



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 02:36 PM GMT

PDB ID : 3A8P
Title : Crystal structure of the Tiam2 PHCCEX domain
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Deposited on : 2009-10-07
Resolution : 2.10 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

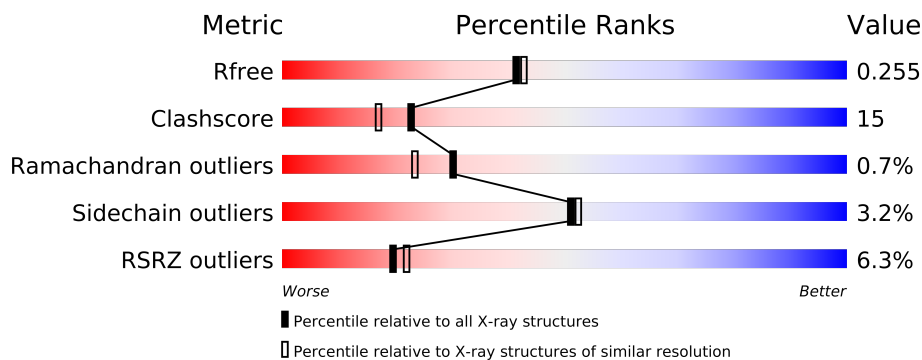
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.10 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	3012 (2.10-2.10)
Clashscore	79885	3649 (2.10-2.10)
Ramachandran outliers	78287	3610 (2.10-2.10)
Sidechain outliers	78261	3611 (2.10-2.10)
RSRZ outliers	66119	3013 (2.10-2.10)

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. 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	263	
1	B	263	
1	C	263	
1	D	263	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7754 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 T-lymphoma invasion and metastasis-inducing protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	232	Total	C	N	O	S	0	0	0
			1856	1184	331	329	12			
1	B	231	Total	C	N	O	S	0	0	0
			1853	1182	330	329	12			
1	C	229	Total	C	N	O	S	0	0	0
			1839	1175	328	324	12			
1	D	228	Total	C	N	O	S	0	0	0
			1823	1163	327	321	12			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	495	GLY	-	EXPRESSION TAG	UNP Q6ZPF3
A	496	PRO	-	EXPRESSION TAG	UNP Q6ZPF3
A	497	LEU	-	EXPRESSION TAG	UNP Q6ZPF3
A	498	GLY	-	EXPRESSION TAG	UNP Q6ZPF3
A	499	SER	-	EXPRESSION TAG	UNP Q6ZPF3
B	495	GLY	-	EXPRESSION TAG	UNP Q6ZPF3
B	496	PRO	-	EXPRESSION TAG	UNP Q6ZPF3
B	497	LEU	-	EXPRESSION TAG	UNP Q6ZPF3
B	498	GLY	-	EXPRESSION TAG	UNP Q6ZPF3
B	499	SER	-	EXPRESSION TAG	UNP Q6ZPF3
C	495	GLY	-	EXPRESSION TAG	UNP Q6ZPF3
C	496	PRO	-	EXPRESSION TAG	UNP Q6ZPF3
C	497	LEU	-	EXPRESSION TAG	UNP Q6ZPF3
C	498	GLY	-	EXPRESSION TAG	UNP Q6ZPF3
C	499	SER	-	EXPRESSION TAG	UNP Q6ZPF3
D	495	GLY	-	EXPRESSION TAG	UNP Q6ZPF3
D	496	PRO	-	EXPRESSION TAG	UNP Q6ZPF3
D	497	LEU	-	EXPRESSION TAG	UNP Q6ZPF3
D	498	GLY	-	EXPRESSION TAG	UNP Q6ZPF3
D	499	SER	-	EXPRESSION TAG	UNP Q6ZPF3

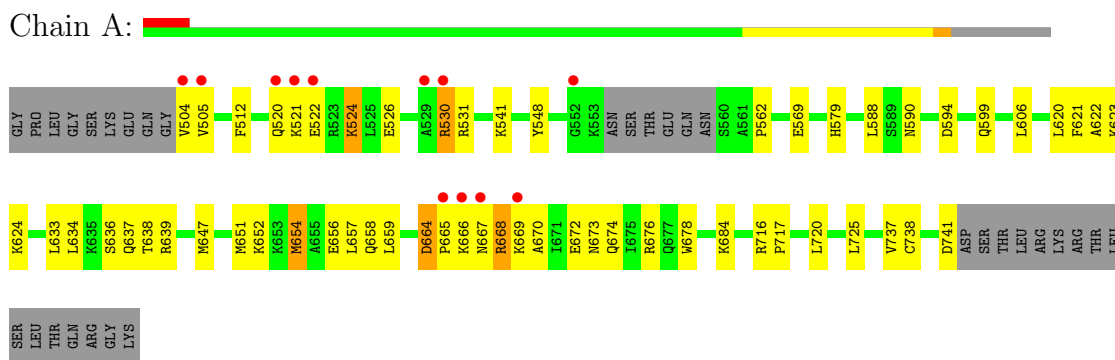
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	126	Total 126	O 126	0	0
2	B	75	Total 75	O 75	0	0
2	C	90	Total 90	O 90	0	0
2	D	92	Total 92	O 92	0	0

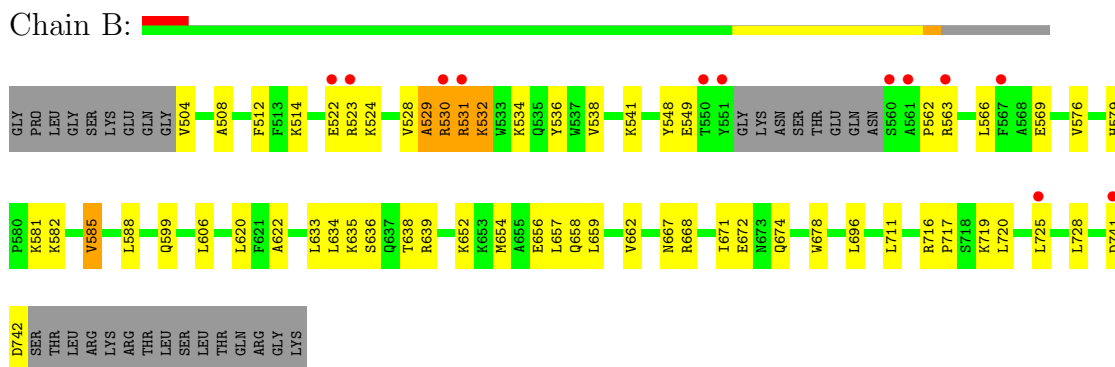
3 Residue-property plots

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.

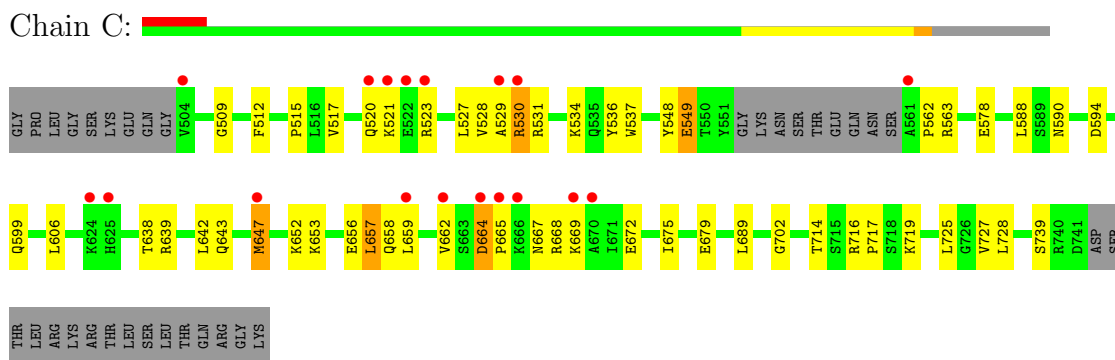
- Molecule 1: T-lymphoma invasion and metastasis-inducing protein 2



- Molecule 1: T-lymphoma invasion and metastasis-inducing protein 2

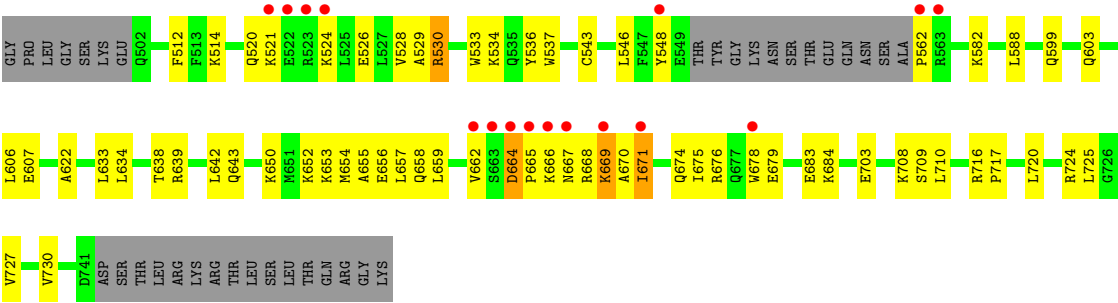


- Molecule 1: T-lymphoma invasion and metastasis-inducing protein 2



- Molecule 1: T-lymphoma invasion and metastasis-inducing protein 2

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	46.69Å 104.79Å 115.97Å 90.00° 80.55° 90.00°	Depositor
Resolution (Å)	29.81 – 2.10 37.76 – 2.08	Depositor EDS
% Data completeness (in resolution range)	98.7 (29.81-2.10) 97.7 (37.76-2.08)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 2.08Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.211 , 0.253 0.212 , 0.255	Depositor DCC
R_{free} test set	6404 reflections (10.10%)	DCC
Wilson B-factor (Å ²)	30.5	Xtriage
Anisotropy	0.546	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 47.5	EDS
Estimated twinning fraction	0.024 for h,-k,h-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	2 of 64541 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7754	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.22% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities 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	0.35	0/1893	0.54	0/2548
1	B	0.34	0/1890	0.55	0/2544
1	C	0.33	0/1876	0.53	0/2525
1	D	0.33	0/1859	0.53	0/2500
All	All	0.34	0/7518	0.54	0/10117

There are no bond length outliers.

There are no bond angle outliers.

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	1856	0	1901	49	0
1	B	1853	0	1901	62	0
1	C	1839	0	1892	49	0
1	D	1823	0	1875	62	0
2	A	126	0	0	2	0
2	B	75	0	0	0	0
2	C	90	0	0	2	0
2	D	92	0	0	3	0
All	All	7754	0	7569	219	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 15.

All (219) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:520:GLN:HG3	1:C:521:LYS:H	1.30	0.94
1:A:664:ASP:HB2	1:A:667:ASN:HB2	1.51	0.93
1:D:650:LYS:O	1:D:653:LYS:HG2	1.78	0.82
1:B:662:VAL:HG13	1:B:668:ARG:NH1	1.95	0.80
1:C:520:GLN:CG	1:C:521:LYS:H	1.94	0.80
1:B:528:VAL:HB	1:B:531:ARG:HH11	1.46	0.79
1:B:514:LYS:HG3	1:B:531:ARG:HD3	1.67	0.77
1:B:514:LYS:CG	1:B:531:ARG:HD3	2.16	0.76
1:B:532:LYS:H	1:B:532:LYS:HD2	1.50	0.75
1:A:590:ASN:HD21	1:A:594:ASP:HB2	1.53	0.72
1:B:725:LEU:HD23	1:B:725:LEU:O	1.90	0.72
1:B:522:GLU:HB3	1:B:524:LYS:HE3	1.70	0.71
1:D:659:LEU:HA	1:D:662:VAL:HG12	1.73	0.71
1:D:530:ARG:H	1:D:530:ARG:NE	1.88	0.71
1:D:684:LYS:HE2	2:D:239:HOH:O	1.89	0.71
1:C:638:THR:O	1:C:642:LEU:HD13	1.91	0.71
1:D:703:GLU:CD	1:D:703:GLU:H	1.93	0.70
1:B:716:ARG:NH1	1:B:720:LEU:HD11	2.07	0.70
1:A:590:ASN:ND2	1:A:594:ASP:HB2	2.07	0.69
1:B:711:LEU:HD22	1:B:728:LEU:HD11	1.76	0.68
1:D:529:ALA:HB3	1:D:530:ARG:NH2	2.09	0.66
1:A:504:VAL:N	1:A:541:LYS:HZ2	1.94	0.66
1:C:669:LYS:HA	1:C:672:GLU:HG2	1.77	0.66
1:C:664:ASP:CB	1:C:667:ASN:HB3	2.26	0.66
1:B:652:LYS:O	1:B:656:GLU:HG3	1.97	0.65
1:C:664:ASP:HB2	1:C:667:ASN:HB3	1.78	0.65
1:C:520:GLN:HG3	1:C:521:LYS:N	2.07	0.64
1:A:654:MET:HE1	1:A:657:LEU:HD23	1.80	0.64
1:D:528:VAL:HG13	1:D:530:ARG:HH11	1.62	0.64
1:D:674:GLN:HG3	1:D:678:TRP:CZ3	2.32	0.63
1:D:634:LEU:O	1:D:638:THR:HG23	1.99	0.63
1:D:524:LYS:HE2	1:D:526:GLU:OE2	1.99	0.63
1:D:725:LEU:HD22	1:D:727:VAL:H	1.65	0.62
1:C:664:ASP:HB2	1:C:667:ASN:HD22	1.65	0.61
1:C:534:LYS:HD3	1:C:536:TYR:OH	2.00	0.61
1:A:716:ARG:O	1:A:720:LEU:HD13	2.01	0.60
1:A:673:ASN:HD22	1:A:676:ARG:NH1	1.99	0.60
1:A:530:ARG:HG2	1:A:530:ARG:O	2.00	0.60
1:B:662:VAL:HG13	1:B:668:ARG:HH11	1.63	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:725:LEU:C	1:D:725:LEU:HD23	2.21	0.60
1:C:520:GLN:CG	1:C:521:LYS:N	2.65	0.59
1:D:671:ILE:O	1:D:675:ILE:HG12	2.02	0.59
1:D:669:LYS:HD3	1:D:670:ALA:H	1.67	0.59
1:D:659:LEU:HA	1:D:662:VAL:CG1	2.33	0.58
1:A:522:GLU:HB3	1:A:524:LYS:HD2	1.85	0.58
1:B:582:LYS:HD3	1:B:599:GLN:NE2	2.18	0.58
1:D:548:TYR:CZ	1:D:562:PRO:HG3	2.39	0.58
1:C:657:LEU:H	1:C:657:LEU:HD22	1.68	0.57
1:A:664:ASP:CB	1:A:667:ASN:HD22	2.17	0.57
1:A:664:ASP:HB3	1:A:667:ASN:HD22	1.70	0.57
1:C:659:LEU:HD13	1:C:668:ARG:HH21	1.69	0.57
1:D:679:GLU:O	1:D:683:GLU:HG3	2.04	0.56
1:A:521:LYS:O	1:A:522:GLU:HB3	2.06	0.56
1:A:504:VAL:HG12	1:A:505:VAL:N	2.21	0.56
1:A:524:LYS:HD3	1:A:524:LYS:H	1.71	0.56
1:B:576:VAL:HB	1:B:585:VAL:CG2	2.36	0.56
1:B:636:SER:HA	1:B:639:ARG:NH1	2.21	0.56
1:B:523:ARG:HH22	1:B:741:ASP:C	2.10	0.55
1:D:709:SER:HB2	2:D:141:HOH:O	2.06	0.55
1:A:673:ASN:HD22	1:A:676:ARG:HH12	1.54	0.55
1:A:654:MET:CE	1:A:654:MET:HA	2.37	0.55
1:A:716:ARG:HB2	1:A:717:PRO:HD3	1.89	0.55
1:D:724:ARG:HH11	1:D:724:ARG:HG3	1.70	0.55
1:B:531:ARG:CD	1:B:531:ARG:H	2.20	0.55
1:B:512:PHE:HB2	1:B:599:GLN:HB3	1.88	0.55
1:C:669:LYS:HA	1:C:672:GLU:CG	2.37	0.54
1:C:725:LEU:HD23	1:C:725:LEU:C	2.26	0.54
1:B:531:ARG:N	1:B:531:ARG:HD2	2.21	0.54
1:D:665:PRO:HA	1:D:668:ARG:HG2	1.89	0.54
1:B:569:GLU:HG3	1:B:620:LEU:HD23	1.89	0.54
1:D:530:ARG:H	1:D:530:ARG:CD	2.19	0.54
1:B:534:LYS:HD3	1:B:536:TYR:OH	2.07	0.54
1:D:676:ARG:HG2	1:D:676:ARG:HH11	1.72	0.54
1:D:671:ILE:O	1:D:671:ILE:HD13	2.08	0.53
1:A:654:MET:HE2	1:A:654:MET:HA	1.89	0.53
1:A:668:ARG:O	1:A:672:GLU:HG3	2.08	0.53
1:B:528:VAL:HB	1:B:531:ARG:NH1	2.19	0.53
1:B:531:ARG:H	1:B:531:ARG:HD2	1.72	0.53
1:A:521:LYS:O	1:A:522:GLU:CB	2.57	0.53
1:C:530:ARG:HD3	1:C:530:ARG:C	2.29	0.52
1:A:579:HIS:O	1:B:719:LYS:HE2	2.09	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:702:GLY:HA3	2:C:278:HOH:O	2.10	0.52
1:C:662:VAL:O	1:C:668:ARG:HD3	2.09	0.52
1:D:603:GLN:O	1:D:607:GLU:HG3	2.09	0.52
1:D:659:LEU:HB2	1:D:671:ILE:HD12	1.90	0.52
1:D:652:LYS:HE3	1:D:679:GLU:HG2	1.92	0.52
1:D:669:LYS:HD3	1:D:670:ALA:N	2.25	0.52
1:D:520:GLN:HG3	1:D:526:GLU:OE1	2.10	0.51
1:B:711:LEU:HD22	1:B:728:LEU:CD1	2.40	0.51
1:D:622:ALA:HA	1:D:633:LEU:HD23	1.92	0.51
1:D:710:LEU:HD23	1:D:730:VAL:HG22	1.91	0.51
1:A:665:PRO:HG2	1:A:666:LYS:H	1.75	0.51
1:B:504:VAL:N	1:B:541:LYS:HE3	2.26	0.51
1:D:671:ILE:C	1:D:671:ILE:HD13	2.30	0.51
1:C:662:VAL:HG13	1:C:664:ASP:OD1	2.10	0.51
1:B:563:ARG:HD2	1:B:563:ARG:N	2.26	0.51
1:C:656:GLU:OE2	1:C:656:GLU:HA	2.11	0.51
1:C:578:GLU:OE2	1:D:708:LYS:HE2	2.11	0.51
1:D:653:LYS:HG3	1:D:654:MET:N	2.26	0.50
1:B:548:TYR:CZ	1:B:562:PRO:HG3	2.46	0.50
1:A:530:ARG:CG	1:A:530:ARG:O	2.59	0.50
1:B:667:ASN:O	1:B:671:ILE:HG13	2.12	0.50
1:A:673:ASN:ND2	1:A:676:ARG:HH12	2.09	0.50
1:B:549:GLU:HA	1:B:549:GLU:OE1	2.12	0.50
1:B:523:ARG:HH22	1:B:742:ASP:N	2.10	0.49
1:B:659:LEU:HA	1:B:662:VAL:HG12	1.94	0.49
1:C:658:GLN:O	1:C:662:VAL:HG23	2.12	0.49
1:C:509:GLY:HA2	1:C:537:TRP:CZ3	2.48	0.49
1:A:684:LYS:HD3	1:A:737:VAL:CG1	2.42	0.49
1:B:529:ALA:O	1:B:531:ARG:N	2.45	0.49
1:D:659:LEU:HD13	1:D:671:ILE:HG23	1.93	0.49
1:C:664:ASP:HB3	1:C:667:ASN:HB3	1.93	0.49
1:B:534:LYS:HG3	1:B:536:TYR:CE2	2.47	0.49
1:C:714:THR:HG22	1:C:719:LYS:HG3	1.92	0.49
1:B:528:VAL:CB	1:B:531:ARG:HH11	2.20	0.49
1:B:716:ARG:HH12	1:B:720:LEU:HD11	1.75	0.49
1:C:725:LEU:HD22	1:C:727:VAL:H	1.77	0.49
1:D:512:PHE:HB2	1:D:599:GLN:HB3	1.95	0.49
1:B:659:LEU:HA	1:B:662:VAL:CG1	2.42	0.49
1:D:662:VAL:O	1:D:662:VAL:HG13	2.13	0.49
1:C:653:LYS:O	1:C:657:LEU:HD22	2.13	0.49
1:A:520:GLN:HG3	1:A:526:GLU:CD	2.32	0.49
1:C:647:MET:HA	1:C:647:MET:CE	2.43	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:652:LYS:O	1:C:656:GLU:HG2	2.13	0.48
1:D:675:ILE:O	1:D:679:GLU:HG3	2.14	0.48
1:B:532:LYS:H	1:B:532:LYS:CD	2.24	0.48
1:D:708:LYS:NZ	2:D:107:HOH:O	2.47	0.48
1:D:548:TYR:CE1	1:D:562:PRO:HG3	2.49	0.47
1:C:725:LEU:HD23	1:C:725:LEU:O	2.14	0.47
1:D:528:VAL:CG1	1:D:530:ARG:HH11	2.28	0.47
1:B:716:ARG:HB3	1:B:717:PRO:HD3	1.95	0.47
1:A:548:TYR:CZ	1:A:562:PRO:HG3	2.49	0.47
1:A:659:LEU:HD11	1:A:668:ARG:HD3	1.97	0.47
1:D:716:ARG:HB3	1:D:717:PRO:HD3	1.95	0.47
1:B:588:LEU:HD23	1:B:588:LEU:C	2.34	0.47
1:A:674:GLN:HE21	1:A:678:TRP:HE1	1.62	0.47
1:B:662:VAL:HG13	1:B:668:ARG:HH12	1.79	0.47
1:D:668:ARG:HA	1:D:671:ILE:HG22	1.97	0.47
1:B:531:ARG:C	1:B:531:ARG:HD2	2.35	0.46
1:D:546:LEU:HD22	1:D:562:PRO:HG2	1.96	0.46
1:A:654:MET:CE	1:A:657:LEU:HD23	2.43	0.46
1:C:659:LEU:O	1:C:659:LEU:HD13	2.15	0.46
1:A:620:LEU:HD12	1:A:623:LYS:HE3	1.97	0.46
1:A:669:LYS:HG3	1:A:670:ALA:N	2.30	0.46
1:B:606:LEU:C	1:B:606:LEU:HD23	2.36	0.46
1:A:652:LYS:O	1:A:656:GLU:HG3	2.15	0.46
1:D:652:LYS:O	1:D:656:GLU:HG3	2.16	0.46
1:C:664:ASP:CB	1:C:667:ASN:HD22	2.28	0.46
1:C:549:GLU:OE2	1:C:563:ARG:NH1	2.49	0.46
1:A:569:GLU:OE2	1:A:624:LYS:HD2	2.15	0.46
1:D:659:LEU:HD13	1:D:671:ILE:CG2	2.46	0.45
1:D:606:LEU:HD23	1:D:606:LEU:C	2.37	0.45
1:C:515:PRO:HB2	1:C:528:VAL:HG21	1.97	0.45
1:C:588:LEU:C	1:C:588:LEU:HD23	2.37	0.45
1:A:667:ASN:HB3	2:A:270:HOH:O	2.16	0.45
1:D:622:ALA:HA	1:D:633:LEU:CD2	2.46	0.45
1:B:532:LYS:N	1:B:532:LYS:HD2	2.27	0.45
1:A:621:PHE:HZ	1:A:637:GLN:HE21	1.65	0.44
1:C:606:LEU:C	1:C:606:LEU:HD23	2.38	0.44
1:A:531:ARG:HA	2:A:311:HOH:O	2.16	0.44
1:A:622:ALA:HA	1:A:633:LEU:CD2	2.48	0.44
1:B:711:LEU:HD13	1:B:728:LEU:HD13	1.98	0.44
1:B:579:HIS:CE1	1:B:581:LYS:HG3	2.53	0.44
1:B:576:VAL:HB	1:B:585:VAL:HG23	1.99	0.44
1:A:512:PHE:HB2	1:A:599:GLN:HB3	1.99	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:665:PRO:C	1:D:667:ASN:H	2.19	0.44
1:A:579:HIS:O	1:B:719:LYS:CE	2.66	0.44
1:C:639:ARG:O	1:C:643:GLN:HG3	2.18	0.44
1:A:634:LEU:O	1:A:638:THR:HG23	2.18	0.44
1:D:537:TRP:HB3	1:D:548:TYR:HB2	2.00	0.43
1:C:728:LEU:HD23	1:C:728:LEU:C	2.38	0.43
1:B:654:MET:HE1	1:B:657:LEU:HD22	2.01	0.43
1:B:634:LEU:HB2	1:B:696:LEU:HD13	1.99	0.43
1:C:668:ARG:HH11	1:C:668:ARG:HG2	1.82	0.43
1:C:590:ASN:ND2	1:C:594:ASP:HB2	2.33	0.43
1:D:520:GLN:HB2	1:D:524:LYS:HG3	2.00	0.43
1:D:725:LEU:HD22	1:D:727:VAL:N	2.32	0.43
1:A:520:GLN:HG2	1:A:526:GLU:HG2	2.01	0.43
1:A:588:LEU:C	1:A:588:LEU:HD23	2.39	0.43
1:D:655:ALA:O	1:D:671:ILE:HD11	2.19	0.43
1:B:674:GLN:HG3	1:B:678:TRP:CZ3	2.54	0.43
1:C:534:LYS:HD3	1:C:536:TYR:CZ	2.54	0.42
1:C:548:TYR:CZ	1:C:562:PRO:HG3	2.53	0.42
1:B:674:GLN:HG3	1:B:678:TRP:CH2	2.55	0.42
1:B:508:ALA:HA	1:B:538:VAL:O	2.19	0.42
1:A:606:LEU:HD23	1:A:606:LEU:C	2.40	0.42
1:C:523:ARG:HG2	1:C:739:SER:O	2.20	0.42
1:C:675:ILE:O	1:C:679:GLU:HG3	2.19	0.42
1:D:534:LYS:HE3	1:D:536:TYR:OH	2.20	0.42
1:D:521:LYS:O	1:D:524:LYS:HG2	2.19	0.42
1:A:738:CYS:HA	1:A:741:ASP:OD2	2.19	0.42
1:C:638:THR:HG23	1:C:689:LEU:HD11	2.01	0.42
1:D:724:ARG:HG3	1:D:724:ARG:NH1	2.35	0.41
1:C:512:PHE:HB2	1:C:599:GLN:HB3	2.02	0.41
1:B:529:ALA:O	1:B:530:ARG:C	2.58	0.41
1:D:716:ARG:O	1:D:720:LEU:HD13	2.20	0.41
1:B:654:MET:CE	1:B:657:LEU:HD22	2.50	0.41
1:B:548:TYR:CE2	1:B:562:PRO:HG3	2.55	0.41
1:D:659:LEU:CA	1:D:662:VAL:HG12	2.46	0.41
1:C:531:ARG:HB3	2:C:314:HOH:O	2.20	0.41
1:B:635:LYS:O	1:B:639:ARG:HG3	2.21	0.41
1:A:654:MET:O	1:A:658:GLN:HG2	2.21	0.41
1:A:530:ARG:HG2	1:A:530:ARG:HH21	1.86	0.41
1:B:579:HIS:HE1	1:B:581:LYS:HG3	1.86	0.41
1:B:531:ARG:N	1:B:531:ARG:CD	2.82	0.41
1:D:668:ARG:HH11	1:D:668:ARG:HG3	1.85	0.41
1:B:634:LEU:O	1:B:638:THR:HG23	2.21	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:716:ARG:HB2	1:C:717:PRO:HD3	2.02	0.41
1:B:522:GLU:HA	1:B:522:GLU:OE2	2.21	0.40
1:B:622:ALA:HA	1:B:633:LEU:HD23	2.03	0.40
1:A:636:SER:HA	1:A:639:ARG:CZ	2.51	0.40
1:A:647:MET:HE3	1:A:651:MET:CG	2.51	0.40
1:C:668:ARG:HD2	1:C:668:ARG:HA	1.91	0.40
1:B:622:ALA:HA	1:B:633:LEU:CD2	2.51	0.40
1:D:639:ARG:O	1:D:643:GLN:HG3	2.21	0.40
1:C:517:VAL:HG22	1:C:527:LEU:HD23	2.02	0.40
1:D:655:ALA:HB3	1:D:675:ILE:HD11	2.04	0.40
1:D:588:LEU:HD23	1:D:588:LEU:C	2.42	0.40
1:D:514:LYS:HE3	1:D:533:TRP:CD1	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	228/263 (87%)	225 (99%)	3 (1%)	0	100	100
1	B	227/263 (86%)	222 (98%)	3 (1%)	2 (1%)	25	17
1	C	225/263 (86%)	214 (95%)	9 (4%)	2 (1%)	25	17
1	D	224/263 (85%)	217 (97%)	5 (2%)	2 (1%)	25	17
All	All	904/1052 (86%)	878 (97%)	20 (2%)	6 (1%)	30	23

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	530	ARG
1	C	529	ALA
1	C	665	PRO
1	B	529	ALA
1	D	664	ASP
1	D	666	LYS

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	204/233 (88%)	198 (97%)	6 (3%)	55	57
1	B	205/233 (88%)	199 (97%)	6 (3%)	55	57
1	C	203/233 (87%)	198 (98%)	5 (2%)	60	63
1	D	201/233 (86%)	192 (96%)	9 (4%)	38	35
All	All	813/932 (87%)	787 (97%)	26 (3%)	51	52

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

Mol	Chain	Res	Type
1	A	524	LYS
1	A	530	ARG
1	A	654	MET
1	A	664	ASP
1	A	668	ARG
1	A	725	LEU
1	B	531	ARG
1	B	532	LYS
1	B	566	LEU
1	B	585	VAL
1	B	658	GLN
1	B	672	GLU
1	C	530	ARG
1	C	549	GLU
1	C	647	MET
1	C	657	LEU
1	C	664	ASP
1	D	530	ARG
1	D	543	CYS
1	D	582	LYS
1	D	642	LEU
1	D	657	LEU
1	D	658	GLN
1	D	664	ASP
1	D	669	LYS
1	D	671	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	574	GLN
1	A	658	GLN
1	A	667	ASN
1	A	673	ASN
1	A	674	GLN
1	A	680	GLN
1	B	574	GLN
1	B	599	GLN
1	B	637	GLN
1	B	658	GLN
1	B	667	ASN
1	B	673	ASN
1	C	637	GLN
1	C	667	ASN
1	C	674	GLN
1	C	680	GLN
1	C	686	HIS
1	D	574	GLN
1	D	599	GLN
1	D	643	GLN
1	D	686	HIS

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, 95th 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	232/263 (88%)	0.22	12 (5%) 26 29	22, 32, 59, 72	0
1	B	231/263 (87%)	0.27	12 (5%) 26 29	20, 37, 66, 80	0
1	C	229/263 (87%)	0.38	18 (7%) 13 14	24, 37, 73, 88	0
1	D	228/263 (86%)	0.37	16 (7%) 16 18	22, 35, 68, 83	0
All	All	920/1052 (87%)	0.31	58 (6%) 19 22	20, 35, 68, 88	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	504	VAL	6.1
1	C	659	LEU	6.0
1	D	665	PRO	5.7
1	D	666	LYS	5.6
1	C	665	PRO	5.1
1	C	669	LYS	4.9
1	B	550	THR	4.7
1	A	666	LYS	4.4
1	D	563	ARG	4.4
1	C	522	GLU	4.1
1	C	530	ARG	3.7
1	C	666	LYS	3.7
1	A	665	PRO	3.7
1	B	551	TYR	3.7
1	C	504	VAL	3.6
1	C	662	VAL	3.5
1	A	530	ARG	3.5
1	D	522	GLU	3.4
1	C	521	LYS	3.4
1	D	521	LYS	3.1
1	A	521	LYS	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	522	GLU	3.0
1	B	741	ASP	3.0
1	D	562	PRO	2.9
1	D	671	ILE	2.9
1	C	625	HIS	2.9
1	A	552	GLY	2.8
1	C	664	ASP	2.8
1	D	523	ARG	2.8
1	B	531	ARG	2.8
1	D	662	VAL	2.7
1	C	529	ALA	2.7
1	D	663	SER	2.7
1	B	522	GLU	2.5
1	C	647	MET	2.5
1	C	520	GLN	2.5
1	D	548	TYR	2.5
1	C	523	ARG	2.5
1	B	561	ALA	2.4
1	D	669	LYS	2.4
1	A	667	ASN	2.3
1	D	678	TRP	2.3
1	B	560	SER	2.3
1	D	664	ASP	2.3
1	A	669	LYS	2.3
1	D	667	ASN	2.3
1	C	561	ALA	2.2
1	A	520	GLN	2.2
1	B	563	ARG	2.2
1	B	567	PHE	2.2
1	B	530	ARG	2.1
1	D	524	LYS	2.1
1	A	529	ALA	2.1
1	C	670	ALA	2.1
1	B	725	LEU	2.1
1	A	505	VAL	2.1
1	C	624	LYS	2.1
1	B	523	ARG	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 ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.