Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 05:42 PM GMT

PDB ID	:	4J7G
Title	:	Crystal structure of EvaA, a 2,3-dehydratase in complex with dTDP-fucose and dTDP-rhamnose
Authors	:	Holden, H.M.; Kubiak, R.L.; Thoden, J.B.
Deposited on	:	2013-02-13
Resolution	:	1.70 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:

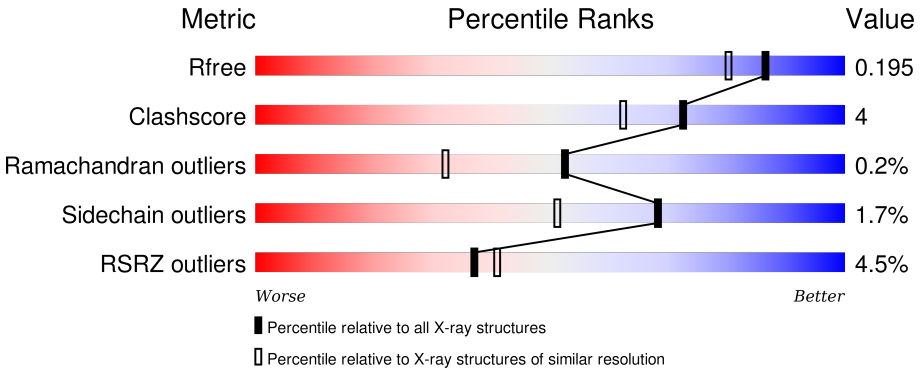
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	3190 (1.70-1.70)
Clashscore	102246	3585 (1.70-1.70)
Ramachandran outliers	100387	3527 (1.70-1.70)
Sidechain outliers	100360	3527 (1.70-1.70)
RSRZ outliers	91569	3200 (1.70-1.70)

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	471	4% 82% 12% • 5%
1	B	471	4% 82% 12% • 5%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7953 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.

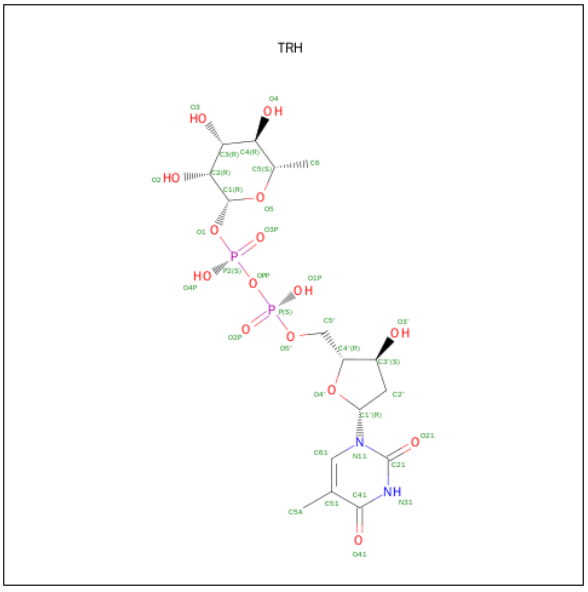
- Molecule 1 is a protein called EvaA 2,3-dehydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	448	Total	C	N	O	S	0	4	0
			3599	2285	644	658	12			
1	B	446	Total	C	N	O	S	0	4	0
			3580	2276	639	654	11			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	381	ALA	ARG	ENGINEERED MUTATION	UNP O52793
B	381	ALA	ARG	ENGINEERED MUTATION	UNP O52793

- Molecule 2 is 2'-DEOXY-THYMIDINE-BETA-L-RHAMNOSE (three-letter code: TRH) (formula: $C_{16}H_{26}N_2O_{15}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			35	16	2	15	2		

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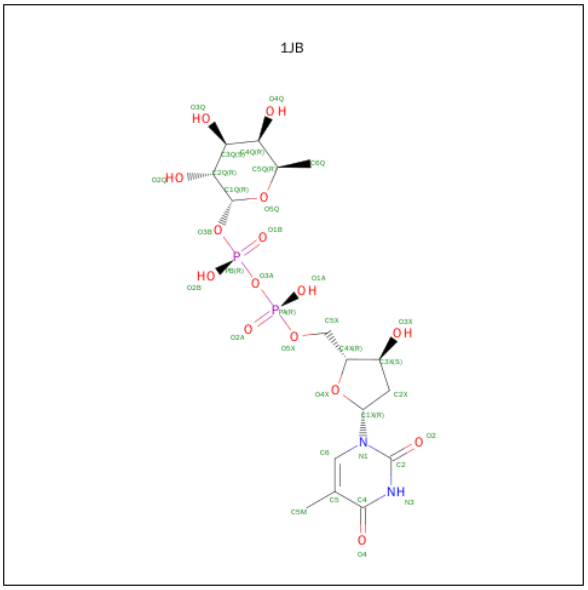
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			35	16	2	15	2		

- Molecule 3 is [[(2R,3S,5R)-5-[5-METHYL-2,4-BIS(OXIDANYLIDENE)PYRIMIDIN-1-YL]-3-OXIDANYL-OXOLAN-2-YL]METHOXY-OXIDANYL-PHOSPHORYL] [(2R,3R,4S,5R,6R)-6-METHYL-3,4,5-TRIS(OXIDANYL)OXAN-2-YL] HYDROGEN PHOSPHATE (three-letter code: 1JB) (formula: $C_{16}H_{26}N_2O_{15}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			35	16	2	15	2		
3	B	1	Total	C	N	O	P	0	0
			35	16	2	15	2		

- Molecule 4 is water.

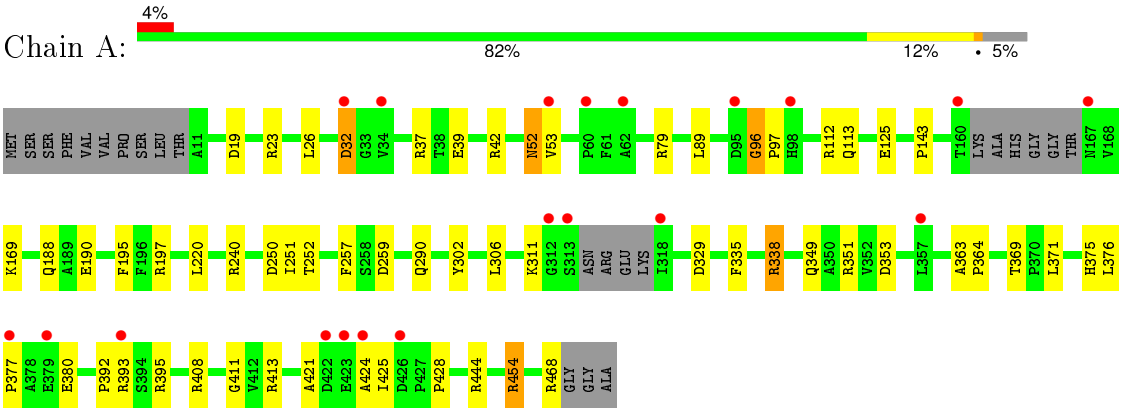
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	352	Total	O	0	0
			352	352		
4	B	282	Total	O	0	0
			282	282		

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.

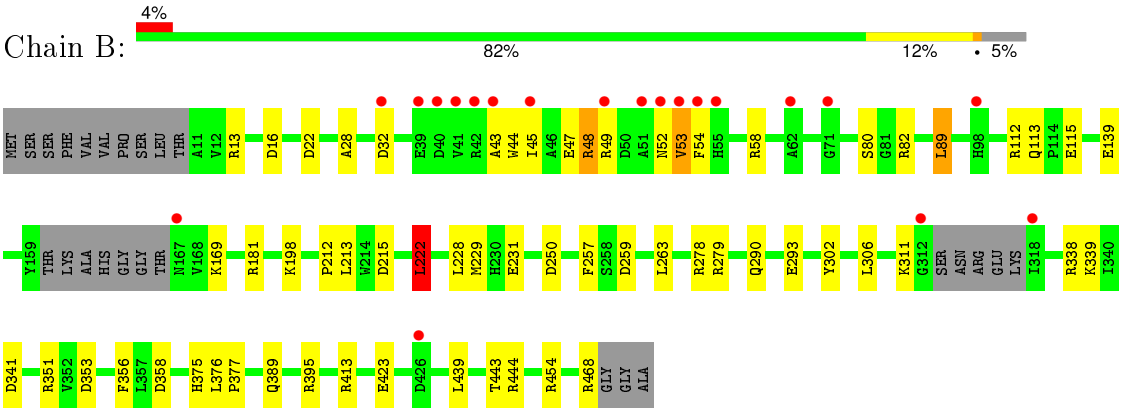
- Molecule 1: EvaA 2,3-dehydratase

Chain A:



- Molecule 1: EvaA 2,3-dehydratase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	105.83Å 108.22Å 110.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	110.21 – 1.70 44.54 – 1.70	Depositor EDS
% Data completeness (in resolution range)	95.3 (110.21-1.70) 95.3 (44.54-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	3.22 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.172 , 0.197 0.171 , 0.195	Depositor DCC
R_{free} test set	6991 reflections (5.53%)	DCC
Wilson B-factor (Å^2)	22.9	Xtriage
Anisotropy	0.045	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 40.9	EDS
Estimated twinning fraction	0.012 for -h,l,k 0.012 for -l,-k,-h 0.014 for k,h,-l 0.000 for k,l,h 0.000 for l,h,k	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 133122 reflections	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	7953	wwPDB-VP
Average B, all atoms (Å^2)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.40% 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: TRH, 1JB

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	0.58	0/3702	1.34	27/5044 (0.5%)
1	B	0.56	0/3686	1.35	37/5029 (0.7%)
All	All	0.57	0/7388	1.35	64/10073 (0.6%)

There are no bond length outliers.

All (64) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	229[A]	MET	CG-SD-CE	-16.77	73.36	100.20
1	B	229[B]	MET	CG-SD-CE	-16.77	73.36	100.20
1	A	468	ARG	NE-CZ-NH2	-15.12	112.74	120.30
1	A	351	ARG	NE-CZ-NH2	-13.30	113.65	120.30
1	A	444	ARG	NE-CZ-NH1	-10.01	115.29	120.30
1	A	250	ASP	CB-CG-OD2	9.80	127.12	118.30
1	B	395	ARG	NE-CZ-NH2	9.73	125.17	120.30
1	B	454	ARG	NE-CZ-NH1	-9.26	115.67	120.30
1	A	112	ARG	NE-CZ-NH2	-8.78	115.91	120.30
1	A	197	ARG	NE-CZ-NH2	-8.50	116.05	120.30
1	A	395	ARG	NE-CZ-NH2	-8.31	116.15	120.30
1	A	468	ARG	NE-CZ-NH1	8.23	124.41	120.30
1	B	278	ARG	NE-CZ-NH1	-8.09	116.26	120.30
1	B	395	ARG	NE-CZ-NH1	-7.95	116.32	120.30
1	A	32	ASP	CB-CG-OD2	-7.95	111.14	118.30
1	B	58	ARG	NE-CZ-NH1	-7.86	116.37	120.30
1	A	444	ARG	NE-CZ-NH2	7.55	124.08	120.30
1	B	13	ARG	NE-CZ-NH1	7.28	123.94	120.30
1	B	181	ARG	NE-CZ-NH1	-7.28	116.66	120.30
1	B	444	ARG	NE-CZ-NH2	7.24	123.92	120.30
1	A	454	ARG	NE-CZ-NH1	-7.18	116.71	120.30
1	A	190	GLU	CA-CB-CG	-7.17	97.64	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	240	ARG	NE-CZ-NH1	-7.16	116.72	120.30
1	B	351	ARG	NE-CZ-NH2	-7.13	116.73	120.30
1	B	16[A]	ASP	CB-CG-OD1	7.10	124.69	118.30
1	B	16[B]	ASP	CB-CG-OD1	7.10	124.69	118.30
1	B	215	ASP	CB-CG-OD2	-6.85	112.14	118.30
1	B	454	ARG	NE-CZ-NH2	6.84	123.72	120.30
1	A	353	ASP	CB-CG-OD1	6.77	124.39	118.30
1	A	197	ARG	NE-CZ-NH1	6.74	123.67	120.30
1	B	22	ASP	CB-CG-OD1	6.68	124.31	118.30
1	B	439	LEU	CB-CG-CD1	-6.66	99.68	111.00
1	A	395	ARG	NE-CZ-NH1	6.57	123.58	120.30
1	B	444	ARG	NE-CZ-NH1	-6.54	117.03	120.30
1	A	257	PHE	CB-CG-CD2	-6.50	116.25	120.80
1	B	259	ASP	CB-CG-OD1	6.43	124.09	118.30
1	B	169	LYS	CD-CE-NZ	-6.32	97.17	111.70
1	A	259	ASP	CB-CG-OD1	6.20	123.88	118.30
1	B	257	PHE	CB-CG-CD2	-6.17	116.48	120.80
1	A	89	LEU	CB-CG-CD1	-6.07	100.68	111.00
1	B	356	PHE	CB-CG-CD2	-5.96	116.63	120.80
1	B	263	LEU	CB-CG-CD1	-5.88	101.01	111.00
1	A	42	ARG	NE-CZ-NH2	5.85	123.22	120.30
1	A	220	LEU	CB-CG-CD2	-5.79	101.16	111.00
1	A	79	ARG	NE-CZ-NH1	5.66	123.13	120.30
1	B	413	ARG	NE-CZ-NH2	5.61	123.10	120.30
1	B	89	LEU	CB-CG-CD1	-5.58	101.51	111.00
1	B	222	LEU	CB-CG-CD1	5.54	120.42	111.00
1	A	96	GLY	N-CA-C	-5.51	99.32	113.10
1	A	19	ASP	CB-CG-OD2	5.51	123.26	118.30
1	B	250	ASP	CB-CG-OD2	5.46	123.21	118.30
1	B	341	ASP	CB-CA-C	-5.44	99.52	110.40
1	B	257	PHE	CB-CG-CD1	5.41	124.59	120.80
1	A	26	LEU	CB-CG-CD1	-5.40	101.81	111.00
1	A	195	PHE	CB-CG-CD2	-5.35	117.05	120.80
1	B	341	ASP	CB-CG-OD1	-5.33	113.50	118.30
1	B	48	ARG	NE-CZ-NH2	-5.28	117.66	120.30
1	A	413	ARG	NE-CZ-NH2	5.26	122.93	120.30
1	B	13	ARG	CB-CG-CD	5.19	125.10	111.60
1	B	358	ASP	CB-CG-OD2	-5.09	113.72	118.30
1	B	443	THR	N-CA-CB	5.08	119.96	110.30
1	B	358	ASP	CB-CG-OD1	5.05	122.85	118.30
1	B	228	LEU	CB-CG-CD1	-5.03	102.44	111.00
1	B	338	ARG	NE-CZ-NH2	5.03	122.81	120.30

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	3599	0	3508	35	0
1	B	3580	0	3484	28	0
2	A	35	0	24	0	0
2	B	35	0	24	0	0
3	A	35	0	25	0	0
3	B	35	0	25	0	0
4	A	352	0	0	3	0
4	B	282	0	0	5	0
All	All	7953	0	7090	61	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 4.

All (61) 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:96:GLY:C	1:A:97:PRO:CA	2.39	0.91
1:A:338[A]:ARG:HH11	1:A:425:ILE:H	1.22	0.85
1:A:113:GLN:NE2	4:A:673:HOH:O	2.06	0.83
1:A:96:GLY:C	1:A:97:PRO:CD	2.54	0.76
1:B:49:ARG:O	1:B:52:ASN:HB3	1.89	0.72
1:A:52:ASN:HD22	1:A:53:VAL:N	1.87	0.72
1:A:371:LEU:O	1:A:371:LEU:HD23	1.91	0.70
1:B:80:SER:OG	1:B:82:ARG:HG2	1.95	0.66
1:B:389:GLN:NE2	4:B:758:HOH:O	2.19	0.66
1:A:113:GLN:HE22	1:A:408:ARG:HH11	1.45	0.65
1:A:376:LEU:HB3	1:A:377:PRO:HD2	1.84	0.59
1:A:23[B]:ARG:NH1	1:A:125:GLU:OE2	2.32	0.58
1:B:375:HIS:H	1:B:375:HIS:CD2	2.22	0.57
1:B:53:VAL:HG12	1:B:54:PHE:N	2.19	0.57
1:B:139:GLU:OE2	4:B:826:HOH:O	2.17	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:113:GLN:OE1	4:B:674:HOH:O	2.17	0.55
1:B:49:ARG:HA	1:B:52:ASN:HB3	1.89	0.54
1:A:338[A]:ARG:NH1	1:A:425:ILE:H	2.01	0.54
1:A:143:PRO:HG3	1:A:411:GLY:HA2	1.91	0.53
1:A:169:LYS:HE3	4:A:951:HOH:O	2.07	0.53
1:B:45:ILE:O	1:B:49:ARG:N	2.33	0.52
1:B:112:ARG:HD2	1:B:279:ARG:HD3	1.91	0.52
1:B:49:ARG:O	1:B:52:ASN:CA	2.58	0.52
1:A:338[B]:ARG:HG2	1:A:421:ALA:HB3	1.93	0.51
1:B:49:ARG:C	1:B:52:ASN:H	2.14	0.51
1:A:349:GLN:HG2	1:A:363:ALA:HB3	1.92	0.51
1:B:231:GLU:OE2	1:B:468:ARG:NH2	2.44	0.51
1:B:49:ARG:O	1:B:52:ASN:N	2.44	0.49
1:A:338[B]:ARG:HD2	1:A:424:ALA:HA	1.94	0.49
1:B:49:ARG:O	1:B:52:ASN:CB	2.58	0.49
1:A:52:ASN:HD22	1:A:53:VAL:H	1.60	0.49
1:B:389:GLN:HA	1:B:389:GLN:HE21	1.78	0.49
1:B:28:ALA:HA	1:B:222:LEU:HB3	1.95	0.48
1:A:52:ASN:C	1:A:52:ASN:HD22	2.14	0.47
1:A:335:PHE:CE1	1:A:364:PRO:HA	2.50	0.47
1:A:251:ILE:HG13	1:A:252:THR:HG23	1.97	0.47
1:A:52:ASN:ND2	1:A:52:ASN:C	2.69	0.46
1:B:112:ARG:HD3	1:B:112:ARG:HH11	1.63	0.45
1:B:115:GLU:O	1:B:198:LYS:NZ	2.49	0.45
1:A:329:ASP:HB3	1:A:369:THR:CG2	2.46	0.45
1:A:37:ARG:NE	1:A:39:GLU:OE1	2.49	0.45
1:B:353:ASP:OD2	4:B:875:HOH:O	2.21	0.45
1:A:96:GLY:CA	1:A:97:PRO:CD	2.95	0.44
1:A:52:ASN:ND2	1:A:53:VAL:N	2.61	0.43
1:A:37:ARG:HH21	1:A:39:GLU:CD	2.22	0.43
1:A:393:ARG:HD3	1:B:213:LEU:HD23	2.01	0.42
1:B:43:ALA:O	1:B:47:GLU:CB	2.67	0.42
1:A:375:HIS:CD2	1:A:375:HIS:H	2.37	0.42
1:A:188:GLN:HB3	1:A:454:ARG:CZ	2.50	0.42
1:A:290:GLN:NE2	4:A:827:HOH:O	2.34	0.42
1:A:392:PRO:HA	1:B:212:PRO:HA	2.02	0.41
1:A:371:LEU:C	1:A:371:LEU:HD23	2.41	0.41
1:B:376:LEU:HB3	1:B:377:PRO:HD2	2.03	0.41
1:A:338[A]:ARG:HH11	1:A:425:ILE:N	2.04	0.41
1:B:44:TRP:O	1:B:48:ARG:HG2	2.21	0.41
1:B:290:GLN:NE2	4:B:796:HOH:O	2.30	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:377:PRO:HG2	1:A:380:GLU:OE1	2.21	0.40
1:B:89:LEU:HA	1:B:89:LEU:HD12	1.89	0.40
1:A:393:ARG:HH11	1:A:393:ARG:HG2	1.87	0.40
1:A:311:LYS:HB2	1:A:311:LYS:HE3	1.92	0.40
1:B:43:ALA:O	1:B:47:GLU:HB2	2.22	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	444/471 (94%)	433 (98%)	10 (2%)	1 (0%)	52	32
1	B	444/471 (94%)	424 (96%)	19 (4%)	1 (0%)	52	32
All	All	888/942 (94%)	857 (96%)	29 (3%)	2 (0%)	52	32

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	302	TYR
1	B	302	TYR

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	383/398 (96%)	377 (98%)	6 (2%)	70	54
1	B	381/398 (96%)	373 (98%)	8 (2%)	61	42
All	All	764/796 (96%)	750 (98%)	14 (2%)	68	49

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	ASP
1	A	52	ASN
1	A	306	LEU
1	A	338[A]	ARG
1	A	338[B]	ARG
1	A	428	PRO
1	B	32	ASP
1	B	53	VAL
1	B	222	LEU
1	B	293	GLU
1	B	306	LEU
1	B	311	LYS
1	B	339	LYS
1	B	423	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	52	ASN
1	A	113	GLN
1	A	135	GLN
1	A	375	HIS
1	B	98	HIS
1	B	135	GLN
1	B	290	GLN
1	B	375	HIS

5.3.3 RNA

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)

4 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	TRH	A	501	-	29,37,37	0.61	0	43,57,57	2.45	5 (11%)
3	1JB	A	502	-	29,37,37	0.68	0	43,57,57	2.11	14 (32%)
2	TRH	B	501	-	29,37,37	0.60	0	43,57,57	2.64	6 (13%)
3	1JB	B	502	-	29,37,37	0.65	0	43,57,57	2.82	19 (44%)

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	TRH	A	501	-	-	0/17/53/53	0/3/3/3
3	1JB	A	502	-	-	0/17/53/53	0/3/3/3
2	TRH	B	501	-	-	0/17/53/53	0/3/3/3
3	1JB	B	502	-	-	0/17/53/53	0/3/3/3

There are no bond length outliers.

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	501	TRH	O5-C1-O1	-8.52	100.12	111.36
2	B	501	TRH	O5-C1-O1	-7.67	101.24	111.36
2	B	501	TRH	C51-C41-N31	-7.00	117.34	125.14
3	B	502	1JB	O5Q-C1Q-O3B	-6.82	102.37	111.36
2	A	501	TRH	C51-C41-N31	-5.77	118.72	125.14
3	B	502	1JB	O4Q-C4Q-C5Q	-5.48	96.99	109.84
3	B	502	1JB	C5-C4-N3	-4.72	119.89	125.14
3	A	502	1JB	C5-C4-N3	-4.41	120.23	125.14
3	A	502	1JB	PA-O3A-PB	-3.89	121.80	132.73
3	A	502	1JB	C2X-C1X-N1	-3.75	105.05	114.16
2	B	501	TRH	C1-O5-C5	-3.65	107.41	113.64
3	A	502	1JB	O5Q-C1Q-O3B	-3.54	106.69	111.36
2	B	501	TRH	C2'-C1'-N11	-3.30	106.14	114.16
3	B	502	1JB	PA-O3A-PB	-3.29	123.50	132.73
3	A	502	1JB	O3X-C3X-C2X	-3.23	100.06	110.74
2	A	501	TRH	O5-C5-C4	-3.17	104.04	109.53
2	A	501	TRH	C1-C2-C3	-2.67	104.70	109.97
3	B	502	1JB	C3Q-C4Q-C5Q	-2.64	105.27	109.72
2	B	501	TRH	O3-C3-C4	-2.62	104.43	110.34
3	B	502	1JB	C2X-C1X-N1	-2.38	108.37	114.16
3	B	502	1JB	O3Q-C3Q-C4Q	-2.25	105.26	110.34
3	B	502	1JB	O3X-C3X-C2X	-2.10	103.81	110.74
3	B	502	1JB	O2B-PB-O3A	-2.09	95.62	105.09
3	B	502	1JB	O2Q-C2Q-C3Q	-2.09	105.64	110.34
3	A	502	1JB	O4Q-C4Q-C5Q	-2.05	105.03	109.84
3	A	502	1JB	O3A-PA-O5X	-2.04	97.52	102.94
3	A	502	1JB	C1Q-C2Q-C3Q	2.15	114.20	109.97
3	B	502	1JB	O4X-C1X-C2X	2.23	110.72	106.27
3	A	502	1JB	O5Q-C5Q-C6Q	2.35	111.78	106.64
3	A	502	1JB	O1A-PA-O3A	2.44	116.18	105.09
3	A	502	1JB	O4X-C1X-N1	2.48	112.00	107.72
3	B	502	1JB	O5Q-C5Q-C6Q	2.50	112.09	106.64
3	B	502	1JB	C2X-C3X-C4X	2.61	108.18	102.77
3	B	502	1JB	O3Q-C3Q-C2Q	2.74	116.51	110.34
3	B	502	1JB	C6Q-C5Q-C4Q	3.32	119.62	113.08
3	A	502	1JB	C2X-C3X-C4X	3.38	109.77	102.77
3	B	502	1JB	C2Q-C3Q-C4Q	3.45	117.24	110.79
3	B	502	1JB	C1Q-O5Q-C5Q	4.21	120.81	113.64
3	A	502	1JB	O5Q-C5Q-C4Q	4.77	117.79	109.53
3	B	502	1JB	O5Q-C5Q-C4Q	5.27	118.66	109.53
3	A	502	1JB	C4-N3-C2	5.50	120.00	115.25
3	B	502	1JB	C4-N3-C2	8.85	122.89	115.25
2	A	501	TRH	C41-N31-C21	10.07	123.95	115.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	501	TRH	C41-N31-C21	10.97	124.73	115.25

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, 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	448/471 (95%)	-0.01	20 (4%)	37	41	13, 23, 47, 64	0
1	B	446/471 (94%)	-0.01	20 (4%)	37	41	14, 24, 48, 72	0
All	All	894/942 (94%)	-0.01	40 (4%)	37	41	13, 24, 48, 72	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	160	THR	5.8
1	B	39	GLU	5.5
1	B	318	ILE	5.5
1	A	318	ILE	5.1
1	B	54	PHE	5.0
1	A	312	GLY	4.8
1	A	422	ASP	4.5
1	B	32	ASP	4.0
1	B	41	VAL	3.9
1	A	313	SER	3.9
1	B	43	ALA	3.6
1	A	426	ASP	3.5
1	A	167	ASN	3.3
1	A	424	ALA	3.2
1	B	71	GLY	3.2
1	A	95	ASP	3.1
1	B	42	ARG	3.0
1	B	167	ASN	3.0
1	A	379	GLU	2.9
1	A	53	VAL	2.8
1	B	62	ALA	2.6
1	B	52	ASN	2.6
1	B	53	VAL	2.6
1	B	312	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	55	HIS	2.5
1	A	60	PRO	2.3
1	A	423	GLU	2.2
1	B	98	HIS	2.2
1	B	51	ALA	2.2
1	A	377	PRO	2.2
1	B	49	ARG	2.2
1	A	357	LEU	2.1
1	A	32	ASP	2.1
1	B	40	ASP	2.1
1	A	98	HIS	2.1
1	A	62	ALA	2.1
1	A	393	ARG	2.1
1	B	45	ILE	2.1
1	B	426	ASP	2.0
1	A	34	VAL	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, 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	1JB	A	502	35/35	0.95	0.09	-0.04	19,27,38,43	0
2	TRH	B	501	35/35	0.97	0.07	-0.26	18,23,26,31	0
3	1JB	B	502	35/35	0.96	0.08	-0.36	20,29,40,46	0
2	TRH	A	501	35/35	0.97	0.06	-0.71	20,22,25,26	0

6.5 Other polymers (i)

There are no such residues in this entry.