



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 03:32 am GMT

PDB ID : 1PD1
Title : Crystal structure of the COPII coat subunit, Sec24, complexed with a peptide containing the Dx_E cargo sorting signal of yeast Sys1 protein
Authors : Mossessova, E.; Bickford, L.C.; Goldberg, J.
Deposited on : 2003-05-18
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 ⓘ symbol.

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

MolProbity : 4.02b-467
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

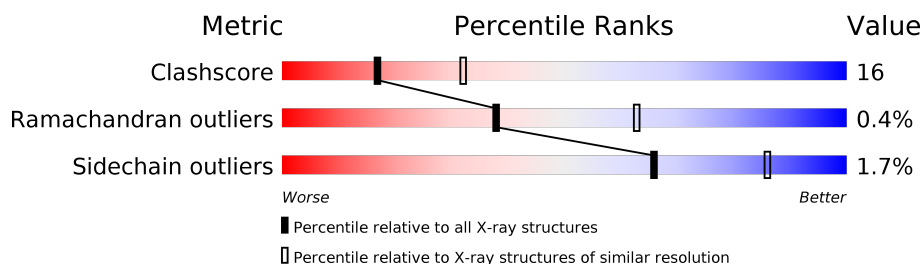
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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))
Clashscore	112137	2895 (2.60-2.60)
Ramachandran outliers	110173	2848 (2.60-2.60)
Sidechain outliers	110143	2848 (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 $\leq 5\%$. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	810	
2	B	9	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6059 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 Protein transport protein Sec24.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	735	Total	C	N	O	S	0	0	0
			5821	3708	993	1083	37			

- Molecule 2 is a protein called DxE cargo sorting signal peptide of yeast Sys1 protein.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	9	Total	C	N	O	0	0	0
			71	44	11	16			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		

- Molecule 4 is water.

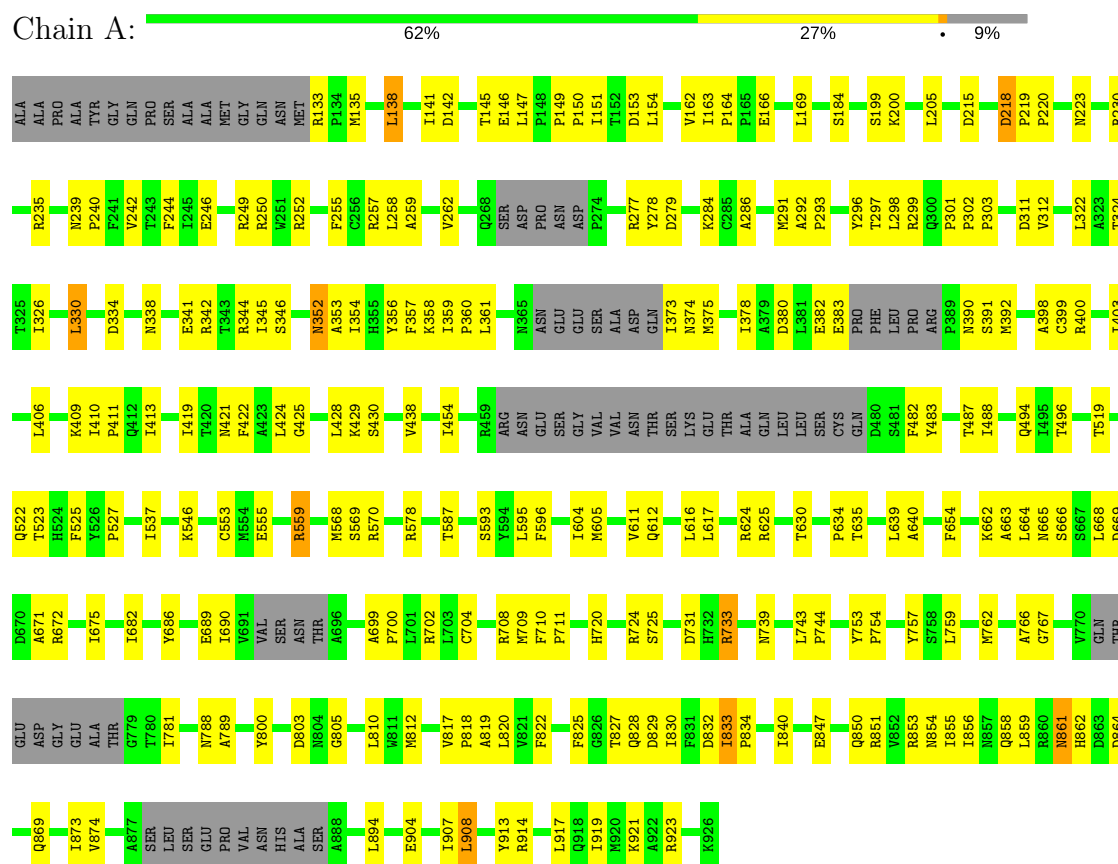
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	165	Total	O	0	0
			165	165		
4	B	1	Total	O	0	0
			1	1		

3 Residue-property plots

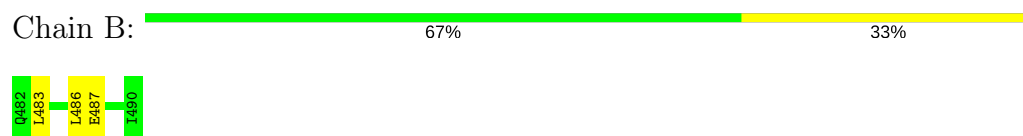
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are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

Note EDS was not executed.

• Molecule 1: Protein transport protein Sec24



• Molecule 2: DxE cargo sorting signal peptide of yeast Sys1 protein



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	93.95 Å 93.95 Å 199.93 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.65 – 2.60	Depositor
% Data completeness (in resolution range)	89.6 (19.65-2.60)	Depositor
R_{merge}	0.03	Depositor
R_{sym}	0.09	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.203 , 0.256	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6059	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	0/5939	0.63	0/8052
2	B	0.36	0/70	0.57	0/91
All	All	0.35	0/6009	0.63	0/8143

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

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	5821	0	5854	185	1
2	B	71	0	70	2	0
3	A	1	0	0	0	0
4	A	165	0	0	4	0
4	B	1	0	0	0	0
All	All	6059	0	5924	185	1

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 16.

All (185) 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:373:ILE:HD12	1:A:409:LYS:HD2	1.30	1.13
1:A:303:PRO:HG2	1:A:344:ARG:HH21	1.27	0.99
1:A:810:LEU:HD22	1:A:840:ILE:HD11	1.59	0.83
1:A:827:THR:HG22	1:A:828:GLN:H	1.45	0.79
1:A:375:MET:HE1	1:A:413:ILE:HG21	1.65	0.76
1:A:297:THR:HG22	1:A:624:ARG:HG2	1.68	0.74
1:A:279:ASP:HA	1:A:284:LYS:HE3	1.70	0.73
1:A:850:GLN:HA	1:A:853:ARG:NH1	2.04	0.72
1:A:215:ASP:O	1:A:219:PRO:HG3	1.90	0.71
1:A:525:PHE:CE2	1:A:527:PRO:HG3	2.26	0.70
1:A:825:PHE:CD2	1:A:834:PRO:HD2	2.27	0.69
1:A:154:LEU:HG	1:A:709:MET:CE	2.24	0.68
1:A:297:THR:CG2	1:A:624:ARG:HG2	2.24	0.67
1:A:759:LEU:O	1:A:762:MET:HB2	1.95	0.67
1:A:142:ASP:OD2	1:A:700:PRO:HB3	1.94	0.66
1:A:424:LEU:HD11	1:A:428:LEU:HD13	1.77	0.66
1:A:380:ASP:OD2	1:A:383:GLU:HB2	1.95	0.66
1:A:151:ILE:HA	1:A:709:MET:CE	2.26	0.66
1:A:429:LYS:HE3	1:A:482:PHE:HE1	1.61	0.66
1:A:154:LEU:HG	1:A:709:MET:HE2	1.79	0.65
1:A:151:ILE:HA	1:A:709:MET:HE1	1.77	0.65
1:A:874:VAL:HG22	1:A:894:LEU:HB3	1.79	0.64
1:A:324:THR:OG1	1:A:537:ILE:HD12	1.98	0.64
1:A:720:HIS:O	1:A:724:ARG:HG3	1.97	0.64
1:A:345:ILE:HG12	1:A:346:SER:N	2.14	0.63
1:A:352:ASN:HD22	1:A:353:ALA:N	1.97	0.63
1:A:488:ILE:HD13	1:A:578:ARG:HH21	1.65	0.61
1:A:825:PHE:CE1	1:A:833:ILE:HD11	2.35	0.61
1:A:390:ASN:O	1:A:391:SER:HB2	2.00	0.61
1:A:494:GLN:HG3	1:A:555:GLU:HG2	1.82	0.61
1:A:425:GLY:HA3	1:A:454:ILE:O	2.01	0.60
1:A:690:ILE:C	1:A:702:ARG:HH22	2.05	0.60
1:A:356:TYR:OH	1:A:430:SER:HB3	2.02	0.59
1:A:639:LEU:HD23	1:A:689:GLU:HB3	1.85	0.58
1:A:303:PRO:HG3	1:A:342:ARG:NH1	2.19	0.58
1:A:357:PHE:CD2	1:A:406:LEU:HD11	2.38	0.58
1:A:302:PRO:HB3	1:A:438:VAL:HG12	1.86	0.57
1:A:724:ARG:O	1:A:733:ARG:NH2	2.36	0.57
1:A:419:ILE:HG22	1:A:421:ASN:H	1.70	0.57
1:A:255:PHE:CZ	1:A:612:GLN:HB2	2.40	0.56
1:A:827:THR:HG22	1:A:828:GLN:N	2.18	0.56
1:A:672:ARG:NH2	1:A:725:SER:HA	2.21	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:138:LEU:CD2	1:A:704:CYS:HB3	2.36	0.55
1:A:829:ASP:HB3	1:A:832:ASP:HB2	1.86	0.55
1:A:322:LEU:O	1:A:326:ILE:HG12	2.06	0.55
1:A:789:ALA:O	1:A:904:GLU:HB2	2.06	0.55
1:A:682:ILE:HD12	1:A:710:PHE:CE2	2.42	0.55
1:A:850:GLN:HA	1:A:853:ARG:HH12	1.69	0.55
1:A:250:ARG:HA	1:A:262:VAL:HG23	1.89	0.54
1:A:345:ILE:HG12	1:A:346:SER:H	1.72	0.54
1:A:662:LYS:O	1:A:666:SER:HB3	2.07	0.54
1:A:919:ILE:O	1:A:923:ARG:HG3	2.07	0.54
1:A:559:ARG:HD2	2:B:487:GLU:OE1	2.08	0.54
1:A:205:LEU:CD1	1:A:611:VAL:HG11	2.37	0.54
1:A:356:TYR:CD2	1:A:392:MET:HE3	2.43	0.54
1:A:424:LEU:CD1	1:A:428:LEU:HD13	2.36	0.54
1:A:373:ILE:HD11	1:A:406:LEU:HA	1.90	0.54
1:A:822:PHE:HA	1:A:827:THR:O	2.08	0.53
1:A:617:LEU:HD12	1:A:617:LEU:C	2.29	0.53
1:A:141:ILE:HD12	1:A:146:GLU:OE1	2.08	0.53
1:A:252:ARG:NH1	1:A:257:ARG:O	2.41	0.53
1:A:767:GLY:HA3	1:A:851:ARG:HB3	1.90	0.53
1:A:356:TYR:CE1	1:A:378:ILE:HD12	2.44	0.53
1:A:488:ILE:HD13	1:A:578:ARG:NH2	2.24	0.53
1:A:399:CYS:O	1:A:403:ILE:HG13	2.09	0.53
1:A:373:ILE:CD1	1:A:409:LYS:HD2	2.20	0.52
1:A:568:MET:CE	1:A:596:PHE:HB3	2.39	0.52
1:A:149:PRO:HG3	1:A:913:TYR:CE2	2.44	0.52
1:A:352:ASN:HD22	1:A:352:ASN:N	2.06	0.52
1:A:147:LEU:HD23	1:A:914:ARG:HG2	1.92	0.52
1:A:373:ILE:HG13	1:A:409:LYS:NZ	2.25	0.52
1:A:595:LEU:HD11	1:A:654:PHE:HA	1.91	0.52
1:A:781:ILE:HG12	1:A:854:ASN:HB3	1.92	0.52
1:A:668:LEU:HD21	1:A:733:ARG:CZ	2.39	0.51
1:A:825:PHE:CD1	1:A:833:ILE:HD11	2.45	0.51
1:A:762:MET:HG2	1:A:766:ALA:HB3	1.93	0.51
1:A:303:PRO:HG3	1:A:342:ARG:CZ	2.40	0.51
1:A:361:LEU:HD22	1:A:398:ALA:HB1	1.93	0.51
1:A:553:CYS:O	1:A:587:THR:HA	2.10	0.51
1:A:855:ILE:O	1:A:859:LEU:HG	2.10	0.51
1:A:334:ASP:CG	1:A:400:ARG:HH12	2.15	0.51
1:A:359:ILE:HD11	1:A:406:LEU:HD23	1.91	0.51
1:A:840:ILE:HD12	1:A:856:ILE:CD1	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:296:TYR:HB3	1:A:616:LEU:HD11	1.93	0.50
1:A:138:LEU:HD12	1:A:640:ALA:HB1	1.93	0.50
1:A:199:SER:O	1:A:200:LYS:HB2	2.11	0.50
1:A:149:PRO:HD2	1:A:788:ASN:ND2	2.26	0.50
1:A:710:PHE:HB3	1:A:711:PRO:HD3	1.94	0.50
1:A:326:ILE:HG22	1:A:330:LEU:HD22	1.93	0.49
1:A:617:LEU:HD11	1:A:625:ARG:HB2	1.93	0.49
1:A:356:TYR:HB3	1:A:392:MET:HE1	1.94	0.49
1:A:817:VAL:HG12	1:A:819:ALA:H	1.77	0.49
1:A:205:LEU:O	1:A:595:LEU:HA	2.13	0.49
1:A:604:ILE:HG21	1:A:635:THR:HG21	1.95	0.48
1:A:593:SER:HB2	1:A:744:PRO:HA	1.94	0.48
1:A:671:ALA:O	1:A:675:ILE:HG13	2.13	0.48
1:A:840:ILE:HD12	1:A:856:ILE:HD12	1.94	0.48
1:A:138:LEU:HD23	1:A:704:CYS:HB3	1.94	0.48
1:A:827:THR:HG22	1:A:829:ASP:H	1.76	0.48
1:A:246:GLU:O	1:A:249:ARG:HG3	2.14	0.48
1:A:142:ASP:CG	1:A:700:PRO:HB3	2.34	0.48
1:A:163:ILE:HD13	1:A:634:PRO:HG3	1.95	0.48
1:A:162:VAL:O	1:A:634:PRO:HG3	2.13	0.48
1:A:546:LYS:HE2	4:A:1030:HOH:O	2.13	0.47
1:A:149:PRO:HG3	1:A:913:TYR:HE2	1.79	0.47
1:A:739:ASN:OD1	1:A:743:LEU:HD13	2.14	0.47
1:A:258:LEU:HD12	1:A:259:ALA:H	1.79	0.47
1:A:496:THR:HA	1:A:519:THR:HB	1.95	0.47
1:A:668:LEU:HB2	4:A:1084:HOH:O	2.13	0.47
1:A:825:PHE:HE2	1:A:873:ILE:HD13	1.77	0.47
1:A:525:PHE:CZ	1:A:731:ASP:HB3	2.50	0.47
1:A:338:ASN:OD1	1:A:341:GLU:HA	2.14	0.47
1:A:352:ASN:H	1:A:352:ASN:HD22	1.63	0.47
1:A:299:ARG:NH2	1:A:438:VAL:O	2.47	0.47
1:A:151:ILE:HG21	1:A:753:TYR:CZ	2.50	0.47
1:A:354:ILE:HD13	1:A:430:SER:OG	2.15	0.46
1:A:483:TYR:O	1:A:487:THR:HG23	2.14	0.46
1:A:223:ASN:O	1:A:291:MET:HG2	2.15	0.46
1:A:663:ALA:C	1:A:665:ASN:H	2.19	0.46
1:A:312:VAL:HG12	1:A:322:LEU:HD13	1.97	0.46
1:A:861:ASN:HD22	1:A:861:ASN:H	1.64	0.46
1:A:352:ASN:H	1:A:352:ASN:ND2	2.13	0.46
1:A:373:ILE:HD13	1:A:406:LEU:HD13	1.97	0.46
1:A:686:TYR:CE1	1:A:690:ILE:HD12	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:LEU:HD11	1:A:611:VAL:HG11	1.97	0.46
1:A:812:MET:HE1	1:A:820:LEU:HG	1.98	0.46
1:A:218:ASP:N	1:A:219:PRO:HD3	2.31	0.45
1:A:352:ASN:HB3	1:A:421:ASN:HB2	1.97	0.45
1:A:166:GLU:CD	1:A:166:GLU:H	2.20	0.45
1:A:759:LEU:HD22	1:A:762:MET:CE	2.46	0.45
1:A:352:ASN:HD22	1:A:353:ALA:H	1.62	0.45
1:A:297:THR:HG22	1:A:624:ARG:CG	2.42	0.45
1:A:373:ILE:HG13	1:A:409:LYS:HZ1	1.80	0.44
1:A:805:GLY:HA3	4:A:995:HOH:O	2.16	0.44
1:A:358:LYS:HB2	1:A:374:ASN:HB2	1.99	0.44
1:A:759:LEU:HD22	1:A:762:MET:HE2	2.00	0.44
1:A:359:ILE:HA	1:A:360:PRO:HD3	1.86	0.44
1:A:153:ASP:OD2	1:A:708:ARG:HD2	2.18	0.44
1:A:164:PRO:HB2	1:A:166:GLU:OE1	2.17	0.44
1:A:669:ASP:HA	1:A:672:ARG:NH1	2.33	0.44
1:A:230:ARG:HD2	1:A:235:ARG:O	2.18	0.44
1:A:862:HIS:HB3	1:A:864:ASP:OD1	2.17	0.43
1:A:708:ARG:O	1:A:708:ARG:HG2	2.18	0.43
1:A:150:PRO:HG2	1:A:708:ARG:HD3	1.99	0.43
1:A:569:SER:O	1:A:570:ARG:HD2	2.18	0.43
1:A:250:ARG:CA	1:A:262:VAL:HG23	2.48	0.43
1:A:356:TYR:HB3	1:A:392:MET:CE	2.49	0.43
1:A:759:LEU:HA	1:A:762:MET:HE2	1.99	0.43
1:A:352:ASN:ND2	1:A:352:ASN:N	2.67	0.43
1:A:711:PRO:HD2	4:A:961:HOH:O	2.18	0.43
1:A:151:ILE:HG21	1:A:753:TYR:OH	2.18	0.43
1:A:757:TYR:O	1:A:800:TYR:HA	2.19	0.43
1:A:301:PRO:HA	1:A:302:PRO:HD3	1.75	0.42
1:A:699:ALA:HB1	1:A:921:LYS:NZ	2.33	0.42
1:A:617:LEU:HD12	1:A:617:LEU:O	2.18	0.42
1:A:133:ARG:HG3	1:A:135:MET:HG2	2.01	0.42
1:A:522:GLN:HG3	1:A:523:THR:N	2.35	0.42
1:A:356:TYR:HE1	1:A:378:ILE:HD12	1.85	0.42
1:A:286:ALA:HB1	1:A:630:THR:O	2.19	0.42
1:A:357:PHE:HA	1:A:374:ASN:O	2.19	0.42
1:A:781:ILE:HD12	1:A:781:ILE:H	1.85	0.41
1:A:279:ASP:HA	1:A:284:LYS:CE	2.45	0.41
1:A:686:TYR:CD1	1:A:690:ILE:HD12	2.55	0.41
1:A:292:ALA:HA	1:A:293:PRO:HD3	1.85	0.41
1:A:311:ASP:HB2	1:A:422:PHE:CE1	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:298:LEU:HG	2:B:486:LEU:HD21	2.02	0.41
1:A:803:ASP:OD1	1:A:869:GLN:OE1	2.39	0.41
1:A:257:ARG:HH22	1:A:605:MET:HB2	1.86	0.41
1:A:184:SER:HA	1:A:205:LEU:HD23	2.03	0.41
1:A:244:PHE:CE1	1:A:277:ARG:HD3	2.56	0.41
1:A:568:MET:HE2	1:A:596:PHE:HB3	2.03	0.41
1:A:913:TYR:CZ	1:A:917:LEU:HD11	2.55	0.41
1:A:219:PRO:HA	1:A:220:PRO:HD3	1.94	0.41
1:A:169:LEU:HD11	1:A:242:VAL:HG22	2.03	0.41
1:A:410:ILE:HB	1:A:411:PRO:HD3	2.02	0.41
1:A:830:ILE:O	1:A:833:ILE:HB	2.21	0.41
1:A:239:ASN:HB2	1:A:240:PRO:CD	2.51	0.41
1:A:303:PRO:HB3	1:A:342:ARG:HB3	2.03	0.41
1:A:278:TYR:O	1:A:284:LYS:HE3	2.22	0.40
1:A:907:ILE:O	1:A:908:LEU:HB2	2.20	0.40
1:A:151:ILE:HA	1:A:709:MET:HE2	2.02	0.40
1:A:352:ASN:ND2	1:A:353:ALA:N	2.66	0.40
1:A:639:LEU:CD2	1:A:689:GLU:HB3	2.49	0.40
1:A:818:PRO:HB3	1:A:828:GLN:HB3	2.03	0.40
1:A:858:GLN:NE2	1:A:861:ASN:HD21	2.19	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:382:GLU:OE2	1:A:382:GLU:OE2[6_555]	1.86	0.34

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	719/810 (89%)	681 (95%)	36 (5%)	2 (0%)	44 70

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	7/9 (78%)	5 (71%)	1 (14%)	1 (14%)	0	0
All	All	726/819 (89%)	686 (94%)	37 (5%)	3 (0%)	38	63

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	664	LEU
1	A	908	LEU
2	B	483	LEU

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	658/720 (91%)	647 (98%)	11 (2%)	66	86
2	B	8/9 (89%)	8 (100%)	0	100	100
All	All	666/729 (91%)	655 (98%)	11 (2%)	66	86

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	138	LEU
1	A	145	THR
1	A	218	ASP
1	A	330	LEU
1	A	352	ASN
1	A	559	ARG
1	A	733	ARG
1	A	754	PRO
1	A	833	ILE
1	A	847	GLU
1	A	861	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	314	GLN
1	A	352	ASN
1	A	412	GLN
1	A	732	HIS
1	A	861	ASN
1	A	869	GLN
1	A	918	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is 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 [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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.