



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 22, 2021 – 01:06 PM EDT

PDB ID : 7MWB
Title : ERAP1 binds peptide C-terminus of a SPF sequence (FKARKF)
Authors : Guo, H.C.; Sui, L.
Deposited on : 2021-05-16
Resolution : 3.20 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.22
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.22

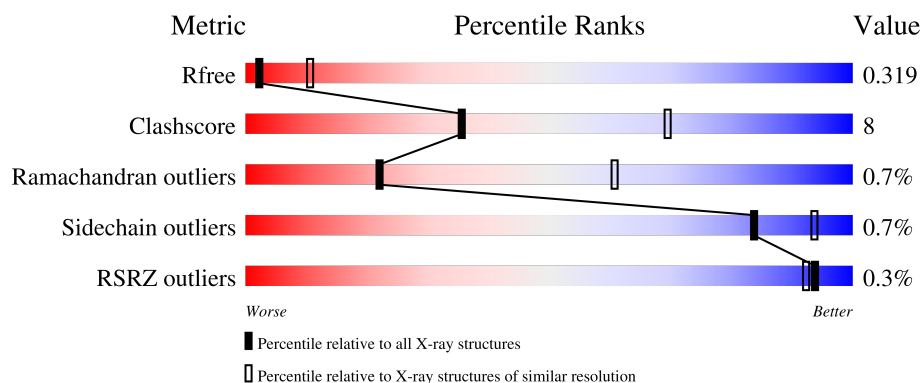
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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}	130704	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)

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 $\leq 5\%$. 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	419	<div> <div></div> <div>79%</div> <div>18%</div> <div>.</div> </div>
1	B	419	<div> <div>81%</div> <div>16%</div> <div>.</div> </div>
1	C	419	<div> <div>78%</div> <div>18%</div> <div>..</div> </div>
1	D	419	<div> <div>80%</div> <div>17%</div> <div>.</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 13373 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 Endoplasmic reticulum aminopeptidase 1,SPF Sequence.

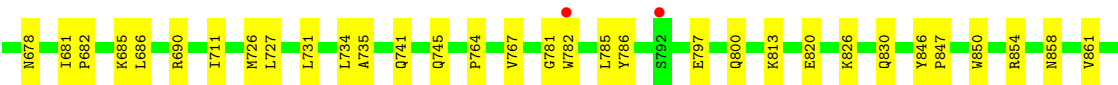
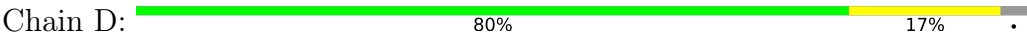
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	407	Total	C	N	O	S	0	0	0
			3320	2135	564	605	16			
1	B	408	Total	C	N	O	S	0	0	0
			3332	2142	565	608	17			
1	C	407	Total	C	N	O	S	0	0	0
			3324	2138	563	606	17			
1	D	406	Total	C	N	O	S	0	0	0
			3313	2131	560	606	16			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	21	Total	O	0	0
			21	21		
2	B	23	Total	O	0	0
			23	23		
2	C	16	Total	O	0	0
			16	16		
2	D	24	Total	O	0	0
			24	24		



● Molecule 1: Endoplasmic reticulum aminopeptidase 1,SPF Sequence



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	58.43Å 69.98Å 121.24Å 90.11° 100.87° 90.15°	Depositor
Resolution (Å)	59.51 – 3.20 59.53 – 3.20	Depositor EDS
% Data completeness (in resolution range)	94.2 (59.51-3.20) 80.0 (59.53-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.97 (at 3.19Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.271 , 0.309 0.271 , 0.319	Depositor DCC
R_{free} test set	1558 reflections (5.30%)	wwPDB-VP
Wilson B-factor (Å ²)	23.5	Xtriage
Anisotropy	0.440	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , -10.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.368 for -h,k,-l	Xtriage
F_o, F_c correlation	0.80	EDS
Total number of atoms	13373	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 21.37 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 7.1077e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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	0.25	0/3391	0.47	0/4572
1	B	0.27	0/3404	0.48	0/4590
1	C	0.27	0/3396	0.49	0/4579
1	D	0.26	0/3384	0.47	0/4563
All	All	0.27	0/13575	0.48	0/18304

There are no bond length outliers.

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	3320	0	3309	64	0
1	B	3332	0	3323	50	0
1	C	3324	0	3317	61	0
1	D	3313	0	3300	55	0
2	A	21	0	0	2	0
2	B	23	0	0	1	0
2	C	16	0	0	1	0
2	D	24	0	0	3	0
All	All	13373	0	13249	214	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 8.

All (214) 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:D:942:PHE:CE1	1:D:944:ALA:HB3	1.80	1.15
1:C:701:ALA:O	1:C:705:ARG:HG3	1.50	1.11
1:D:942:PHE:HE1	1:D:944:ALA:HB3	1.13	1.02
1:D:532:LEU:HG	1:D:547:GLU:OE2	1.62	1.00
1:C:936:GLU:O	1:C:940:ARG:HG3	1.63	0.98
1:A:806:CYS:SG	1:A:838:LEU:HD11	2.12	0.90
1:A:942:PHE:CE1	1:A:945:ARG:HB2	2.15	0.82
1:D:681:ILE:O	1:D:685:LYS:HG2	1.80	0.81
1:D:782:TRP:CZ3	1:D:813:LYS:HD2	2.16	0.79
1:D:550:MET:HA	2:D:1006:HOH:O	1.84	0.77
1:A:546:GLN:NE2	1:A:582:LEU:CD1	2.49	0.75
1:D:942:PHE:CD1	1:D:944:ALA:HB3	2.22	0.75
1:B:879:THR:HG22	1:B:915:ILE:HD11	1.70	0.71
1:C:879:THR:HG22	1:C:915:ILE:HD11	1.72	0.71
1:D:942:PHE:CE1	1:D:944:ALA:CB	2.70	0.69
1:A:571:THR:HG22	1:A:598:ILE:HG13	1.75	0.69
1:B:943:LYS:NZ	1:B:946:LYS:HD2	2.08	0.68
1:A:943:LYS:HB2	1:D:678:ASN:OD1	1.93	0.68
1:C:711:ILE:HD13	1:C:735:ALA:HB2	1.75	0.68
1:D:942:PHE:CD1	1:D:944:ALA:CB	2.77	0.68
1:B:733:LEU:HD21	1:B:803:PHE:HE2	1.59	0.67
1:D:538:ARG:HG3	1:D:541:ASN:HB3	1.77	0.67
1:B:678:ASN:HA	1:C:943:LYS:HG2	1.75	0.66
1:B:943:LYS:HG3	1:B:943:LYS:O	1.97	0.65
1:C:942:PHE:CD1	1:C:944:ALA:HB3	2.32	0.65
1:B:624:LEU:HD12	1:B:660:LEU:HD21	1.79	0.65
1:A:806:CYS:SG	1:A:838:LEU:CD1	2.86	0.64
1:B:726:MET:CE	1:C:942:PHE:HB3	2.28	0.63
1:C:936:GLU:OE2	1:C:940:ARG:HD2	1.99	0.62
1:A:623:LEU:HD21	1:A:630:ALA:HB3	1.80	0.62
1:A:562:LEU:HD21	1:A:585:LYS:HD2	1.82	0.62
1:C:822:PHE:HE1	1:C:860:LEU:HD23	1.65	0.62
1:A:689:LYS:HE2	1:A:881:GLN:HB2	1.81	0.62
1:B:708:ARG:NH1	1:B:740:TYR:OH	2.32	0.61
1:A:850:TRP:CE2	1:A:854:ARG:HD2	2.35	0.61
1:A:942:PHE:HE1	1:A:945:ARG:HB2	1.62	0.61
1:B:818:LEU:HD21	1:B:839:ILE:HG13	1.82	0.60
1:C:939:GLU:HA	1:C:942:PHE:CD2	2.37	0.60
1:C:737:VAL:HG13	1:C:807:ARG:HE	1.66	0.60
1:C:748:GLU:HG3	1:C:752:ARG:HH21	1.66	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:939:GLU:O	1:C:942:PHE:CE2	2.55	0.60
1:A:537:VAL:HG22	1:A:542:VAL:HG13	1.84	0.59
1:C:600:PHE:N	1:C:612:TYR:OH	2.35	0.59
1:C:939:GLU:HA	1:C:942:PHE:HD2	1.67	0.59
1:A:546:GLN:NE2	1:A:565:VAL:HG23	2.18	0.59
1:A:938:LEU:HD23	1:D:726:MET:HG2	1.83	0.59
1:D:547:GLU:O	1:D:547:GLU:HG2	2.02	0.59
1:B:571:THR:HG22	1:B:598:ILE:HG13	1.85	0.59
1:B:734:LEU:HB2	1:C:947:PHE:CE2	2.37	0.59
1:D:573:LYS:NZ	1:D:594:GLU:O	2.30	0.59
1:A:546:GLN:HE22	1:A:565:VAL:HG23	1.67	0.58
1:D:866:LEU:HD22	1:D:905:LEU:HD21	1.86	0.58
1:C:942:PHE:HD1	1:C:944:ALA:HB3	1.67	0.58
1:D:600:PHE:N	1:D:612:TYR:OH	2.35	0.58
1:A:600:PHE:N	1:A:612:TYR:OH	2.31	0.58
1:D:563:TRP:CE2	1:D:565:VAL:HG22	2.39	0.57
1:A:546:GLN:HE21	1:A:582:LEU:HD12	1.69	0.57
1:A:726:MET:CE	1:A:730:GLN:HG2	2.35	0.56
1:B:769:LEU:HG	1:B:800:GLN:HG2	1.86	0.56
1:A:546:GLN:NE2	1:A:582:LEU:HD12	2.21	0.56
1:A:806:CYS:CB	1:A:838:LEU:CD1	2.82	0.56
1:B:733:LEU:HD13	1:C:947:PHE:CD2	2.41	0.56
1:C:748:GLU:HG3	1:C:752:ARG:NH2	2.21	0.55
1:D:764:PRO:HB2	1:D:767:VAL:HG12	1.87	0.55
1:B:860:LEU:HD11	1:B:874:MET:HE3	1.89	0.55
1:A:946:LYS:HD2	1:D:685:LYS:HG3	1.89	0.55
1:B:825:ASP:OD1	1:B:825:ASP:N	2.40	0.55
1:A:563:TRP:CE2	1:A:565:VAL:HG22	2.41	0.55
1:A:705:ARG:CZ	1:A:936:GLU:OE2	2.53	0.55
1:B:533:ILE:HD11	1:B:565:VAL:HG11	1.89	0.55
1:B:755:LYS:HG3	1:B:784:PHE:CD2	2.41	0.55
1:C:825:ASP:N	1:C:825:ASP:OD1	2.40	0.54
1:C:685:LYS:HD3	1:C:881:GLN:HE22	1.73	0.54
1:A:897:SER:HB3	2:A:1012:HOH:O	2.07	0.54
1:B:568:THR:OG1	1:B:601:ASN:OD1	2.25	0.54
1:B:945:ARG:HB3	2:B:1017:HOH:O	2.08	0.53
1:C:939:GLU:O	1:C:942:PHE:HE2	1.91	0.53
1:A:755:LYS:HG2	1:A:784:PHE:CE2	2.43	0.53
1:D:690:ARG:HD3	1:D:884:THR:HG22	1.90	0.53
1:A:776:ALA:HB2	1:A:785:LEU:HD23	1.90	0.53
1:D:943:LYS:HG3	1:D:943:LYS:O	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:947:PHE:HE2	1:D:734:LEU:HB2	1.74	0.52
1:B:781:GLY:O	1:B:785:LEU:HB2	2.09	0.52
1:A:825:ASP:N	1:A:825:ASP:OD1	2.41	0.52
1:D:858:ASN:O	1:D:861:VAL:HG22	2.11	0.51
1:A:944:ALA:O	1:A:946:LYS:HG2	2.10	0.51
1:B:726:MET:HE1	1:C:942:PHE:HB3	1.93	0.51
1:D:531:PRO:HA	1:D:548:HIS:HA	1.93	0.51
1:A:811:LYS:HE2	1:A:848:LEU:HD11	1.93	0.51
1:C:733:LEU:HD21	1:C:803:PHE:CE2	2.45	0.50
1:A:546:GLN:HE21	1:A:582:LEU:CD1	2.21	0.50
1:B:658:LEU:O	1:B:661:SER:OG	2.28	0.50
1:C:571:THR:HG22	1:C:598:ILE:HG13	1.94	0.50
1:A:781:GLY:O	1:A:785:LEU:HB2	2.12	0.50
1:B:682:PRO:HG3	1:C:946:LYS:HZ2	1.76	0.50
1:A:727:LEU:O	1:A:731:LEU:HB2	2.12	0.50
1:C:769:LEU:HG	1:C:800:GLN:HG2	1.93	0.50
1:D:942:PHE:HD1	1:D:944:ALA:CB	2.24	0.50
1:B:677:LEU:HD11	1:B:707:LEU:HD11	1.94	0.49
1:C:942:PHE:HD1	1:C:944:ALA:CB	2.25	0.49
1:D:674:PHE:O	1:D:678:ASN:N	2.45	0.49
1:A:942:PHE:CD1	1:A:942:PHE:C	2.86	0.49
1:C:533:ILE:HD11	1:C:565:VAL:HG11	1.93	0.49
1:C:922:MET:HE3	1:C:926:PHE:HB2	1.93	0.49
1:A:658:LEU:O	1:A:661:SER:OG	2.26	0.49
1:A:821:SER:O	1:A:829:THR:HA	2.13	0.49
1:B:833:PRO:HA	1:B:874:MET:SD	2.53	0.49
1:D:942:PHE:HD1	1:D:944:ALA:HB2	1.77	0.49
1:D:546:GLN:NE2	1:D:562:LEU:HD13	2.28	0.49
1:B:601:ASN:OD1	1:B:601:ASN:O	2.31	0.49
1:D:924:LYS:HB3	2:D:1021:HOH:O	2.12	0.49
1:A:755:LYS:HG2	1:A:784:PHE:CD2	2.48	0.48
1:D:727:LEU:O	1:D:731:LEU:HB2	2.12	0.48
1:A:858:ASN:O	1:A:861:VAL:HG22	2.13	0.48
1:A:702:PHE:HD1	1:A:705:ARG:HH11	1.61	0.48
1:A:945:ARG:O	1:A:945:ARG:HG2	2.13	0.48
1:A:789:TYR:CE1	1:A:798:LYS:HG3	2.49	0.48
1:A:822:PHE:HE2	1:A:859:LYS:HE2	1.78	0.48
1:D:781:GLY:O	1:D:785:LEU:HB2	2.13	0.48
1:B:942:PHE:CD1	1:B:944:ALA:HB3	2.49	0.48
1:A:726:MET:HE2	1:A:730:GLN:HG2	1.95	0.48
1:A:633:SER:HB3	2:A:1001:HOH:O	2.12	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:624:LEU:HD12	1:C:660:LEU:HD11	1.95	0.47
1:A:546:GLN:NE2	1:A:565:VAL:CG2	2.77	0.47
1:A:674:PHE:O	1:A:678:ASN:N	2.46	0.47
1:B:711:ILE:HD13	1:B:735:ALA:HB2	1.96	0.47
1:C:529:GLY:N	1:C:549:TYR:O	2.47	0.47
1:A:822:PHE:CE2	1:A:859:LYS:HE2	2.49	0.47
1:B:705:ARG:NH2	1:B:936:GLU:OE2	2.47	0.47
1:B:581:LEU:HD23	1:B:583:LYS:HE2	1.95	0.47
1:B:733:LEU:CD1	1:C:947:PHE:CD2	2.97	0.47
1:B:853:LEU:HD12	1:B:878:THR:HG21	1.96	0.47
1:D:741:GLN:HG3	1:D:745:GLN:HG3	1.97	0.47
1:A:533:ILE:HD13	1:A:546:GLN:HA	1.97	0.47
1:A:833:PRO:HG3	1:A:874:MET:HE2	1.97	0.47
1:B:644:PHE:HB2	1:B:679:GLU:HG2	1.97	0.47
1:C:568:THR:OG1	1:C:577:VAL:HG13	2.15	0.47
1:C:855:LYS:HD3	2:C:1014:HOH:O	2.14	0.46
1:D:786:TYR:OH	1:D:826:LYS:HB3	2.15	0.46
1:D:850:TRP:CE2	1:D:854:ARG:HD2	2.50	0.46
1:D:682:PRO:O	1:D:686:LEU:HD13	2.16	0.46
1:B:734:LEU:HA	1:C:947:PHE:CD2	2.51	0.46
1:C:834:GLN:O	1:C:834:GLN:HG3	2.15	0.46
1:C:898:LEU:HB2	1:C:903:SER:HB2	1.98	0.46
1:B:755:LYS:NZ	1:B:780:GLU:HB3	2.31	0.46
1:D:595:VAL:HG11	1:D:598:ILE:HG12	1.97	0.46
1:C:785:LEU:HD12	1:C:785:LEU:HA	1.79	0.46
1:B:898:LEU:HB2	1:B:903:SER:HB2	1.98	0.46
1:B:764:PRO:HB2	1:B:767:VAL:HG12	1.97	0.45
1:D:591:LEU:HD13	1:D:595:VAL:HG21	1.97	0.45
1:A:662:LEU:HA	1:A:706:LEU:HD21	1.98	0.45
1:A:537:VAL:HG13	1:A:542:VAL:HG22	1.99	0.45
1:A:737:VAL:HG13	1:A:807:ARG:HE	1.82	0.45
1:A:710:LEU:HD21	1:A:727:LEU:HD21	1.97	0.45
1:A:857:TRP:O	1:A:861:VAL:HG13	2.17	0.45
1:D:563:TRP:CZ2	1:D:565:VAL:HG13	2.52	0.45
1:C:851:GLN:OE1	1:C:854:ARG:NH2	2.50	0.45
1:D:533:ILE:HD11	1:D:565:VAL:HG11	1.98	0.45
1:C:658:LEU:O	1:C:661:SER:OG	2.28	0.44
1:B:942:PHE:HD1	1:B:944:ALA:HB3	1.82	0.44
1:D:797:GLU:HA	1:D:800:GLN:OE1	2.17	0.44
1:C:831:GLU:N	1:C:831:GLU:OE2	2.51	0.44
1:C:943:LYS:HD3	1:C:943:LYS:HA	1.46	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:695:VAL:HG21	1:C:926:PHE:HE1	1.82	0.44
1:B:755:LYS:HA	1:B:784:PHE:HE2	1.83	0.44
1:D:610:VAL:H	1:D:642:ASN:ND2	2.16	0.44
1:D:671:MET:HB2	1:D:672:PRO:HD3	2.00	0.43
1:D:532:LEU:CG	1:D:547:GLU:OE2	2.51	0.43
1:A:898:LEU:HB2	1:A:903:SER:HB2	2.00	0.43
1:B:720:GLY:CA	1:B:724:GLU:OE1	2.67	0.43
1:C:857:TRP:CZ2	1:C:866:LEU:CD2	3.02	0.42
1:A:542:VAL:HB	1:A:589:LEU:HD23	2.01	0.42
1:A:563:TRP:CZ2	1:A:565:VAL:HG13	2.54	0.42
1:A:816:TRP:CZ2	1:A:820:GLU:HG3	2.54	0.42
1:D:609:ILE:HA	1:D:642:ASN:OD1	2.19	0.42
1:A:946:LYS:HB2	1:D:685:LYS:HE2	2.01	0.42
1:B:726:MET:HE3	1:C:942:PHE:HB3	1.99	0.42
1:B:857:TRP:O	1:B:861:VAL:HG23	2.19	0.42
1:D:846:TYR:CG	1:D:847:PRO:HD3	2.54	0.42
1:C:895:PHE:HA	1:C:898:LEU:HD23	2.01	0.42
1:A:623:LEU:HA	1:D:830:GLN:OE1	2.19	0.42
1:A:733:LEU:HD21	1:A:803:PHE:CE2	2.55	0.42
1:A:838:LEU:HA	1:A:841:ARG:HB2	2.01	0.42
1:C:702:PHE:HA	1:C:705:ARG:HD2	2.02	0.42
1:C:741:GLN:HB3	1:C:742:PRO:HD3	2.02	0.42
1:D:741:GLN:HG3	2:D:1011:HOH:O	2.20	0.42
1:D:854:ARG:HA	1:D:894:PHE:CZ	2.54	0.42
1:A:658:LEU:HD11	1:A:929:ILE:HG12	2.01	0.41
1:B:755:LYS:HA	1:B:784:PHE:CE2	2.55	0.41
1:C:571:THR:H	1:C:574:SER:HB3	1.86	0.41
1:C:846:TYR:CG	1:C:847:PRO:HD3	2.55	0.41
1:C:751:PHE:O	1:C:755:LYS:N	2.46	0.41
1:B:563:TRP:CE2	1:B:565:VAL:HG22	2.56	0.41
1:B:755:LYS:HZ1	1:B:780:GLU:HB3	1.84	0.41
1:C:751:PHE:HB2	1:C:771:VAL:HG13	2.02	0.41
1:C:822:PHE:CE1	1:C:860:LEU:HD23	2.50	0.41
1:C:693:ASN:O	1:C:697:THR:HG22	2.21	0.41
1:B:720:GLY:HA3	1:B:724:GLU:OE1	2.20	0.41
1:B:866:LEU:HD21	1:B:905:LEU:HD11	2.02	0.41
1:C:613:GLU:HG2	1:C:614:ASP:H	1.86	0.41
1:C:683:MET:HB3	1:C:683:MET:HE2	1.74	0.41
1:C:883:SER:HB3	1:C:918:ASN:HB3	2.02	0.41
1:D:711:ILE:HD13	1:D:735:ALA:HB2	2.03	0.41
1:D:571:THR:HG22	1:D:598:ILE:CD1	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:673:VAL:O	1:B:677:LEU:HD13	2.21	0.40
1:B:908:VAL:O	1:B:911:THR:HG22	2.21	0.40
1:D:546:GLN:NE2	1:D:561:TYR:O	2.37	0.40
1:C:533:ILE:CD1	1:C:567:LEU:HD21	2.52	0.40
1:D:846:TYR:CD2	1:D:847:PRO:HD3	2.57	0.40
1:C:674:PHE:O	1:C:678:ASN:N	2.55	0.40
1:B:572:SER:HB2	1:B:597:TRP:CD1	2.56	0.40
1:A:946:LYS:HD2	1:D:685:LYS:HE2	2.03	0.40
1:C:840:GLY:HA2	1:C:846:TYR:HA	2.03	0.40
1:D:820:GLU:CG	1:D:826:LYS:HB2	2.51	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%)	372 (93%)	25 (6%)	2 (0%)	29	67
1	B	402/419 (96%)	373 (93%)	26 (6%)	3 (1%)	22	61
1	C	401/419 (96%)	377 (94%)	21 (5%)	3 (1%)	22	61
1	D	398/419 (95%)	370 (93%)	25 (6%)	3 (1%)	19	58
All	All	1600/1676 (96%)	1492 (93%)	97 (6%)	11 (1%)	22	61

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	606	GLY
1	B	602	VAL
1	D	946	LYS
1	A	944	ALA
1	B	942	PHE

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Mol	Chain	Res	Type
1	D	602	VAL
1	B	943	LYS
1	C	942	PHE
1	C	865	GLU
1	D	603	GLY
1	C	592	PRO

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	365/379 (96%)	364 (100%)	1 (0%)	92	96
1	B	367/379 (97%)	364 (99%)	3 (1%)	81	93
1	C	366/379 (97%)	362 (99%)	4 (1%)	73	88
1	D	365/379 (96%)	363 (100%)	2 (0%)	88	95
All	All	1463/1516 (96%)	1453 (99%)	10 (1%)	84	94

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	663	TYR
1	B	663	TYR
1	B	702	PHE
1	B	940	ARG
1	C	579	ARG
1	C	663	TYR
1	C	702	PHE
1	C	943	LYS
1	D	546	GLN
1	D	663	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	546	GLN
1	C	578	HIS
1	D	745	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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, 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	407/419 (97%)	-0.16	3 (0%) 87 81	22, 36, 56, 85	1 (0%)
1	B	408/419 (97%)	-0.26	0 100 100	18, 29, 48, 65	0
1	C	407/419 (97%)	-0.34	0 100 100	14, 27, 46, 70	1 (0%)
1	D	406/419 (96%)	-0.23	2 (0%) 91 86	16, 32, 58, 78	2 (0%)
All	All	1628/1676 (97%)	-0.25	5 (0%) 94 92	14, 31, 54, 85	4 (0%)

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	568	THR	2.5
1	A	794	SER	2.5
1	A	792	SER	2.4
1	D	792	SER	2.1
1	D	782	TRP	2.1

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 monosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.