Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 06:50 AM GMT

PDB ID	:	2YK0
Title	:	Structure of the N-terminal NTS-DBL1-alpha and CIDR-gamma double domain of the PfEMP1 protein from Plasmodium falciparum varO strain.
Authors	:	Lewit-Bentley, A.; Juillerat, A.; Vigan-Womas, I.; Guillotte, M.; Hessel, A.; Raynal, B.; Mercereau-Puijalon, O.; Bentley, G.A.
Deposited on	:	2011-05-25
Resolution	:	2.80 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

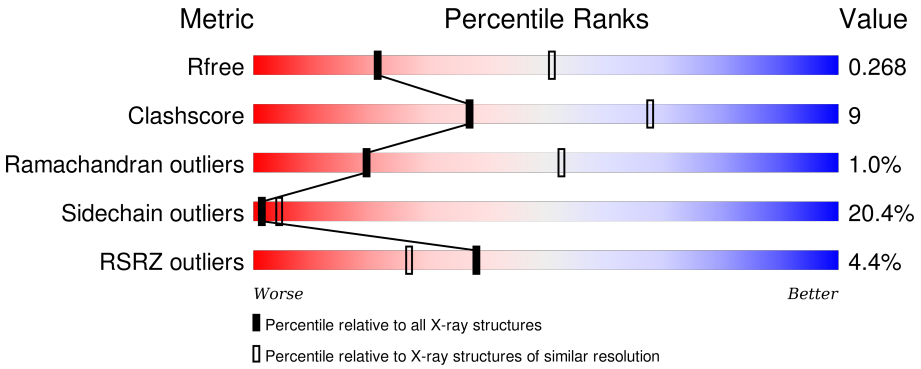
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 2.80 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	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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 <=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	790	4% 61% 22% 5% 12%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SO4	A	1719	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5728 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 ERYTHROCYTE MEMBRANE PROTEIN 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	698	Total	C	N	O	S	0	4	0
			5684	3570	1012	1066	36			

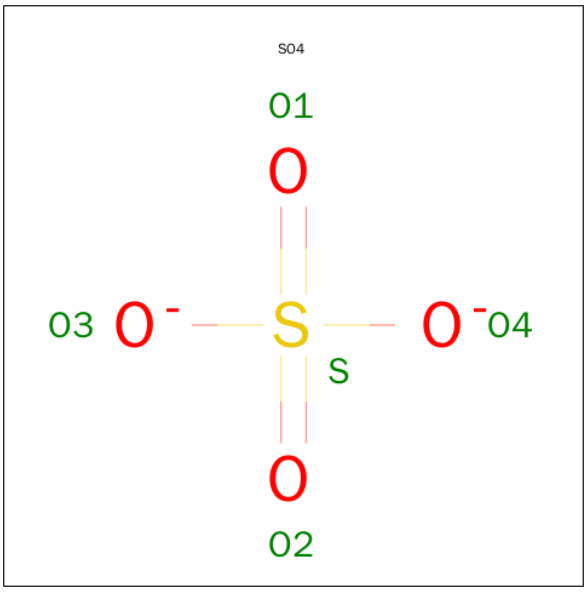
There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	ILE	-	EXPRESSION TAG	UNP B7T1P0
A	-1	SER	-	EXPRESSION TAG	UNP B7T1P0
A	0	GLU	-	EXPRESSION TAG	UNP B7T1P0
A	1	PHE	-	EXPRESSION TAG	UNP B7T1P0
A	10	ALA	THR	ENGINEERED MUTATION	UNP B7T1P0
A	125	ALA	SER	ENGINEERED MUTATION	UNP B7T1P0
A	128	ALA	THR	ENGINEERED MUTATION	UNP B7T1P0
A	154	ALA	THR	ENGINEERED MUTATION	UNP B7T1P0
A	186	ALA	SER	ENGINEERED MUTATION	UNP B7T1P0
A	250	ALA	SER	ENGINEERED MUTATION	UNP B7T1P0
A	360	ALA	SER	ENGINEERED MUTATION	UNP B7T1P0
A	657	ALA	SER	ENGINEERED MUTATION	UNP B7T1P0

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Mg	0	0
			2	2		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0

- Molecule 4 is water.

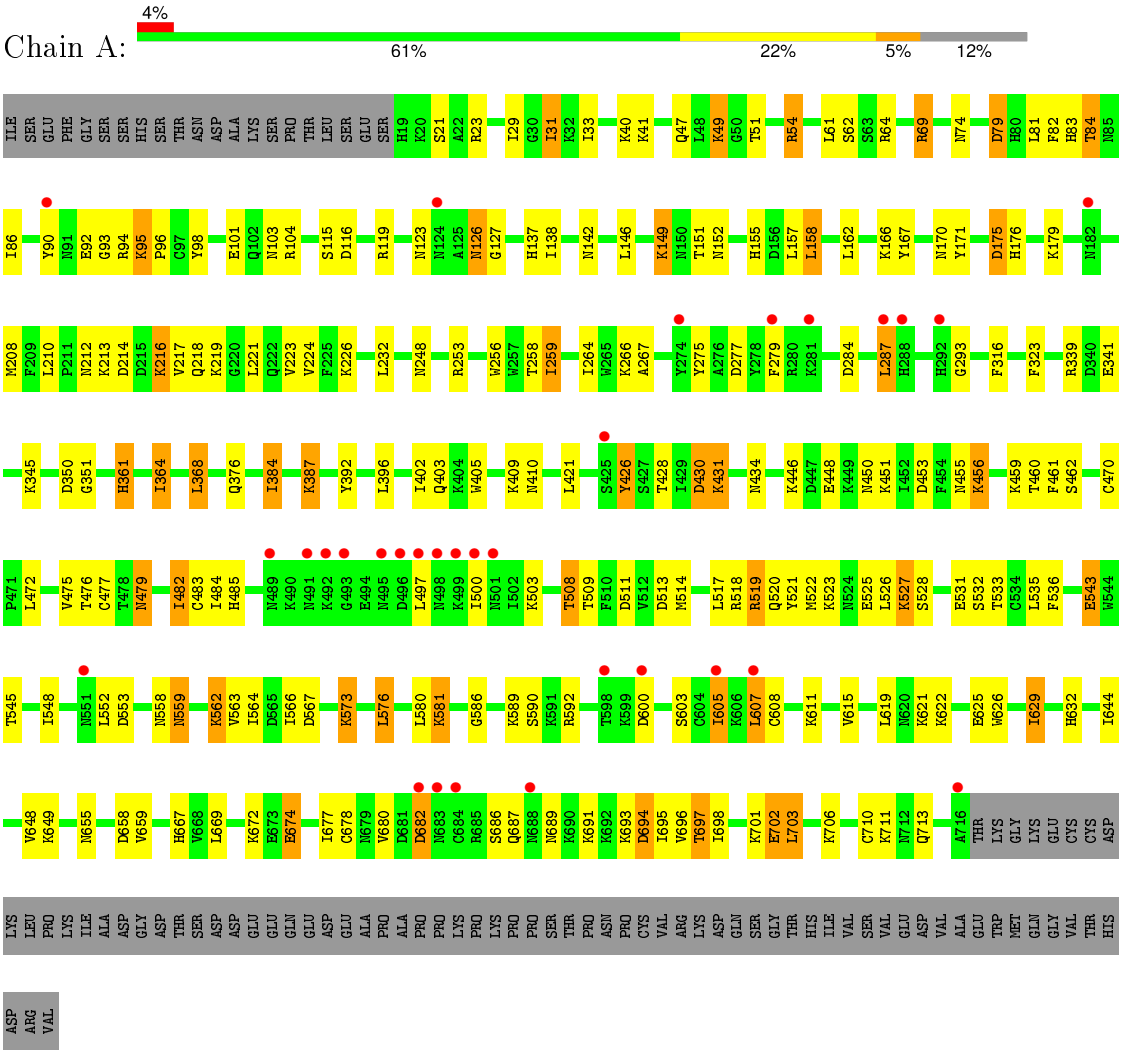
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	32	Total O 32 32	0	0

3 Residue-property plots (i)

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: ERYTHROCYTE MEMBRANE PROTEIN 1

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	157.91Å 144.36Å 75.84Å 90.00° 102.89° 90.00°	Depositor
Resolution (Å)	40.23 – 2.80 40.09 – 2.80	Depositor EDS
% Data completeness (in resolution range)	96.8 (40.23-2.80) 96.8 (40.09-2.80)	Depositor EDS
R_{merge}	0.25	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.53 (at 2.81Å)	Xtriage
Refinement program	BUSTER 2.9.3	Depositor
R, R_{free}	0.213 , 0.255 0.234 , 0.268	Depositor DCC
R_{free} test set	2007 reflections (5.36%)	DCC
Wilson B-factor (Å^2)	51.8	Xtriage
Anisotropy	0.757	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 88.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.44$, $< L^2 > = 0.27$	Xtriage
Outliers	0 of 39458 reflections	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	5728	wwPDB-VP
Average B, all atoms (Å^2)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.73% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, SO4

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	0.49	0/5824	0.76	0/7833

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	5684	0	5469	105	0
2	A	2	0	0	0	0
3	A	10	0	0	0	0
4	A	32	0	0	2	0
All	All	5728	0	5469	105	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 9.

All (105) 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:626:TRP:HZ3	1:A:697:THR:HG22	1.21	1.04
1:A:626:TRP:CZ3	1:A:697:THR:HG22	2.02	0.94
1:A:508:THR:CG2	1:A:632:HIS:HD2	1.88	0.85
1:A:54:ARG:HG2	1:A:54:ARG:HH11	1.40	0.85
1:A:626:TRP:HH2	1:A:694:ASP:HB2	1.42	0.84
1:A:508:THR:CG2	1:A:632:HIS:CD2	2.64	0.80
1:A:51:THR:H	1:A:155:HIS:HD2	1.27	0.79
1:A:364:ILE:HA	1:A:573:LYS:HG3	1.65	0.79
1:A:79:ASP:H	1:A:83:HIS:HD2	1.32	0.77
1:A:84:THR:HG23	1:A:86:ILE:H	1.52	0.73
1:A:79:ASP:H	1:A:83:HIS:CD2	2.07	0.71
1:A:519:ARG:HH21	1:A:532:SER:HB3	1.58	0.69
1:A:626:TRP:HH2	1:A:694:ASP:CB	2.06	0.69
1:A:533:THR:HG23	1:A:536:PHE:H	1.57	0.67
1:A:508:THR:HG23	1:A:632:HIS:CD2	2.30	0.66
1:A:508:THR:HG22	1:A:632:HIS:HD2	1.63	0.63
1:A:84:THR:HG21	1:A:94:ARG:HB3	1.80	0.63
1:A:51:THR:H	1:A:155:HIS:CD2	2.11	0.63
1:A:548:ILE:HG12	1:A:553:ASP:HB3	1.82	0.62
1:A:277:ASP:OD1	1:A:293:GLY:HA2	1.99	0.61
1:A:384:ILE:HD12	1:A:461:PHE:HA	1.82	0.61
1:A:453:ASP:H	1:A:460:THR:HG21	1.66	0.60
1:A:176:HIS:O	1:A:179:LYS:HB2	2.01	0.60
1:A:674:GLU:HB3	1:A:693:LYS:HD3	1.83	0.59
1:A:626:TRP:CH2	1:A:694:ASP:HB2	2.31	0.59
1:A:323:PHE:HD1	1:A:387:LYS:HG2	1.66	0.59
1:A:543:GLU:H	1:A:559:ASN:ND2	2.01	0.58
1:A:104:ARG:HD2	1:A:137:HIS:HB2	1.85	0.57
1:A:51:THR:N	1:A:155:HIS:HD2	2.00	0.57
1:A:576:LEU:HD22	1:A:580:LEU:HD11	1.87	0.56
1:A:648:VAL:HG12	1:A:695:ILE:HD12	1.88	0.55
1:A:430:ASP:HA	1:A:455:ASN:HD21	1.71	0.55
1:A:697:THR:O	1:A:701:LYS:HD3	2.08	0.53
1:A:519:ARG:HA	1:A:526:LEU:HD12	1.91	0.53
1:A:79:ASP:N	1:A:83:HIS:HD2	2.05	0.53
1:A:462:SER:HA	1:A:518:ARG:HH12	1.74	0.53
1:A:47:GLN:NE2	1:A:232:LEU:HA	2.24	0.52
1:A:559:ASN:H	1:A:559:ASN:HD22	1.57	0.52
1:A:157:LEU:HD13	1:A:224:VAL:HG11	1.93	0.51
1:A:426:TYR:HA	1:A:431:LYS:HB2	1.93	0.51
1:A:477:CYS:HA	1:A:483:CYS:HA	1.93	0.50
1:A:149:LYS:H	1:A:149:LYS:HE3	1.75	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:667:HIS:CD2	1:A:667:HIS:H	2.30	0.50
1:A:49:LYS:HD2	1:A:259:ILE:HG22	1.93	0.50
1:A:217:VAL:O	1:A:221:LEU:HB2	2.12	0.49
1:A:287:LEU:H	1:A:287:LEU:HD22	1.77	0.49
1:A:90:TYR:HD1	1:A:98:TYR:CE1	2.30	0.49
1:A:626:TRP:HZ3	1:A:697:THR:CG2	2.09	0.49
1:A:475:VAL:HG12	1:A:485:HIS:HA	1.94	0.49
1:A:90:TYR:CD1	1:A:98:TYR:CE1	3.01	0.48
1:A:96:PRO:HG2	1:A:167:TYR:CG	2.48	0.48
1:A:351:GLY:O	1:A:368:LEU:HB2	2.12	0.48
1:A:562:LYS:HE3	1:A:563:VAL:HG23	1.94	0.48
1:A:341:GLU:HB3	1:A:482:ILE:HA	1.96	0.48
1:A:253:ARG:NH2	4:A:2018:HOH:O	2.41	0.48
1:A:518:ARG:HG3	4:A:2027:HOH:O	2.13	0.48
1:A:519:ARG:HH12	1:A:527:LYS:HZ3	1.60	0.48
1:A:126:ASN:HB3	1:A:396:LEU:CD1	2.44	0.47
1:A:93:GLY:HA3	1:A:95:LYS:HE3	1.95	0.47
1:A:536:PHE:HZ	1:A:586:GLY:HA2	1.79	0.47
1:A:47:GLN:HE22	1:A:232:LEU:HA	1.78	0.47
1:A:166:LYS:HA	1:A:267:ALA:O	2.14	0.47
1:A:23:ARG:NH2	1:A:279:PHE:O	2.47	0.47
1:A:626:TRP:CH2	1:A:696:VAL:HB	2.51	0.46
1:A:54:ARG:HG2	1:A:54:ARG:NH1	2.19	0.46
1:A:79:ASP:OD1	1:A:81:LEU:HB2	2.16	0.46
1:A:608:CYS:HA	1:A:611:LYS:HD2	1.97	0.46
1:A:29:ILE:O	1:A:33:ILE:HG12	2.16	0.45
1:A:702:GLU:HG3	1:A:703:LEU:N	2.31	0.45
1:A:543:GLU:HG3	1:A:558:ASN:HB2	1.99	0.44
1:A:219:LYS:O	1:A:223:VAL:HG23	2.17	0.44
1:A:655:ASN:O	1:A:659:VAL:HG23	2.17	0.44
1:A:479:ASN:H	1:A:479:ASN:ND2	2.16	0.44
1:A:430:ASP:HB2	1:A:431:LYS:HE3	2.00	0.44
1:A:511:ASP:OD2	1:A:621:LYS:NZ	2.45	0.44
1:A:69[B]:ARG:HD3	1:A:82:PHE:HD1	1.83	0.43
1:A:54:ARG:HH11	1:A:54:ARG:CG	2.18	0.43
1:A:323:PHE:CD1	1:A:387:LYS:HG2	2.49	0.43
1:A:361:HIS:CD2	1:A:470:CYS:HB2	2.54	0.43
1:A:119:ARG:HD2	1:A:208:MET:HA	1.99	0.43
1:A:127:GLY:HA2	1:A:396:LEU:HB2	2.01	0.43
1:A:622:LYS:HA	1:A:625:GLU:HB2	2.01	0.43
1:A:453:ASP:OD2	1:A:456:LYS:HB2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:LEU:HD23	1:A:256:TRP:CE2	2.54	0.42
1:A:581:LYS:HA	1:A:581:LYS:HD2	1.91	0.42
1:A:69[A]:ARG:HG3	1:A:82:PHE:HB3	2.01	0.42
1:A:605:ILE:HG22	1:A:607:LEU:HG	2.02	0.42
1:A:69[B]:ARG:HD3	1:A:82:PHE:CD1	2.54	0.42
1:A:103:ASN:ND2	1:A:175:ASP:OD2	2.52	0.42
1:A:431:LYS:HG2	1:A:431:LYS:H	1.61	0.42
1:A:158:LEU:O	1:A:162:LEU:HG	2.20	0.42
1:A:686:SER:HB2	1:A:689:ASN:HB2	2.02	0.42
1:A:31:ILE:H	1:A:31:ILE:HG13	1.70	0.42
1:A:361:HIS:HD2	1:A:470:CYS:HB2	1.84	0.41
1:A:682:ASP:OD1	1:A:682:ASP:N	2.49	0.41
1:A:92:GLU:O	1:A:94:ARG:HD3	2.21	0.41
1:A:590:SER:HB3	1:A:615:VAL:HG13	2.03	0.41
1:A:626:TRP:HD1	1:A:629:ILE:HG12	1.85	0.41
1:A:392:TYR:CZ	1:A:396:LEU:HD11	2.56	0.41
1:A:216:LYS:NZ	1:A:216:LYS:H	2.19	0.41
1:A:171:TYR:C	1:A:171:TYR:CD1	2.94	0.41
1:A:158:LEU:HG	1:A:162:LEU:HG	2.03	0.41
1:A:461:PHE:O	1:A:518:ARG:NH1	2.54	0.41
1:A:644:ILE:HG23	1:A:694:ASP:OD2	2.21	0.40
1:A:562:LYS:HD3	1:A:562:LYS:H	1.86	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	700/790 (89%)	637 (91%)	56 (8%)	7 (1%)	19 52

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	523	LYS
1	A	175	ASP
1	A	527	LYS
1	A	687	GLN
1	A	213	LYS
1	A	522	MET
1	A	600	ASP

5.3.2 Protein sidechains (i)

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	618/720 (86%)	490 (79%)	128 (21%)	1 4

All (128) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	SER
1	A	31	ILE
1	A	40	LYS
1	A	41	LYS
1	A	49	LYS
1	A	54	ARG
1	A	61	LEU
1	A	62	SER
1	A	64	ARG
1	A	69[A]	ARG
1	A	69[B]	ARG
1	A	74	ASN
1	A	79	ASP
1	A	84	THR
1	A	95	LYS
1	A	101	GLU
1	A	115	SER
1	A	116	ASP
1	A	123	ASN
1	A	126	ASN

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Mol	Chain	Res	Type
1	A	138	ILE
1	A	142	ASN
1	A	146	LEU
1	A	149	LYS
1	A	151	THR
1	A	152	ASN
1	A	158	LEU
1	A	170	ASN
1	A	210	LEU
1	A	212	ASN
1	A	214	ASP
1	A	216	LYS
1	A	218	GLN
1	A	226	LYS
1	A	248	ASN
1	A	258	THR
1	A	259	ILE
1	A	264	ILE
1	A	266	LYS
1	A	275	TYR
1	A	284	ASP
1	A	287	LEU
1	A	316	PHE
1	A	339	ARG
1	A	345	LYS
1	A	350	ASP
1	A	361	HIS
1	A	364	ILE
1	A	368	LEU
1	A	376	GLN
1	A	384	ILE
1	A	387	LYS
1	A	402	ILE
1	A	403	GLN
1	A	405	TRP
1	A	409	LYS
1	A	410	ASN
1	A	421	LEU
1	A	426	TYR
1	A	428	THR
1	A	430	ASP
1	A	431	LYS

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Mol	Chain	Res	Type
1	A	434	ASN
1	A	446	LYS
1	A	448	GLU
1	A	450	ASN
1	A	451	LYS
1	A	456	LYS
1	A	459	LYS
1	A	472	LEU
1	A	476	THR
1	A	479	ASN
1	A	482	ILE
1	A	484	ILE
1	A	497	LEU
1	A	500	ILE
1	A	503	LYS
1	A	508	THR
1	A	509	THR
1	A	513	ASP
1	A	514	MET
1	A	517	LEU
1	A	519	ARG
1	A	520	GLN
1	A	521	TYR
1	A	525[A]	GLU
1	A	525[B]	GLU
1	A	528	SER
1	A	531[A]	GLU
1	A	531[B]	GLU
1	A	535	LEU
1	A	543	GLU
1	A	545	THR
1	A	552	LEU
1	A	559	ASN
1	A	562	LYS
1	A	564	ILE
1	A	566	ILE
1	A	567	ASP
1	A	573	LYS
1	A	576	LEU
1	A	581	LYS
1	A	589	LYS
1	A	592	ARG

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Mol	Chain	Res	Type
1	A	603	SER
1	A	605	ILE
1	A	607	LEU
1	A	619	LEU
1	A	629	ILE
1	A	649	LYS
1	A	658	ASP
1	A	669	LEU
1	A	672	LYS
1	A	674	GLU
1	A	677	ILE
1	A	678	CYS
1	A	680	VAL
1	A	682	ASP
1	A	691	LYS
1	A	694	ASP
1	A	697	THR
1	A	698	ILE
1	A	702	GLU
1	A	703	LEU
1	A	706	LYS
1	A	710	CYS
1	A	711	LYS
1	A	713	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	GLN
1	A	74	ASN
1	A	83	HIS
1	A	150	ASN
1	A	155	HIS
1	A	160	ASN
1	A	170	ASN
1	A	174	ASN
1	A	248	ASN
1	A	263	GLN
1	A	352	HIS
1	A	362	ASN
1	A	363	GLN
1	A	410	ASN

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Mol	Chain	Res	Type
1	A	455	ASN
1	A	479	ASN
1	A	559	ASN
1	A	632	HIS
1	A	638	HIS
1	A	667	HIS

5.3.3 RNA

There are no RNA molecules 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

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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	$\#\|Z\| > 2$
3	SO4	A	1719	-	4,4,4	0.17	0	6,6,6	0.09	0
3	SO4	A	1720	-	4,4,4	0.34	0	6,6,6	0.06	0

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
3	SO4	A	1719	-	-	0/0/0/0	0/0/0/0
3	SO4	A	1720	-	-	0/0/0/0	0/0/0/0

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

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, 95^{th} 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	698/790 (88%)	0.19	31 (4%) 38 26	30, 70, 132, 147	5 (0%)

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	496	ASP	5.4
1	A	495	ASN	5.4
1	A	716	ALA	5.1
1	A	493	GLY	4.5
1	A	90	TYR	4.0
1	A	292	HIS	3.5
1	A	598	THR	3.5
1	A	124	ASN	3.5
1	A	497	LEU	3.4
1	A	489	ASN	3.3
1	A	688	ASN	3.3
1	A	600	ASP	3.3
1	A	683	ASN	3.2
1	A	492	LYS	2.8
1	A	498	ASN	2.8
1	A	607	LEU	2.8
1	A	274	TYR	2.8
1	A	279	PHE	2.7
1	A	682	ASP	2.6
1	A	491	ASN	2.6
1	A	684	CYS	2.5
1	A	499	LYS	2.4
1	A	281	LYS	2.3
1	A	551	ASN	2.3
1	A	501	ASN	2.3
1	A	500	ILE	2.3
1	A	182	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	605	ILE	2.2
1	A	288	HIS	2.2
1	A	287	LEU	2.2
1	A	425	SER	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 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, 95^{th} 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
3	SO4	A	1719	5/5	0.73	0.34	2.95	134,138,139,139	3
2	MG	A	1717	1/1	0.77	0.16	-0.56	60,60,60,60	0
2	MG	A	1718	1/1	0.91	0.09	-	43,43,43,43	0
3	SO4	A	1720	5/5	0.91	0.47	-	75,79,80,81	5

6.5 Other polymers (i)

There are no such residues in this entry.