



Full wwPDB X-ray Structure Validation Report

Mar 1, 2014 – 12:21 AM GMT

PDB ID : 2FU3
Title : Crystal structure of gephyrin E-domain
Authors : Kim, E.Y.; Schindelin, H.
Deposited on : 2006-01-25
Resolution : 2.70 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>

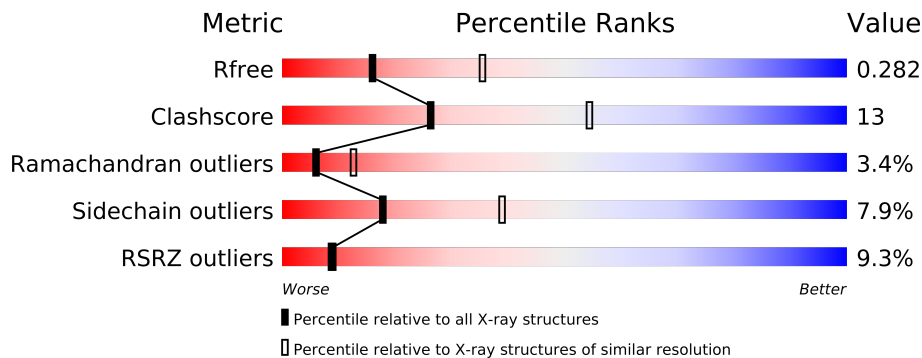
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.70 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	1557 (2.70-2.70)
Clashscore	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)
RSRZ outliers	66119	1559 (2.70-2.70)

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	419	
1	B	419	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6257 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 gephyrin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	405	Total	C	N	O	S	0	0	0
			3054	1930	531	572	21			
1	B	419	Total	C	N	O	S	0	0	0
			3137	1978	548	591	20			

- Molecule 2 is water.

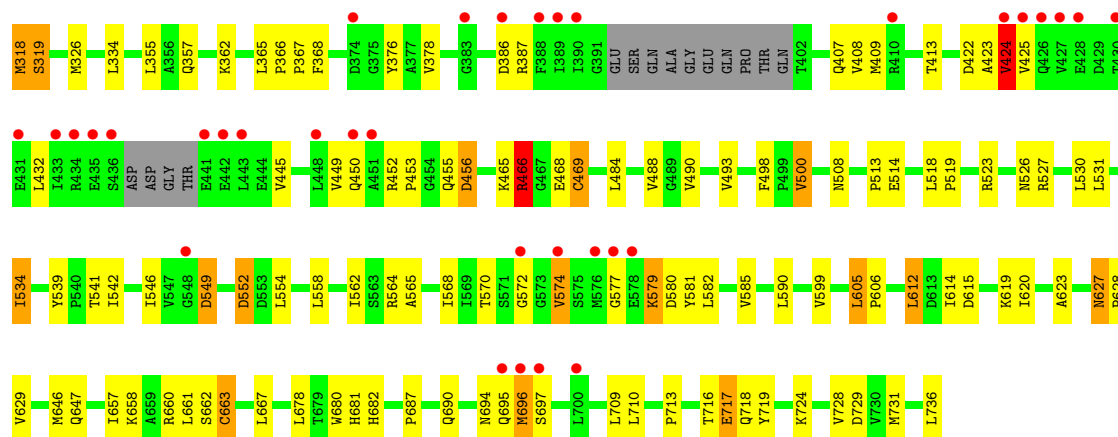
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	37	Total	O	0	0
			37	37		
2	B	29	Total	O	0	0
			29	29		

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 ($\text{RSRZ} > 2$). 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.

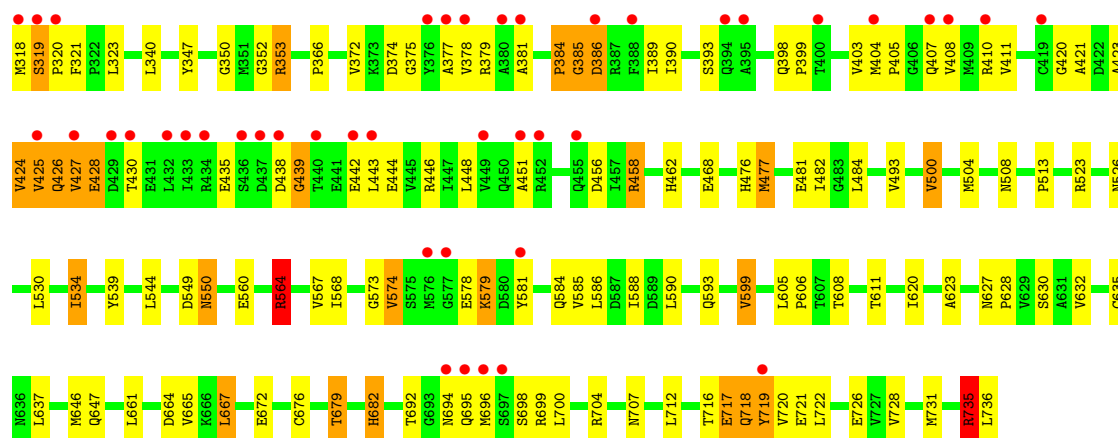
• Molecule 1: gephyrin

Chain A: 



• Molecule 1: gephyrin

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	108.82Å 156.19Å 51.36Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.70 39.92 – 2.38	Depositor EDS
% Data completeness (in resolution range)	99.9 (20.00-2.70) 96.9 (39.92-2.38)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.11 (at 2.37Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.202 , 0.273 0.218 , 0.282	Depositor DCC
R_{free} test set	1266 reflections (5.37%)	DCC
Wilson B-factor (Å ²)	58.7	Xtriage
Anisotropy	0.278	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 80.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 35018 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6257	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.08% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality i

5.1 Standard geometry i

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	1.03	4/3112 (0.1%)	0.76	1/4233 (0.0%)
1	B	1.03	6/3198 (0.2%)	0.81	2/4355 (0.0%)
All	All	1.03	10/6310 (0.2%)	0.78	3/8588 (0.0%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	676	CYS	CB-SG	-6.58	1.71	1.82
1	B	635	CYS	CB-SG	-6.56	1.71	1.82
1	A	514	GLU	CB-CG	6.28	1.64	1.52
1	A	469	CYS	CB-SG	-6.01	1.72	1.82
1	B	647	GLN	CB-CG	-5.53	1.37	1.52
1	A	663	CYS	CB-SG	-5.52	1.72	1.81
1	A	424	VAL	CA-CB	5.34	1.66	1.54
1	B	500	VAL	CB-CG1	-5.28	1.41	1.52
1	B	574	VAL	CA-CB	5.09	1.65	1.54
1	B	468	GLU	CG-CD	5.07	1.59	1.51

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	735	ARG	NE-CZ-NH1	6.96	123.78	120.30
1	A	466	ARG	NE-CZ-NH2	5.52	123.06	120.30
1	B	564	ARG	NE-CZ-NH1	5.03	122.81	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 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	3054	0	3072	79	0
1	B	3137	0	3127	91	0
2	A	37	0	0	1	0
2	B	29	0	0	1	0
All	All	6257	0	6199	165	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 13.

All (165) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:526:ASN:HD21	1:B:628:PRO:HA	1.40	0.86
1:A:362:LYS:O	1:A:466:ARG:HD2	1.76	0.86
1:A:449:VAL:HG12	1:A:450:GLN:H	1.41	0.85
1:A:627:ASN:HD21	1:A:695:GLN:HE21	1.22	0.84
1:B:424:VAL:O	1:B:425:VAL:HG23	1.81	0.79
1:A:378:VAL:HG21	1:A:425:VAL:HG21	1.66	0.76
1:B:534:ILE:CD1	1:B:539:TYR:HB2	2.19	0.71
1:B:627:ASN:HD22	1:B:630:SER:H	1.36	0.71
1:A:629:VAL:HG21	1:A:695:GLN:HA	1.71	0.70
1:A:627:ASN:ND2	1:A:695:GLN:HE21	1.90	0.69
1:B:424:VAL:CG2	1:B:456:ASP:HB2	2.22	0.69
1:B:423:ALA:O	1:B:424:VAL:HB	1.92	0.68
1:B:718:GLN:O	1:B:719:TYR:C	2.33	0.67
1:B:378:VAL:O	1:B:423:ALA:HA	1.95	0.66
1:B:425:VAL:HG12	1:B:426:GLN:N	2.11	0.64
1:A:534:ILE:CD1	1:A:539:TYR:HB2	2.28	0.64
1:B:584:GLN:HG3	1:B:588:ILE:HD12	1.81	0.63
1:B:375:GLY:HA3	1:B:425:VAL:HB	1.81	0.62
1:B:379:ARG:HG2	1:B:421:ALA:HB2	1.80	0.62
1:A:562:ILE:O	1:A:619:LYS:NZ	2.31	0.62
1:A:466:ARG:HG3	1:A:466:ARG:HH21	1.66	0.61
1:A:423:ALA:N	2:A:34:HOH:O	2.32	0.61
1:B:534:ILE:HD12	1:B:539:TYR:HB2	1.83	0.60
1:A:562:ILE:N	1:A:568:ILE:HD11	2.17	0.59
1:B:442:GLU:O	1:B:444:GLU:N	2.32	0.59
1:A:526:ASN:O	1:A:527:ARG:C	2.41	0.58
1:B:513:PRO:HA	1:B:523:ARG:HD3	1.86	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:389:ILE:HD12	1:B:442:GLU:O	2.04	0.58
1:B:564:ARG:HH11	1:B:564:ARG:HG2	1.68	0.57
1:A:552:ASP:OD2	1:A:552:ASP:N	2.36	0.57
1:A:449:VAL:HG12	1:A:450:GLN:N	2.16	0.57
1:A:612:LEU:HD13	1:A:614:ILE:HD11	1.86	0.57
1:B:424:VAL:HG21	1:B:456:ASP:HB2	1.87	0.56
1:A:581:TYR:O	1:A:585:VAL:HG23	2.05	0.56
1:B:372:VAL:HB	1:B:456:ASP:HB3	1.86	0.56
1:B:735:ARG:HB2	1:B:735:ARG:HH11	1.70	0.56
1:A:534:ILE:HD12	1:A:539:TYR:HB2	1.87	0.55
1:A:500:VAL:HG22	1:A:565:ALA:HA	1.88	0.55
1:A:696:MET:CE	1:B:462:HIS:CE1	2.89	0.55
1:A:574:VAL:HA	1:A:579:LYS:HB2	1.87	0.54
1:A:680:TRP:CE2	1:A:687:PRO:HB3	2.42	0.54
1:A:424:VAL:HG13	1:A:425:VAL:H	1.73	0.54
1:B:347:TYR:O	1:B:482:ILE:HD12	2.09	0.53
1:A:376:TYR:HB3	1:A:408:VAL:HG13	1.91	0.53
1:B:573:GLY:HA2	1:B:579:LYS:HB3	1.90	0.53
1:B:323:LEU:HD21	1:B:605:LEU:HD11	1.90	0.52
1:B:526:ASN:ND2	1:B:628:PRO:HA	2.19	0.52
1:B:692:THR:HA	1:B:700:LEU:HD13	1.91	0.52
1:B:728:VAL:HG13	2:B:39:HOH:O	2.09	0.51
1:B:404:MET:HB3	1:B:405:PRO:HD2	1.93	0.51
1:B:716:THR:HG22	1:B:718:GLN:H	1.75	0.51
1:B:679:THR:HB	1:B:704:ARG:HH22	1.75	0.51
1:B:398:GLN:OE1	1:B:399:PRO:HD2	2.11	0.51
1:A:508:ASN:HD21	1:A:549:ASP:H	1.58	0.51
1:A:526:ASN:HD21	1:A:628:PRO:HA	1.76	0.51
1:B:679:THR:HB	1:B:704:ARG:NH2	2.25	0.51
1:B:477:MET:HG3	1:B:481:GLU:HB2	1.93	0.51
1:B:735:ARG:CG	1:B:735:ARG:HH11	2.24	0.50
1:B:377:ALA:HA	1:B:423:ALA:CB	2.41	0.50
1:A:542:ILE:HD13	1:A:564:ARG:HG2	1.92	0.50
1:A:513:PRO:HA	1:A:523:ARG:HD3	1.93	0.50
1:A:424:VAL:HG12	1:A:455:GLN:HE21	1.77	0.49
1:A:627:ASN:ND2	1:A:629:VAL:H	2.10	0.49
1:A:605:LEU:C	1:A:605:LEU:HD13	2.32	0.49
1:A:530:LEU:O	1:A:534:ILE:HG23	2.12	0.49
1:A:657:ILE:HD11	1:A:678:LEU:HD13	1.95	0.49
1:A:570:THR:HG22	1:A:623:ALA:HA	1.94	0.49
1:B:340:LEU:HG	1:B:646:MET:HG2	1.94	0.49
1:B:661:LEU:HG	1:B:728:VAL:HG21	1.94	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:658:LYS:NZ	1:A:729:ASP:OD1	2.44	0.49
1:B:550:ASN:HD22	1:B:550:ASN:N	2.11	0.49
1:B:667:LEU:HD13	1:B:721:GLU:HA	1.95	0.48
1:B:477:MET:HG3	1:B:481:GLU:CB	2.44	0.48
1:A:736:LEU:CD1	1:B:736:LEU:HD12	2.43	0.48
1:B:581:TYR:O	1:B:585:VAL:HG23	2.13	0.48
1:B:378:VAL:HG23	1:B:425:VAL:HG23	1.96	0.48
1:B:508:ASN:HD21	1:B:549:ASP:H	1.62	0.48
1:A:627:ASN:HD22	1:A:629:VAL:H	1.62	0.48
1:A:424:VAL:HG12	1:A:455:GLN:NE2	2.29	0.47
1:A:680:TRP:CD2	1:A:687:PRO:HB3	2.49	0.47
1:B:404:MET:HB2	1:B:407:GLN:OE1	2.15	0.47
1:B:424:VAL:HG23	1:B:456:ASP:HB2	1.97	0.47
1:A:407:GLN:O	1:A:408:VAL:HG23	2.15	0.47
1:A:488:VAL:HG23	1:A:490:VAL:HG23	1.96	0.47
1:A:542:ILE:CD1	1:A:564:ARG:HG2	2.45	0.47
1:B:424:VAL:HG13	1:B:425:VAL:N	2.30	0.47
1:B:379:ARG:C	1:B:381:ALA:H	2.17	0.46
1:B:599:VAL:O	1:B:605:LEU:HA	2.15	0.46
1:A:527:ARG:HH21	1:A:546:ILE:HD11	1.79	0.46
1:B:389:ILE:HA	1:B:408:VAL:O	2.15	0.46
1:A:713:PRO:HG3	1:A:719:TYR:HE1	1.79	0.46
1:A:424:VAL:HG22	1:A:425:VAL:N	2.31	0.46
1:B:586:LEU:HD23	1:B:590:LEU:HD12	1.98	0.46
1:A:554:LEU:CD1	1:A:582:LEU:HD13	2.46	0.46
1:B:352:GLY:O	1:B:476:HIS:CD2	2.68	0.45
1:B:377:ALA:HA	1:B:423:ALA:HB1	1.97	0.45
1:A:663:CYS:HA	1:A:724:LYS:HB2	1.98	0.45
1:A:627:ASN:HD22	1:A:629:VAL:N	2.14	0.45
1:B:375:GLY:O	1:B:410:ARG:HD2	2.17	0.45
1:A:696:MET:HE1	1:B:462:HIS:CE1	2.52	0.45
1:A:386:ASP:C	1:A:387:ARG:HG3	2.37	0.45
1:A:452:ARG:HB3	1:A:455:GLN:NE2	2.32	0.45
1:B:318:MET:O	1:B:319:SER:CB	2.64	0.45
1:B:608:THR:HB	1:B:623:ALA:HB3	1.99	0.44
1:B:593:GLN:NE2	1:B:611:THR:OG1	2.45	0.44
1:B:425:VAL:O	1:B:426:GLN:CB	2.65	0.44
1:B:393:SER:HB3	1:B:411:VAL:HG23	1.99	0.44
1:B:717:GLU:OE1	1:B:717:GLU:N	2.51	0.44
1:A:518:LEU:HB3	1:A:519:PRO:HD2	1.98	0.44
1:A:716:THR:HG22	1:A:717:GLU:N	2.31	0.44
1:B:426:GLN:O	1:B:428:GLU:N	2.51	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:504:MET:HB3	1:B:544:LEU:HB2	1.99	0.44
1:A:680:TRP:CZ2	1:A:687:PRO:HG3	2.52	0.44
1:B:605:LEU:HB3	1:B:606:PRO:HD3	1.99	0.44
1:B:726:GLU:O	1:B:728:VAL:HG23	2.17	0.44
1:A:620:ILE:HD11	1:A:646:MET:CE	2.48	0.44
1:B:665:VAL:O	1:B:665:VAL:HG23	2.18	0.44
1:B:425:VAL:O	1:B:426:GLN:HB2	2.17	0.43
1:A:629:VAL:HG12	1:A:629:VAL:O	2.18	0.43
1:B:319:SER:O	1:B:321:PHE:N	2.51	0.43
1:B:564:ARG:HH11	1:B:564:ARG:CG	2.31	0.43
1:B:319:SER:C	1:B:321:PHE:N	2.70	0.43
1:B:350:GLY:O	1:B:353:ARG:HB2	2.18	0.43
1:B:403:VAL:HB	1:B:420:GLY:HA3	1.99	0.43
1:B:568:ILE:N	1:B:568:ILE:HD12	2.33	0.43
1:B:672:GLU:HB2	1:B:712:LEU:HB2	2.00	0.43
1:B:438:ASP:O	1:B:439:GLY:C	2.57	0.42
1:B:735:ARG:CB	1:B:735:ARG:HH11	2.31	0.42
1:A:646:MET:HE3	1:A:646:MET:HB2	1.93	0.42
1:A:526:ASN:HD22	1:A:526:ASN:N	2.17	0.42
1:B:735:ARG:HG3	1:B:735:ARG:HH11	1.84	0.42
1:A:468:GLU:HG2	1:A:469:CYS:N	2.33	0.42
1:A:424:VAL:HA	1:A:456:ASP:HB2	2.01	0.42
1:B:620:ILE:CD1	1:B:646:MET:HE1	2.49	0.42
1:A:709:LEU:HB2	1:A:731:MET:HB3	2.01	0.42
1:A:558:LEU:HD13	1:A:590:LEU:CD1	2.48	0.42
1:A:710:LEU:HD11	1:A:728:VAL:HG21	2.00	0.42
1:A:661:LEU:HD21	1:A:710:LEU:HD21	2.02	0.42
1:A:612:LEU:CD1	1:A:614:ILE:HD11	2.48	0.42
1:A:580:ASP:O	1:A:581:TYR:C	2.58	0.42
1:A:696:MET:HE3	1:B:462:HIS:CE1	2.55	0.42
1:A:318:MET:O	1:A:319:SER:CB	2.68	0.42
1:A:681:HIS:O	1:A:682:HIS:C	2.58	0.42
1:A:620:ILE:CD1	1:A:646:MET:HE2	2.50	0.42
1:A:365:LEU:HD12	1:A:366:PRO:HA	2.01	0.42
1:A:367:PRO:HG2	1:A:368:PHE:CD2	2.55	0.41
1:B:366:PRO:HG3	1:B:458:ARG:HB3	2.02	0.41
1:A:605:LEU:HB3	1:A:606:PRO:HD3	2.03	0.41
1:A:579:LYS:O	1:A:580:ASP:C	2.58	0.41
1:A:662:SER:OG	1:A:690:GLN:NE2	2.54	0.41
1:B:567:VAL:HG22	1:B:620:ILE:CG1	2.50	0.41
1:B:386:ASP:N	1:B:386:ASP:OD2	2.43	0.41
1:B:530:LEU:O	1:B:534:ILE:HG23	2.20	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:430:THR:HA	1:B:446:ARG:O	2.21	0.41
1:B:637:LEU:HD23	1:B:731:MET:CE	2.50	0.41
1:A:376:TYR:HA	1:A:409:MET:O	2.21	0.41
1:A:736:LEU:HD11	1:B:736:LEU:HD12	2.03	0.40
1:A:432:LEU:HA	1:A:445:VAL:HA	2.03	0.40
1:A:531:LEU:CD2	1:A:541:THR:HB	2.51	0.40
1:A:355:LEU:CD2	1:A:493:VAL:HG11	2.51	0.40
1:B:384:PRO:O	1:B:385:GLY:O	2.39	0.40
1:B:435:GLU:HA	1:B:442:GLU:HA	2.03	0.40
1:B:718:GLN:HG3	1:B:719:TYR:N	2.37	0.40
1:B:390:ILE:HG23	1:B:407:GLN:HB3	2.04	0.40
1:B:374:ASP:OD1	1:B:427:VAL:HG23	2.21	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	399/419 (95%)	365 (92%)	23 (6%)	11 (3%)	8	18
1	B	417/419 (100%)	364 (87%)	36 (9%)	17 (4%)	4	9
All	All	816/838 (97%)	729 (89%)	59 (7%)	28 (3%)	6	12

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	319	SER
1	A	413	THR
1	A	456	ASP
1	B	319	SER
1	B	385	GLY
1	B	424	VAL
1	B	425	VAL
1	B	443	LEU

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Mol	Chain	Res	Type
1	A	422	ASP
1	A	574	VAL
1	A	577	GLY
1	A	696	MET
1	A	697	SER
1	B	426	GLN
1	B	719	TYR
1	B	428	GLU
1	B	439	GLY
1	B	448	LEU
1	B	682	HIS
1	B	696	MET
1	B	698	SER
1	A	424	VAL
1	A	453	PRO
1	B	320	PRO
1	B	384	PRO
1	B	451	ALA
1	A	572	GLY
1	B	427	VAL

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	334/356 (94%)	310 (93%)	24 (7%)	21	45
1	B	338/356 (95%)	309 (91%)	29 (9%)	15	33
All	All	672/712 (94%)	619 (92%)	53 (8%)	18	39

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	318	MET
1	A	326	MET
1	A	334	LEU
1	A	357	GLN
1	A	465	LYS

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Mol	Chain	Res	Type
1	A	466	ARG
1	A	484	LEU
1	A	498	PHE
1	A	500	VAL
1	A	534	ILE
1	A	549	ASP
1	A	552	ASP
1	A	579	LYS
1	A	599	VAL
1	A	605	LEU
1	A	612	LEU
1	A	615	ASP
1	A	627	ASN
1	A	647	GLN
1	A	660	ARG
1	A	667	LEU
1	A	694	ASN
1	A	717	GLU
1	A	718	GLN
1	B	353	ARG
1	B	386	ASP
1	B	458	ARG
1	B	477	MET
1	B	484	LEU
1	B	493	VAL
1	B	500	VAL
1	B	534	ILE
1	B	550	ASN
1	B	560	GLU
1	B	564	ARG
1	B	574	VAL
1	B	578	GLU
1	B	579	LYS
1	B	599	VAL
1	B	632	VAL
1	B	664	ASP
1	B	667	LEU
1	B	679	THR
1	B	682	HIS
1	B	694	ASN
1	B	695	GLN
1	B	699	ARG

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Mol	Chain	Res	Type
1	B	707	ASN
1	B	717	GLU
1	B	718	GLN
1	B	720	VAL
1	B	722	LEU
1	B	735	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	455	GLN
1	A	508	ASN
1	A	526	ASN
1	A	556	ASN
1	A	627	ASN
1	A	690	GLN
1	B	476	HIS
1	B	508	ASN
1	B	526	ASN
1	B	550	ASN
1	B	556	ASN
1	B	627	ASN
1	B	707	ASN
1	B	723	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	405/419 (96%)	0.13	34 (8%) 11 12	49, 64, 71, 79	0
1	B	419/419 (100%)	0.22	43 (10%) 7 7	55, 65, 71, 74	0
All	All	824/838 (98%)	0.18	77 (9%) 9 9	49, 65, 71, 79	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	430	THR	7.7
1	A	695	GLN	7.1
1	B	318	MET	6.7
1	B	576	MET	6.0
1	B	388	PHE	6.0
1	B	696	MET	5.5
1	A	425	VAL	5.2
1	B	432	LEU	5.2
1	A	697	SER	5.2
1	A	386	ASP	4.8
1	A	433	ILE	4.7
1	B	381	ALA	4.3
1	A	436	SER	4.3
1	A	428	GLU	4.3
1	B	378	VAL	4.2
1	A	431	GLU	4.1
1	A	435	GLU	4.0
1	B	400	THR	4.0
1	A	443	LEU	3.9
1	B	376	TYR	3.9
1	A	390	ILE	3.8
1	B	695	GLN	3.6
1	A	383	GLY	3.6
1	A	430	THR	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	437	ASP	3.6
1	A	426	GLN	3.6
1	A	576	MET	3.5
1	B	434	ARG	3.5
1	B	407	GLN	3.4
1	A	442	GLU	3.3
1	B	719	TYR	3.3
1	B	408	VAL	3.3
1	A	388	PHE	3.2
1	A	427	VAL	3.2
1	A	448	LEU	3.2
1	A	574	VAL	3.1
1	A	696	MET	3.0
1	A	451	ALA	3.0
1	B	451	ALA	3.0
1	B	455	GLN	3.0
1	B	377	ALA	2.9
1	B	419	CYS	2.8
1	B	440	THR	2.8
1	B	436	SER	2.8
1	B	427	VAL	2.8
1	A	577	GLY	2.8
1	A	450	GLN	2.8
1	A	548	GLY	2.7
1	B	395	ALA	2.7
1	A	424	VAL	2.7
1	B	319	SER	2.7
1	B	577	GLY	2.6
1	B	429	ASP	2.5
1	A	578	GLU	2.5
1	A	441	GLU	2.5
1	B	433	ILE	2.5
1	B	443	LEU	2.5
1	B	386	ASP	2.4
1	B	442	GLU	2.4
1	B	438	ASP	2.4
1	B	452	ARG	2.3
1	A	434	ARG	2.3
1	B	694	ASN	2.3
1	B	394	GLN	2.3
1	B	404	MET	2.3
1	B	449	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	374	ASP	2.2
1	B	425	VAL	2.2
1	A	700	LEU	2.2
1	A	389	ILE	2.1
1	B	697	SER	2.1
1	B	380	ALA	2.1
1	B	581	TYR	2.1
1	A	572	GLY	2.0
1	A	410	ARG	2.0
1	B	410	ARG	2.0
1	B	320	PRO	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.