



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:11 AM GMT

PDB ID : 2ZX5
Title : alpha-L-fucosidase complexed with inhibitor, F10
Authors : Wu, H.-J.; Ko, T.-P.; Ho, C.-W.; Lin, C.-H.; Wang, A.H.-J.
Deposited on : 2008-12-19
Resolution : 2.65 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

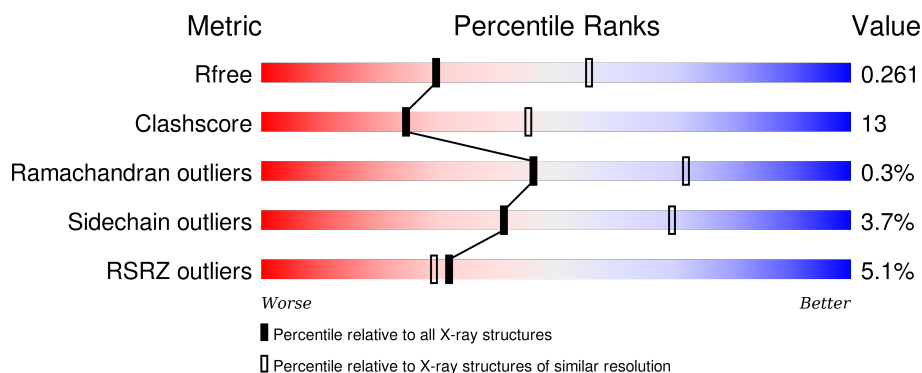
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 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	3152 (2.70-2.62)
Clashscore	102246	3524 (2.70-2.62)
Ramachandran outliers	100387	3469 (2.70-2.62)
Sidechain outliers	100360	3469 (2.70-2.62)
RSRZ outliers	91569	3161 (2.70-2.62)

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	455	<div> <div>5%</div> <div>71%</div> <div>24%</div> <div>• •</div> </div>
1	B	455	<div> <div>5%</div> <div>67%</div> <div>27%</div> <div>• •</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7667 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.

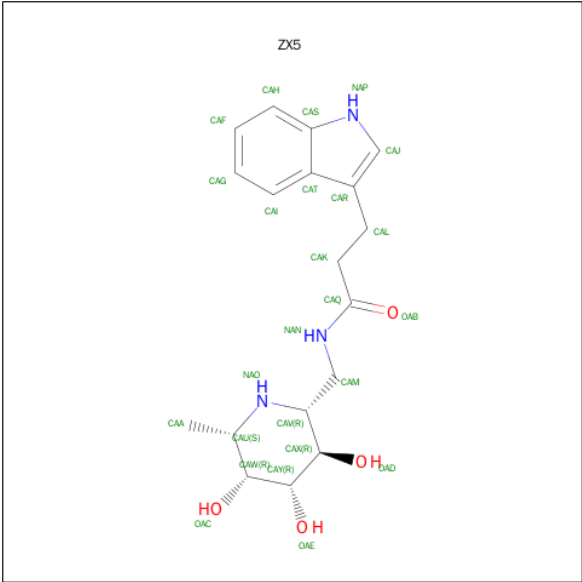
- Molecule 1 is a protein called Alpha-L-fucosidase, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	441	Total	C	N	O	S	0	0	0
			3639	2372	599	660	8			
1	B	441	Total	C	N	O	S	0	0	0
			3639	2372	599	660	8			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	450	HIS	-	EXPRESSION TAG	UNP Q9WYE2
A	451	HIS	-	EXPRESSION TAG	UNP Q9WYE2
A	452	HIS	-	EXPRESSION TAG	UNP Q9WYE2
A	453	HIS	-	EXPRESSION TAG	UNP Q9WYE2
A	454	HIS	-	EXPRESSION TAG	UNP Q9WYE2
A	455	HIS	-	EXPRESSION TAG	UNP Q9WYE2
B	450	HIS	-	EXPRESSION TAG	UNP Q9WYE2
B	451	HIS	-	EXPRESSION TAG	UNP Q9WYE2
B	452	HIS	-	EXPRESSION TAG	UNP Q9WYE2
B	453	HIS	-	EXPRESSION TAG	UNP Q9WYE2
B	454	HIS	-	EXPRESSION TAG	UNP Q9WYE2
B	455	HIS	-	EXPRESSION TAG	UNP Q9WYE2

- Molecule 2 is 3-(1H-INDOL-3-YL)-N-([(2R,3R,4R,5R,6S)-3,4,5-TRIHYDROXY-6-METHYLPIPERIDIN-2-YL]METHYL)PROPANAMIDE (three-letter code: ZX5) (formula: C₁₈H₂₅N₃O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			25	18	3	4		
2	B	1	Total	C	N	O	0	0
			25	18	3	4		

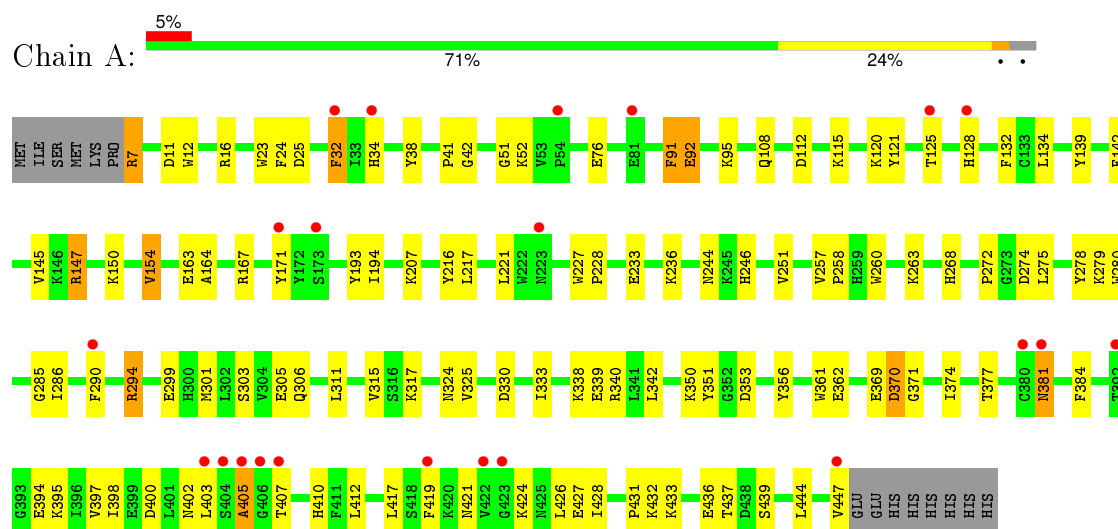
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	174	Total	O	0	0
			174	174		
3	B	165	Total	O	0	0
			165	165		

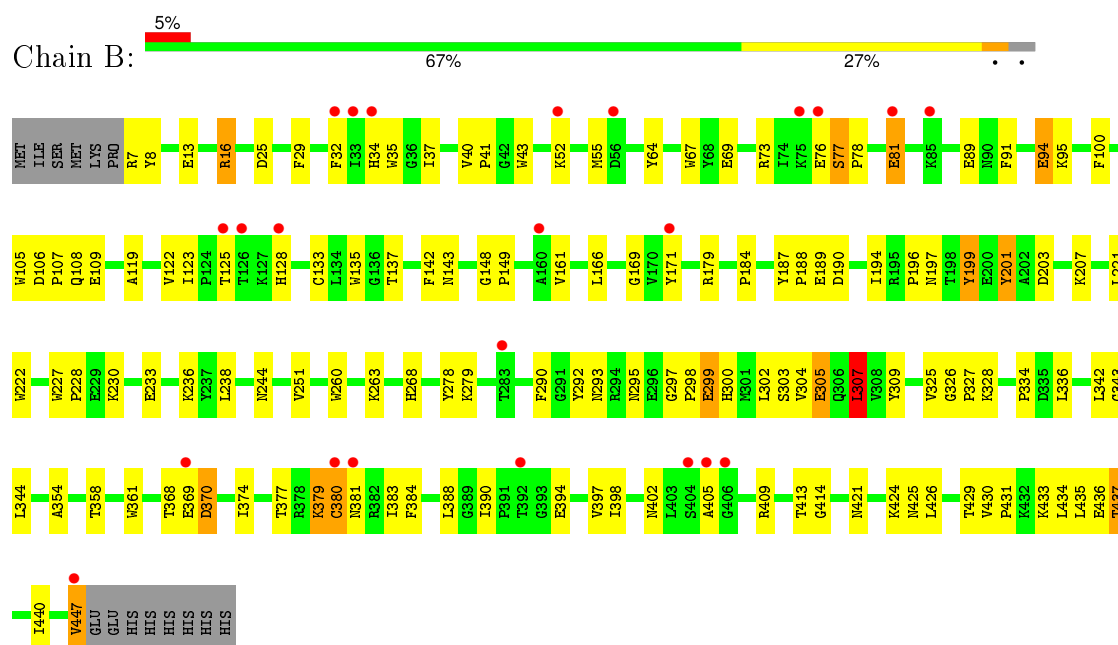
3 Residue-property plots

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: Alpha-L-fucosidase, putative



- Molecule 1: Alpha-L-fucosidase, putative



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	180.52Å 180.52Å 169.27Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.65 29.37 – 2.65	Depositor EDS
% Data completeness (in resolution range)	99.4 (30.00-2.65) 99.4 (29.37-2.65)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.18 (at 2.64Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.191 , 0.256 0.194 , 0.261	Depositor DCC
R_{free} test set	1521 reflections (4.96%)	DCC
Wilson B-factor (Å ²)	47.0	Xtriage
Anisotropy	0.077	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 61.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.54$, $\langle L^2 \rangle = 0.38$	Xtriage
Outliers	0 of 30649 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7667	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.39% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZX5

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	0.96	1/3759 (0.0%)	0.93	3/5108 (0.1%)
1	B	0.95	2/3759 (0.1%)	0.93	3/5108 (0.1%)
All	All	0.96	3/7518 (0.0%)	0.93	6/10216 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	94	GLU	CG-CD	7.17	1.62	1.51
1	B	305	GLU	CG-CD	5.19	1.59	1.51
1	A	154	VAL	CA-CB	5.15	1.65	1.54

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	380	CYS	CA-CB-SG	-6.22	102.81	114.00
1	B	77	SER	N-CA-C	-6.11	94.51	111.00
1	A	330	ASP	CB-CG-OD2	6.02	123.72	118.30
1	A	147	ARG	NE-CZ-NH1	-5.82	117.39	120.30
1	A	371	GLY	N-CA-C	5.24	126.20	113.10
1	B	307	LEU	CB-CG-CD1	-5.10	102.33	111.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	199	TYR	Sidechain
1	B	201	TYR	Sidechain
1	B	309	TYR	Sidechain

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	3639	0	3506	92	0
1	B	3639	0	3506	99	0
2	A	25	0	25	2	0
2	B	25	0	25	4	0
3	A	174	0	0	1	0
3	B	165	0	0	2	0
All	All	7667	0	7062	186	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 13.

All (186) 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:B:106:ASP:OD2	1:B:109:GLU:HG3	1.69	0.93
1:B:297:GLY:H	1:B:300:HIS:HD2	1.22	0.87
1:A:217:LEU:HD11	1:A:246:HIS:HB2	1.57	0.85
1:B:409:ARG:NH1	1:B:414:GLY:O	2.11	0.82
1:B:297:GLY:H	1:B:300:HIS:CD2	2.02	0.77
1:B:433:LYS:HG3	1:B:434:LEU:HD23	1.71	0.72
1:B:161:VAL:HG13	1:B:166:LEU:HB2	1.73	0.71
1:B:233:GLU:OE1	1:B:236:LYS:HE3	1.91	0.71
1:B:190:ASP:HB3	1:B:194:ILE:HD12	1.73	0.70
1:B:227:TRP:CG	1:B:228:PRO:HD2	2.28	0.69
1:B:25:ASP:HA	1:B:279:LYS:HE2	1.76	0.68
1:A:290:PHE:HA	1:A:324:ASN:ND2	2.09	0.67
1:B:297:GLY:N	1:B:300:HIS:HD2	1.92	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:430:VAL:HG12	1:B:435:LEU:HG	1.79	0.65
1:B:125:THR:HA	1:B:171:TYR:HB3	1.77	0.65
1:A:108:GLN:HE22	1:A:163:GLU:HB2	1.63	0.64
1:B:370:ASP:OD1	3:B:595:HOH:O	2.15	0.64
1:B:413:THR:HG21	1:B:434:LEU:HD22	1.78	0.63
1:A:395:LYS:HE3	1:A:427:GLU:OE1	1.99	0.63
1:B:388:LEU:N	1:B:388:LEU:HD23	2.14	0.62
1:B:402:ASN:HA	1:B:421:ASN:OD1	1.99	0.62
1:A:91:PHE:CE2	1:A:95:LYS:HB3	2.36	0.61
1:A:121:TYR:HB3	1:A:167:ARG:HB2	1.83	0.61
1:A:400:ASP:HA	1:A:424:LYS:O	2.00	0.61
1:A:405:ALA:HB1	1:A:447:VAL:C	2.21	0.60
1:B:328:LYS:HE3	1:B:334:PRO:HG3	1.84	0.60
1:B:263:LYS:HD3	1:B:278:TYR:CE1	2.36	0.60
1:A:236:LYS:HE2	3:A:612:HOH:O	2.00	0.59
1:B:238:LEU:HD23	1:B:238:LEU:C	2.22	0.59
2:B:902:ZX5:HAI	2:B:902:ZX5:NAN	2.18	0.59
1:A:34:HIS:CG	1:A:290:PHE:HB3	2.37	0.58
1:A:381:ASN:HD22	1:A:381:ASN:N	2.01	0.58
1:A:405:ALA:HB1	1:A:447:VAL:O	2.04	0.58
1:B:405:ALA:HB1	1:B:447:VAL:O	2.03	0.58
1:A:134:LEU:O	1:A:154:VAL:HG23	2.03	0.58
1:A:12:TRP:CZ2	1:A:236:LYS:HG2	2.39	0.57
1:B:43:TRP:CD1	1:B:295:ASN:ND2	2.72	0.57
1:B:434:LEU:O	1:B:437:THR:HG22	2.03	0.57
1:B:369:GLU:OE2	1:B:397:VAL:HG11	2.04	0.57
1:A:260:TRP:CE2	1:B:16:ARG:HG2	2.40	0.57
1:A:433:LYS:O	1:A:437:THR:HG23	2.05	0.57
1:A:285:GLY:HA2	1:A:324:ASN:HB3	1.87	0.56
1:B:307:LEU:HD13	1:B:344:LEU:HD22	1.86	0.56
1:A:263:LYS:HD3	1:A:278:TYR:CE1	2.41	0.55
1:A:260:TRP:CZ2	1:B:16:ARG:HG2	2.41	0.55
1:B:227:TRP:CD2	1:B:228:PRO:HD2	2.42	0.54
1:A:125:THR:HA	1:A:171:TYR:HB3	1.87	0.54
1:A:274:ASP:CG	1:A:275:LEU:H	2.10	0.54
1:B:201:TYR:CE2	1:B:228:PRO:HG3	2.43	0.54
2:B:902:ZX5:CAI	2:B:902:ZX5:NAN	2.70	0.54
1:B:73:ARG:HB2	1:B:184:PRO:HB3	1.90	0.53
1:B:358:THR:HG22	1:B:379:LYS:HD2	1.90	0.53
1:A:274:ASP:OD2	1:A:275:LEU:N	2.41	0.53
1:B:379:LYS:HB2	1:B:384:PHE:HE1	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:419:PHE:HB2	1:A:428:ILE:HG12	1.89	0.53
1:B:34:HIS:CG	1:B:290:PHE:HB3	2.43	0.53
1:B:233:GLU:OE1	1:B:236:LYS:CE	2.57	0.53
1:A:305:GLU:CD	1:A:305:GLU:H	2.12	0.52
1:B:394:GLU:O	1:B:429:THR:HA	2.09	0.52
1:B:91:PHE:CE1	1:B:95:LYS:HD2	2.44	0.52
1:A:52:LYS:HG2	1:A:268:HIS:CD2	2.44	0.52
1:A:16:ARG:HD2	1:B:16:ARG:HH21	1.74	0.52
1:A:7:ARG:HB2	1:A:244:ASN:OD1	2.10	0.52
1:A:275:LEU:HD23	1:A:317:LYS:HA	1.92	0.51
1:A:16:ARG:HH11	1:B:16:ARG:CZ	2.22	0.51
1:B:142:PHE:CE2	1:B:179:ARG:HD2	2.45	0.51
1:A:92:GLU:OE2	1:A:95:LYS:HD2	2.10	0.51
1:A:128:HIS:CE1	2:A:901:ZX5:OAE	2.64	0.51
1:A:381:ASN:HD22	1:A:381:ASN:H	1.57	0.51
1:B:55:MET:HB3	1:B:188:PRO:HG2	1.92	0.51
1:A:403:LEU:HG	1:A:426:LEU:HD22	1.93	0.51
1:B:302:LEU:HD12	1:B:307:LEU:HD23	1.92	0.50
1:B:221:LEU:HB2	1:B:251:VAL:HG12	1.93	0.50
1:B:361:TRP:CE2	1:B:383:ILE:HD13	2.46	0.50
1:A:25:ASP:HA	1:A:279:LYS:HE2	1.93	0.50
1:A:301:MET:HE2	1:A:340:ARG:HD2	1.94	0.50
1:A:272:PRO:HG3	1:A:280:TRP:CZ2	2.47	0.50
1:B:171:TYR:CD2	1:B:171:TYR:C	2.86	0.50
1:B:161:VAL:CG1	1:B:166:LEU:HB2	2.40	0.49
1:B:135:TRP:CE2	1:B:137:THR:HB	2.46	0.49
1:A:374:ILE:HD12	1:A:398:ILE:HG23	1.94	0.49
1:A:311:LEU:O	1:A:315:VAL:HG23	2.12	0.49
1:A:91:PHE:C	1:A:91:PHE:CD2	2.86	0.49
1:A:16:ARG:HG2	1:B:260:TRP:CE2	2.48	0.49
1:A:369:GLU:HB2	1:A:397:VAL:HB	1.94	0.49
1:A:377:THR:OG1	1:A:384:PHE:HB2	2.13	0.48
1:B:122:VAL:HG23	1:B:166:LEU:HD12	1.96	0.48
1:A:115:LYS:NZ	1:A:164:ALA:O	2.43	0.48
1:B:40:VAL:O	1:B:41:PRO:C	2.51	0.48
1:B:78:PRO:O	1:B:81:GLU:HB2	2.14	0.47
1:B:100:PHE:O	1:B:149:PRO:HD3	2.14	0.47
1:A:120:LYS:HE2	1:A:356:TYR:CE2	2.50	0.47
1:B:37:ILE:HG22	1:B:100:PHE:CD2	2.50	0.47
1:B:433:LYS:HG3	1:B:434:LEU:CD2	2.43	0.47
1:A:11:ASP:HA	1:A:236:LYS:HD2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:305:GLU:HG3	1:B:440:ILE:HG12	1.97	0.47
1:A:216:TYR:O	1:A:217:LEU:C	2.53	0.46
1:A:370:ASP:OD2	1:A:370:ASP:O	2.33	0.46
1:A:42:GLY:O	1:A:294:ARG:HG2	2.14	0.46
1:B:179:ARG:HH21	1:B:179:ARG:HG2	1.81	0.46
1:A:145:VAL:O	1:A:150:LYS:HA	2.16	0.46
2:B:902:ZX5:HNAN	2:B:902:ZX5:HAI	1.79	0.46
1:A:147:ARG:HA	1:A:150:LYS:HE2	1.96	0.46
1:A:303:SER:OG	1:A:306:GLN:HG3	2.15	0.46
1:B:292:TYR:HB2	1:B:326:GLY:O	2.15	0.45
1:B:69:GLU:O	1:B:73:ARG:HG3	2.16	0.45
1:A:402:ASN:ND2	1:A:421:ASN:HB3	2.30	0.45
1:B:105:TRP:CZ2	1:B:107:PRO:HB3	2.52	0.45
1:B:91:PHE:CD1	1:B:95:LYS:HD2	2.51	0.45
1:B:433:LYS:HG3	1:B:434:LEU:N	2.31	0.45
1:B:368:THR:HG23	1:B:374:ILE:HD11	1.99	0.45
1:B:77:SER:O	1:B:78:PRO:C	2.54	0.45
1:B:302:LEU:CD1	1:B:307:LEU:HD23	2.47	0.45
1:B:325:VAL:O	1:B:327:PRO:HD3	2.16	0.45
1:B:201:TYR:HE2	1:B:228:PRO:HG3	1.80	0.45
1:B:89:GLU:C	1:B:91:PHE:H	2.20	0.45
1:A:402:ASN:HD22	1:A:402:ASN:HA	1.65	0.45
1:A:419:PHE:CB	1:A:428:ILE:HG12	2.47	0.45
1:B:94:GLU:H	1:B:94:GLU:CD	2.16	0.45
1:B:354:ALA:HB2	1:B:384:PHE:CD2	2.53	0.44
1:A:221:LEU:HB2	1:A:251:VAL:HG12	1.99	0.44
1:B:429:THR:O	1:B:431:PRO:HD3	2.17	0.44
1:B:29:PHE:CD1	1:B:119:ALA:HA	2.52	0.44
1:A:217:LEU:CD1	1:A:246:HIS:HB2	2.38	0.44
1:B:303:SER:O	1:B:304:VAL:C	2.54	0.44
1:A:350:LYS:HG2	1:A:351:TYR:CE2	2.53	0.44
1:A:16:ARG:HD2	1:B:16:ARG:NH2	2.32	0.44
1:B:377:THR:OG1	1:B:384:PHE:HB2	2.18	0.43
1:A:134:LEU:C	1:A:154:VAL:HG23	2.39	0.43
1:B:35:TRP:CD1	1:B:327:PRO:HB2	2.53	0.43
1:A:233:GLU:CD	1:A:236:LYS:HE3	2.39	0.43
1:A:139:TYR:CE2	1:A:207:LYS:HE2	2.54	0.43
1:B:203:ASP:O	1:B:207:LYS:HG3	2.17	0.43
1:A:108:GLN:NE2	1:A:163:GLU:HB2	2.31	0.43
1:B:293:ASN:OD1	1:B:295:ASN:HB2	2.18	0.43
1:B:424:LYS:HG3	1:B:425:ASN:OD1	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:410:HIS:HA	1:A:444:LEU:HD23	2.00	0.43
1:A:51:GLY:HA3	1:A:268:HIS:HB2	2.00	0.43
1:B:123:ILE:HG12	1:B:169:GLY:HA3	2.01	0.43
1:A:257:VAL:HB	1:A:258:PRO:HD2	2.01	0.43
1:B:298:PRO:N	1:B:299:GLU:OE1	2.52	0.43
1:A:227:TRP:CG	1:A:228:PRO:HD2	2.54	0.43
1:B:108:GLN:HG2	3:B:593:HOH:O	2.18	0.43
1:A:394:GLU:OE2	1:A:431:PRO:HA	2.19	0.42
1:B:368:THR:HA	1:B:397:VAL:O	2.19	0.42
1:B:16:ARG:HH11	1:B:16:ARG:CG	2.32	0.42
1:A:23:TRP:CD2	1:A:24:PHE:N	2.88	0.42
1:B:199:TYR:CD2	1:B:230:LYS:HD3	2.55	0.42
1:B:421:ASN:OD1	1:B:426:LEU:HD13	2.19	0.42
1:A:286:ILE:HG22	1:A:301:MET:CE	2.50	0.42
1:B:133:CYS:O	1:B:143:ASN:HA	2.20	0.42
1:B:52:LYS:HG3	1:B:268:HIS:CD2	2.55	0.42
1:B:64:TYR:O	1:B:67:TRP:HB2	2.19	0.42
1:B:196:PRO:O	1:B:197:ASN:HB2	2.19	0.42
1:A:417:LEU:HD11	1:A:444:LEU:HD22	2.01	0.42
1:A:381:ASN:ND2	1:A:381:ASN:H	2.17	0.42
1:A:38:TYR:O	1:A:41:PRO:HG2	2.19	0.42
1:B:434:LEU:O	1:B:435:LEU:C	2.57	0.42
1:B:187:TYR:HB3	1:B:189:GLU:OE1	2.19	0.42
1:A:121:TYR:HA	1:A:167:ARG:O	2.20	0.42
1:A:132:PHE:HB2	1:A:142:PHE:CE2	2.54	0.42
1:B:8:TYR:H	1:B:244:ASN:HD21	1.66	0.42
1:A:257:VAL:HB	1:A:258:PRO:CD	2.50	0.41
1:B:128:HIS:CE1	2:B:902:ZX5:OAE	2.73	0.41
1:A:128:HIS:HE1	2:A:901:ZX5:OAE	2.03	0.41
1:A:432:LYS:HG3	1:A:436:GLU:HG3	2.02	0.41
1:A:338:LYS:HB2	1:A:338:LYS:HE3	1.70	0.41
1:B:297:GLY:C	1:B:299:GLU:OE1	2.59	0.41
1:A:233:GLU:OE1	1:A:236:LYS:HE3	2.20	0.41
1:B:336:LEU:HD23	1:B:336:LEU:HA	1.71	0.41
1:A:32:PHE:HB3	1:A:324:ASN:HA	2.02	0.41
1:B:398:ILE:HD12	1:B:426:LEU:HD23	2.02	0.41
1:A:16:ARG:CG	1:B:260:TRP:CE2	3.03	0.41
1:A:16:ARG:HG3	1:B:260:TRP:CZ2	2.56	0.41
1:A:325:VAL:HG11	1:A:333:ILE:HD13	2.02	0.41
1:B:342:LEU:O	1:B:343:GLY:C	2.59	0.41
1:A:108:GLN:OE1	1:A:108:GLN:HA	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:263:LYS:HD3	1:A:278:TYR:CD1	2.56	0.41
1:A:342:LEU:HD23	1:A:342:LEU:HA	1.85	0.41
1:A:444:LEU:HD23	1:A:444:LEU:HA	1.87	0.40
1:A:361:TRP:CG	1:A:362:GLU:N	2.89	0.40
1:B:13:GLU:O	1:B:16:ARG:HB3	2.21	0.40
1:A:193:TYR:CE1	1:A:194:ILE:HG13	2.56	0.40
1:A:405:ALA:O	1:A:407:THR:N	2.51	0.40
1:A:410:HIS:CE1	1:A:412:LEU:HB3	2.56	0.40
1:A:303:SER:HB2	1:A:305:GLU:OE1	2.21	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	439/455 (96%)	411 (94%)	27 (6%)	1 (0%)	52	77
1	B	439/455 (96%)	409 (93%)	28 (6%)	2 (0%)	34	59
All	All	878/910 (96%)	820 (93%)	55 (6%)	3 (0%)	46	72

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	405	ALA
1	B	148	GLY
1	B	76	GLU

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	381/395 (96%)	368 (97%)	13 (3%)	44	72
1	B	381/395 (96%)	366 (96%)	15 (4%)	39	67
All	All	762/790 (96%)	734 (96%)	28 (4%)	41	69

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ARG
1	A	32	PHE
1	A	76	GLU
1	A	91	PHE
1	A	92	GLU
1	A	112	ASP
1	A	294	ARG
1	A	299	GLU
1	A	339	GLU
1	A	353	ASP
1	A	370	ASP
1	A	381	ASN
1	A	439	SER
1	B	7	ARG
1	B	16	ARG
1	B	32	PHE
1	B	81	GLU
1	B	222	TRP
1	B	299	GLU
1	B	307	LEU
1	B	370	ASP
1	B	379	LYS
1	B	380	CYS
1	B	381	ASN
1	B	390	ILE
1	B	436	GLU
1	B	437	THR
1	B	447	VAL

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	61	GLN
1	A	108	GLN
1	A	270	ASN
1	A	381	ASN
1	A	402	ASN
1	B	300	HIS
1	B	402	ASN

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 ⓘ

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	ZX5	A	901	-	26,27,27	0.96	1 (3%)	24,38,38	0.62	0
2	ZX5	B	902	-	26,27,27	0.97	1 (3%)	24,38,38	1.22	3 (12%)

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	ZX5	A	901	-	-	0/9/30/30	0/3/3/3
2	ZX5	B	902	-	-	0/9/30/30	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	901	ZX5	CAJ-CAR	-2.75	1.33	1.38
2	B	902	ZX5	CAJ-CAR	-2.72	1.33	1.38

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	902	ZX5	CAV-CAM-NAN	-2.98	104.55	112.84
2	B	902	ZX5	CAL-CAR-CAJ	-2.59	123.52	127.88
2	B	902	ZX5	CAK-CAQ-NAN	-2.20	112.63	116.46

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	901	ZX5	2	0
2	B	902	ZX5	4	0

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, 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	441/455 (96%)	0.14	22 (4%) 32 30	22, 39, 66, 82	0
1	B	441/455 (96%)	0.21	23 (5%) 31 28	25, 41, 68, 79	0
All	All	882/910 (96%)	0.17	45 (5%) 32 29	22, 40, 68, 82	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	406	GLY	4.7
1	A	392	THR	4.0
1	A	381	ASN	3.9
1	A	406	GLY	3.9
1	A	171	TYR	3.7
1	B	125	THR	3.6
1	A	404	SER	3.5
1	B	171	TYR	3.5
1	A	405	ALA	3.5
1	B	34	HIS	3.4
1	B	405	ALA	3.3
1	B	404	SER	3.1
1	B	381	ASN	3.1
1	A	403	LEU	2.9
1	A	422	VAL	2.9
1	B	56	ASP	2.9
1	A	125	THR	2.8
1	B	75	LYS	2.8
1	A	34	HIS	2.8
1	A	447	VAL	2.8
1	A	380	CYS	2.7
1	B	380	CYS	2.6
1	B	126	THR	2.6
1	B	81	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	173	SER	2.5
1	A	81	GLU	2.5
1	A	54	PRO	2.5
1	A	423	GLY	2.5
1	B	128	HIS	2.4
1	A	32	PHE	2.3
1	B	32	PHE	2.3
1	B	76	GLU	2.3
1	B	369	GLU	2.3
1	B	392	THR	2.3
1	B	85	LYS	2.3
1	B	52	LYS	2.3
1	A	128	HIS	2.2
1	B	447	VAL	2.2
1	B	160	ALA	2.1
1	B	283	THR	2.1
1	A	290	PHE	2.1
1	A	419	PHE	2.1
1	A	407	THR	2.0
1	A	223	ASN	2.0
1	B	33	ILE	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	ZX5	A	901	25/25	0.81	0.35	1.35	41,54,72,73	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	ZX5	B	902	25/25	0.80	0.29	0.52	32,44,64,65	0

6.5 Other polymers [i](#)

There are no such residues in this entry.