Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 06:13 AM GMT

PDB ID	:	2WCS
Title	:	CRYSTAL STRUCTURE OF DEBRANCHING ENZYME FROM NOSTOC PUNCTIFORME (NPDE)
Authors	:	Dumbrepatil, A.B.; Choi, J.H.; Nam, S.H.; Park, K.H.; Woo, E.J.
Deposited on	:	2009-03-16
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:

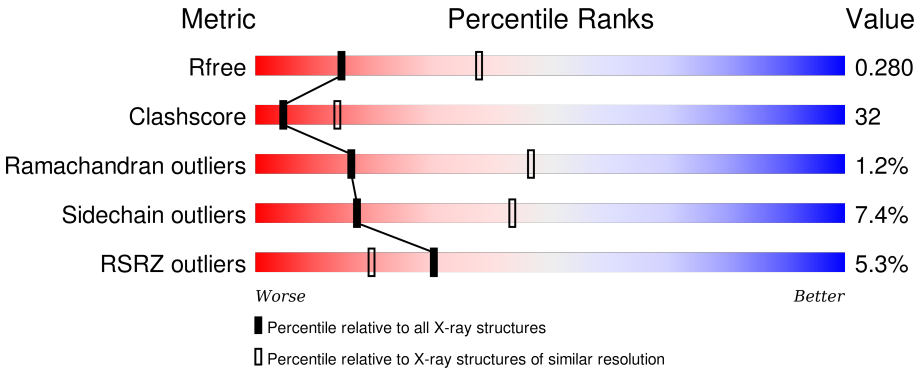
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)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

1 Overall quality at a glance (i)

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	488	5% 53% 40% 5% •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 3968 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 AMYLASE, CATALYTIC REGION.

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

- Molecule 2 is water.

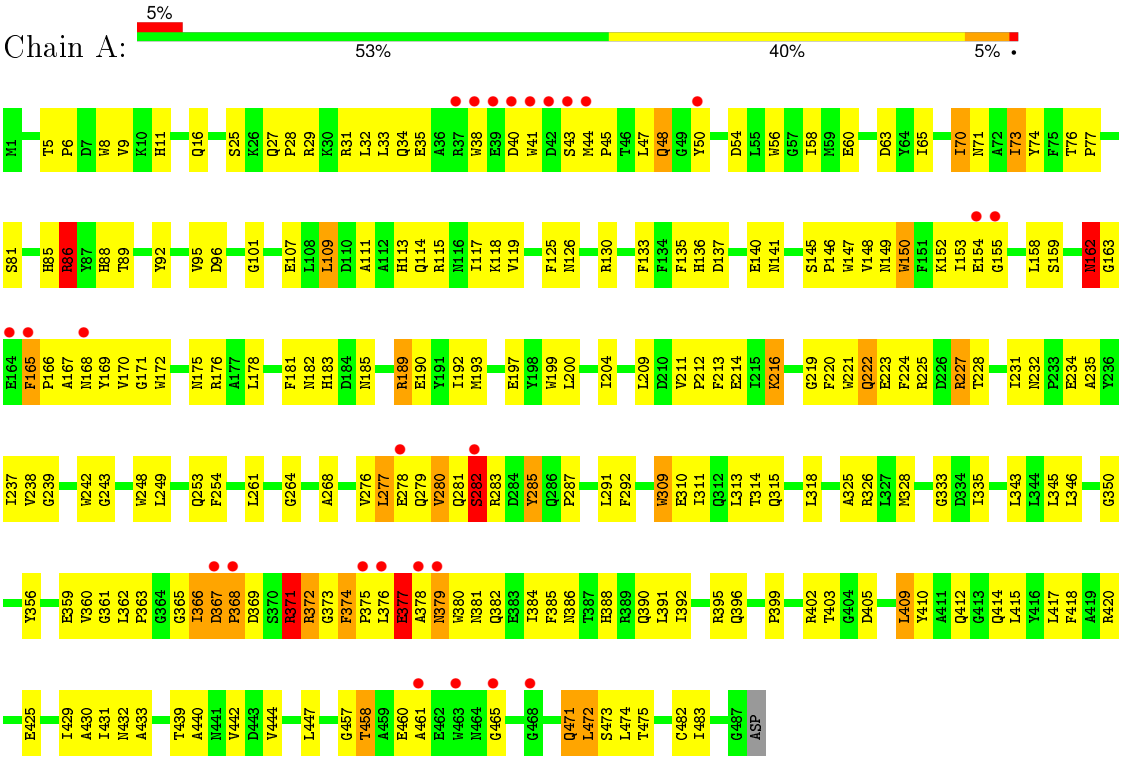
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	46	46	46	0	0

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: ALPHA AMYLASE, CATALYTIC REGION

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α, β, γ	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
$< I/\sigma(I) >$ [1]	9.44 (at 2.80Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.224 , 0.284 0.224 , 0.280	Depositor DCC
R_{free} test set	785 reflections (5.07%)	DCC
Wilson B-factor (Å^2)	48.2	Xtriage
Anisotropy	0.729	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 62.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	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^2)	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.*

[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

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 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	3922	0	3726	243	0
2	A	46	0	0	1	0
All	All	3968	0	3726	243	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 32.

All (243) 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: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	Interatomic distance (Å)	Clash overlap (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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:167:ALA:HB1	1:A:169:TYR:CE2	2.35	0.61
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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 (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	482/488 (99%)	415 (86%)	61 (13%)	6 (1%)	16 47

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 (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	406/409 (99%)	376 (93%)	30 (7%)	17 43

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
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 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

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^{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	487/488 (99%)	0.11	26 (5%) 30 20	24, 49, 105, 133	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	376	LEU	30.6
1	A	42	ASP	4.7
1	A	155	GLY	4.4
1	A	461	ALA	4.1
1	A	50	TYR	3.9
1	A	38	TRP	3.3
1	A	463	TRP	3.3
1	A	43	SER	3.1
1	A	379	ASN	2.9
1	A	465	GLY	2.7
1	A	368	PRO	2.6
1	A	282	SER	2.6
1	A	278	GLU	2.6
1	A	41	TRP	2.6
1	A	44	MET	2.6
1	A	168	ASN	2.6
1	A	165	PHE	2.6
1	A	39	GLU	2.6
1	A	375	PRO	2.5
1	A	40	ASP	2.4
1	A	37	ARG	2.3
1	A	367	ASP	2.2
1	A	378	ALA	2.2
1	A	468	GLY	2.1
1	A	154	GLU	2.1
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.