:A:216:LEU:O	1:A:220:LEU:HD13	2.03	0.59
1:A:124:ASP:O	1:A:127:ARG:HB3	2.03	0.59
1:B:201:ALA:O	1:B:202:GLU:HB2	2.01	0.59
1:B:51:VAL:HG13	1:B:55:ARG:NH1	2.17	0.59
1:B:192:ALA:CB	1:B:225:LEU:HD11	2.33	0.59
1:B:124:ASP:O	1:B:127:ARG:HB3	2.03	0.59
1:B:216:LEU:O	1:B:220:LEU:HD13	2.02	0.59
1:A:45:SER:HG	1:A:117:PHE:HZ	1.49	0.58
1:A:169:GLY:HA3	2:P:2:LEU:HD11	1.84	0.58
1:A:192:ALA:CB	1:A:225:LEU:HD11	2.33	0.58
1:B:162:PRO:C	1:B:164:HIS:H	2.07	0.58
1:B:169:GLY:HA3	2:Q:2:LEU:HD11	1.84	0.58
1:B:101:LEU:HD21	1:B:121:MET:HB3	1.86	0.58
1:B:213:ASP:O	1:B:217:ILE:HG13	2.04	0.58
1:A:8:GLN:HE21	1:B:78:MET:HE2	1.69	0.58
1:A:30:THR:C	1:A:32:GLN:H	2.07	0.57
1:B:73:GLU:O	1:B:76:GLN:HB3	2.04	0.57
1:A:101:LEU:HD21	1:A:121:MET:HB3	1.86	0.57
1:A:139:LYS:HA	1:A:142:VAL:HG12	1.86	0.57
1:B:87:GLU:HG2	1:B:132:VAL:HG13	1.87	0.57
1:A:38:ASN:HA	1:A:41:ARG:HB3	1.87	0.57
1:A:163:THR:O	1:A:163:THR:HG22	2.05	0.57
1:A:162:PRO:C	1:A:164:HIS:H	2.07	0.56
1:A:73:GLU:O	1:A:76:GLN:HB3	2.04	0.56
1:A:213:ASP:O	1:A:217:ILE:HG13	2.04	0.56
1:B:38:ASN:HA	1:B:41:ARG:HB3	1.87	0.56
1:A:129:LEU:HB3	1:A:141:ILE:HG21	1.87	0.56
1:B:212:LYS:O	1:B:215:THR:HB	2.06	0.56
1:A:87:GLU:HG2	1:A:132:VAL:HG13	1.87	0.56
1:B:30:THR:C	1:B:32:GLN:H	2.07	0.56
1:A:150:GLN:O	1:A:150:GLN:HG3	2.05	0.56
1:B:163:THR:HG22	1:B:163:THR:O	2.05	0.56
1:B:139:LYS:HA	1:B:142:VAL:HG12	1.86	0.56
1:A:212:LYS:O	1:A:215:THR:HB	2.06	0.56
1:B:129:LEU:HB3	1:B:141:ILE:HG21	1.87	0.55
1:B:150:GLN:O	1:B:150:GLN:HG3	2.05	0.55
1:B:59:TRP:NE1	1:B:87:GLU:HG3	2.22	0.55
1:B:170:LEU:HD12	1:B:174:PHE:HB2	1.88	0.54
1:A:170:LEU:HD12	1:A:174:PHE:HB2	1.88	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:59:TRP:NE1	1:A:87:GLU:HG3	2.22	0.53
1:A:124:ASP:HA	1:A:149:TYR:OH	2.08	0.53
1:A:162:PRO:O	1:A:168:LEU:HD13	2.08	0.53
1:B:168:LEU:HB3	1:B:217:ILE:HG21	1.90	0.53
1:B:124:ASP:HA	1:B:149:TYR:OH	2.08	0.53
1:A:168:LEU:HB3	1:A:217:ILE:HG21	1.90	0.53
1:B:169:GLY:O	1:B:172:LEU:HB3	2.08	0.53
1:B:162:PRO:O	1:B:168:LEU:HD13	2.08	0.53
1:B:116:VAL:HG21	1:B:156:SER:HB3	1.91	0.52
1:B:87:GLU:CD	1:B:132:VAL:HG13	2.30	0.52
1:A:126:TYR:CD1	1:A:144:GLN:HB3	2.45	0.52
1:B:126:TYR:CE1	1:B:144:GLN:HB3	2.45	0.52
1:B:87:GLU:HG2	1:B:132:VAL:CG1	2.40	0.52
1:A:87:GLU:CD	1:A:132:VAL:HG13	2.30	0.52
1:A:87:GLU:HG2	1:A:132:VAL:CG1	2.40	0.52
1:A:116:VAL:HG21	1:A:156:SER:HB3	1.91	0.52
1:A:61:VAL:HG13	1:A:62:VAL:H	1.75	0.52
1:B:61:VAL:HG13	1:B:62:VAL:H	1.75	0.52
1:A:169:GLY:O	1:A:172:LEU:HB3	2.09	0.52
1:A:126:TYR:CE1	1:A:144:GLN:HB3	2.45	0.52
1:B:87:GLU:CG	1:B:132:VAL:HG13	2.40	0.52
1:A:125:TYR:HA	1:A:128:TYR:CD2	2.45	0.52
1:A:87:GLU:CG	1:A:132:VAL:HG13	2.40	0.52
1:B:126:TYR:CD1	1:B:144:GLN:HB3	2.45	0.52
1:B:109:ALA:HB1	1:B:114:SER:HB2	1.93	0.51
1:A:98:LEU:HD11	1:A:126:TYR:CD2	2.46	0.51
1:B:98:LEU:HD11	1:B:126:TYR:CD2	2.46	0.51
1:A:14:GLU:HB2	1:A:22:MET:SD	2.51	0.51
1:B:14:GLU:HB2	1:B:22:MET:SD	2.51	0.50
1:A:109:ALA:HB1	1:A:114:SER:HB2	1.93	0.50
1:B:125:TYR:HA	1:B:128:TYR:CD2	2.45	0.50
1:A:8:GLN:CD	1:B:75:LYS:HD2	2.31	0.50
1:B:14:GLU:C	1:B:16:ALA:H	2.15	0.50
1:B:153:PHE:O	1:B:157:LYS:HB2	2.12	0.50
1:A:14:GLU:C	1:A:16:ALA:H	2.15	0.49
1:A:192:ALA:O	1:A:196:PHE:HB3	2.13	0.49
1:A:178:TYR:HD1	1:A:184:SER:HB3	1.77	0.49
1:A:76:GLN:O	1:A:79:ALA:HB3	2.13	0.49
1:A:152:ALA:HA	1:A:155:ILE:HB	1.94	0.49
1:B:76:GLN:O	1:B:79:ALA:HB3	2.13	0.49
1:B:162:PRO:O	1:B:168:LEU:HD22	2.13	0.48
1:B:173:ASN:HA	1:B:176:VAL:HG12	1.95	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:8:GLN:HE21	1:B:78:MET:CE	2.26	0.48
1:A:45:SER:O	1:A:49:LYS:HB3	2.14	0.48
1:A:220:LEU:H	1:A:220:LEU:CD1	2.27	0.48
1:B:178:TYR:HD1	1:B:184:SER:HB3	1.77	0.48
1:B:63:SER:O	1:B:67:GLN:N	2.45	0.48
1:B:45:SER:O	1:B:49:LYS:HB3	2.14	0.48
1:B:152:ALA:HA	1:B:155:ILE:HB	1.94	0.48
1:B:213:ASP:HA	1:B:216:LEU:HD12	1.96	0.48
1:A:153:PHE:O	1:A:157:LYS:HB2	2.12	0.48
1:B:23:ALA:HB1	1:B:48:TYR:CE2	2.49	0.48
1:B:223:ASP:O	1:B:226:THR:HB	2.14	0.48
1:B:221:LEU:O	1:B:224:ASN:HB2	2.14	0.48
1:B:172:LEU:O	1:B:175:SER:HB3	2.14	0.48
1:B:192:ALA:O	1:B:196:PHE:HB3	2.13	0.47
1:A:211:TYR:CD1	1:A:212:LYS:N	2.82	0.47
1:A:63:SER:O	1:A:67:GLN:N	2.45	0.47
1:A:23:ALA:HB1	1:A:48:TYR:CE2	2.49	0.47
1:A:213:ASP:HA	1:A:216:LEU:HD12	1.96	0.47
1:A:117:PHE:HB2	1:A:166:ILE:HD11	1.96	0.47
1:A:162:PRO:O	1:A:168:LEU:HD22	2.13	0.47
1:B:61:VAL:HG13	1:B:62:VAL:N	2.29	0.47
1:A:221:LEU:O	1:A:224:ASN:HB2	2.14	0.47
1:B:220:LEU:CD1	1:B:220:LEU:H	2.27	0.47
1:A:173:ASN:HA	1:A:176:VAL:HG12	1.95	0.47
1:B:117:PHE:HB2	1:B:166:ILE:HD11	1.96	0.47
1:A:172:LEU:O	1:A:175:SER:HB3	2.14	0.47
1:B:211:TYR:CD1	1:B:212:LYS:N	2.82	0.47
2:P:1:TRP:CD1	2:P:2:LEU:N	2.80	0.47
1:A:217:ILE:O	1:A:221:LEU:CD2	2.61	0.47
2:Q:1:TRP:CD1	2:Q:2:LEU:N	2.80	0.46
1:A:220:LEU:N	1:A:220:LEU:CD1	2.78	0.46
1:B:220:LEU:N	1:B:220:LEU:CD1	2.78	0.46
1:A:223:ASP:O	1:A:226:THR:HB	2.14	0.46
1:A:61:VAL:HG13	1:A:62:VAL:N	2.29	0.46
1:A:169:GLY:HA3	2:P:2:LEU:CD1	2.46	0.46
1:B:217:ILE:O	1:B:221:LEU:CD2	2.61	0.46
1:A:82:TYR:O	1:A:86:ILE:HG13	2.16	0.45
1:B:31:GLU:HA	1:B:104:PHE:CZ	2.51	0.45
1:B:145:SER:HB3	1:B:177:PHE:CE1	2.51	0.45
2:Q:2:LEU:HG	2:Q:3:ASP:N	2.31	0.45
1:A:31:GLU:HA	1:A:104:PHE:CZ	2.51	0.45
1:A:27:LYS:HA	1:A:100:LEU:HD21	1.98	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:116:VAL:HA	1:A:152:ALA:CB	2.47	0.45
1:B:92:ASP:O	1:B:96:ASP:HB2	2.17	0.45
1:B:82:TYR:O	1:B:86:ILE:HG13	2.16	0.45
1:B:37:SER:HB3	1:B:40:GLU:HG3	1.98	0.45
1:B:101:LEU:HD21	1:B:121:MET:CG	2.44	0.45
1:A:37:SER:HB3	1:A:40:GLU:HG3	1.98	0.45
1:A:145:SER:HB3	1:A:177:PHE:CE1	2.51	0.45
1:A:1:MET:O	1:A:2:ASP:HB3	2.17	0.45
1:A:92:ASP:O	1:A:96:ASP:HB2	2.16	0.45
1:A:119:LEU:HD13	1:A:155:ILE:HD12	1.98	0.45
1:B:27:LYS:HA	1:B:100:LEU:HD21	1.98	0.45
1:A:101:LEU:HD12	1:A:101:LEU:HA	1.75	0.44
1:A:211:TYR:CD1	1:A:211:TYR:C	2.90	0.44
1:A:101:LEU:HD21	1:A:121:MET:CB	2.47	0.44
1:A:101:LEU:HD21	1:A:121:MET:CG	2.44	0.44
1:B:171:ALA:O	1:B:175:SER:HB2	2.18	0.44
1:B:1:MET:O	1:B:2:ASP:HB3	2.17	0.44
1:B:211:TYR:CD1	1:B:211:TYR:C	2.90	0.44
1:A:30:THR:HG22	1:A:36:LEU:HD21	2.00	0.44
2:P:2:LEU:HG	2:P:3:ASP:N	2.31	0.44
1:A:142:VAL:HG23	1:A:181:ILE:CG2	2.48	0.44
1:A:119:LEU:HD22	1:A:151:GLU:HB3	1.99	0.44
1:B:45:SER:OG	1:B:117:PHE:HZ	2.01	0.44
1:A:171:ALA:O	1:A:175:SER:HB2	2.18	0.44
1:A:211:TYR:HD1	1:A:211:TYR:C	2.21	0.44
1:B:142:VAL:HG23	1:B:181:ILE:CG2	2.48	0.44
1:B:119:LEU:HD22	1:B:151:GLU:HB3	1.98	0.44
1:A:46:VAL:HG13	1:A:47:ALA:N	2.33	0.44
1:B:119:LEU:HD13	1:B:155:ILE:HD12	1.98	0.44
1:B:46:VAL:HG13	1:B:47:ALA:N	2.33	0.44
1:A:59:TRP:CE2	1:A:132:VAL:HG22	2.52	0.44
1:B:169:GLY:HA3	2:Q:2:LEU:CD1	2.46	0.43
1:B:116:VAL:HA	1:B:152:ALA:CB	2.47	0.43
1:B:59:TRP:CE2	1:B:132:VAL:HG22	2.52	0.43
1:B:101:LEU:HD21	1:B:121:MET:CB	2.47	0.43
1:A:174:PHE:CD1	1:A:174:PHE:O	2.72	0.43
1:B:33:GLY:O	1:B:34:ALA:O	2.37	0.43
1:B:96:ASP:O	1:B:100:LEU:CD1	2.67	0.43
1:B:30:THR:HG22	1:B:36:LEU:HD21	2.00	0.43
1:A:12:LEU:HG	1:B:82:TYR:HD2	1.82	0.43
1:B:142:VAL:O	1:B:177:PHE:HZ	2.02	0.43
1:A:142:VAL:O	1:A:177:PHE:HZ	2.02	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:178:TYR:O	1:A:183:ASN:N	2.52	0.43
1:A:96:ASP:O	1:A:100:LEU:CD1	2.67	0.43
1:B:211:TYR:HD1	1:B:211:TYR:C	2.21	0.42
1:A:220:LEU:H	1:A:220:LEU:HD13	1.85	0.42
1:B:119:LEU:CD2	1:B:151:GLU:HB3	2.49	0.42
1:A:119:LEU:CD2	1:A:151:GLU:HB3	2.49	0.42
1:B:174:PHE:CD1	1:B:174:PHE:O	2.72	0.42
1:A:196:PHE:O	1:A:199:ALA:HB3	2.19	0.42
1:B:196:PHE:O	1:B:199:ALA:HB3	2.19	0.42
1:A:174:PHE:HD1	1:A:174:PHE:O	2.02	0.42
1:B:85:LYS:O	1:B:89:GLU:HG3	2.20	0.42
1:A:46:VAL:CG1	1:A:47:ALA:N	2.83	0.42
1:B:221:LEU:O	1:B:225:LEU:HD13	2.20	0.42
1:B:174:PHE:HD1	1:B:174:PHE:O	2.02	0.42
1:A:162:PRO:C	1:A:164:HIS:N	2.72	0.42
1:A:162:PRO:HG3	1:A:167:ARG:CZ	2.50	0.42
1:B:162:PRO:HG3	1:B:167:ARG:CZ	2.50	0.42
1:B:179:TYR:CD1	1:B:179:TYR:C	2.93	0.42
1:B:220:LEU:HD13	1:B:220:LEU:H	1.85	0.42
1:A:33:GLY:O	1:A:34:ALA:O	2.37	0.42
1:A:164:HIS:HA	1:A:165:PRO:HD2	1.76	0.42
1:A:221:LEU:O	1:A:225:LEU:HD13	2.20	0.42
1:B:165:PRO:CB	1:B:217:ILE:HD11	2.50	0.42
1:B:126:TYR:HD1	1:B:141:ILE:HG23	1.85	0.42
1:B:46:VAL:CG1	1:B:47:ALA:N	2.83	0.41
1:A:165:PRO:CB	1:A:217:ILE:HD11	2.50	0.41
1:B:178:TYR:O	1:B:183:ASN:N	2.52	0.41
1:A:126:TYR:HD1	1:A:141:ILE:HG23	1.85	0.41
1:A:45:SER:OG	1:A:117:PHE:HZ	2.01	0.41
1:A:179:TYR:C	1:A:179:TYR:CD1	2.93	0.41
1:A:221:LEU:O	1:A:225:LEU:CD1	2.68	0.41
1:B:221:LEU:O	1:B:225:LEU:CD1	2.68	0.41
1:B:219:GLN:O	1:B:222:ARG:HB3	2.21	0.41
1:A:85:LYS:O	1:A:89:GLU:HG3	2.20	0.41
1:B:162:PRO:C	1:B:164:HIS:N	2.72	0.41
1:B:210:SER:O	1:B:211:TYR:HB2	2.21	0.41
1:A:210:SER:O	1:A:211:TYR:HB2	2.21	0.41
1:A:12:LEU:HG	1:B:82:TYR:CD2	2.56	0.41
1:A:165:PRO:HB3	1:A:217:ILE:CD1	2.51	0.41
1:B:165:PRO:HB3	1:B:217:ILE:CD1	2.51	0.41
1:B:127:ARG:O	1:B:130:ALA:HB3	2.21	0.41
1:B:38:ASN:O	1:B:41:ARG:HB3	2.21	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:120:LYS:HB3	1:A:120:LYS:HE2	1.94	0.41
1:B:105:LEU:HD12	1:B:105:LEU:HA	1.93	0.40
1:A:225:LEU:N	1:A:225:LEU:HD12	2.37	0.40
1:A:27:LYS:HA	1:A:100:LEU:CD2	2.51	0.40
1:B:101:LEU:HA	1:B:101:LEU:HD12	1.76	0.40
1:B:31:GLU:HA	1:B:104:PHE:CE2	2.57	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	210/245 (86%)	152 (72%)	48 (23%)	10 (5%)	4	31
1	B	210/245 (86%)	152 (72%)	48 (23%)	10 (5%)	4	31
2	P	3/20 (15%)	2 (67%)	1 (33%)	0	100	100
2	Q	3/20 (15%)	2 (67%)	1 (33%)	0	100	100
All	All	426/530 (80%)	308 (72%)	98 (23%)	20 (5%)	4	31

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	34	ALA
1	A	166	ILE
1	A	211	TYR
1	B	34	ALA
1	B	166	ILE
1	B	211	TYR
1	A	201	ALA
1	A	212	LYS
1	B	201	ALA
1	B	212	LYS
1	A	17	GLU
1	A	35	GLU

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Mol	Chain	Res	Type
1	B	17	GLU
1	B	35	GLU
1	A	113	GLU
1	A	163	THR
1	A	197	ASP
1	B	113	GLU
1	B	163	THR
1	B	197	ASP

### 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	187/209 (90%)	157 (84%)	30 (16%)	3	19
1	B	187/209 (90%)	157 (84%)	30 (16%)	3	19
2	P	5/19 (26%)	5 (100%)	0	100	100
2	Q	5/19 (26%)	5 (100%)	0	100	100
All	All	384/456 (84%)	324 (84%)	60 (16%)	4	20

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	22	MET
1	A	29	VAL
1	A	30	THR
1	A	45	SER
1	A	49	LYS
1	A	78	MET
1	A	81	GLU
1	A	82	TYR
1	A	90	LEU
1	A	96	ASP
1	A	98	LEU
1	A	101	LEU
1	A	105	LEU

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Mol	Chain	Res	Type
1	A	110	SER
1	A	114	SER
1	A	119	LEU
1	A	136	ASP
1	A	150	GLN
1	A	160	MET
1	A	168	LEU
1	A	179	TYR
1	A	183	ASN
1	A	191	LEU
1	A	211	TYR
1	A	213	ASP
1	A	218	MET
1	A	221	LEU
1	A	222	ARG
1	A	223	ASP
1	B	1	MET
1	B	22	MET
1	B	29	VAL
1	B	30	THR
1	B	45	SER
1	B	49	LYS
1	B	78	MET
1	B	81	GLU
1	B	82	TYR
1	B	90	LEU
1	B	96	ASP
1	B	98	LEU
1	B	101	LEU
1	B	105	LEU
1	B	110	SER
1	B	114	SER
1	B	119	LEU
1	B	136	ASP
1	B	150	GLN
1	B	160	MET
1	B	168	LEU
1	B	179	TYR
1	B	183	ASN
1	B	191	LEU
1	B	211	TYR
1	B	213	ASP

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Mol	Chain	Res	Type
1	B	218	MET
1	B	221	LEU
1	B	222	ARG
1	B	223	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	32	GLN
1	A	67	GLN
1	A	76	GLN
1	A	111	GLN
1	A	146	GLN
1	A	164	HIS
1	A	183	ASN
1	B	8	GLN
1	B	32	GLN
1	B	67	GLN
1	B	76	GLN
1	B	111	GLN
1	B	146	GLN
1	B	164	HIS
1	B	183	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains 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<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	216/245 (88%)	0.47	4 (1%) 64 27	16, 54, 85, 94	0
1	B	216/245 (88%)	0.43	5 (2%) 57 22	16, 54, 85, 94	0
2	P	5/20 (25%)	1.31	1 (20%) 2 2	73, 77, 82, 83	0
2	Q	5/20 (25%)	0.91	1 (20%) 2 2	73, 77, 82, 83	0
All	All	442/530 (83%)	0.46	11 (2%) 54 20	16, 54, 85, 94	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	73	GLU	4.2
1	B	112	ALA	3.7
1	A	184	SER	3.1
1	B	116	VAL	2.4
1	B	212	LYS	2.4
2	P	5	GLU	2.4
1	A	185	PRO	2.3
1	A	73	GLU	2.3
1	A	182	LEU	2.2
1	B	1	MET	2.1
2	Q	4	LEU	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 ⓘ

There are no ligands in this entry.

## 6.5 Other polymers ⓘ

There are no such residues in this entry.