



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:41 AM GMT

PDB ID : 2I7N
Title : Crystal structure of human PANK1 alpha: the catalytic core domain in complex with AcCoA
Authors : Hong, B.S.; Wang, L.; Tempel, W.; Loppnau, P.; Allali-Hassani, A.; Arrow-smith, C.H.; Edwards, A.M.; Sundstrom, M.; Weigelt, J.; Bochkarev, A.; Park, H.W.
Deposited on : 2006-08-31
Resolution : 1.90 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
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

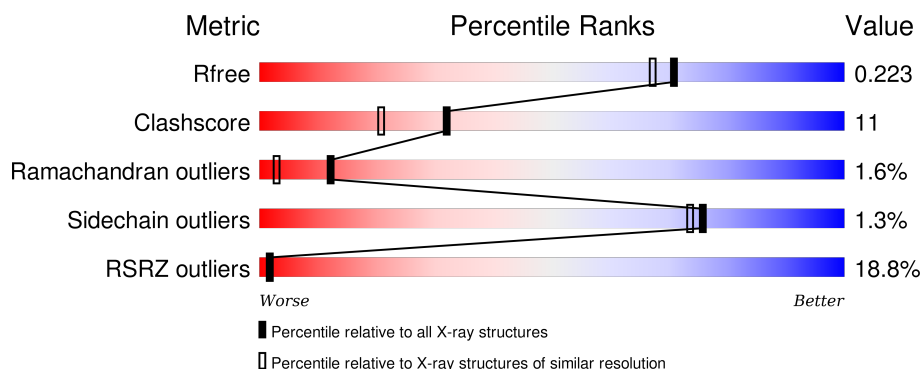
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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.

Mol	Chain	Length	Quality of chain
1	A	360	<div> <div>21%</div> <div>68%</div> <div>21%</div> <div>10%</div> </div>
1	B	360	<div> <div>14%</div> <div>71%</div> <div>20%</div> <div>8%</div> </div>

2 Entry composition [i](#)

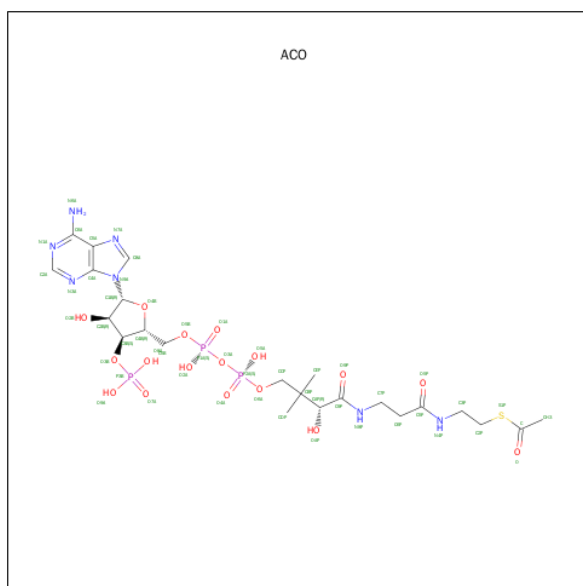
There are 3 unique types of molecules in this entry. The entry contains 5453 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Pantothenate kinase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	325	Total	C	N	O	S	0	0	0
			2538	1628	419	468	23			
1	B	333	Total	C	N	O	S	0	0	0
			2605	1671	431	479	24			

- Molecule 2 is ACETYL COENZYME *A (three-letter code: ACO) (formula: $C_{23}H_{38}N_7O_{17}P_3S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P S	0	0
			29	13	2	11	2 1		
2	B	1	Total	C	N	O	P S	0	0
			29	13	2	11	2 1		

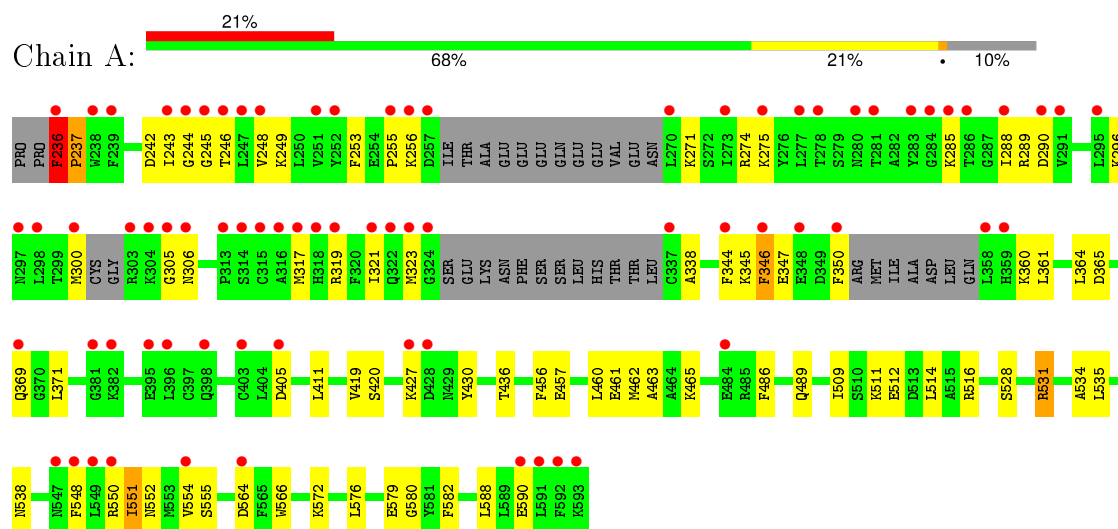
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	92	Total 92	O 92	0	0
3	B	160	Total 160	O 160	0	0

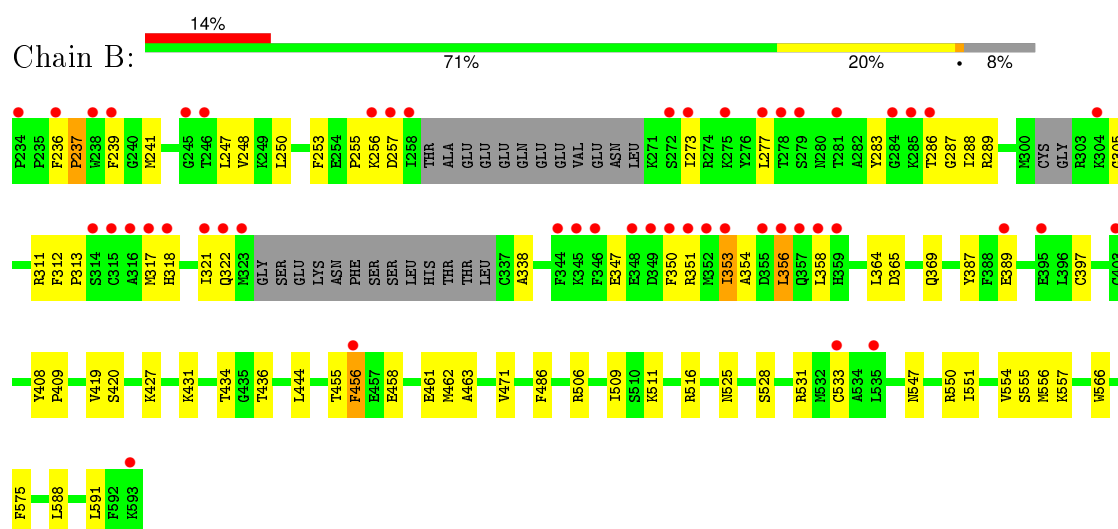
3 Residue-property plots

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 ($RSRZ > 2$). 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: Pantothenate kinase 1



• Molecule 1: Pantothenate kinase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, α , β , γ	92.99 Å 92.99 Å 197.78 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.46 – 1.90 39.46 – 1.90	Depositor EDS
% Data completeness (in resolution range)	96.7 (39.46-1.90) 96.8 (39.46-1.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.93 (at 1.89 Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.222 , 0.243 0.223 , 0.223	Depositor DCC
R_{free} test set	3699 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	24.5	Xtriage
Anisotropy	0.011	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 36.6	EDS
Estimated twinning fraction	0.108 for h,-h-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 73382 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5453	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.28% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

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: ACO

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	0.36	0/2586	0.58	2/3471 (0.1%)
1	B	0.39	0/2656	0.71	3/3568 (0.1%)
All	All	0.37	0/5242	0.65	5/7039 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	236	PHE	C-N-CD	-20.45	75.60	120.60
1	B	236	PHE	C-N-CA	13.64	179.29	122.00
1	A	236	PHE	C-N-CD	-11.72	94.82	120.60
1	A	551	ILE	N-CA-C	5.43	125.67	111.00
1	B	237	PRO	CA-N-CD	-5.33	104.04	111.50

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	2538	0	2528	60	0
1	B	2605	0	2602	52	0
2	A	29	0	23	0	0
2	B	29	0	23	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	92	0	0	3	0
3	B	160	0	0	6	0
All	All	5453	0	5176	110	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 11.

All (110) 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:550:ARG:HG3	1:A:551:ILE:HG13	1.54	0.89
1:A:457:GLU:O	1:A:461:GLU:HG2	1.74	0.86
1:A:552:ASN:OD1	1:A:554:VAL:HG12	1.83	0.78
1:B:239:PHE:HD2	1:B:250:LEU:HD11	1.55	0.69
1:A:344:PHE:HA	1:A:347:GLU:HB2	1.75	0.69
1:B:256:LYS:HG3	1:B:305:GLY:HA2	1.74	0.69
1:A:296:LYS:HG2	1:A:306:ASN:ND2	2.08	0.68
1:B:351:ARG:HB2	1:B:356:LEU:HB2	1.76	0.68
1:A:285:LYS:HD3	1:A:319:ARG:HH12	1.61	0.66
1:B:551:ILE:N	1:B:551:ILE:HD12	2.10	0.65
1:B:462:MET:HE1	1:B:511:LYS:HB3	1.80	0.63
1:B:365:ASP:O	1:B:369:GLN:HG2	1.99	0.62
1:B:431:LYS:HG2	3:B:994:HOH:O	2.00	0.60
1:B:353:ILE:HD12	1:B:353:ILE:N	2.17	0.60
1:B:286:THR:HG22	1:B:313:PRO:HD2	1.83	0.60
1:A:457:GLU:HG3	1:A:551:ILE:HD13	1.84	0.60
1:A:456:PHE:CE2	1:A:460:LEU:HD11	2.37	0.60
1:B:273:ILE:HD12	1:B:273:ILE:N	2.16	0.60
1:B:347:GLU:HG3	1:B:356:LEU:HD11	1.85	0.59
1:A:285:LYS:HD3	1:A:319:ARG:NH1	2.17	0.58
1:B:321:ILE:HD11	1:B:350:PHE:CE1	2.38	0.58
1:A:317:MET:O	1:A:321:ILE:HG12	2.04	0.57
1:B:247:LEU:HD23	1:B:248:VAL:N	2.18	0.57
1:B:273:ILE:HD12	1:B:273:ILE:H	1.68	0.57
1:B:463:ALA:O	1:B:516:ARG:HD3	2.05	0.56
1:B:356:LEU:HG	1:B:358:LEU:H	1.70	0.55
1:A:344:PHE:HD1	1:A:347:GLU:HG3	1.73	0.54
1:A:554:VAL:HG13	1:A:555:SER:N	2.22	0.54
1:A:531:ARG:NH1	1:A:535:LEU:HG	2.22	0.54
1:A:243:ILE:HD12	1:A:317:MET:HE1	1.89	0.54
1:A:365:ASP:OD1	1:A:430:TYR:OH	2.26	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:419:VAL:HB	1:A:436:THR:CG2	2.39	0.53
1:B:338:ALA:HB3	1:B:358:LEU:HD11	1.91	0.53
1:A:564:ASP:OD1	1:A:572:LYS:HD3	2.08	0.53
1:A:509:ILE:HD12	1:A:514:LEU:HD21	1.90	0.53
1:A:538:ASN:ND2	3:A:870:HOH:O	2.41	0.52
1:B:287:GLY:HA2	1:B:312:PHE:HB3	1.90	0.52
1:A:419:VAL:HB	1:A:436:THR:HG22	1.91	0.52
1:B:317:MET:O	1:B:321:ILE:HG12	2.10	0.52
1:B:528:SER:HB2	1:B:566:TRP:CE2	2.44	0.51
1:A:528:SER:HB2	1:A:566:TRP:CE2	2.46	0.50
1:A:364:LEU:HD22	1:B:486:PHE:HZ	1.77	0.50
1:B:462:MET:CE	1:B:511:LYS:HB3	2.42	0.50
1:A:456:PHE:HE2	1:A:460:LEU:HD11	1.76	0.49
1:A:288:ILE:HG22	1:A:289:ARG:N	2.28	0.48
1:B:347:GLU:O	1:B:356:LEU:HD13	2.13	0.48
1:A:243:ILE:HD12	1:A:317:MET:CE	2.44	0.48
1:A:465:LYS:HE2	1:A:512:GLU:OE2	2.14	0.48
1:A:456:PHE:HZ	1:A:548:PHE:HD1	1.62	0.48
1:B:387:TYR:CE1	1:B:389:GLU:HG2	2.49	0.48
1:B:427:LYS:HE2	3:B:965:HOH:O	2.12	0.47
1:B:247:LEU:HD21	1:B:311:ARG:HB2	1.95	0.47
1:B:444:LEU:HD22	1:B:456:PHE:CD1	2.49	0.47
1:A:296:LYS:HG2	1:A:306:ASN:HD21	1.78	0.47
1:A:344:PHE:CD1	1:A:347:GLU:HG3	2.50	0.47
1:A:253:PHE:O	1:A:255:PRO:HD3	2.15	0.47
1:A:256:LYS:HG2	1:A:305:GLY:HA2	1.97	0.47
1:B:588:LEU:HD12	1:B:591:LEU:HD12	1.95	0.46
1:A:319:ARG:O	1:A:323:MET:HG3	2.16	0.46
1:A:345:LYS:C	1:A:347:GLU:H	2.19	0.46
1:A:344:PHE:CE2	1:A:360:LYS:HD3	2.51	0.46
1:B:318:HIS:O	1:B:322:GLN:HG3	2.16	0.46
1:A:249:LYS:HD3	1:A:249:LYS:N	2.31	0.46
1:A:579:GLU:HA	1:A:582:PHE:CE2	2.51	0.46
1:B:397:CYS:SG	1:B:557:LYS:HE2	2.56	0.45
1:A:271:LYS:O	1:A:275:LYS:HG3	2.16	0.45
1:B:531:ARG:NH1	3:B:956:HOH:O	2.49	0.45
1:A:236:PHE:N	1:A:236:PHE:CD1	2.85	0.45
1:A:236:PHE:HA	1:A:237:PRO:HD2	1.52	0.45
1:B:554:VAL:HG13	1:B:555:SER:N	2.31	0.45
1:A:486:PHE:HZ	1:B:364:LEU:HD22	1.81	0.44
1:A:338:ALA:O	1:A:360:LYS:HA	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:351:ARG:CB	1:B:356:LEU:HB2	2.46	0.44
1:A:248:VAL:HG11	1:A:317:MET:HE1	1.99	0.44
1:A:371:LEU:HG	1:A:576:LEU:CD1	2.46	0.44
1:B:506:ARG:O	1:B:509:ILE:HG12	2.17	0.44
1:B:277:LEU:HA	1:B:283:TYR:HE2	1.82	0.44
1:A:242:ASP:HB3	1:A:249:LYS:HE2	2.00	0.44
1:B:283:TYR:HE1	1:B:289:ARG:HB2	1.82	0.44
1:B:419:VAL:HB	1:B:436:THR:HG22	1.99	0.44
1:A:462:MET:CE	1:A:511:LYS:HB3	2.48	0.44
1:B:591:LEU:HD21	3:B:985:HOH:O	2.16	0.43
1:B:556:MET:HG2	1:B:575:PHE:CE2	2.53	0.43
1:B:547:ASN:HA	1:B:550:ARG:CD	2.49	0.43
1:A:249:LYS:HG2	1:A:580:GLY:O	2.18	0.43
1:A:411:LEU:HD11	1:A:534:ALA:HA	2.01	0.43
1:B:551:ILE:CD1	1:B:551:ILE:N	2.80	0.43
1:B:461:GLU:HG3	3:B:971:HOH:O	2.17	0.43
1:B:241:MET:HG2	1:B:250:LEU:HD13	1.99	0.43
1:B:547:ASN:HA	1:B:550:ARG:HD2	2.00	0.43
1:B:353:ILE:H	1:B:353:ILE:HD12	1.83	0.43
1:A:300:MET:HE1	1:A:588:LEU:HD21	2.01	0.43
1:B:288:ILE:HD12	1:B:311:ARG:HH12	1.83	0.43
1:A:489:GLN:HG2	3:A:856:HOH:O	2.18	0.42
1:A:253:PHE:HB2	1:A:300:MET:HE1	2.01	0.42
1:A:427:LYS:HG2	3:A:862:HOH:O	2.18	0.42
1:B:525:ASN:HB3	3:B:916:HOH:O	2.19	0.42
1:A:345:LYS:HD3	1:A:346:PHE:CE1	2.54	0.42
1:B:434:THR:HG21	1:B:533:CYS:SG	2.60	0.42
1:A:248:VAL:C	1:A:249:LYS:HD3	2.40	0.42
1:A:361:LEU:HD21	1:A:590:GLU:HG2	2.01	0.42
1:A:365:ASP:O	1:A:369:GLN:HG3	2.21	0.41
1:A:461:GLU:O	1:A:465:LYS:HD3	2.21	0.41
1:A:244:GLY:O	1:A:246:THR:N	2.54	0.41
1:B:408:TYR:HA	1:B:409:PRO:HA	1.85	0.41
1:A:463:ALA:O	1:A:516:ARG:HD3	2.20	0.41
1:B:455:THR:OG1	1:B:458:GLU:HG3	2.21	0.41
1:B:253:PHE:O	1:B:255:PRO:HD3	2.22	0.40
1:A:248:VAL:HG21	1:A:317:MET:HE2	2.02	0.40
1:A:274:ARG:HG2	1:A:274:ARG:HH11	1.87	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	315/360 (88%)	298 (95%)	12 (4%)	5 (2%)	12	3
1	B	325/360 (90%)	310 (95%)	10 (3%)	5 (2%)	13	3
All	All	640/720 (89%)	608 (95%)	22 (3%)	10 (2%)	12	3

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	237	PRO
1	B	237	PRO
1	B	356	LEU
1	A	245	GLY
1	A	290	ASP
1	A	405	ASP
1	A	346	PHE
1	B	257	ASP
1	B	354	ALA
1	B	353	ILE

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	273/305 (90%)	269 (98%)	4 (2%)	72	69
1	B	281/305 (92%)	278 (99%)	3 (1%)	80	79
All	All	554/610 (91%)	547 (99%)	7 (1%)	76	73

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	236	PHE
1	A	350	PHE
1	A	420	SER
1	A	531	ARG
1	B	420	SER
1	B	456	PHE
1	B	471	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	429	ASN
1	A	538	ASN
1	B	292	HIS
1	B	499	ASN
1	B	538	ASN

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](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	ACO	A	796	-	23,28,53	1.50	4 (17%)	33,40,79	1.75	6 (18%)
2	ACO	B	896	-	23,28,53	1.48	4 (17%)	33,40,79	1.79	6 (18%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ACO	A	796	-	-	0/36/36/67	0/0/0/3
2	ACO	B	896	-	-	0/36/36/67	0/0/0/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	896	ACO	C-S1P	-3.29	1.52	1.74
2	A	796	ACO	C-S1P	-3.00	1.54	1.74
2	A	796	ACO	O5P-C5P	2.00	1.27	1.23
2	B	896	ACO	C9P-N8P	2.13	1.38	1.33
2	B	896	ACO	O-C	2.15	1.32	1.21
2	A	796	ACO	O-C	2.43	1.34	1.21
2	A	796	ACO	CH3-C	3.40	1.65	1.50
2	B	896	ACO	CH3-C	3.47	1.66	1.50

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	796	ACO	O-C-CH3	-4.48	101.46	122.83
2	B	896	ACO	O-C-CH3	-4.22	102.69	122.83
2	A	796	ACO	C2P-C3P-N4P	-3.02	106.31	112.36
2	B	896	ACO	C3P-C2P-S1P	-2.73	104.06	111.36
2	B	896	ACO	C2P-C3P-N4P	-2.50	107.36	112.36
2	A	796	ACO	C3P-C2P-S1P	-2.14	105.64	111.36
2	B	896	ACO	O2A-P1A-O1A	2.33	118.07	110.58
2	A	796	ACO	O2A-P1A-O3A	2.34	115.69	105.09
2	A	796	ACO	O2A-P1A-O1A	2.40	118.32	110.58
2	B	896	ACO	O2A-P1A-O3A	2.49	116.41	105.09
2	A	796	ACO	C3P-N4P-C5P	5.17	132.97	122.79
2	B	896	ACO	C3P-N4P-C5P	5.50	133.60	122.79

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 ⓘ

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	325/360 (90%)	1.30	75 (23%) 1 1	14, 38, 75, 83	0
1	B	333/360 (92%)	0.99	49 (14%) 3 3	12, 28, 72, 82	0
All	All	658/720 (91%)	1.14	124 (18%) 2 2	12, 33, 74, 83	0

All (124) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	350	PHE	16.7
1	B	318	HIS	10.0
1	A	321	ILE	8.9
1	B	356	LEU	8.9
1	B	352	MET	8.7
1	A	238	TRP	7.6
1	B	285	LYS	7.5
1	A	286	THR	7.3
1	B	258	ILE	7.0
1	A	236	PHE	6.6
1	B	353	ILE	6.5
1	A	246	THR	6.1
1	B	321	ILE	5.7
1	A	281	THR	5.7
1	B	351	ARG	5.7
1	A	318	HIS	5.6
1	B	236	PHE	5.5
1	B	286	THR	5.4
1	A	247	LEU	5.4
1	B	355	ASP	5.3
1	A	315	CYS	5.1
1	B	357	GLN	5.1
1	A	398	GLN	5.0
1	A	319	ARG	5.0

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Mol	Chain	Res	Type	RSRZ
1	A	322	GLN	5.0
1	B	346	PHE	4.9
1	B	273	ILE	4.9
1	A	244	GLY	4.8
1	A	593	LYS	4.8
1	A	358	LEU	4.7
1	A	273	ILE	4.6
1	B	257	ASP	4.5
1	A	403	CYS	4.4
1	A	592	PHE	4.4
1	B	317	MET	4.4
1	A	248	VAL	4.3
1	A	288	ILE	4.2
1	A	297	ASN	4.2
1	B	278	THR	4.1
1	B	348	GLU	4.1
1	A	550	ARG	4.0
1	A	239	PHE	4.0
1	A	303	ARG	3.9
1	A	314	SER	3.9
1	B	350	PHE	3.8
1	B	239	PHE	3.8
1	A	291	VAL	3.7
1	A	280	ASN	3.7
1	B	395	GLU	3.7
1	A	245	GLY	3.7
1	A	270	LEU	3.6
1	B	284	GLY	3.6
1	A	344	PHE	3.6
1	A	405	ASP	3.6
1	B	359	HIS	3.4
1	A	277	LEU	3.4
1	B	323	MET	3.4
1	B	344	PHE	3.4
1	B	304	LYS	3.4
1	B	593	LYS	3.4
1	A	317	MET	3.4
1	B	234	PRO	3.3
1	A	359	HIS	3.3
1	A	298	LEU	3.2
1	B	358	LEU	3.2
1	A	256	LYS	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	547	ASN	3.2
1	A	285	LYS	3.2
1	B	272	SER	3.2
1	A	300	MET	3.1
1	A	382	LYS	3.1
1	B	315	CYS	3.1
1	B	314	SER	3.1
1	A	252	TYR	3.0
1	A	275	LYS	3.0
1	A	395	GLU	3.0
1	B	345	LYS	3.0
1	B	246	THR	2.9
1	A	324	GLY	2.9
1	A	427	LYS	2.9
1	A	313	PRO	2.8
1	B	322	GLN	2.8
1	B	403	CYS	2.8
1	A	283	TYR	2.8
1	B	245	GLY	2.8
1	A	348	GLU	2.8
1	B	316	ALA	2.8
1	A	278	THR	2.8
1	B	238	TRP	2.8
1	A	323	MET	2.7
1	B	279	SER	2.7
1	A	381	GLY	2.7
1	A	590	GLU	2.6
1	B	349	ASP	2.6
1	A	290	ASP	2.6
1	B	277	LEU	2.5
1	A	257	ASP	2.5
1	B	533	CYS	2.5
1	A	591	LEU	2.5
1	A	346	PHE	2.5
1	A	548	PHE	2.5
1	A	305	GLY	2.5
1	A	396	LEU	2.5
1	A	337	CYS	2.4
1	A	284	GLY	2.4
1	B	275	LYS	2.4
1	A	243	ILE	2.4
1	B	256	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	316	ALA	2.3
1	B	535	LEU	2.3
1	B	281	THR	2.3
1	A	251	VAL	2.3
1	A	484	GLU	2.2
1	A	369	GLN	2.2
1	B	456	PHE	2.2
1	A	428	ASP	2.1
1	A	564	ASP	2.1
1	A	554	VAL	2.1
1	A	306	ASN	2.1
1	A	295	LEU	2.1
1	A	549	LEU	2.1
1	A	255	PRO	2.0
1	A	304	LYS	2.0
1	B	389	GLU	2.0

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 carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	ACO	B	896	29/51	0.92	0.14	1.13	16,20,37,39	0
2	ACO	A	796	29/51	0.93	0.14	0.30	15,20,43,45	0

6.5 Other polymers

There are no such residues in this entry.