



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2014 – 03:32 AM GMT

PDB ID : 2WCS  
Title : CRYSTAL STRUCTURE OF DEBRANCHING ENZYME FROM NOSTOC PUNCTIFORME (NPDE)  
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Deposited on : 2009-03-16  
Resolution : 2.80 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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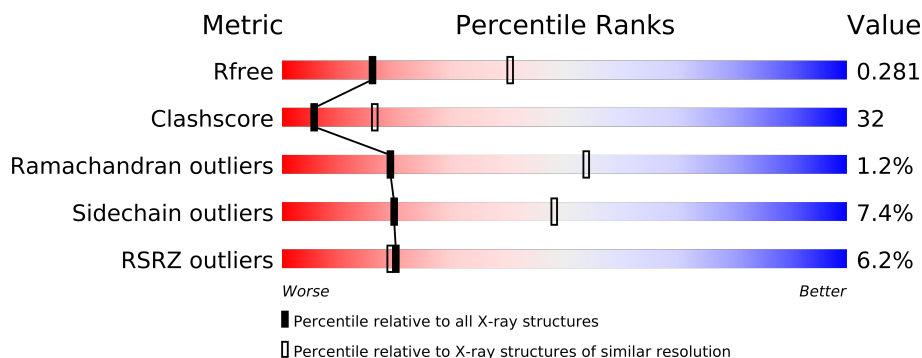
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

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))
$R_{free}$	66092	1799 (2.80-2.80)
Clashscore	79885	2295 (2.80-2.80)
Ramachandran outliers	78287	2252 (2.80-2.80)
Sidechain outliers	78261	2254 (2.80-2.80)
RSRZ outliers	66119	1802 (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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	488	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3968 atoms, of which 0 are hydrogen and 0 are deuterium.

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 AMYLASE, CATALYTIC REGION.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	487	Total	C	N	O	S	0	0	2
			3922	2526	662	727	7			

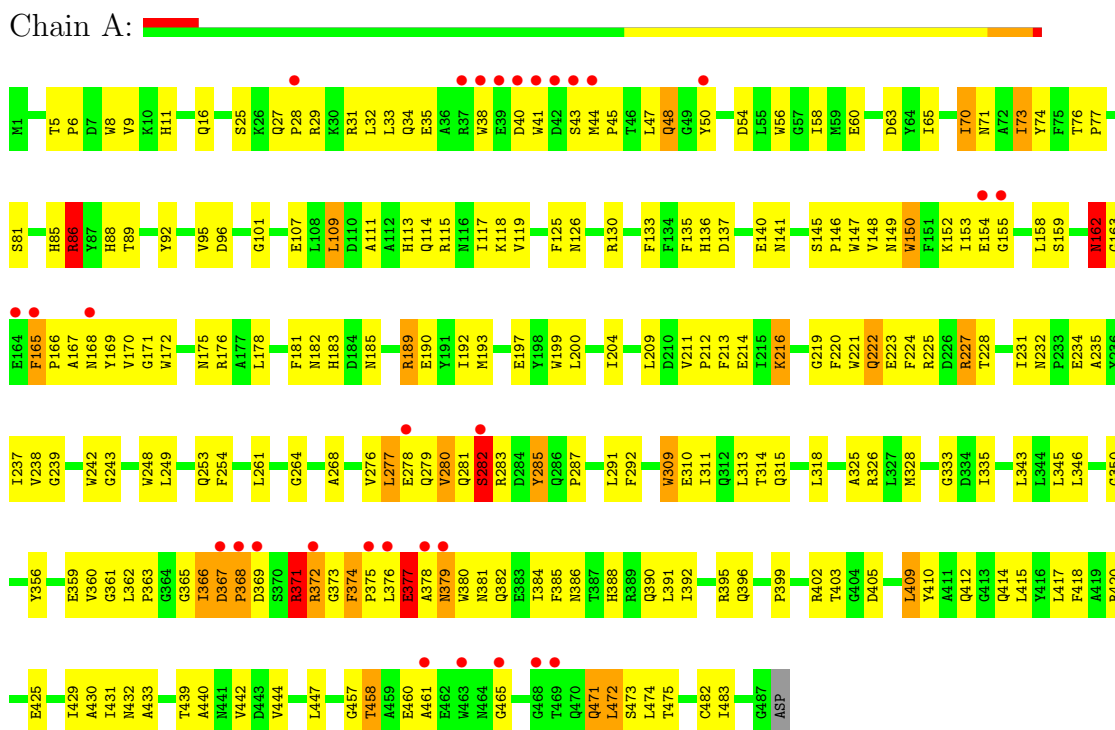
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	46	Total	O	0	0
			46	46		

### 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 AMYLASE, CATALYTIC REGION



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	93.73Å 93.73Å 257.04Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.88 – 2.80 29.88 – 2.80	Depositor EDS
% Data completeness (in resolution range)	94.2 (29.88-2.80) 94.3 (29.88-2.80)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	9.44 (at 2.80Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, $R_{free}$	0.224 , 0.284 0.224 , 0.281	Depositor DCC
$R_{free}$ test set	785 reflections (5.07%)	DCC
Wilson B-factor (Å <sup>2</sup> )	48.2	Xtriage
Anisotropy	0.729	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 40.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 17252 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	3968	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.64% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

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	0.32	0/4042	0.53	0/5515

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	A	2	9

There are no bond length outliers.

There are no bond angle outliers.

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	281	GLN	CA
1	A	379	ASN	CA

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	166	PRO	Peptide
1	A	280	VAL	Peptide
1	A	282	SER	Peptide
1	A	366	ILE	Peptide
1	A	367	ASP	Peptide
1	A	369	ASP	Peptide
1	A	371	ARG	Peptide
1	A	374	PHE	Peptide
1	A	377	GLU	Peptide

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3922	0	3726	243	0
2	A	46	0	0	1	0
All	All	3968	0	3726	243	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 32.

All (243) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:377:GLU:O	1:A:379:ASN:HA	1.45	1.17
1:A:155:GLY:H	1:A:168:ASN:HB3	1.06	1.11
1:A:367:ASP:HB3	1:A:368:PRO:HD3	1.31	1.09
1:A:193:MET:HE2	1:A:227:ARG:HD3	1.40	1.01
1:A:367:ASP:CG	1:A:368:PRO:HB3	1.82	1.00
1:A:362:LEU:HD11	1:A:373:GLY:HA3	1.46	0.96
1:A:367:ASP:OD2	1:A:368:PRO:HB3	1.66	0.95
1:A:343:LEU:HD11	1:A:483:ILE:HD11	1.48	0.94
1:A:360:VAL:HG13	1:A:380:TRP:HA	1.47	0.94
1:A:228:THR:HG21	1:A:237:ILE:HD11	1.48	0.93
1:A:31:ARG:HB2	1:A:31:ARG:NH2	1.83	0.93
1:A:155:GLY:N	1:A:168:ASN:HB3	1.84	0.92
1:A:379:ASN:O	1:A:380:TRP:HD1	1.55	0.88
1:A:361:GLY:N	1:A:384:ILE:HD11	1.91	0.85
1:A:73:ILE:HD11	1:A:119:VAL:HG22	1.60	0.84
1:A:367:ASP:HB3	1:A:368:PRO:CD	2.08	0.83
1:A:31:ARG:HB2	1:A:31:ARG:HH21	1.41	0.83
1:A:76:THR:HB	1:A:77:PRO:HD2	1.59	0.83
1:A:189:ARG:O	1:A:193:MET:HG2	1.79	0.82
1:A:54:ASP:OD1	1:A:56:TRP:HB2	1.79	0.81
1:A:11:HIS:HD2	1:A:402:ARG:HE	1.28	0.81
1:A:461:ALA:HB3	1:A:475:THR:N	1.95	0.80
1:A:65:ILE:HG22	1:A:70:ILE:CG1	2.12	0.80
1:A:366:ILE:CG2	1:A:367:ASP:H	1.95	0.79
1:A:280:VAL:HG12	1:A:282:SER:N	1.99	0.78

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:360:VAL:HG13	1:A:380:TRP:CA	2.15	0.77
1:A:154:GLU:N	1:A:168:ASN:O	2.18	0.76
1:A:366:ILE:HG23	1:A:367:ASP:H	1.50	0.76
1:A:366:ILE:HG23	1:A:367:ASP:N	2.01	0.75
1:A:76:THR:HB	1:A:77:PRO:CD	2.17	0.74
1:A:367:ASP:OD2	1:A:368:PRO:CB	2.36	0.74
1:A:279:GLN:HB3	1:A:325:ALA:CB	2.19	0.73
1:A:360:VAL:CG1	1:A:380:TRP:HA	2.17	0.73
1:A:171:GLY:HA3	1:A:175:ASN:O	1.89	0.73
1:A:377:GLU:C	1:A:379:ASN:HA	2.08	0.73
1:A:27:GLN:HB2	1:A:28:PRO:HD2	1.70	0.72
1:A:32:LEU:HB3	1:A:47:LEU:HD21	1.71	0.72
1:A:159:SER:OG	1:A:162:ASN:HB3	1.90	0.71
1:A:153:ILE:HD11	1:A:158:LEU:HD21	1.73	0.70
1:A:472:LEU:HD23	1:A:473:SER:N	2.06	0.70
1:A:377:GLU:O	1:A:379:ASN:CA	2.32	0.70
1:A:376:LEU:O	1:A:377:GLU:N	2.25	0.69
1:A:163:GLY:HA2	1:A:176:ARG:NH1	2.07	0.69
1:A:5:THR:HB	1:A:314:THR:HG21	1.74	0.69
1:A:461:ALA:O	1:A:474:LEU:HA	1.93	0.69
1:A:276:VAL:HG12	1:A:278:GLU:H	1.57	0.68
1:A:43:SER:O	1:A:45:PRO:HD3	1.94	0.67
1:A:281:GLN:H	1:A:366:ILE:HD11	1.58	0.67
1:A:366:ILE:CG2	1:A:367:ASP:N	2.55	0.67
1:A:172:TRP:HZ3	1:A:214:GLU:HG3	1.60	0.66
1:A:414:GLN:HG3	1:A:433:ALA:HB3	1.77	0.66
1:A:277:LEU:HD12	1:A:277:LEU:O	1.95	0.66
1:A:219:GLY:O	1:A:222:GLN:HG2	1.95	0.66
1:A:193:MET:CE	1:A:224:PHE:HA	2.26	0.66
1:A:279:GLN:HB3	1:A:325:ALA:HB2	1.77	0.66
1:A:197:GLU:HG2	1:A:231:ILE:HD11	1.78	0.65
1:A:193:MET:HE3	1:A:224:PHE:HA	1.79	0.65
1:A:362:LEU:HD23	1:A:363:PRO:O	1.97	0.65
1:A:379:ASN:N	1:A:379:ASN:OD1	2.30	0.65
1:A:360:VAL:O	1:A:360:VAL:CG1	2.46	0.63
1:A:292:PHE:HA	1:A:414:GLN:HE22	1.62	0.62
1:A:429:ILE:HD12	1:A:483:ILE:HG13	1.82	0.62
1:A:310:GLU:HG2	2:A:2034:HOH:O	1.99	0.62
1:A:461:ALA:HB3	1:A:475:THR:H	1.65	0.61
1:A:41:TRP:HA	1:A:44:MET:HE2	1.80	0.61
1:A:16:GLN:HB2	1:A:74:TYR:CZ	2.36	0.61
1:A:167:ALA:HB1	1:A:169:TYR:CE2	2.35	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:379:ASN:O	1:A:380:TRP:CD1	2.46	0.61
1:A:147:TRP:O	1:A:150:TRP:HB2	2.01	0.60
1:A:65:ILE:HG22	1:A:70:ILE:HG13	1.84	0.60
1:A:133:PHE:O	1:A:136:HIS:HB3	2.01	0.60
1:A:54:ASP:O	1:A:58:ILE:HD13	2.01	0.60
1:A:211:VAL:O	1:A:211:VAL:HG12	2.02	0.59
1:A:182:ASN:ND2	1:A:185:ASN:HB2	2.16	0.59
1:A:418:PHE:CE1	1:A:429:ILE:HB	2.36	0.59
1:A:73:ILE:CD1	1:A:119:VAL:HG22	2.30	0.59
1:A:155:GLY:H	1:A:168:ASN:CB	1.99	0.58
1:A:41:TRP:NE1	1:A:48:GLN:HA	2.18	0.58
1:A:460:GLU:CD	1:A:461:ALA:H	2.06	0.58
1:A:165:PHE:N	1:A:165:PHE:CD1	2.69	0.58
1:A:249:LEU:HD13	1:A:311:ILE:HG21	1.86	0.58
1:A:71:ASN:O	1:A:117:ILE:HA	2.03	0.58
1:A:309:TRP:CZ3	1:A:313:LEU:HD11	2.39	0.58
1:A:417:LEU:HG	1:A:430:ALA:HB2	1.87	0.57
1:A:152:LYS:HE3	1:A:181:PHE:O	2.03	0.57
1:A:211:VAL:HG13	1:A:213:PHE:CZ	2.40	0.57
1:A:285:TYR:CD2	1:A:285:TYR:N	2.71	0.57
1:A:367:ASP:OD2	1:A:368:PRO:CA	2.53	0.57
1:A:167:ALA:N	1:A:176:ARG:NH2	2.53	0.57
1:A:277:LEU:CD1	1:A:277:LEU:O	2.53	0.57
1:A:109:LEU:HD22	1:A:113:HIS:CE1	2.40	0.56
1:A:70:ILE:HD12	1:A:70:ILE:C	2.25	0.56
1:A:71:ASN:C	1:A:117:ILE:HG23	2.26	0.56
1:A:384:ILE:HG13	1:A:385:PHE:N	2.20	0.56
1:A:379:ASN:C	1:A:380:TRP:CD1	2.79	0.56
1:A:32:LEU:H	1:A:47:LEU:HD23	1.70	0.56
1:A:382:GLN:O	1:A:386:ASN:ND2	2.39	0.56
1:A:280:VAL:HG12	1:A:282:SER:HB3	1.88	0.56
1:A:153:ILE:HG23	1:A:153:ILE:O	2.06	0.56
1:A:211:VAL:HG13	1:A:213:PHE:CE1	2.42	0.55
1:A:211:VAL:CG1	1:A:214:GLU:HG2	2.37	0.55
1:A:172:TRP:CZ3	1:A:214:GLU:HG3	2.40	0.55
1:A:165:PHE:N	1:A:165:PHE:HD1	2.05	0.55
1:A:149:ASN:O	1:A:185:ASN:ND2	2.40	0.55
1:A:367:ASP:CB	1:A:368:PRO:CD	2.82	0.54
1:A:461:ALA:O	1:A:474:LEU:HD23	2.07	0.54
1:A:356:TYR:CD1	1:A:375:PRO:HG3	2.42	0.54
1:A:248:TRP:HA	1:A:253:GLN:NE2	2.23	0.54
1:A:439:THR:HA	1:A:474:LEU:O	2.07	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:283:ARG:O	1:A:283:ARG:HD2	2.07	0.54
1:A:367:ASP:OD2	1:A:368:PRO:HA	2.08	0.54
1:A:211:VAL:HG12	1:A:214:GLU:HG2	1.89	0.54
1:A:360:VAL:HG13	1:A:380:TRP:CB	2.37	0.53
1:A:377:GLU:CD	1:A:377:GLU:N	2.62	0.53
1:A:379:ASN:C	1:A:380:TRP:HD1	2.09	0.53
1:A:280:VAL:HG12	1:A:282:SER:CB	2.39	0.53
1:A:345:LEU:HD13	1:A:346:LEU:HD23	1.91	0.53
1:A:25:SER:HB2	1:A:60:GLU:OE2	2.09	0.53
1:A:73:ILE:HD13	1:A:73:ILE:H	1.74	0.53
1:A:417:LEU:HD21	1:A:474:LEU:HD12	1.91	0.52
1:A:209:LEU:O	1:A:212:PRO:HD3	2.09	0.52
1:A:461:ALA:CB	1:A:475:THR:HB	2.39	0.52
1:A:65:ILE:CG2	1:A:70:ILE:HD11	2.39	0.52
1:A:405:ASP:O	1:A:420:ARG:HA	2.09	0.52
1:A:367:ASP:CG	1:A:368:PRO:CB	2.66	0.52
1:A:56:TRP:HZ3	1:A:107:GLU:OE2	1.92	0.52
1:A:309:TRP:CE3	1:A:313:LEU:HD11	2.44	0.52
1:A:346:LEU:HD12	1:A:391:LEU:CB	2.38	0.52
1:A:280:VAL:HG12	1:A:282:SER:CA	2.39	0.52
1:A:58:ILE:H	1:A:58:ILE:CD1	2.22	0.52
1:A:457:GLY:HA3	1:A:482:CYS:HA	1.90	0.52
1:A:472:LEU:HD23	1:A:473:SER:H	1.71	0.51
1:A:6:PRO:HB2	1:A:9:VAL:HG23	1.92	0.51
1:A:234:GLU:OE1	1:A:234:GLU:HA	2.10	0.51
1:A:335:ILE:HD12	1:A:335:ILE:N	2.26	0.51
1:A:285:TYR:N	1:A:285:TYR:HD2	2.08	0.51
1:A:356:TYR:HD1	1:A:375:PRO:HG3	1.74	0.51
1:A:197:GLU:HG2	1:A:231:ILE:CD1	2.41	0.51
1:A:343:LEU:HD11	1:A:483:ILE:CD1	2.32	0.51
1:A:372:ARG:HG2	1:A:373:GLY:N	2.24	0.50
1:A:58:ILE:H	1:A:58:ILE:HD12	1.77	0.50
1:A:162:ASN:C	1:A:162:ASN:HD22	2.14	0.50
1:A:197:GLU:HA	1:A:231:ILE:HD12	1.94	0.50
1:A:223:GLU:HG2	1:A:227:ARG:HD2	1.94	0.50
1:A:65:ILE:HG22	1:A:70:ILE:CD1	2.41	0.50
1:A:65:ILE:HG22	1:A:70:ILE:HD11	1.93	0.50
1:A:292:PHE:HA	1:A:414:GLN:NE2	2.27	0.49
1:A:285:TYR:H	1:A:285:TYR:HD2	1.60	0.49
1:A:58:ILE:N	1:A:58:ILE:HD12	2.27	0.48
1:A:346:LEU:HD12	1:A:391:LEU:HB2	1.94	0.48
1:A:178:LEU:N	1:A:178:LEU:HD22	2.28	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:199:TRP:O	1:A:204:ILE:HG12	2.12	0.48
1:A:410:TYR:CZ	1:A:412:GLN:HB2	2.48	0.48
1:A:309:TRP:HH2	1:A:403:THR:O	1.97	0.48
1:A:181:PHE:HB2	1:A:183:HIS:CE1	2.49	0.48
1:A:200:LEU:HD11	1:A:228:THR:HG23	1.96	0.47
1:A:31:ARG:CZ	1:A:31:ARG:HB2	2.44	0.47
1:A:279:GLN:NE2	1:A:365:GLY:HA2	2.29	0.47
1:A:209:LEU:HD12	1:A:254:PHE:CE1	2.50	0.47
1:A:264:GLY:O	1:A:287:PRO:HA	2.14	0.47
1:A:81:SER:OG	1:A:86:ARG:HA	2.14	0.47
1:A:197:GLU:CG	1:A:231:ILE:HD11	2.44	0.47
1:A:378:ALA:HA	1:A:380:TRP:CD1	2.50	0.47
1:A:381:ASN:HB3	1:A:384:ILE:HG12	1.97	0.47
1:A:444:VAL:HG23	1:A:444:VAL:O	2.15	0.47
1:A:415:LEU:HD23	1:A:440:ALA:HB2	1.97	0.47
1:A:399:PRO:HD2	1:A:425:GLU:OE1	2.15	0.47
1:A:76:THR:CB	1:A:77:PRO:CD	2.83	0.46
1:A:360:VAL:O	1:A:360:VAL:HG12	2.14	0.46
1:A:130:ARG:HA	1:A:135:PHE:CD2	2.51	0.46
1:A:126:ASN:HA	1:A:183:HIS:CE1	2.51	0.46
1:A:268:ALA:HB3	1:A:291:LEU:HD12	1.96	0.46
1:A:85:HIS:HB3	1:A:88:HIS:CD2	2.50	0.46
1:A:318:LEU:HD23	1:A:318:LEU:N	2.31	0.46
1:A:141:ASN:HB2	1:A:145:SER:HB2	1.97	0.45
1:A:34:GLN:HG3	1:A:35:GLU:HG2	1.98	0.45
1:A:111:ALA:O	1:A:114:GLN:HG2	2.16	0.45
1:A:362:LEU:HD13	1:A:375:PRO:HB3	1.98	0.45
1:A:371:ARG:HG2	1:A:371:ARG:H	1.38	0.45
1:A:190:GLU:OE2	1:A:227:ARG:NH2	2.49	0.45
1:A:11:HIS:CD2	1:A:402:ARG:HE	2.19	0.45
1:A:86:ARG:NH2	1:A:96:ASP:OD2	2.50	0.45
1:A:154:GLU:HB2	1:A:168:ASN:HB3	1.97	0.45
1:A:5:THR:HG21	1:A:311:ILE:HD13	1.98	0.45
1:A:350:GLY:O	1:A:395:ARG:NH2	2.49	0.45
1:A:167:ALA:H	1:A:176:ARG:NH2	2.15	0.45
1:A:216:LYS:HE2	1:A:216:LYS:HB3	1.86	0.45
1:A:359:GLU:O	1:A:385:PHE:HA	2.16	0.45
1:A:328:MET:HG2	1:A:361:GLY:O	2.16	0.44
1:A:16:GLN:NE2	1:A:74:TYR:OH	2.49	0.44
1:A:209:LEU:HB2	1:A:239:GLY:HA2	1.99	0.44
1:A:209:LEU:HD13	1:A:221:TRP:CZ3	2.53	0.44
1:A:130:ARG:HA	1:A:135:PHE:HD2	1.83	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:415:LEU:HD12	1:A:431:ILE:O	2.17	0.44
1:A:243:GLY:HA2	1:A:261:LEU:HD21	1.97	0.44
1:A:225:ARG:HD2	1:A:253:GLN:O	2.18	0.44
1:A:386:ASN:O	1:A:390:GLN:HB2	2.18	0.44
1:A:95:VAL:HB	1:A:101:GLY:C	2.38	0.44
1:A:65:ILE:HG22	1:A:70:ILE:HG12	1.96	0.44
1:A:376:LEU:O	1:A:380:TRP:NE1	2.51	0.43
1:A:172:TRP:CZ3	1:A:211:VAL:HG11	2.53	0.43
1:A:211:VAL:HG12	1:A:214:GLU:CG	2.48	0.43
1:A:346:LEU:HD12	1:A:391:LEU:HB3	2.00	0.43
1:A:125:PHE:HB3	1:A:192:ILE:HD13	2.00	0.43
1:A:360:VAL:O	1:A:360:VAL:HG13	2.19	0.43
1:A:282:SER:O	1:A:282:SER:OG	2.30	0.43
1:A:197:GLU:HA	1:A:231:ILE:CD1	2.48	0.43
1:A:63:ASP:OD2	1:A:115:ARG:NH2	2.48	0.43
1:A:461:ALA:O	1:A:473:SER:O	2.37	0.43
1:A:73:ILE:HD13	1:A:118:LYS:O	2.18	0.43
1:A:311:ILE:O	1:A:315:GLN:HG2	2.19	0.43
1:A:154:GLU:HB2	1:A:168:ASN:CA	2.49	0.43
1:A:361:GLY:H	1:A:384:ILE:HD11	1.78	0.43
1:A:465:GLY:HA3	1:A:471:GLN:OE1	2.19	0.43
1:A:472:LEU:CD2	1:A:473:SER:N	2.78	0.42
1:A:461:ALA:HB2	1:A:475:THR:HB	2.01	0.42
1:A:458:THR:HG23	1:A:458:THR:O	2.18	0.42
1:A:33:LEU:H	1:A:33:LEU:HD22	1.84	0.42
1:A:409:LEU:HB3	1:A:442:VAL:HG11	2.02	0.42
1:A:65:ILE:CG2	1:A:70:ILE:CG1	2.93	0.42
1:A:71:ASN:O	1:A:117:ILE:HG23	2.20	0.42
1:A:333:GLY:O	1:A:335:ILE:HD12	2.20	0.42
1:A:326:ARG:CZ	1:A:362:LEU:HD22	2.50	0.42
1:A:356:TYR:O	1:A:356:TYR:CG	2.73	0.41
1:A:152:LYS:HB2	1:A:170:VAL:HB	2.02	0.41
1:A:89:THR:HG21	1:A:92:TYR:CE2	2.55	0.41
1:A:211:VAL:CG1	1:A:214:GLU:CG	2.98	0.41
1:A:232:ASN:HB3	1:A:235:ALA:HB2	2.01	0.41
1:A:417:LEU:HG	1:A:430:ALA:CB	2.50	0.41
1:A:420:ARG:HH11	1:A:420:ARG:HG3	1.85	0.41
1:A:137:ASP:OD2	1:A:147:TRP:HZ3	2.03	0.41
1:A:392:ILE:O	1:A:396:GLN:HG2	2.21	0.41
1:A:359:GLU:HA	1:A:388:HIS:CD2	2.55	0.41
1:A:193:MET:HB3	1:A:227:ARG:NH1	2.36	0.41
1:A:8:TRP:CZ3	1:A:118:LYS:HG3	2.56	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:461:ALA:HB3	1:A:475:THR:HB	2.03	0.41
1:A:31:ARG:NE	1:A:50:TYR:CD2	2.89	0.40
1:A:5:THR:CB	1:A:314:THR:HG21	2.46	0.40
1:A:65:ILE:H	1:A:65:ILE:HG12	1.55	0.40
1:A:148:VAL:HG12	1:A:153:ILE:HG21	2.03	0.40
1:A:145:SER:HA	1:A:146:PRO:HD3	1.82	0.40
1:A:366:ILE:HG22	1:A:367:ASP:H	1.79	0.40
1:A:40:ASP:O	1:A:43:SER:HB2	2.22	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	482/488 (99%)	415 (86%)	61 (13%)	6 (1%)	19	54

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	162	ASN
1	A	374	PHE
1	A	368	PRO
1	A	86	ARG
1	A	220	PHE
1	A	458	THR

### 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	406/409 (99%)	376 (93%)	30 (7%)	20	48

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	ARG
1	A	38	TRP
1	A	48	GLN
1	A	70	ILE
1	A	73	ILE
1	A	86	ARG
1	A	109	LEU
1	A	140	GLU
1	A	150	TRP
1	A	162	ASN
1	A	165	PHE
1	A	189	ARG
1	A	216	LYS
1	A	222	GLN
1	A	227	ARG
1	A	238	VAL
1	A	242	TRP
1	A	277	LEU
1	A	282	SER
1	A	285	TYR
1	A	309	TRP
1	A	371	ARG
1	A	372	ARG
1	A	377	GLU
1	A	379	ASN
1	A	409	LEU
1	A	432	ASN
1	A	447	LEU
1	A	471	GLN
1	A	472	LEU

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	2	GLN
1	A	11	HIS
1	A	16	GLN

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Mol	Chain	Res	Type
1	A	48	GLN
1	A	80	GLN
1	A	88	HIS
1	A	114	GLN
1	A	126	ASN
1	A	127	HIS
1	A	141	ASN
1	A	162	ASN
1	A	182	ASN
1	A	222	GLN
1	A	247	GLN
1	A	386	ASN
1	A	388	HIS
1	A	412	GLN
1	A	414	GLN
1	A	432	ASN
1	A	470	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no ligands in this entry.

### 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	487/488 (99%)	0.11	30 (6%) 20 19	24, 49, 105, 133	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	376	LEU	30.9
1	A	42	ASP	4.6
1	A	461	ALA	4.2
1	A	155	GLY	4.1
1	A	50	TYR	3.8
1	A	463	TRP	3.7
1	A	38	TRP	3.5
1	A	43	SER	3.1
1	A	465	GLY	3.0
1	A	368	PRO	2.8
1	A	379	ASN	2.8
1	A	278	GLU	2.6
1	A	375	PRO	2.6
1	A	165	PHE	2.6
1	A	44	MET	2.6
1	A	37	ARG	2.6
1	A	41	TRP	2.5
1	A	282	SER	2.4
1	A	168	ASN	2.4
1	A	378	ALA	2.4
1	A	39	GLU	2.3
1	A	468	GLY	2.3
1	A	40	ASP	2.2
1	A	367	ASP	2.2
1	A	469	THR	2.2
1	A	372	ARG	2.1
1	A	369	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	154	GLU	2.1
1	A	28	PRO	2.0
1	A	164	GLU	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

There are no ligands in this entry.

## 6.5 Other polymers ⓘ

There are no such residues in this entry.