



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:56 AM UTC

PDB ID : 1UZJ / pdb_00001uzj
Title : Integrin binding cbEGF22-TB4-cbEGF33 fragment of human fibrillin-1, holo form.
Authors : Lee, S.S.J.; Knott, V.; Harlos, K.; Handford, P.A.; Stuart, D.I.
Deposited on : 2004-03-12
Resolution : 2.25 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

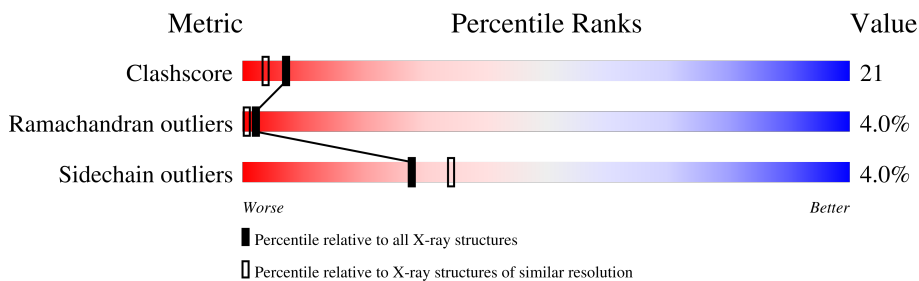
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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))
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	162	65% 27% 8% .
1	B	162	69% 25% 5% .
1	C	162	60% 33% 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3806 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 FIBRILLIN-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	162	Total	C	N	O	S	0	0	0
			1195	725	202	247	21			
1	B	162	Total	C	N	O	S	0	0	0
			1195	725	202	247	21			
1	C	162	Total	C	N	O	S	0	0	0
			1195	725	202	247	21			

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Ca	0	0
			2	2		
2	B	2	Total	Ca	0	0
			2	2		
2	C	2	Total	Ca	0	0
			2	2		

- Molecule 3 is water.

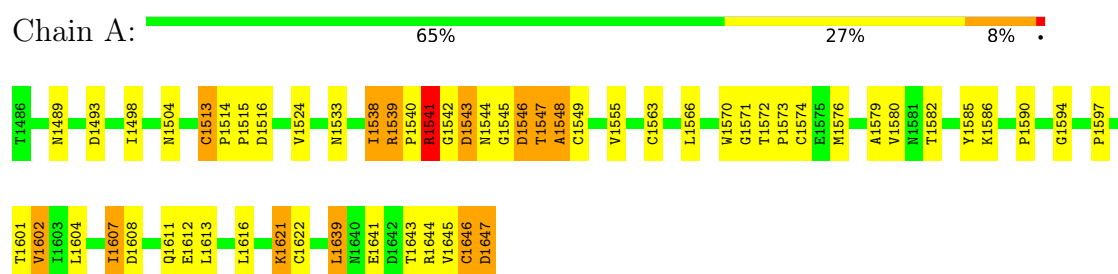
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	81	Total	O	0	0
			81	81		
3	B	79	Total	O	0	0
			79	79		
3	C	55	Total	O	0	0
			55	55		

3 Residue-property plots

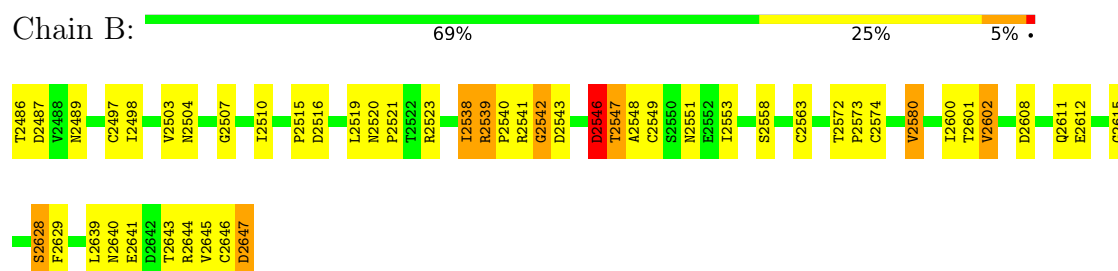
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

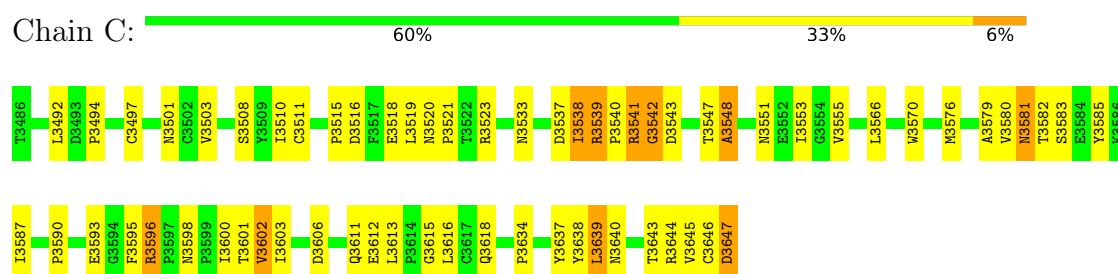
• Molecule 1: FIBRILLIN-1



• Molecule 1: FIBRILLIN-1



• Molecule 1: FIBRILLIN-1



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	43.10 Å 105.60 Å 65.00 Å 90.00° 104.20° 90.00°	Depositor
Resolution (Å)	30.00 – 2.25	Depositor
% Data completeness (in resolution range)	96.2 (30.00-2.25)	Depositor
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.247 , 0.318	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3806	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

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:
CA

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	0.62	0/1218	1.10	6/1659 (0.4%)
1	B	0.57	0/1218	1.15	10/1659 (0.6%)
1	C	0.50	0/1218	1.02	6/1659 (0.4%)
All	All	0.57	0/3654	1.09	22/4977 (0.4%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	2546	ASP	N-CA-C	-7.24	98.03	107.73
1	A	1493	ASP	CA-C-N	6.66	127.20	119.47
1	A	1493	ASP	C-N-CA	6.66	127.20	119.47
1	B	2612	GLU	N-CA-C	6.51	120.44	112.23
1	C	3598	ASN	CA-C-N	6.24	126.71	119.47
1	C	3598	ASN	C-N-CA	6.24	126.71	119.47
1	B	2520	ASN	CA-C-N	6.10	126.28	119.32
1	B	2520	ASN	C-N-CA	6.10	126.28	119.32
1	C	3542	GLY	N-CA-C	6.08	118.64	110.43
1	C	3570	TRP	N-CA-C	5.68	118.03	109.23
1	B	2497	CYS	N-CA-C	5.65	118.09	111.02
1	C	3611	GLN	N-CA-C	5.51	119.18	112.23
1	A	1524	VAL	N-CA-C	-5.50	107.92	113.47
1	C	3600	ILE	N-CA-C	5.40	116.80	111.67
1	B	2558	SER	N-CA-C	-5.39	103.42	110.53
1	B	2547	THR	N-CA-C	5.29	117.43	109.23
1	B	2640	ASN	N-CA-C	-5.21	101.65	109.15
1	A	1563	CYS	N-CA-C	5.13	117.77	111.82
1	B	2563	CYS	N-CA-C	5.10	117.73	111.82
1	A	1513	CYS	N-CA-C	5.09	116.91	109.84
1	A	1607	ILE	N-CA-C	-5.07	100.47	108.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	2580	VAL	N-CA-C	5.04	119.17	112.76

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-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 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	1195	0	1100	47	0
1	B	1195	0	1100	39	0
1	C	1195	0	1100	59	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
3	A	81	0	0	4	0
3	B	79	0	0	4	0
3	C	55	0	0	4	0
All	All	3806	0	3300	143	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 21.

All (143) 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:1538:ILE:HG23	1:A:1539:ARG:H	1.37	0.88
1:C:3497:CYS:HB3	3:C:2018:HOH:O	1.78	0.83
1:A:1548:ALA:HB3	1:A:1574:CYS:SG	2.19	0.83
1:C:3540:PRO:HD2	1:C:3548:ALA:HB1	1.59	0.81
1:C:3511:CYS:HB3	3:C:2018:HOH:O	1.81	0.79
1:A:1572:THR:HA	1:A:1573:PRO:C	2.08	0.78
1:C:3618:GLN:HG3	1:C:3645:VAL:HG22	1.67	0.77
1:A:1611:GLN:NE2	1:B:2641:GLU:HB3	2.00	0.77
1:A:1538:ILE:HD13	1:A:1549:CYS:SG	2.25	0.76
1:C:3541:ARG:HD3	1:C:3541:ARG:H	1.50	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:3539:ARG:HD3	1:C:3540:PRO:HD3	1.68	0.74
1:A:1641:GLU:HG2	3:A:2076:HOH:O	1.87	0.74
1:B:2503:VAL:HB	1:B:2510:ILE:HG12	1.69	0.74
1:B:2538:ILE:HA	3:B:2043:HOH:O	1.87	0.73
1:C:3596:ARG:O	1:C:3596:ARG:HD2	1.89	0.72
1:A:1541:ARG:HD2	1:A:1541:ARG:H	1.53	0.71
1:A:1646:CYS:O	1:A:1647:ASP:HB2	1.91	0.71
1:C:3542:GLY:O	1:C:3547:THR:HA	1.90	0.70
1:B:2498:ILE:HD12	1:B:2602:VAL:HG12	1.73	0.69
1:B:2503:VAL:HB	1:B:2510:ILE:CG1	2.24	0.67
1:B:2519:LEU:HD23	1:B:2523:ARG:HG2	1.78	0.66
1:C:3538:ILE:HA	1:C:3548:ALA:O	1.96	0.65
1:A:1539:ARG:HD2	1:A:1540:PRO:HD2	1.77	0.65
1:B:2538:ILE:HD13	1:B:2549:CYS:SG	2.36	0.65
1:A:1590:PRO:HA	3:A:2047:HOH:O	1.96	0.65
1:B:2601:THR:O	1:B:2602:VAL:HB	1.98	0.64
1:C:3580:VAL:HG12	1:C:3585:TYR:CE2	2.33	0.63
1:C:3615:GLY:HA2	1:C:3618:GLN:NE2	2.15	0.62
1:C:3515:PRO:O	1:C:3516:ASP:HB2	2.00	0.61
1:C:3510:ILE:HD12	1:C:3510:ILE:C	2.25	0.61
1:A:1541:ARG:HD2	1:A:1541:ARG:N	2.14	0.61
1:B:2602:VAL:HG12	1:B:2602:VAL:O	2.00	0.60
1:C:3587:ILE:O	1:C:3590:PRO:HD3	2.01	0.60
1:B:2538:ILE:HD12	3:B:2043:HOH:O	2.01	0.60
1:A:1590:PRO:HG2	1:A:1604:LEU:HD21	1.83	0.59
1:C:3618:GLN:HB2	1:C:3645:VAL:HA	1.84	0.58
1:C:3646:CYS:O	1:C:3647:ASP:HB2	2.03	0.58
1:B:2608:ASP:CG	1:B:2611:GLN:HG3	2.29	0.57
1:C:3538:ILE:HD13	1:C:3576:MET:HG3	1.87	0.56
1:A:1538:ILE:HG23	1:A:1539:ARG:N	2.16	0.56
1:C:3580:VAL:O	1:C:3580:VAL:HG23	2.05	0.56
1:A:1566:LEU:HD12	1:A:1616:LEU:HD13	1.88	0.56
1:C:3643:THR:O	1:C:3644:ARG:HB2	2.06	0.55
1:A:1571:GLY:O	1:A:1574:CYS:HA	2.06	0.55
1:A:1542:GLY:O	1:A:1543:ASP:HB2	2.05	0.55
1:A:1572:THR:CA	1:A:1573:PRO:C	2.78	0.55
1:B:2521:PRO:HG2	1:B:2600:ILE:HG23	1.87	0.55
1:B:2643:THR:O	1:B:2644:ARG:HB2	2.06	0.55
1:C:3593:GLU:HG2	3:C:2045:HOH:O	2.06	0.54
1:B:2541:ARG:HB3	3:B:2003:HOH:O	2.07	0.54
1:C:3640:ASN:OD1	1:C:3643:THR:HG23	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1645:VAL:O	1:A:1647:ASP:N	2.41	0.54
1:A:1515:PRO:O	1:A:1516:ASP:HB2	2.07	0.53
1:A:1533:ASN:OD1	1:A:1555:VAL:HA	2.08	0.53
1:A:1621:LYS:HD3	1:A:1622:CYS:N	2.24	0.53
1:C:3583:SER:O	1:C:3587:ILE:HG12	2.08	0.53
1:A:1586:LYS:HD2	3:A:2047:HOH:O	2.07	0.53
1:B:2515:PRO:O	1:B:2516:ASP:HB2	2.09	0.53
1:A:1611:GLN:NE2	1:B:2641:GLU:CB	2.71	0.52
1:A:1601:THR:C	1:A:1602:VAL:HG23	2.34	0.51
1:A:1498:ILE:HB	1:A:1602:VAL:HG11	1.93	0.51
1:B:2628:SER:OG	1:B:2629:PHE:N	2.43	0.51
1:A:1643:THR:O	1:A:1644:ARG:HB2	2.10	0.51
1:A:1616:LEU:O	1:A:1644:ARG:HD3	2.11	0.51
1:B:2608:ASP:OD1	1:B:2611:GLN:HG3	2.11	0.51
1:B:2489:ASN:HA	1:B:2504:ASN:OD1	2.11	0.50
1:C:3618:GLN:CG	1:C:3645:VAL:HG22	2.39	0.50
1:C:3638:TYR:O	1:C:3647:ASP:N	2.45	0.50
1:A:1540:PRO:O	1:A:1541:ARG:C	2.55	0.50
1:A:1590:PRO:HG3	1:A:1597:PRO:HG3	1.94	0.50
1:C:3566:LEU:HD12	1:C:3616:LEU:HD13	1.93	0.50
1:C:3612:GLU:O	1:C:3613:LEU:HD23	2.12	0.50
1:A:1579:ALA:HB3	1:A:1582:THR:HG23	1.94	0.50
1:B:2538:ILE:O	1:B:2542:GLY:HA2	2.11	0.49
1:B:2580:VAL:HG22	3:B:2056:HOH:O	2.11	0.49
1:C:3537:ASP:OD1	1:C:3551:ASN:OD1	2.30	0.49
1:C:3538:ILE:HA	1:C:3548:ALA:HB3	1.94	0.49
1:C:3639:LEU:HD23	1:C:3640:ASN:N	2.28	0.49
1:B:2572:THR:HA	1:B:2573:PRO:C	2.36	0.49
1:C:3540:PRO:O	1:C:3548:ALA:HB2	2.13	0.48
1:C:3602:VAL:HG12	1:C:3602:VAL:O	2.14	0.48
1:A:1572:THR:HA	1:A:1574:CYS:N	2.29	0.48
1:B:2543:ASP:HB3	1:B:2546:ASP:O	2.13	0.48
1:A:1580:VAL:HG12	1:A:1585:TYR:CE2	2.49	0.48
1:B:2519:LEU:HG	1:B:2523:ARG:HA	1.95	0.48
1:B:2539:ARG:C	1:B:2541:ARG:N	2.71	0.47
1:B:2646:CYS:O	1:B:2647:ASP:HB2	2.12	0.47
1:A:1547:THR:HG23	1:A:1548:ALA:N	2.29	0.47
1:A:1616:LEU:O	1:A:1644:ARG:CD	2.63	0.47
1:B:2539:ARG:C	1:B:2541:ARG:H	2.23	0.47
1:C:3595:PHE:CD2	1:C:3606:ASP:HA	2.49	0.47
1:C:3510:ILE:HD12	1:C:3510:ILE:O	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:3601:THR:OG1	1:C:3603:ILE:HG13	2.15	0.46
1:C:3581:ASN:ND2	3:C:2040:HOH:O	2.47	0.46
1:C:3580:VAL:HG12	1:C:3585:TYR:CZ	2.51	0.46
1:C:3580:VAL:O	1:C:3581:ASN:HB3	2.15	0.46
1:C:3634:PRO:HG2	1:C:3637:TYR:CD2	2.50	0.46
1:A:1594:GLY:C	1:A:1607:ILE:HD12	2.41	0.46
1:C:3515:PRO:O	1:C:3516:ASP:CB	2.64	0.46
1:C:3596:ARG:CD	1:C:3596:ARG:C	2.88	0.46
1:C:3618:GLN:HG3	1:C:3645:VAL:CG2	2.42	0.46
1:A:1576:MET:HG2	3:A:2041:HOH:O	2.16	0.45
1:B:2602:VAL:O	1:B:2602:VAL:CG1	2.64	0.45
1:C:3492:LEU:O	1:C:3494:PRO:HD3	2.17	0.45
1:C:3596:ARG:HD2	1:C:3596:ARG:C	2.41	0.45
1:C:3616:LEU:O	1:C:3644:ARG:HD3	2.15	0.45
1:C:3533:ASN:OD1	1:C:3555:VAL:HA	2.17	0.45
1:C:3640:ASN:HB3	1:C:3643:THR:OG1	2.16	0.45
1:A:1639:LEU:CD2	1:A:1639:LEU:C	2.90	0.45
1:A:1546:ASP:O	1:A:1547:THR:HB	2.16	0.44
1:C:3547:THR:O	1:C:3548:ALA:HB2	2.17	0.44
1:C:3637:TYR:O	1:C:3638:TYR:CD1	2.70	0.44
1:B:2521:PRO:HD2	1:B:2600:ILE:O	2.18	0.44
1:C:3553:ILE:HA	1:C:3615:GLY:HA3	1.99	0.43
1:B:2628:SER:O	1:B:2629:PHE:HB3	2.18	0.43
1:B:2574:CYS:O	1:B:2574:CYS:SG	2.77	0.43
1:B:2547:THR:HG22	1:B:2548:ALA:N	2.33	0.43
1:A:1489:ASN:OD1	1:A:1504:ASN:HB2	2.18	0.43
1:B:2539:ARG:HB3	1:B:2540:PRO:HD3	1.99	0.43
1:A:1539:ARG:HB3	1:A:1541:ARG:HD3	2.00	0.43
1:C:3520:ASN:HB2	1:C:3521:PRO:HD2	2.00	0.43
1:C:3579:ALA:HB3	1:C:3582:THR:HG23	2.00	0.43
1:C:3539:ARG:N	1:C:3540:PRO:CD	2.82	0.43
1:C:3581:ASN:CG	1:C:3582:THR:N	2.77	0.43
1:C:3618:GLN:CB	1:C:3645:VAL:HG22	2.49	0.43
1:A:1570:TRP:CG	1:A:1571:GLY:N	2.86	0.42
1:C:3646:CYS:O	1:C:3647:ASP:CB	2.67	0.42
1:B:2553:ILE:HA	1:B:2615:GLY:HA3	2.01	0.42
1:A:1580:VAL:HA	1:A:1585:TYR:CD2	2.55	0.42
1:B:2542:GLY:HA2	1:B:2547:THR:HG21	2.01	0.42
1:A:1612:GLU:O	1:A:1613:LEU:HD23	2.20	0.41
1:A:1608:ASP:OD1	1:A:1608:ASP:C	2.63	0.41
1:B:2498:ILE:HD12	1:B:2602:VAL:CG1	2.45	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2645:VAL:O	1:B:2647:ASP:N	2.53	0.41
1:C:3640:ASN:HB3	1:C:3643:THR:HG1	1.86	0.41
1:C:3503:VAL:HB	1:C:3510:ILE:HD11	2.03	0.41
1:A:1544:ASN:HB3	1:A:1545:GLY:H	1.49	0.41
1:A:1546:ASP:OD2	1:A:1546:ASP:N	2.52	0.41
1:B:2487:ASP:OD2	1:B:2507:GLY:N	2.40	0.41
1:B:2538:ILE:HA	1:B:2538:ILE:HD12	1.95	0.41
1:C:3492:LEU:HD23	1:C:3492:LEU:HA	1.96	0.41
1:C:3519:LEU:CD2	1:C:3523:ARG:HG2	2.51	0.40
1:A:1513:CYS:HA	1:A:1514:PRO:HD3	1.78	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	160/162 (99%)	137 (86%)	16 (10%)	7 (4%)	2	0
1	B	160/162 (99%)	139 (87%)	15 (9%)	6 (4%)	2	1
1	C	160/162 (99%)	135 (84%)	19 (12%)	6 (4%)	2	1
All	All	480/486 (99%)	411 (86%)	50 (10%)	19 (4%)	2	1

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1602	VAL
1	C	3548	ALA
1	A	1541	ARG
1	A	1543	ASP
1	A	1646	CYS
1	B	2602	VAL
1	C	3538	ILE

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Mol	Chain	Res	Type
1	C	3543	ASP
1	C	3581	ASN
1	B	2542	GLY
1	B	2551	ASN
1	A	1548	ALA
1	A	1538	ILE
1	B	2628	SER
1	C	3508	SER
1	B	2538	ILE
1	B	2539	ARG
1	A	1539	ARG
1	C	3602	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	140/140 (100%)	134 (96%)	6 (4%)	26	31
1	B	140/140 (100%)	136 (97%)	4 (3%)	37	47
1	C	140/140 (100%)	133 (95%)	7 (5%)	22	24
All	All	420/420 (100%)	403 (96%)	17 (4%)	28	34

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1541	ARG
1	A	1546	ASP
1	A	1547	THR
1	A	1621	LYS
1	A	1639	LEU
1	A	1647	ASP
1	B	2486	THR
1	B	2546	ASP
1	B	2639	LEU
1	B	2647	ASP

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Mol	Chain	Res	Type
1	C	3501	ASN
1	C	3518	GLU
1	C	3539	ARG
1	C	3541	ARG
1	C	3596	ARG
1	C	3639	LEU
1	C	3647	ASP

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

Mol	Chain	Res	Type
1	A	1551	ASN
1	A	1618	GLN
1	A	1630	GLN
1	B	2551	ASN
1	B	2611	GLN
1	B	2630	GLN
1	C	3581	ASN
1	C	3618	GLN
1	C	3630	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

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](#)

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section is therefore empty.

6.4 Ligands [i](#)

EDS was not executed - this section is therefore empty.

6.5 Other polymers [i](#)

EDS was not executed - this section is therefore empty.