Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 10:11 PM GMT

PDB ID	:	1SJ8
Title	:	Crystal Structure of talin residues 482-789
Authors	:	Papagrigoriou, E.; Gingras, A.R.; Barsukov, I.L.; Critchley, D.R.; Emsley, J.
Deposited on	:	2004-03-03
Resolution	:	2.60 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
A user guide is available at
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

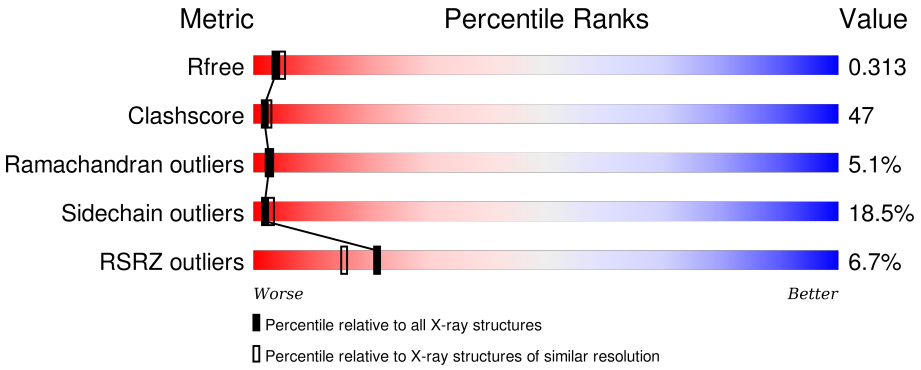
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	308	6% 40% 40% 12%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 2199 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 Talin 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	297	2113	1291	373	440	9	0	0	0

- Molecule 2 is water.

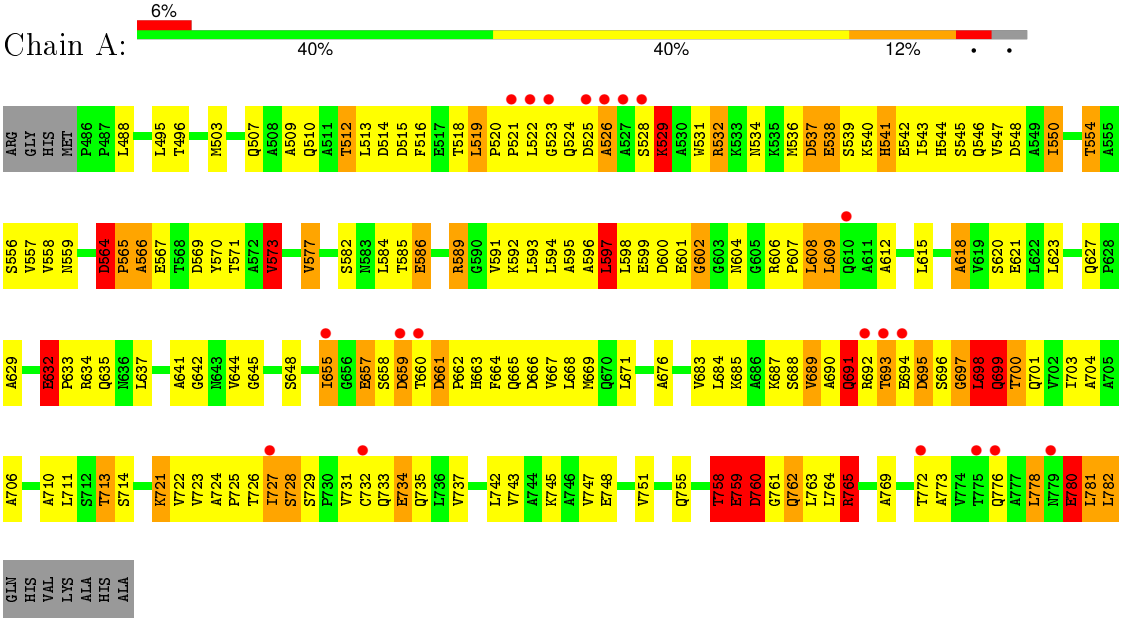
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	86	86	86	0	0

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: Talin 1

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α, β, γ	105.81Å 105.81Å 173.79Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.60 28.97 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-2.60) 99.7 (28.97-2.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	2.60 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5	Depositor
R, R_{free}	0.258 , 0.297 0.293 , 0.313	Depositor DCC
R_{free} test set	578 reflections (4.93%)	DCC
Wilson B-factor (Å^2)	56.7	Xtriage
Anisotropy	0.514	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 58.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Outliers	0 of 13179 reflections	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	2199	wwPDB-VP
Average B, all atoms (Å^2)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.89% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\| > 5$	RMSZ	$\#\|Z\| > 5$
1	A	1.41	14/2131 (0.7%)	1.47	31/2902 (1.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	3	2

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	714	SER	CB-OG	-7.01	1.33	1.42
1	A	621	GLU	CD-OE2	6.77	1.33	1.25
1	A	550	ILE	C-O	-6.42	1.11	1.23
1	A	573	VAL	CA-CB	6.37	1.68	1.54
1	A	618	ALA	CA-CB	-6.05	1.39	1.52
1	A	586	GLU	CD-OE1	5.99	1.32	1.25
1	A	748	GLU	CD-OE1	5.76	1.31	1.25
1	A	748	GLU	CG-CD	5.72	1.60	1.51
1	A	586	GLU	CD-OE2	5.56	1.31	1.25
1	A	699	GLN	CB-CG	-5.46	1.37	1.52
1	A	503	MET	CG-SD	-5.40	1.67	1.81
1	A	632	GLU	CD-OE1	5.34	1.31	1.25
1	A	745	LYS	CD-CE	5.22	1.64	1.51
1	A	710	ALA	CA-CB	5.15	1.63	1.52

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	698	LEU	O-C-N	-10.48	105.94	122.70
1	A	697	GLY	C-N-CA	-9.93	96.88	121.70

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	515	ASP	CB-CG-OD2	9.25	126.62	118.30
1	A	698	LEU	CB-CA-C	8.93	127.16	110.20
1	A	659	ASP	CB-CA-C	-8.62	93.15	110.40
1	A	698	LEU	CA-C-N	8.37	135.62	117.20
1	A	760	ASP	CB-CG-OD2	8.10	125.59	118.30
1	A	514	ASP	CB-CG-OD2	7.74	125.26	118.30
1	A	699	GLN	N-CA-C	7.05	130.05	111.00
1	A	659	ASP	C-N-CA	-6.98	104.25	121.70
1	A	569	ASP	CB-CG-OD2	6.74	124.37	118.30
1	A	564	ASP	CB-CG-OD2	6.69	124.33	118.30
1	A	693	THR	CA-C-N	-6.62	102.64	117.20
1	A	537	ASP	CB-CG-OD2	6.50	124.15	118.30
1	A	608	LEU	CB-CG-CD1	-6.25	100.37	111.00
1	A	697	GLY	N-CA-C	-6.20	97.59	113.10
1	A	780	GLU	CB-CA-C	6.13	122.65	110.40
1	A	658	SER	N-CA-C	6.06	127.35	111.00
1	A	690	ALA	C-N-CA	-6.02	106.66	121.70
1	A	597	LEU	CA-CB-CG	5.82	128.69	115.30
1	A	782	LEU	CA-CB-CG	-5.69	102.22	115.30
1	A	689	VAL	C-N-CA	-5.57	107.78	121.70
1	A	698	LEU	C-N-CA	5.47	135.37	121.70
1	A	691	GLN	N-CA-CB	5.42	120.35	110.60
1	A	657	GLU	C-N-CA	5.38	135.14	121.70
1	A	620	SER	CB-CA-C	-5.26	100.10	110.10
1	A	782	LEU	CB-CG-CD1	-5.26	102.06	111.00
1	A	666	ASP	CB-CG-OD2	5.19	122.97	118.30
1	A	582	SER	CB-CA-C	5.18	119.94	110.10
1	A	609	LEU	CB-CG-CD2	-5.11	102.32	111.00
1	A	765	ARG	CB-CA-C	5.04	120.47	110.40

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	695	ASP	CA
1	A	698	LEU	CA
1	A	699	GLN	CA

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	698	LEU	Mainchain,Peptide

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	2113	0	2127	198	0
2	A	86	0	0	3	0
All	All	2199	0	2127	198	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 47.

All (198) 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:769:ALA:O	1:A:772:THR:HG22	1.31	1.24
1:A:598:LEU:HD12	1:A:598:LEU:H	1.23	0.98
1:A:692:ARG:HD3	1:A:762:GLN:HE22	1.29	0.97
1:A:664:PHE:O	1:A:667:VAL:HG12	1.67	0.94
1:A:541:HIS:HD2	1:A:542:GLU:N	1.66	0.93
1:A:598:LEU:HD12	1:A:598:LEU:N	1.81	0.93
1:A:761:GLY:HA2	1:A:764:LEU:HD13	1.51	0.92
1:A:632:GLU:HB3	1:A:633:PRO:O	1.74	0.87
1:A:759:GLU:O	1:A:760:ASP:HB2	1.72	0.87
1:A:755:GLN:O	1:A:758:THR:HG23	1.74	0.87
1:A:697:GLY:O	1:A:698:LEU:C	2.03	0.87
1:A:726:THR:O	1:A:728:SER:N	2.07	0.87
1:A:541:HIS:CD2	1:A:542:GLU:N	2.42	0.86
1:A:769:ALA:O	1:A:772:THR:CG2	2.23	0.85
1:A:660:THR:O	1:A:661:ASP:C	2.14	0.85
1:A:726:THR:O	1:A:727:ILE:HG13	1.77	0.84
1:A:559:ASN:ND2	1:A:735:GLN:HE21	1.77	0.83
1:A:726:THR:HG22	1:A:727:ILE:N	1.94	0.83
1:A:496:THR:HG23	1:A:627:GLN:HE22	1.46	0.81
1:A:554:THR:HG21	1:A:644:VAL:HG11	1.62	0.81
1:A:721:LYS:C	1:A:721:LYS:HD3	2.00	0.80
1:A:693:THR:HB	1:A:699:GLN:HB2	1.64	0.79
1:A:532:ARG:O	1:A:536:MET:HB2	1.82	0.79
1:A:726:THR:CG2	1:A:729:SER:HB3	2.17	0.75
1:A:727:ILE:HD12	1:A:727:ILE:C	2.08	0.74

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:559:ASN:HD22	1:A:735:GLN:HE21	1.32	0.74
1:A:698:LEU:CB	1:A:701:GLN:H	2.00	0.74
1:A:529:LYS:HA	1:A:532:ARG:NE	2.04	0.73
1:A:726:THR:HG21	1:A:729:SER:HB3	1.68	0.73
1:A:564:ASP:OD1	1:A:564:ASP:N	2.21	0.73
1:A:539:SER:O	1:A:543:ILE:HG12	1.88	0.72
1:A:676:ALA:HA	1:A:713:THR:HG21	1.73	0.71
1:A:721:LYS:O	1:A:721:LYS:HD3	1.90	0.71
1:A:565:PRO:O	1:A:566:ALA:HB2	1.91	0.71
1:A:523:GLY:O	1:A:532:ARG:NH2	2.23	0.70
1:A:660:THR:HG22	1:A:664:PHE:HD2	1.55	0.70
1:A:755:GLN:CA	1:A:758:THR:HG23	2.22	0.70
1:A:698:LEU:HB2	1:A:701:GLN:N	2.06	0.70
1:A:598:LEU:H	1:A:598:LEU:CD1	2.02	0.69
1:A:769:ALA:C	1:A:772:THR:HG22	2.11	0.69
1:A:697:GLY:C	1:A:698:LEU:HD23	2.12	0.69
1:A:565:PRO:O	1:A:566:ALA:CB	2.39	0.69
1:A:559:ASN:HD22	1:A:735:GLN:NE2	1.91	0.68
1:A:697:GLY:CA	1:A:698:LEU:HD23	2.24	0.68
1:A:641:ALA:O	1:A:722:VAL:HG22	1.93	0.68
1:A:512:THR:HG23	1:A:592:LYS:HD2	1.74	0.68
1:A:684:LEU:HD23	1:A:684:LEU:C	2.14	0.67
1:A:541:HIS:HD2	1:A:542:GLU:H	1.40	0.67
1:A:698:LEU:HB2	1:A:701:GLN:H	1.57	0.67
1:A:657:GLU:HA	1:A:657:GLU:OE2	1.93	0.67
1:A:703:ILE:HG13	1:A:704:ALA:N	2.10	0.67
1:A:758:THR:HG22	1:A:763:LEU:HD11	1.77	0.66
1:A:697:GLY:HA2	1:A:698:LEU:HD23	1.78	0.66
1:A:737:VAL:HG12	1:A:781:LEU:HD13	1.78	0.66
1:A:529:LYS:HA	1:A:532:ARG:CZ	2.25	0.66
1:A:585:THR:HG22	1:A:589:ARG:NH1	2.11	0.66
1:A:698:LEU:HD23	1:A:698:LEU:N	2.10	0.65
1:A:726:THR:HG21	1:A:732:CYS:SG	2.36	0.65
1:A:532:ARG:H	1:A:532:ARG:HD2	1.61	0.65
1:A:541:HIS:CD2	1:A:541:HIS:C	2.69	0.65
1:A:546:GLN:O	1:A:550:ILE:HG12	1.96	0.65
1:A:755:GLN:HA	1:A:758:THR:HG23	1.79	0.65
1:A:698:LEU:HB2	1:A:701:GLN:CB	2.27	0.65
1:A:698:LEU:CB	1:A:701:GLN:N	2.60	0.65
1:A:726:THR:C	1:A:728:SER:H	1.99	0.64
1:A:697:GLY:O	1:A:699:GLN:N	2.31	0.64

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:698:LEU:HD22	1:A:700:THR:HG23	1.80	0.64
1:A:655:ILE:HG13	1:A:657:GLU:HB2	1.79	0.64
1:A:661:ASP:O	1:A:664:PHE:HB3	1.97	0.64
1:A:762:GLN:HA	1:A:765:ARG:HD2	1.80	0.63
1:A:516:PHE:CZ	1:A:606:ARG:HD2	2.34	0.63
1:A:520:PRO:O	1:A:520:PRO:HG2	1.98	0.62
1:A:559:ASN:ND2	1:A:735:GLN:NE2	2.45	0.61
1:A:698:LEU:HB3	1:A:700:THR:HG23	1.81	0.61
1:A:683:VAL:HG22	1:A:706:ALA:O	2.01	0.61
1:A:519:LEU:HD21	1:A:597:LEU:HD13	1.81	0.61
1:A:698:LEU:HB3	1:A:700:THR:N	2.15	0.61
1:A:759:GLU:O	1:A:760:ASP:CB	2.48	0.60
1:A:645:GLY:HA3	1:A:722:VAL:HA	1.82	0.60
1:A:755:GLN:O	1:A:758:THR:CG2	2.49	0.60
1:A:550:ILE:O	1:A:554:THR:HG23	2.01	0.60
1:A:755:GLN:C	1:A:758:THR:HG23	2.22	0.59
1:A:594:LEU:O	1:A:598:LEU:CD1	2.50	0.59
1:A:595:ALA:HB2	1:A:608:LEU:HB3	1.84	0.59
1:A:723:VAL:O	1:A:727:ILE:HG23	2.03	0.58
1:A:536:MET:HG3	1:A:594:LEU:CD1	2.34	0.58
1:A:693:THR:CB	1:A:699:GLN:HB2	2.32	0.58
1:A:637:LEU:C	1:A:637:LEU:HD23	2.24	0.58
1:A:685:LYS:HD2	1:A:769:ALA:HB2	1.87	0.57
1:A:698:LEU:HB3	1:A:700:THR:H	1.69	0.57
1:A:762:GLN:O	1:A:765:ARG:HG2	2.05	0.57
1:A:692:ARG:O	1:A:694:GLU:N	2.37	0.56
1:A:536:MET:O	1:A:540:LYS:HB2	2.06	0.56
1:A:778:LEU:HD22	1:A:778:LEU:O	2.06	0.56
1:A:509:ALA:O	1:A:512:THR:HB	2.06	0.56
1:A:542:GLU:O	1:A:543:ILE:C	2.41	0.55
1:A:520:PRO:O	1:A:521:PRO:C	2.42	0.55
1:A:634:ARG:HG2	1:A:711:LEU:HD11	1.89	0.55
1:A:545:SER:O	1:A:548:ASP:HB2	2.05	0.55
1:A:665:GLN:O	1:A:669:MET:HG2	2.07	0.54
1:A:593:LEU:HD13	1:A:597:LEU:CD2	2.38	0.54
1:A:528:SER:O	1:A:531:TRP:N	2.37	0.54
1:A:543:ILE:HD12	1:A:591:VAL:HG23	1.91	0.53
1:A:593:LEU:HD13	1:A:593:LEU:C	2.29	0.53
1:A:525:ASP:O	1:A:526:ALA:C	2.47	0.53
1:A:632:GLU:CB	1:A:633:PRO:CA	2.86	0.53
1:A:534:ASN:O	1:A:538:GLU:HG2	2.09	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:595:ALA:HA	1:A:598:LEU:HD13	1.91	0.53
1:A:573:VAL:O	1:A:577:VAL:HG13	2.08	0.52
1:A:733:GLN:O	1:A:734:GLU:C	2.45	0.52
1:A:671:LEU:HB3	1:A:780:GLU:HB2	1.91	0.52
1:A:564:ASP:HB2	1:A:567:GLU:OE1	2.10	0.52
1:A:687:LYS:O	1:A:691:GLN:HG3	2.10	0.52
1:A:723:VAL:HB	1:A:732:CYS:HB3	1.92	0.52
1:A:667:VAL:O	1:A:671:LEU:HB2	2.10	0.52
1:A:525:ASP:O	1:A:526:ALA:O	2.27	0.51
1:A:529:LYS:O	1:A:529:LYS:HD3	2.10	0.51
1:A:541:HIS:CE1	2:A:87:HOH:O	2.63	0.51
1:A:546:GLN:NE2	1:A:586:GLU:HB3	2.26	0.51
1:A:761:GLY:HA2	1:A:764:LEU:CD1	2.33	0.51
1:A:724:ALA:N	1:A:725:PRO:HD2	2.26	0.50
1:A:495:LEU:HD11	1:A:573:VAL:HG22	1.93	0.50
1:A:772:THR:HG23	1:A:773:ALA:N	2.26	0.50
1:A:743:VAL:O	1:A:747:VAL:HG23	2.12	0.50
1:A:547:VAL:CG1	1:A:648:SER:HB3	2.42	0.50
1:A:684:LEU:HD23	1:A:685:LYS:N	2.27	0.50
1:A:692:ARG:HH12	1:A:760:ASP:CG	2.15	0.50
1:A:554:THR:O	1:A:558:VAL:HG23	2.13	0.49
1:A:758:THR:HG22	1:A:763:LEU:CD1	2.42	0.49
1:A:513:LEU:HD13	1:A:612:ALA:HB3	1.94	0.49
1:A:696:SER:C	1:A:697:GLY:O	2.45	0.49
1:A:662:PRO:C	1:A:664:PHE:H	2.16	0.49
1:A:541:HIS:ND1	2:A:87:HOH:O	2.35	0.48
1:A:540:LYS:O	1:A:657:GLU:HG2	2.14	0.48
1:A:660:THR:CG2	1:A:727:ILE:HD11	2.44	0.48
1:A:698:LEU:HB2	1:A:701:GLN:HB2	1.96	0.48
1:A:660:THR:CG2	1:A:664:PHE:HD2	2.23	0.48
1:A:698:LEU:HB3	1:A:701:GLN:N	2.29	0.47
1:A:513:LEU:HD13	1:A:612:ALA:CB	2.44	0.47
1:A:532:ARG:HG2	1:A:601:GLU:CD	2.35	0.47
1:A:692:ARG:HD3	1:A:762:GLN:NE2	2.13	0.47
1:A:698:LEU:HB3	1:A:701:GLN:H	1.78	0.47
1:A:594:LEU:O	1:A:598:LEU:HD12	2.14	0.47
1:A:600:ASP:O	1:A:601:GLU:C	2.50	0.47
1:A:609:LEU:N	1:A:609:LEU:HD22	2.31	0.46
1:A:632:GLU:HB3	1:A:633:PRO:C	2.35	0.46
1:A:601:GLU:O	1:A:602:GLY:O	2.34	0.46
1:A:759:GLU:H	1:A:763:LEU:HD12	1.81	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:726:THR:C	1:A:727:ILE:HG13	2.35	0.45
1:A:735:GLN:HA	1:A:735:GLN:NE2	2.30	0.45
1:A:660:THR:HG21	1:A:727:ILE:HD11	1.99	0.45
1:A:676:ALA:HA	1:A:713:THR:CG2	2.44	0.45
1:A:724:ALA:O	1:A:726:THR:N	2.49	0.45
1:A:724:ALA:C	1:A:726:THR:H	2.19	0.45
1:A:723:VAL:CG1	1:A:732:CYS:HB3	2.46	0.45
1:A:518:THR:HG23	1:A:518:THR:O	2.16	0.45
1:A:608:LEU:HA	1:A:608:LEU:HD12	1.58	0.45
1:A:699:GLN:O	1:A:699:GLN:HG2	2.11	0.45
1:A:495:LEU:HD22	1:A:570:TYR:HD1	1.82	0.45
1:A:660:THR:O	1:A:661:ASP:O	2.34	0.45
1:A:632:GLU:CB	1:A:633:PRO:HA	2.47	0.44
1:A:564:ASP:HA	1:A:565:PRO:HD2	1.86	0.44
1:A:599:GLU:OE1	1:A:604:ASN:HA	2.18	0.44
1:A:516:PHE:HB3	1:A:596:ALA:HB2	2.00	0.44
1:A:607:PRO:HG3	2:A:99:HOH:O	2.17	0.43
1:A:634:ARG:CG	1:A:711:LEU:HD11	2.47	0.43
1:A:615:LEU:O	1:A:618:ALA:HB3	2.19	0.43
1:A:660:THR:CG2	1:A:664:PHE:CD2	3.01	0.43
1:A:544:HIS:CD2	1:A:657:GLU:HG3	2.54	0.43
1:A:726:THR:C	1:A:728:SER:N	2.64	0.43
1:A:635:GLN:HG3	1:A:635:GLN:O	2.18	0.43
1:A:684:LEU:O	1:A:687:LYS:N	2.47	0.42
1:A:762:GLN:HA	1:A:765:ARG:CD	2.48	0.42
1:A:727:ILE:HD12	1:A:728:SER:N	2.34	0.42
1:A:782:LEU:HD13	1:A:782:LEU:HA	1.22	0.42
1:A:632:GLU:HB2	1:A:633:PRO:HA	2.02	0.42
1:A:781:LEU:HA	1:A:781:LEU:HD22	1.89	0.42
1:A:488:LEU:HD11	1:A:629:ALA:HB2	2.01	0.42
1:A:688:SER:O	1:A:689:VAL:C	2.58	0.42
1:A:693:THR:HB	1:A:699:GLN:CB	2.44	0.42
1:A:727:ILE:CD1	1:A:727:ILE:C	2.77	0.42
1:A:698:LEU:HD22	1:A:700:THR:CG2	2.47	0.42
1:A:593:LEU:O	1:A:597:LEU:HD22	2.19	0.42
1:A:586:GLU:HG2	1:A:589:ARG:NH2	2.35	0.42
1:A:532:ARG:H	1:A:532:ARG:CD	2.25	0.42
1:A:662:PRO:O	1:A:664:PHE:N	2.51	0.42
1:A:724:ALA:C	1:A:726:THR:N	2.72	0.42
1:A:543:ILE:CD1	1:A:591:VAL:HG23	2.50	0.42
1:A:585:THR:HG22	1:A:589:ARG:HH12	1.83	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:661:ASP:HA	1:A:662:PRO:HD2	1.81	0.41
1:A:523:GLY:C	1:A:525:ASP:H	2.24	0.41
1:A:556:SER:O	1:A:557:VAL:C	2.58	0.41
1:A:655:ILE:HD13	1:A:655:ILE:HG21	1.76	0.41
1:A:642:GLY:HA2	1:A:722:VAL:HG23	2.03	0.41
1:A:606:ARG:N	1:A:607:PRO:CD	2.84	0.41
1:A:772:THR:CG2	1:A:773:ALA:N	2.84	0.40
1:A:727:ILE:O	1:A:733:GLN:NE2	2.54	0.40
1:A:727:ILE:HD12	1:A:727:ILE:O	2.20	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	295/308 (96%)	247 (84%)	33 (11%)	15 (5%)	2 3

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	526	ALA
1	A	659	ASP
1	A	699	GLN
1	A	727	ILE
1	A	758	THR
1	A	760	ASP
1	A	565	PRO
1	A	566	ALA
1	A	695	ASP
1	A	632	GLU
1	A	529	LYS
1	A	661	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	602	GLY
1	A	655	ILE
1	A	759	GLU

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	222/230 (96%)	181 (82%)	41 (18%)	2 3

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	507	GLN
1	A	510	GLN
1	A	512	THR
1	A	519	LEU
1	A	522	LEU
1	A	524	GLN
1	A	529	LYS
1	A	532	ARG
1	A	537	ASP
1	A	538	GLU
1	A	541	HIS
1	A	554	THR
1	A	564	ASP
1	A	571	THR
1	A	573	VAL
1	A	577	VAL
1	A	584	LEU
1	A	589	ARG
1	A	597	LEU
1	A	623	LEU
1	A	663	HIS
1	A	668	LEU
1	A	691	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	695	ASP
1	A	698	LEU
1	A	700	THR
1	A	713	THR
1	A	721	LYS
1	A	728	SER
1	A	731	VAL
1	A	734	GLU
1	A	742	LEU
1	A	751	VAL
1	A	758	THR
1	A	759	GLU
1	A	762	GLN
1	A	765	ARG
1	A	776	GLN
1	A	778	LEU
1	A	780	GLU
1	A	781	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	492	GLN
1	A	510	GLN
1	A	541	HIS
1	A	544	HIS
1	A	627	GLN
1	A	653	GLN
1	A	733	GLN
1	A	735	GLN
1	A	762	GLN
1	A	776	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	297/308 (96%)	0.24	20 (6%)	21	15	18, 41, 45, 66	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	660	THR	4.9
1	A	523	GLY	4.7
1	A	727	ILE	3.8
1	A	527	ALA	3.8
1	A	693	THR	3.7
1	A	776	GLN	3.0
1	A	659	ASP	3.0
1	A	526	ALA	3.0
1	A	522	LEU	2.9
1	A	521	PRO	2.7
1	A	692	ARG	2.7
1	A	779	ASN	2.6
1	A	528	SER	2.6
1	A	525	ASP	2.5
1	A	732	CYS	2.4
1	A	694	GLU	2.4
1	A	772	THR	2.3
1	A	655	ILE	2.3
1	A	775	THR	2.3
1	A	610	GLN	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.