



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:36 AM GMT

PDB ID : 2RFD  
Title : Crystal structure of the complex between the EGFR kinase domain and a Mig6 peptide  
Authors : Zhang, X.; Pickin, K.A.; Bose, R.; Jura, N.; Cole, P.A.; Kuriyan, J.  
Deposited on : 2007-09-28  
Resolution : 3.60 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

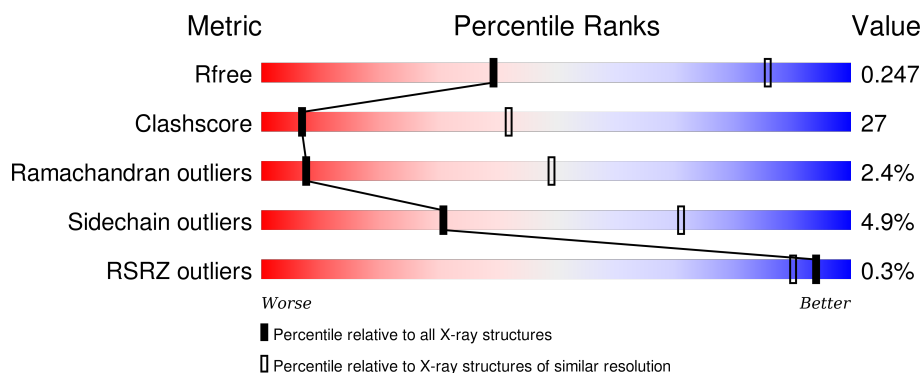
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.60 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	1408 (3.80-3.40)
Clashscore	102246	1010 (3.74-3.46)
Ramachandran outliers	100387	1007 (3.76-3.44)
Sidechain outliers	100360	1007 (3.76-3.44)
RSRZ outliers	91569	1003 (3.78-3.42)

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	324	<div> <div>46%</div> <div>39%</div> <div>•</div> <div>11%</div> </div>
1	B	324	<div> <div>44%</div> <div>42%</div> <div>•</div> <div>12%</div> </div>
2	C	25	<div> <div>52%</div> <div>12%</div> <div>36%</div> </div>
2	D	25	<div> <div>32%</div> <div>24%</div> <div>•</div> <div>40%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4745 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 Epidermal growth factor receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	288	Total	C	N	O	S	0	0	0
			2260	1458	375	410	17			
1	B	286	Total	C	N	O	S	0	0	0
			2242	1450	370	406	16			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	675	GLY	-	EXPRESSION TAG	UNP P00533
A	676	ALA	-	EXPRESSION TAG	UNP P00533
A	677	MET	-	EXPRESSION TAG	UNP P00533
A	799	GLU	LYS	ENGINEERED	UNP P00533
B	675	GLY	-	EXPRESSION TAG	UNP P00533
B	676	ALA	-	EXPRESSION TAG	UNP P00533
B	677	MET	-	EXPRESSION TAG	UNP P00533
B	799	GLU	LYS	ENGINEERED	UNP P00533

- Molecule 2 is a protein called ERBB receptor feedback inhibitor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	16	Total	C	N	O	S	0	0	0
			122	79	18	24	1			
2	D	15	Total	C	N	O	S	0	0	0
			116	76	17	22	1			

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).

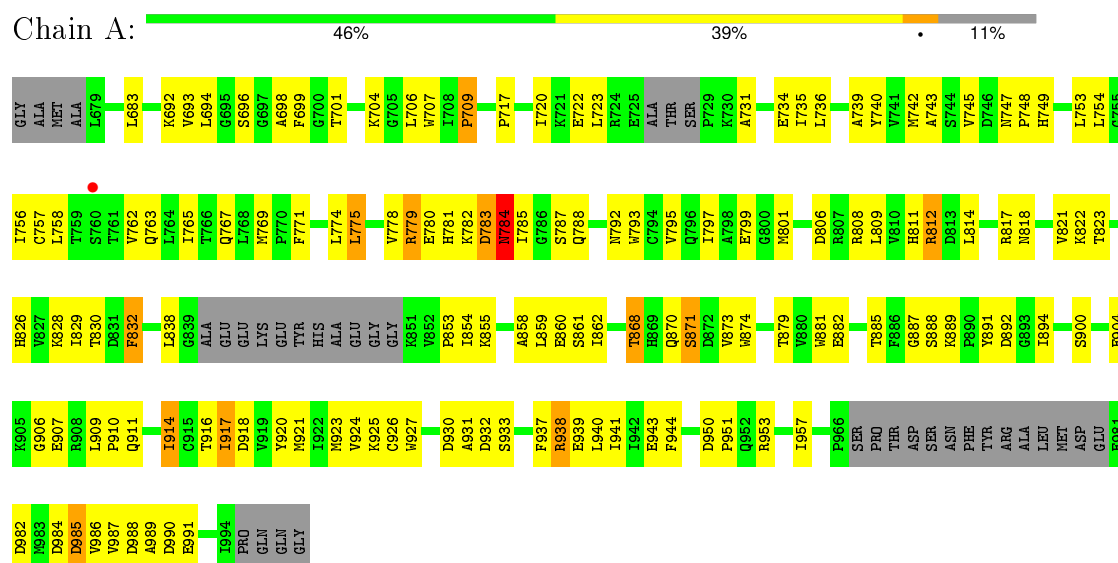


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		

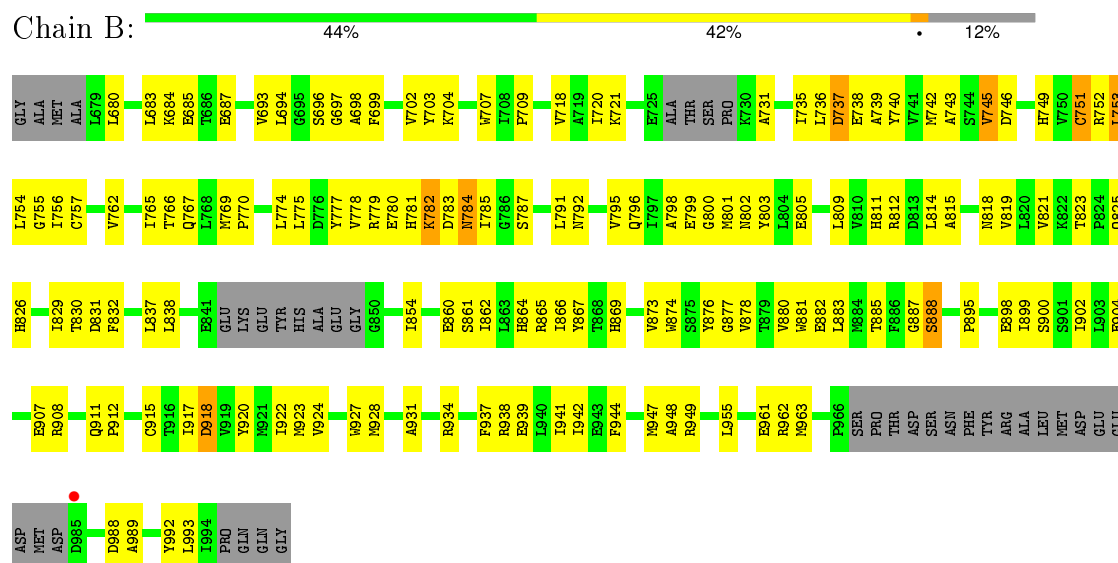
### 3 Residue-property plots

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.

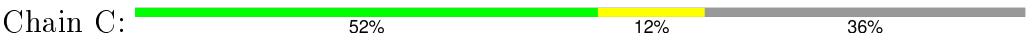
#### • Molecule 1: Epidermal growth factor receptor



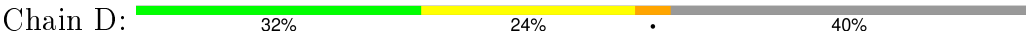
#### • Molecule 1: Epidermal growth factor receptor



#### • Molecule 2: ERBB receptor feedback inhibitor 1



- Molecule 2: ERBB receptor feedback inhibitor 1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	143.96Å 143.96Å 237.85Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.01 – 3.60 49.00 – 3.60	Depositor EDS
% Data completeness (in resolution range)	93.8 (49.01-3.60) 94.3 (49.00-3.60)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	6.73 (at 3.57Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.234 , 0.278 0.217 , 0.247	Depositor DCC
$R_{free}$ test set	798 reflections (4.85%)	DCC
Wilson B-factor (Å <sup>2</sup> )	82.1	Xtriage
Anisotropy	0.209	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 64.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 17232 reflections	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	4745	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 4.05% 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: SO4

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	0.47	2/2306 (0.1%)	0.65	0/3129
1	B	0.45	0/2288	0.66	0/3105
2	C	0.45	0/127	0.65	0/174
2	D	0.42	0/121	0.57	0/166
All	All	0.46	2/4842 (0.0%)	0.65	0/6574

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	943	GLU	CD-OE2	5.31	1.31	1.25
1	A	943	GLU	CG-CD	5.20	1.59	1.51

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2260	0	2260	136	0
1	B	2242	0	2243	125	0
2	C	122	0	115	7	0
2	D	116	0	110	8	0
3	A	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4745	0	4728	253	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 27.

All (253) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:907:GLU:HB3	2:C:349:THR:CG2	1.98	0.93
1:A:891:TYR:HB3	1:A:894:ILE:HD12	1.55	0.88
1:A:907:GLU:HB3	2:C:349:THR:HG21	1.55	0.86
1:A:787:SER:HB3	1:A:951:PRO:HB2	1.58	0.85
1:B:752:ARG:HH12	1:B:754:LEU:HD23	1.43	0.83
1:A:987:VAL:HG13	1:A:991:GLU:HB2	1.61	0.82
1:B:895:PRO:HD2	1:B:898:GLU:OE2	1.81	0.80
1:A:696:SER:HB3	1:A:701:THR:HG23	1.62	0.79
1:A:937:PHE:HA	1:A:940:LEU:HD12	1.63	0.78
1:A:795:VAL:O	1:A:799:GLU:HG3	1.83	0.77
1:B:753:LEU:HD12	1:B:754:LEU:N	2.00	0.77
1:A:683:LEU:HD12	1:A:765:ILE:HG13	1.69	0.74
1:B:899:ILE:HD13	1:B:902:ILE:HD11	1.70	0.73
1:A:787:SER:CB	1:A:951:PRO:HB2	2.18	0.73
1:A:723:LEU:HD22	1:A:838:LEU:HD11	1.70	0.72
1:A:722:GLU:HG3	1:A:763:GLN:HG2	1.73	0.71
1:B:753:LEU:HD12	1:B:754:LEU:H	1.53	0.70
1:A:984:ASP:O	1:A:985:ASP:HB2	1.89	0.70
1:B:699:PHE:HB3	1:B:838:LEU:CD1	2.22	0.69
1:A:788:GLN:HE21	1:A:951:PRO:HG3	1.58	0.68
1:B:907:GLU:OE1	2:D:349:THR:HG21	1.93	0.68
1:B:795:VAL:O	1:B:799:GLU:HG3	1.93	0.68
1:A:881:TRP:HA	1:A:923:MET:HE1	1.76	0.68
1:B:736:LEU:HB3	1:B:740:TYR:CE1	2.29	0.68
1:A:780:GLU:HG2	1:A:780:GLU:O	1.92	0.67
1:A:758:LEU:HD21	1:A:762:VAL:HG23	1.77	0.67
1:A:694:LEU:HD21	1:A:704:LYS:HB2	1.75	0.67
1:B:947:MET:O	1:B:949:ARG:N	2.29	0.66
1:A:706:LEU:HD23	1:A:717:PRO:HA	1.77	0.66
1:B:938:ARG:HG3	1:B:939:GLU:N	2.10	0.65
1:A:757:CYS:HB3	1:A:763:GLN:HB2	1.77	0.65
1:A:906:GLY:HA3	2:C:359:VAL:HG22	1.79	0.65
1:B:775:LEU:O	1:B:779:ARG:HG3	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:694:LEU:HD11	1:B:704:LYS:CB	2.27	0.64
1:B:860:GLU:HG2	1:B:861:SER:N	2.12	0.64
1:A:707:TRP:O	1:A:709:PRO:HD3	1.98	0.63
1:A:758:LEU:CD2	1:A:762:VAL:HG23	2.28	0.63
1:A:900:SER:O	1:A:904:GLU:HG3	1.99	0.63
1:A:775:LEU:O	1:A:775:LEU:HD12	1.99	0.62
1:B:829:ILE:HG22	1:B:830:THR:N	2.14	0.62
1:B:947:MET:C	1:B:949:ARG:H	2.03	0.62
1:B:837:LEU:O	1:B:837:LEU:HD23	2.00	0.61
1:B:683:LEU:HD12	1:B:765:ILE:HG13	1.82	0.61
1:B:811:HIS:O	1:B:812:ARG:HB2	1.99	0.61
1:B:922:ILE:HG21	1:B:944:PHE:CE2	2.36	0.60
1:B:877:GLY:HA3	1:B:927:TRP:HE1	1.67	0.60
1:B:801:MET:CE	1:B:814:LEU:HD22	2.31	0.60
1:B:885:THR:HB	1:B:888:SER:OG	2.02	0.59
1:B:911:GLN:HG3	1:B:920:TYR:CG	2.37	0.59
1:B:799:GLU:HA	1:B:941:ILE:HD11	1.85	0.59
1:B:928:MET:HE3	2:D:357:LYS:HB3	1.84	0.59
1:A:868:THR:HG22	1:B:938:ARG:HD3	1.85	0.58
1:A:694:LEU:HD11	1:A:704:LYS:HB2	1.86	0.58
1:B:756:ILE:HD11	1:B:762:VAL:HG23	1.84	0.58
1:A:723:LEU:HB2	1:A:762:VAL:CG1	2.33	0.58
1:A:868:THR:HG23	1:B:938:ARG:HB2	1.86	0.58
1:A:930:ASP:HB3	1:B:949:ARG:NH2	2.18	0.58
1:A:911:GLN:HA	1:A:920:TYR:CD1	2.39	0.58
1:A:911:GLN:HA	1:A:920:TYR:CE1	2.38	0.57
1:A:870:GLN:HG3	1:B:942:ILE:HD13	1.84	0.57
1:A:736:LEU:HG	1:A:740:TYR:CE1	2.40	0.57
1:A:917:ILE:HG23	1:A:918:ASP:N	2.19	0.57
1:A:731:ALA:O	1:A:735:ILE:HG12	2.04	0.57
1:A:736:LEU:HG	1:A:740:TYR:HE1	1.69	0.57
1:B:854:ILE:HG23	1:B:862:ILE:CD1	2.35	0.56
1:B:791:LEU:HG	1:B:955:LEU:HD12	1.88	0.56
1:B:752:ARG:NH1	1:B:754:LEU:HD23	2.15	0.56
1:A:881:TRP:HA	1:A:923:MET:CE	2.35	0.56
1:B:805:GLU:HA	1:B:869:HIS:CE1	2.40	0.56
1:A:683:LEU:CD1	1:A:765:ILE:HG13	2.34	0.56
1:B:693:VAL:HG22	1:B:703:TYR:CE2	2.41	0.55
1:A:811:HIS:O	1:A:812:ARG:HB2	2.06	0.55
1:B:756:ILE:HD11	1:B:762:VAL:CG2	2.37	0.54
1:B:815:ALA:O	1:B:819:VAL:HG23	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:818:ASN:O	1:A:830:THR:HG22	2.07	0.54
1:B:860:GLU:HG2	1:B:861:SER:H	1.72	0.53
1:B:829:ILE:HG22	1:B:830:THR:H	1.74	0.53
1:A:826:HIS:C	1:A:826:HIS:CD2	2.83	0.53
1:A:756:ILE:HD11	1:A:758:LEU:HD21	1.92	0.52
1:A:754:LEU:HD11	1:A:767:GLN:N	2.25	0.52
1:A:987:VAL:CG1	1:A:991:GLU:HB2	2.34	0.52
1:B:736:LEU:O	1:B:737:ASP:C	2.47	0.52
1:A:778:VAL:HA	1:A:785:ILE:HD11	1.92	0.52
1:A:808:ARG:NH1	1:B:939:GLU:OE1	2.43	0.51
1:B:864:HIS:HB2	1:B:866:ILE:HG13	1.92	0.51
1:B:800:GLY:O	1:B:803:TYR:HB3	2.09	0.51
1:A:736:LEU:O	1:A:740:TYR:HD1	1.94	0.51
1:A:699:PHE:CB	1:A:838:LEU:HD13	2.40	0.51
1:B:924:VAL:HG11	2:D:352:PHE:CE2	2.46	0.51
1:B:693:VAL:HG22	1:B:703:TYR:CD2	2.46	0.51
1:B:684:LYS:HB2	1:B:687:GLU:HG3	1.93	0.51
1:A:914:ILE:HD12	1:A:957:ILE:HG12	1.93	0.51
1:B:912:PRO:HB2	1:B:915:CYS:SG	2.51	0.51
1:B:809:LEU:HD13	1:B:832:PHE:CZ	2.46	0.50
1:A:988:ASP:O	1:A:991:GLU:N	2.44	0.50
1:A:720:ILE:HG23	1:A:765:ILE:HD13	1.93	0.50
1:A:720:ILE:HG23	1:A:765:ILE:CD1	2.42	0.50
1:A:986:VAL:O	1:A:986:VAL:HG12	2.10	0.50
1:B:781:HIS:O	1:B:782:LYS:C	2.50	0.50
1:B:801:MET:HE1	1:B:814:LEU:HD22	1.94	0.50
1:B:696:SER:OG	1:B:697:GLY:N	2.44	0.50
1:B:928:MET:CE	2:D:357:LYS:HB3	2.42	0.50
1:B:854:ILE:HG23	1:B:862:ILE:HD13	1.94	0.49
1:B:812:ARG:CG	1:B:867:TYR:CD2	2.95	0.49
1:B:756:ILE:HG12	1:B:757:CYS:N	2.27	0.49
2:D:354:PRO:O	2:D:356:PRO:HD3	2.12	0.49
1:A:808:ARG:NH1	1:B:939:GLU:OE2	2.45	0.49
1:B:864:HIS:O	1:B:865:ARG:C	2.50	0.49
1:A:774:LEU:HD11	1:A:793:TRP:CZ3	2.48	0.49
1:A:885:THR:OG1	1:A:888:SER:HB3	2.13	0.49
1:A:781:HIS:O	1:A:782:LYS:C	2.52	0.48
1:B:887:GLY:O	1:B:888:SER:C	2.51	0.48
1:A:723:LEU:HD12	1:A:762:VAL:CG1	2.43	0.48
1:A:906:GLY:HA3	2:C:359:VAL:CG2	2.43	0.48
1:A:871:SER:O	1:A:874:TRP:HB3	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:799:GLU:HA	1:A:941:ILE:HD11	1.94	0.48
1:B:680:LEU:HD13	1:B:743:ALA:HB2	1.94	0.48
1:B:774:LEU:HD23	1:B:883:LEU:HD21	1.95	0.48
1:A:809:LEU:HD22	1:A:832:PHE:HZ	1.78	0.48
1:B:720:ILE:HD12	1:B:720:ILE:N	2.28	0.48
1:B:787:SER:HB2	1:B:955:LEU:HB2	1.94	0.48
1:B:826:HIS:C	1:B:826:HIS:CD2	2.86	0.48
1:A:881:TRP:CD2	1:A:909:LEU:HD13	2.49	0.48
1:B:992:TYR:HD1	1:B:993:LEU:HG	1.78	0.48
1:A:775:LEU:C	1:A:775:LEU:HD12	2.34	0.48
1:B:931:ALA:HA	1:B:934:ARG:CZ	2.44	0.47
1:A:923:MET:O	1:A:926:CYS:HB2	2.15	0.47
1:B:778:VAL:HA	1:B:785:ILE:HD11	1.97	0.47
1:A:987:VAL:HG13	1:A:991:GLU:CB	2.39	0.47
1:A:696:SER:HB3	1:A:701:THR:CG2	2.40	0.47
1:B:736:LEU:O	1:B:739:ALA:N	2.47	0.47
1:B:718:VAL:HG23	1:B:720:ILE:CD1	2.44	0.47
1:B:812:ARG:HG2	1:B:867:TYR:CD2	2.48	0.47
1:B:774:LEU:O	1:B:778:VAL:HG13	2.14	0.47
1:B:878:VAL:O	1:B:882:GLU:HG3	2.15	0.47
1:A:881:TRP:CZ2	1:A:885:THR:HG21	2.49	0.47
1:B:775:LEU:C	1:B:777:TYR:H	2.17	0.47
1:A:907:GLU:CB	2:C:349:THR:HG21	2.35	0.47
1:A:783:ASP:O	1:A:784:ASN:HB2	2.15	0.47
1:A:747:ASN:ND2	1:A:748:PRO:HD2	2.30	0.47
1:A:870:GLN:HA	1:A:870:GLN:OE1	2.14	0.47
1:A:911:GLN:HG3	1:A:920:TYR:CG	2.50	0.47
1:A:723:LEU:HD12	1:A:762:VAL:HG13	1.96	0.46
1:A:938:ARG:HE	1:A:938:ARG:HB3	1.63	0.46
1:A:868:THR:CG2	1:B:938:ARG:HD3	2.45	0.46
1:A:692:LYS:HG2	1:A:693:VAL:N	2.29	0.46
1:A:921:MET:O	1:A:925:LYS:HB2	2.16	0.46
2:D:349:THR:HG23	2:D:350:GLN:N	2.29	0.46
1:B:873:VAL:O	1:B:876:TYR:HB3	2.16	0.46
1:A:950:ASP:OD1	1:A:953:ARG:HB2	2.15	0.46
1:A:909:LEU:HB2	1:A:927:TRP:CZ3	2.51	0.46
1:B:735:ILE:CD1	1:B:837:LEU:HD21	2.46	0.46
1:B:702:VAL:HG22	1:B:721:LYS:HE3	1.98	0.45
1:A:988:ASP:HB3	1:A:991:GLU:HG3	1.99	0.45
1:A:734:GLU:O	1:A:735:ILE:C	2.53	0.45
1:A:932:ASP:OD1	1:B:949:ARG:NH2	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:916:THR:O	1:A:917:ILE:C	2.54	0.45
1:A:771:PHE:HB2	1:A:821:VAL:HB	1.99	0.45
1:A:860:GLU:HG2	1:A:861:SER:N	2.30	0.45
1:B:799:GLU:HA	1:B:941:ILE:CD1	2.45	0.45
1:A:881:TRP:CA	1:A:923:MET:CE	2.94	0.45
1:B:720:ILE:HG13	1:B:765:ILE:HD12	1.98	0.45
1:B:900:SER:O	1:B:904:GLU:HG3	2.16	0.45
1:B:751:CYS:HB2	1:B:829:ILE:O	2.16	0.45
1:B:780:GLU:CG	1:B:780:GLU:O	2.65	0.45
1:B:745:VAL:O	1:B:745:VAL:HG23	2.17	0.45
1:A:788:GLN:O	1:A:792:ASN:ND2	2.50	0.45
1:B:752:ARG:NH1	1:B:752:ARG:HG2	2.32	0.45
1:B:947:MET:C	1:B:949:ARG:N	2.67	0.45
1:A:769:MET:HE3	1:A:822:LYS:CA	2.47	0.45
1:A:858:ALA:HA	1:A:874:TRP:CD2	2.52	0.44
1:A:879:THR:O	1:A:882:GLU:HB2	2.18	0.44
1:B:707:TRP:CD1	1:B:707:TRP:C	2.90	0.44
1:A:754:LEU:CD1	1:A:767:GLN:N	2.80	0.44
1:A:745:VAL:HG23	1:A:745:VAL:O	2.17	0.44
1:B:767:GLN:HG3	1:B:989:ALA:HB3	1.99	0.44
1:B:881:TRP:HA	1:B:923:MET:HE1	2.00	0.44
1:A:855:LYS:HD3	1:A:891:TYR:HB2	2.00	0.44
1:A:795:VAL:HG22	1:A:944:PHE:HB3	2.00	0.44
1:B:791:LEU:HG	1:B:955:LEU:CD1	2.48	0.44
1:A:870:GLN:CG	1:B:942:ILE:HD13	2.49	0.43
1:B:829:ILE:CG2	1:B:830:THR:N	2.82	0.43
1:A:938:ARG:HG2	1:A:939:GLU:N	2.34	0.43
1:A:909:LEU:HD22	1:A:910:PRO:HD2	2.00	0.43
1:A:916:THR:O	1:A:918:ASP:N	2.52	0.43
1:A:858:ALA:HA	1:A:874:TRP:CE2	2.53	0.43
1:A:775:LEU:O	1:A:779:ARG:CG	2.66	0.43
1:B:683:LEU:CD1	1:B:765:ILE:HG13	2.48	0.43
1:B:865:ARG:HB2	1:B:865:ARG:HE	1.68	0.43
1:A:774:LEU:HD11	1:A:793:TRP:CE3	2.53	0.43
1:A:742:MET:O	1:A:745:VAL:HG22	2.18	0.43
1:B:731:ALA:O	1:B:735:ILE:HG12	2.19	0.43
1:B:881:TRP:HA	1:B:923:MET:CE	2.48	0.43
1:A:817:ARG:HH22	1:A:853:PRO:HB3	1.83	0.43
1:A:756:ILE:HG12	1:A:757:CYS:N	2.34	0.43
1:B:699:PHE:HB3	1:B:838:LEU:HD12	2.00	0.43
1:A:780:GLU:O	1:A:781:HIS:CG	2.72	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:854:ILE:HG23	1:B:862:ILE:HD11	2.01	0.43
1:B:769:MET:HA	1:B:770:PRO:HD2	1.75	0.42
1:A:769:MET:HE2	1:A:828:LYS:HD2	2.01	0.42
1:A:787:SER:HB2	1:A:951:PRO:HB2	2.01	0.42
1:A:696:SER:HA	1:A:701:THR:HA	2.02	0.42
1:B:899:ILE:O	1:B:902:ILE:HG12	2.19	0.42
1:B:818:ASN:O	1:B:830:THR:HG22	2.19	0.42
1:B:798:ALA:O	1:B:799:GLU:C	2.58	0.42
1:B:938:ARG:HG3	1:B:939:GLU:H	1.81	0.42
1:A:911:GLN:HG3	1:A:920:TYR:HB2	2.01	0.42
1:A:793:TRP:O	1:A:797:ILE:HG13	2.20	0.42
1:A:723:LEU:HB2	1:A:762:VAL:HG13	2.00	0.42
1:A:988:ASP:O	1:A:989:ALA:C	2.58	0.42
1:B:699:PHE:HB3	1:B:838:LEU:HD11	2.01	0.42
1:A:860:GLU:N	1:A:860:GLU:OE1	2.52	0.42
1:A:788:GLN:NE2	1:A:788:GLN:HA	2.35	0.42
1:A:788:GLN:NE2	1:A:951:PRO:HG3	2.29	0.42
1:A:870:GLN:HE21	1:A:932:ASP:HA	1.84	0.42
1:B:720:ILE:HG13	1:B:765:ILE:CD1	2.50	0.42
1:B:880:VAL:HG12	1:B:923:MET:HE3	2.02	0.42
1:A:873:VAL:HG13	1:A:940:LEU:HD11	2.02	0.42
1:A:762:VAL:HG13	1:A:762:VAL:O	2.19	0.42
1:A:909:LEU:CD2	1:A:910:PRO:HD2	2.50	0.42
1:B:920:TYR:HH	2:D:352:PHE:HZ	1.65	0.42
1:B:924:VAL:O	2:D:358:TYR:HE2	2.03	0.42
1:A:739:ALA:O	1:A:740:TYR:C	2.56	0.42
1:A:906:GLY:HA2	2:C:358:TYR:O	2.20	0.41
1:B:718:VAL:HG23	1:B:720:ILE:HD11	2.02	0.41
1:A:749:HIS:O	1:A:829:ILE:N	2.44	0.41
1:B:821:VAL:HG13	1:B:823:THR:O	2.20	0.41
1:A:775:LEU:O	1:A:779:ARG:HG2	2.19	0.41
1:B:908:ARG:HG2	1:B:908:ARG:HH11	1.85	0.41
1:A:854:ILE:HA	1:A:854:ILE:HD13	1.88	0.41
1:B:749:HIS:ND1	1:B:796:GLN:HB3	2.35	0.41
1:B:753:LEU:HD11	1:B:755:GLY:O	2.21	0.41
1:A:779:ARG:HD3	1:A:887:GLY:HA3	2.02	0.41
1:B:799:GLU:O	1:B:802:ASN:HB3	2.19	0.41
1:A:889:LYS:HB2	1:A:892:ASP:HB2	2.03	0.41
1:A:811:HIS:O	1:A:812:ARG:CB	2.69	0.41
1:B:873:VAL:HG23	1:B:937:PHE:CE1	2.56	0.41
1:B:961:GLU:O	1:B:963:MET:N	2.44	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:874:TRP:CE3	1:B:927:TRP:HA	2.56	0.41
1:B:707:TRP:CE2	1:B:709:PRO:HG3	2.56	0.41
1:B:881:TRP:N	1:B:923:MET:HE2	2.35	0.41
1:A:932:ASP:HB3	1:B:942:ILE:CG2	2.51	0.41
1:B:703:TYR:N	1:B:703:TYR:CD1	2.88	0.41
1:B:738:GLU:O	1:B:742:MET:HG3	2.21	0.41
1:A:801:MET:SD	1:A:814:LEU:HD22	2.61	0.40
1:B:917:ILE:HG23	1:B:918:ASP:N	2.36	0.40
1:A:743:ALA:HB2	1:A:753:LEU:HB3	2.01	0.40
1:A:917:ILE:CG2	1:A:918:ASP:N	2.85	0.40
1:A:924:VAL:O	2:C:358:TYR:HE2	2.03	0.40
1:B:830:THR:O	1:B:831:ASP:HB2	2.22	0.40
1:A:931:ALA:C	1:A:933:SER:H	2.24	0.40
1:A:859:LEU:O	1:A:862:ILE:N	2.54	0.40
1:B:860:GLU:C	1:B:862:ILE:N	2.75	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	280/324 (86%)	232 (83%)	41 (15%)	7 (2%)	7	48
1	B	278/324 (86%)	233 (84%)	38 (14%)	7 (2%)	7	48
2	C	14/25 (56%)	13 (93%)	1 (7%)	0	100	100
2	D	13/25 (52%)	11 (85%)	2 (15%)	0	100	100
All	All	585/698 (84%)	489 (84%)	82 (14%)	14 (2%)	7	49

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	698	ALA
1	A	784	ASN
1	B	698	ALA
1	B	745	VAL
1	B	784	ASN
1	A	917	ILE
1	A	982	ASP
1	B	948	ALA
1	B	962	ARG
1	B	888	SER
1	A	709	PRO
1	A	812	ARG
1	B	782	LYS
1	A	985	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	244/284 (86%)	232 (95%)	12 (5%)	31	71
1	B	241/284 (85%)	229 (95%)	12 (5%)	30	70
2	C	15/22 (68%)	15 (100%)	0	100	100
2	D	14/22 (64%)	13 (93%)	1 (7%)	18	59
All	All	514/612 (84%)	489 (95%)	25 (5%)	31	71

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	775	LEU
1	A	779	ARG
1	A	783	ASP
1	A	784	ASN
1	A	806	ASP
1	A	823	THR
1	A	832	PHE
1	A	868	THR

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Mol	Chain	Res	Type
1	A	871	SER
1	A	914	ILE
1	A	938	ARG
1	A	990	ASP
1	B	685	GLU
1	B	737	ASP
1	B	746	ASP
1	B	751	CYS
1	B	753	LEU
1	B	766	THR
1	B	783	ASP
1	B	784	ASN
1	B	792	ASN
1	B	825	GLN
1	B	918	ASP
1	B	988	ASP
2	D	349	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	781	HIS
1	A	788	GLN
1	A	792	ASN
1	A	952	GLN
1	A	958	GLN
1	B	749	HIS
1	B	767	GLN
1	B	826	HIS
1	B	869	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	SO4	A	1	-	4,4,4	0.21	0	6,6,6	0.12	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SO4	A	1	-	-	0/0/0/0	0/0/0/0

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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	288/324 (88%)	-0.09	1 (0%) 94 90	18, 52, 104, 140	0
1	B	286/324 (88%)	-0.23	1 (0%) 94 90	15, 49, 113, 144	0
2	C	16/25 (64%)	-0.09	0 100 100	32, 50, 74, 78	0
2	D	15/25 (60%)	0.02	0 100 100	51, 60, 76, 81	0
All	All	605/698 (86%)	-0.15	2 (0%) 94 90	15, 51, 104, 144	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	985	ASP	2.8
1	A	760	SER	2.2

### 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](#)

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( $\text{\AA}^2$ )	Q<0.9
3	SO4	A	1	5/5	0.93	0.23	0.04	76,77,77,77	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.