



Full wwPDB X-ray Structure Validation Report

May 30, 2014 – 01:46 AM EDT

PDB ID : 4C5U
Title : Structural Investigations into the Stereochemistry and Activity of a Phenylalanine-2,3-Aminomutase from *Taxus chinensis*
Authors : Wybenga, G.G.; Szymanski, W.; Wu, B.; Feringa, B.L.; Janssen, D.B.; Dijkstra, B.W.
Deposited on : 2013-09-16
Resolution : 2.19 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

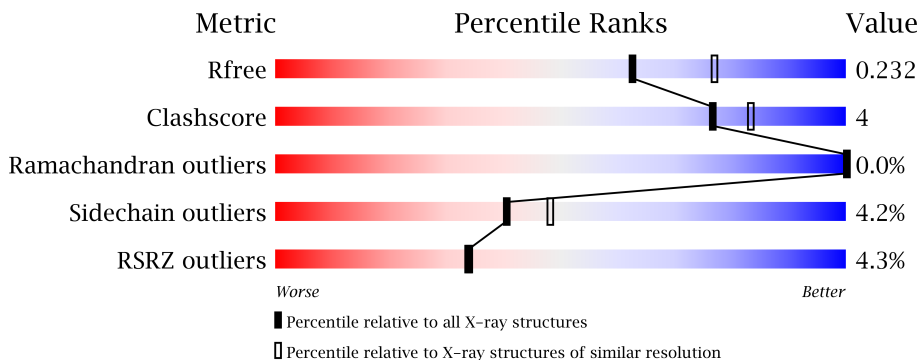
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.16 November 2013
Xtriage (Phenix) : dev-1439
EDS : stable23161
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) : stable23161

1 Overall quality at a glance

The reported resolution of this entry is 2.19 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	2938 (2.20-2.20)
Clashscore	79885	3751 (2.20-2.20)
Ramachandran outliers	78287	3681 (2.20-2.20)
Sidechain outliers	78261	3682 (2.20-2.20)
RSRZ outliers	66119	2939 (2.20-2.20)

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. 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	707	
1	B	707	
1	C	707	
1	D	707	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 19779 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 PHENYLALANINE AMMONIA-LYASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	641	Total	C	N	O	S	0	0	0
			4962	3147	855	937	23			
1	B	636	Total	C	N	O	S	0	0	0
			4922	3120	847	932	23			
1	C	639	Total	C	N	O	S	0	0	0
			4941	3133	853	932	23			
1	D	636	Total	C	N	O	S	0	0	0
			4917	3121	848	925	23			

There are 4 discrepancies between the modelled and reference sequences:

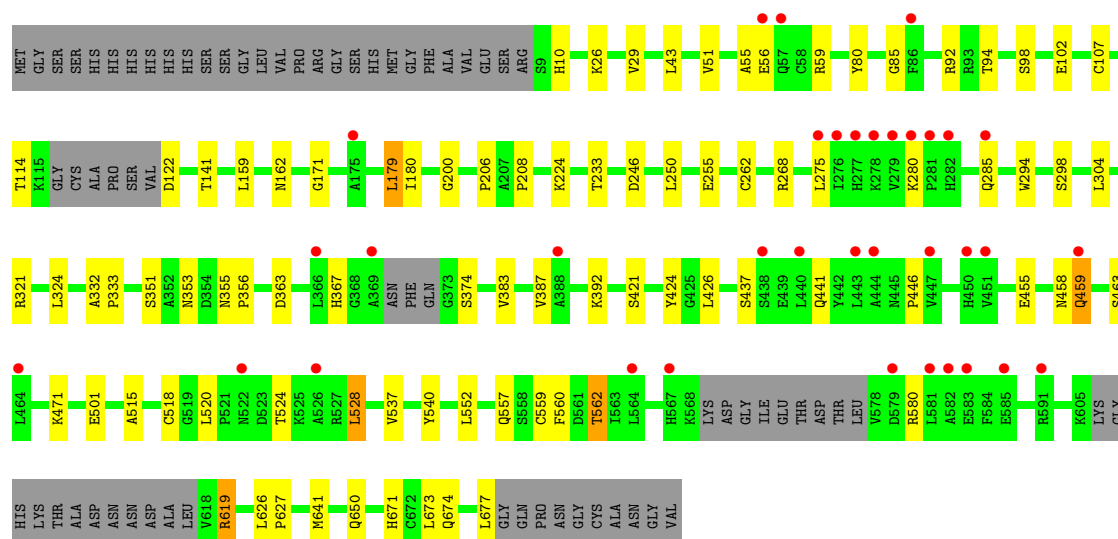
Chain	Residue	Modelled	Actual	Comment	Reference
A	322	ALA	TYR	ENGINEERED MUTATION	UNP Q68G84
B	322	ALA	TYR	ENGINEERED MUTATION	UNP Q68G84
C	322	ALA	TYR	ENGINEERED MUTATION	UNP Q68G84
D	322	ALA	TYR	ENGINEERED MUTATION	UNP Q68G84

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	15	Total	O	0	0
			15	15		
2	B	6	Total	O	0	0
			6	6		
2	C	10	Total	O	0	0
			10	10		
2	D	6	Total	O	0	0
			6	6		

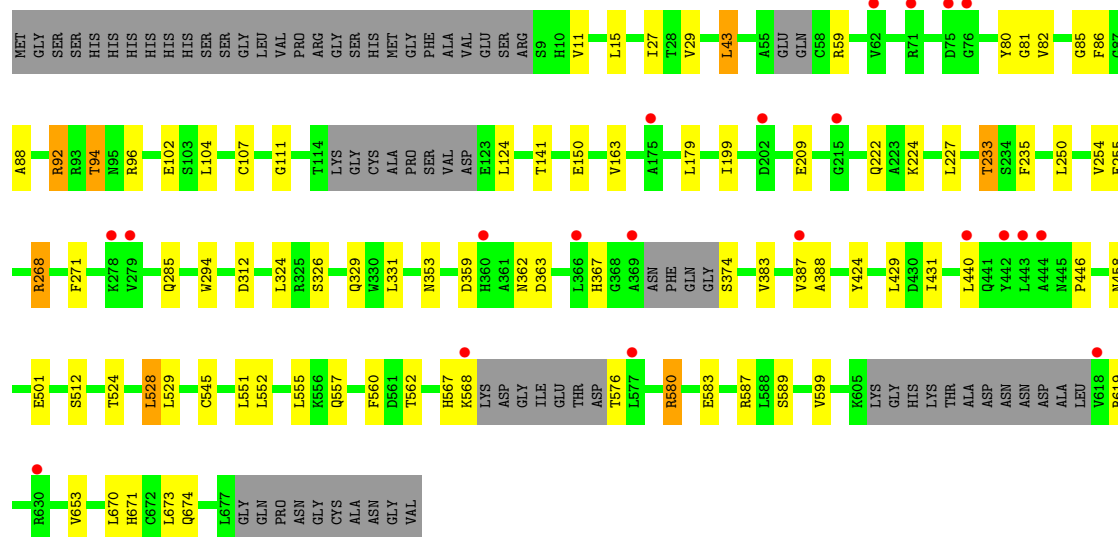
• Molecule 1: PHENYLALANINE AMMONIA-LYASE

Chain C:



• Molecule 1: PHENYLALANINE AMMONIA-LYASE

Chain D:



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	99.48Å 147.27Å 99.77Å 90.00° 99.80° 90.00°	Depositor
Resolution (Å)	48.38 – 2.19 48.38 – 2.19	Depositor EDS
% Data completeness (in resolution range)	94.3 (48.38-2.19) 98.1 (48.38-2.19)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.83 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.189 , 0.224 0.194 , 0.232	Depositor DCC
R_{free} test set	7176 reflections (5.33%)	DCC
Wilson B-factor (Å ²)	37.4	Xtriage
Anisotropy	0.069	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 26.2	EDS
Estimated twinning fraction	0.865 for H, K, L 0.135 for L, -K, H 0.000 for l,-k,h	Xtriage
Reported twinning fraction	0.865 for H, K, L 0.135 for L, -K, H	Depositor
L-test for twinning	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Outliers	1 of 141777 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	19779	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

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

¹Intensities 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	0.57	1/5045 (0.0%)	0.69	1/6843 (0.0%)
1	B	0.54	2/5002 (0.0%)	0.65	1/6782 (0.0%)
1	C	0.55	1/5024 (0.0%)	0.67	1/6813 (0.0%)
1	D	0.55	1/4999 (0.0%)	0.68	0/6780
All	All	0.56	5/20070 (0.0%)	0.67	3/27218 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	294	TRP	CD2-CE2	5.71	1.48	1.41
1	B	68	TRP	CD2-CE2	5.47	1.48	1.41
1	B	294	TRP	CD2-CE2	5.26	1.47	1.41
1	A	68	TRP	CD2-CE2	5.07	1.47	1.41
1	D	294	TRP	CD2-CE2	5.06	1.47	1.41

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	619	ARG	NE-CZ-NH2	-5.38	117.61	120.30
1	B	619	ARG	NE-CZ-NH2	-5.38	117.61	120.30
1	C	321	ARG	NE-CZ-NH2	-5.37	117.62	120.30

There are no chirality outliers.

There are no planarity outliers.

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	4962	0	5044	50	0
1	B	4922	0	4993	40	0
1	C	4941	0	5021	45	0
1	D	4917	0	5004	40	0
2	A	15	0	0	0	0
2	B	6	0	0	0	0
2	C	10	0	0	0	0
2	D	6	0	0	0	0
All	All	19779	0	20062	151	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 4.

All (151) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:29:VAL:HG12	1:A:141:THR:HG21	1.58	0.85
1:D:580:ARG:HH11	1:D:580:ARG:CG	2.01	0.74
1:A:80:TYR:HB3	1:A:367:HIS:HD2	1.52	0.73
1:B:671:HIS:O	1:B:674:GLN:HG2	1.89	0.72
1:D:29:VAL:HG12	1:D:141:THR:HG21	1.71	0.71
1:B:446:PRO:HD3	1:D:446:PRO:HD3	1.72	0.71
1:A:557:GLN:NE2	1:D:557:GLN:OE1	2.21	0.71
1:C:501:GLU:OE1	1:C:619:ARG:HD2	1.91	0.70
1:B:557:GLN:NE2	1:C:557:GLN:OE1	2.23	0.69
1:A:650:GLN:HG3	1:B:111:GLY:O	1.93	0.68
1:A:80:TYR:HB3	1:A:367:HIS:CD2	2.30	0.66
1:C:224:LYS:HE2	1:C:356:PRO:HD2	1.78	0.65
1:A:43:LEU:HD22	1:A:134:LEU:HD22	1.79	0.64
1:B:224:LYS:HE2	1:B:356:PRO:HD2	1.80	0.63
1:C:515:ALA:HA	1:C:520:LEU:HD12	1.82	0.61
1:D:163:VAL:HG22	1:D:199:ILE:HG12	1.83	0.61
1:B:29:VAL:HG12	1:B:141:THR:HG21	1.82	0.60
1:C:304:LEU:HD11	1:C:619:ARG:HD3	1.84	0.60
1:B:233:THR:HG21	1:B:373:GLY:N	2.15	0.60
1:C:437:SER:O	1:C:441:GLN:HG2	2.02	0.59
1:B:123:GLU:OE1	1:B:166:LYS:HE3	2.05	0.57
1:D:580:ARG:HH11	1:D:580:ARG:HG3	1.68	0.57
1:C:524:THR:HG22	1:C:528:LEU:HD22	1.87	0.56
1:C:29:VAL:HG12	1:C:141:THR:HG21	1.86	0.56
1:C:455:GLU:HB3	1:C:459:GLN:HG3	1.88	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:157:LYS:HD2	1:A:216:LEU:HD21	1.88	0.55
1:C:171:GLY:O	1:C:180:ILE:HD12	2.06	0.55
1:A:374:SER:HA	1:A:463:SER:HB3	1.89	0.54
1:D:82:VAL:HG22	1:D:224:LYS:HB2	1.87	0.54
1:A:501:GLU:OE1	1:A:619:ARG:HD2	2.06	0.54
1:A:446:PRO:HD3	1:C:446:PRO:HD3	1.90	0.54
1:D:671:HIS:O	1:D:674:GLN:HG2	2.08	0.54
1:D:501:GLU:OE1	1:D:619:ARG:HD2	2.08	0.53
1:A:255:GLU:HG2	1:A:331:LEU:HD13	1.91	0.52
1:A:351:SER:HB2	1:C:280:LYS:HA	1.91	0.52
1:A:88:ALA:HB2	1:A:458:ASN:HB2	1.90	0.52
1:A:424:TYR:OH	1:B:85:GLY:HA2	2.10	0.52
1:A:81:GLY:HA3	1:A:227:LEU:HD22	1.92	0.52
1:A:551:LEU:O	1:A:555:LEU:HG	2.10	0.52
1:D:233:THR:HG23	1:D:233:THR:O	2.10	0.51
1:A:321:ARG:HG2	1:C:458:ASN:HA	1.93	0.50
1:C:179:LEU:HD23	1:D:431:ILE:HD13	1.94	0.50
1:A:29:VAL:CG1	1:A:141:THR:HG21	2.36	0.50
1:B:334:LEU:HD22	1:B:386:ALA:HA	1.93	0.50
1:B:327:SER:OG	1:B:328:PRO:HD3	2.12	0.50
1:D:94:THR:HB	1:D:96:ARG:H	1.77	0.50
1:B:265:ILE:HG21	1:B:324:LEU:HD12	1.94	0.49
1:B:388:ALA:HA	1:B:440:LEU:HG	1.94	0.49
1:C:114:THR:OG1	1:C:122:ASP:HB2	2.12	0.49
1:C:80:TYR:HB3	1:C:367:HIS:CD2	2.48	0.49
1:B:15:LEU:HD21	1:B:673:LEU:HD23	1.95	0.49
1:C:285:GLN:HB2	1:C:332:ALA:HB2	1.95	0.49
1:A:332:ALA:HB3	1:A:333:PRO:CD	2.42	0.49
1:C:374:SER:HA	1:C:463:SER:HB2	1.95	0.49
1:B:437:SER:O	1:B:441:GLN:HG2	2.13	0.48
1:D:88:ALA:HB2	1:D:458:ASN:HB2	1.94	0.48
1:C:10:HIS:NE2	1:C:255:GLU:OE1	2.47	0.48
1:A:545:CYS:SG	1:A:589:SER:HA	2.54	0.48
1:A:334:LEU:HD22	1:A:386:ALA:HA	1.95	0.48
1:C:671:HIS:O	1:C:674:GLN:HG2	2.14	0.48
1:C:332:ALA:HB3	1:C:333:PRO:HD3	1.94	0.48
1:B:27:ILE:CD1	1:B:43:LEU:HB2	2.43	0.48
1:B:329:GLN:O	1:D:374:SER:OG	2.27	0.48
1:A:437:SER:O	1:A:441:GLN:HG2	2.13	0.47
1:D:27:ILE:CD1	1:D:43:LEU:HB2	2.44	0.47
1:A:267:GLY:HA3	1:A:324:LEU:HD11	1.96	0.47
1:A:560:PHE:HB2	1:D:560:PHE:HB2	1.96	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:424:TYR:OH	1:D:85:GLY:HA2	2.14	0.47
1:A:579:ASP:OD1	1:D:567:HIS:NE2	2.42	0.47
1:B:501:GLU:OE1	1:B:619:ARG:HD2	2.14	0.47
1:D:429:LEU:HD21	1:D:653:VAL:HG21	1.96	0.47
1:B:131:SER:HB2	1:B:238:ALA:HB1	1.97	0.47
1:A:388:ALA:HA	1:A:440:LEU:HG	1.97	0.47
1:C:383:VAL:O	1:C:387:VAL:HG23	2.14	0.46
1:A:567:HIS:O	1:A:568:LYS:HD3	2.15	0.46
1:B:24:VAL:HG21	1:B:39:HIS:HD2	1.80	0.46
1:A:94:THR:HG23	1:B:421:SER:OG	2.16	0.46
1:D:81:GLY:HA3	1:D:227:LEU:HD22	1.98	0.46
1:B:59:ARG:HD2	1:B:63:GLU:OE1	2.16	0.46
1:A:419:ASP:OD2	1:B:96:ARG:NH1	2.47	0.46
1:D:86:PHE:CD2	1:D:104:LEU:HD13	2.51	0.46
1:A:596:MET:O	1:A:599:VAL:HG12	2.16	0.46
1:C:304:LEU:CD1	1:C:619:ARG:HD3	2.45	0.46
1:A:27:ILE:CD1	1:A:43:LEU:HB2	2.47	0.45
1:A:181:PRO:HB2	1:A:235:PHE:CD2	2.51	0.45
1:C:518:CYS:O	1:C:580:ARG:HD3	2.16	0.45
1:D:268:ARG:O	1:D:271:PHE:HD1	1.99	0.45
1:B:181:PRO:HB2	1:B:235:PHE:CD2	2.51	0.45
1:B:369:ALA:CB	1:D:326:SER:HB3	2.46	0.45
1:A:500:GLU:HG3	1:A:537:VAL:HG12	1.97	0.45
1:C:246:ASP:O	1:C:250:LEU:HG	2.16	0.45
1:C:559:CYS:O	1:C:562:THR:HG22	2.16	0.45
1:D:359:ASP:OD2	1:D:362:ASN:HB2	2.17	0.45
1:A:80:TYR:CB	1:A:367:HIS:CD2	3.00	0.45
1:D:285:GLN:CD	1:D:329:GLN:HG3	2.38	0.45
1:A:163:VAL:HG22	1:A:199:ILE:HG23	1.99	0.44
1:A:224:LYS:O	1:A:224:LYS:HG2	2.18	0.44
1:A:374:SER:HA	1:A:463:SER:CB	2.47	0.44
1:B:560:PHE:CE1	1:C:559:CYS:HB3	2.52	0.44
1:D:383:VAL:O	1:D:387:VAL:HG23	2.17	0.44
1:C:332:ALA:HB3	1:C:333:PRO:CD	2.47	0.44
1:C:51:VAL:HG21	1:C:159:LEU:HD13	2.00	0.44
1:D:11:VAL:O	1:D:15:LEU:HG	2.18	0.44
1:B:560:PHE:HB2	1:C:560:PHE:HB2	2.00	0.44
1:C:537:VAL:HA	1:C:540:TYR:CE2	2.53	0.43
1:C:421:SER:HB3	1:D:92:ARG:NH1	2.33	0.43
1:B:27:ILE:HD12	1:B:43:LEU:HB2	2.01	0.43
1:A:111:GLY:O	1:B:650:GLN:HG3	2.17	0.43
1:B:268:ARG:O	1:B:271:PHE:HD1	2.01	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:206:PRO:HB2	1:C:208:PRO:HD2	1.99	0.43
1:B:309:TYR:O	1:B:313:LYS:HG3	2.19	0.43
1:D:524:THR:HG22	1:D:528:LEU:HD22	2.01	0.43
1:C:426:LEU:HD11	1:C:641:MET:SD	2.58	0.43
1:A:131:SER:CB	1:A:238:ALA:HB1	2.49	0.43
1:D:388:ALA:HA	1:D:440:LEU:HG	2.01	0.43
1:D:512:SER:HB2	1:D:529:LEU:HD21	2.00	0.43
1:A:82:VAL:HG22	1:A:224:LYS:HB2	2.01	0.43
1:D:545:CYS:SG	1:D:589:SER:HA	2.58	0.43
1:A:80:TYR:CB	1:A:367:HIS:HD2	2.27	0.42
1:B:626:LEU:HB3	1:B:627:PRO:HD3	2.01	0.42
1:C:355:ASN:OD1	1:C:356:PRO:HA	2.19	0.42
1:C:55:ALA:O	1:C:59:ARG:HB3	2.20	0.42
1:A:233:THR:O	1:A:233:THR:HG23	2.20	0.42
1:B:307:GLU:O	1:B:311:ILE:HG12	2.19	0.42
1:A:279:VAL:O	1:C:351:SER:HB2	2.18	0.42
1:B:54:GLU:OE1	1:B:54:GLU:HA	2.20	0.41
1:A:332:ALA:O	1:A:336:GLN:HG3	2.20	0.41
1:B:82:VAL:HG22	1:B:224:LYS:HB2	2.01	0.41
1:D:255:GLU:HG2	1:D:331:LEU:HD13	2.03	0.41
1:C:262:CYS:HB2	1:C:324:LEU:HD21	2.02	0.41
1:C:85:GLY:HA2	1:D:424:TYR:OH	2.20	0.41
1:D:124:LEU:HD13	1:D:235:PHE:CE1	2.54	0.41
1:A:626:LEU:HB3	1:A:627:PRO:HD3	2.01	0.41
1:D:250:LEU:O	1:D:254:VAL:HG23	2.21	0.41
1:D:583:GLU:O	1:D:587:ARG:HG3	2.20	0.41
1:A:51:VAL:HG21	1:A:159:LEU:HD13	2.02	0.41
1:A:280:LYS:HA	1:C:351:SER:HB2	2.03	0.41
1:C:626:LEU:N	1:C:627:PRO:HD2	2.36	0.41
1:B:72:LYS:HD3	1:B:72:LYS:HA	1.91	0.41
1:B:94:THR:HG22	1:B:96:ARG:H	1.85	0.41
1:C:471:LYS:HA	1:C:471:LYS:HD3	1.90	0.41
1:A:33:THR:HA	1:A:34:PRO:HD3	1.87	0.41
1:B:162:ASN:ND2	1:B:200:GLY:HA2	2.36	0.41
1:D:80:TYR:HB3	1:D:367:HIS:CD2	2.56	0.41
1:D:551:LEU:O	1:D:555:LEU:HG	2.21	0.40
1:C:650:GLN:HG3	1:D:111:GLY:O	2.22	0.40
1:A:65:CYS:O	1:A:69:VAL:HG23	2.20	0.40
1:B:304:LEU:HD11	1:B:619:ARG:HD3	2.04	0.40
1:C:162:ASN:ND2	1:C:200:GLY:HA2	2.37	0.40
1:A:224:LYS:HE2	1:A:356:PRO:HD2	2.04	0.40
1:B:380:MET:CE	1:B:380:MET:HA	2.51	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	631/707 (89%)	616 (98%)	14 (2%)	1 (0%)	56	62
1	B	622/707 (88%)	603 (97%)	19 (3%)	0	100	100
1	C	629/707 (89%)	609 (97%)	20 (3%)	0	100	100
1	D	624/707 (88%)	610 (98%)	14 (2%)	0	100	100
All	All	2506/2828 (89%)	2438 (97%)	67 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	175	ALA

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	543/592 (92%)	516 (95%)	27 (5%)	34	39
1	B	538/592 (91%)	522 (97%)	16 (3%)	53	64
1	C	539/592 (91%)	516 (96%)	23 (4%)	40	47
1	D	537/592 (91%)	512 (95%)	25 (5%)	36	42
All	All	2157/2368 (91%)	2066 (96%)	91 (4%)	40	48

All (91) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	LEU
1	A	57	GLN
1	A	59	ARG
1	A	84	THR
1	A	92	ARG
1	A	94	THR
1	A	102	GLU
1	A	107	CYS
1	A	115	LYS
1	A	162	ASN
1	A	166	LYS
1	A	179	LEU
1	A	202	ASP
1	A	209	GLU
1	A	233	THR
1	A	253	LEU
1	A	268	ARG
1	A	353	ASN
1	A	364	ARG
1	A	374	SER
1	A	528	LEU
1	A	562	THR
1	A	568	LYS
1	A	575	ASP
1	A	672	CYS
1	A	673	LEU
1	A	674	GLN
1	B	43	LEU
1	B	59	ARG
1	B	94	THR
1	B	102	GLU
1	B	107	CYS
1	B	179	LEU
1	B	233	THR
1	B	268	ARG
1	B	312	ASP
1	B	374	SER
1	B	392	LYS
1	B	528	LEU
1	B	562	THR
1	B	605	LYS
1	B	618	VAL
1	B	673	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	26	LYS
1	C	43	LEU
1	C	56	GLU
1	C	92	ARG
1	C	94	THR
1	C	98	SER
1	C	102	GLU
1	C	107	CYS
1	C	179	LEU
1	C	233	THR
1	C	268	ARG
1	C	275	LEU
1	C	298	SER
1	C	353	ASN
1	C	363	ASP
1	C	392	LYS
1	C	459	GLN
1	C	528	LEU
1	C	552	LEU
1	C	562	THR
1	C	619	ARG
1	C	673	LEU
1	C	677	LEU
1	D	43	LEU
1	D	59	ARG
1	D	92	ARG
1	D	94	THR
1	D	102	GLU
1	D	107	CYS
1	D	150	GLU
1	D	179	LEU
1	D	209	GLU
1	D	222	GLN
1	D	233	THR
1	D	268	ARG
1	D	312	ASP
1	D	324	LEU
1	D	353	ASN
1	D	363	ASP
1	D	528	LEU
1	D	552	LEU
1	D	562	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	568	LYS
1	D	576	THR
1	D	580	ARG
1	D	599	VAL
1	D	670	LEU
1	D	673	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	336	GLN
1	D	674	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, 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	641/707 (90%)	0.36	28 (4%)	33 33	20, 37, 61, 92	0
1	B	636/707 (89%)	0.34	27 (4%)	35 35	20, 40, 66, 81	0
1	C	639/707 (90%)	0.39	35 (5%)	24 24	19, 39, 72, 93	0
1	D	636/707 (89%)	0.37	21 (3%)	44 45	20, 39, 64, 85	0
All	All	2552/2828 (90%)	0.36	111 (4%)	34 34	19, 39, 66, 93	0

All (111) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	276	ILE	6.6
1	C	275	LEU	6.0
1	C	279	VAL	5.8
1	B	276	ILE	5.6
1	C	369	ALA	4.8
1	C	278	LYS	4.8
1	A	32	THR	4.1
1	C	281	PRO	4.0
1	C	277	HIS	3.9
1	A	367	HIS	3.8
1	A	276	ILE	3.8
1	B	564	LEU	3.6
1	A	363	ASP	3.5
1	C	579	ASP	3.5
1	C	57	GLN	3.4
1	B	522	ASN	3.4
1	D	568	LYS	3.3
1	A	121	VAL	3.2
1	D	440	LEU	3.1
1	A	590	ASP	3.1
1	C	280	LYS	3.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	D	577	LEU	3.1
1	A	275	LEU	3.0
1	D	279	VAL	3.0
1	D	215	GLY	3.0
1	A	86	PHE	3.0
1	B	443	LEU	3.0
1	A	444	ALA	3.0
1	A	591	ARG	2.9
1	C	585	GLU	2.9
1	B	562	THR	2.8
1	C	443	LEU	2.8
1	A	57	GLN	2.8
1	D	369	ALA	2.8
1	C	56	GLU	2.8
1	A	366	LEU	2.7
1	B	529	LEU	2.7
1	B	444	ALA	2.7
1	D	444	ALA	2.7
1	B	584	PHE	2.7
1	A	587	ARG	2.7
1	D	360	HIS	2.7
1	A	581	LEU	2.7
1	B	447	VAL	2.7
1	B	565	ALA	2.7
1	C	567	HIS	2.6
1	D	366	LEU	2.6
1	C	581	LEU	2.6
1	B	566	LEU	2.6
1	C	282	HIS	2.6
1	C	583	GLU	2.6
1	C	366	LEU	2.5
1	C	522	ASN	2.5
1	C	526	ALA	2.5
1	C	451	VAL	2.5
1	D	71	ARG	2.5
1	A	359	ASP	2.5
1	A	364	ARG	2.5
1	B	521	PRO	2.5
1	D	443	LEU	2.5
1	A	443	LEU	2.5
1	C	175	ALA	2.4
1	C	438	SER	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	440	LEU	2.4
1	D	630	ARG	2.4
1	D	75	ASP	2.4
1	D	202	ASP	2.4
1	C	582	ALA	2.4
1	D	175	ALA	2.4
1	C	447	VAL	2.4
1	D	76	GLY	2.4
1	B	449	THR	2.4
1	C	464	LEU	2.3
1	B	442	TYR	2.3
1	A	578	VAL	2.3
1	C	444	ALA	2.3
1	C	564	LEU	2.3
1	A	279	VAL	2.3
1	B	563	ILE	2.3
1	C	591	ARG	2.2
1	A	281	PRO	2.2
1	B	451	VAL	2.2
1	B	277	HIS	2.2
1	A	358	ILE	2.2
1	B	440	LEU	2.2
1	B	581	LEU	2.2
1	C	450	HIS	2.2
1	B	510	VAL	2.2
1	D	618	VAL	2.2
1	D	442	TYR	2.1
1	A	677	LEU	2.1
1	C	86	PHE	2.1
1	A	217	ARG	2.1
1	B	364	ARG	2.1
1	B	576	THR	2.1
1	C	388	ALA	2.1
1	D	62	VAL	2.1
1	B	603	TYR	2.1
1	D	387	VAL	2.1
1	B	454	ALA	2.1
1	B	203	VAL	2.1
1	D	278	LYS	2.1
1	A	448	THR	2.0
1	B	446	PRO	2.0
1	A	373	GLY	2.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	389	GLY	2.0
1	C	459	GLN	2.0
1	B	559	CYS	2.0
1	A	579	ASP	2.0
1	A	442	TYR	2.0
1	C	285	GLN	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.