



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 01:53 AM GMT

PDB ID : 4L5T
Title : Crystal structure of the tetrameric p202 HIN2
Authors : Yin, Q.; Tian, Y.; Wu, H.
Deposited on : 2013-06-11
Resolution : 3.40 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>

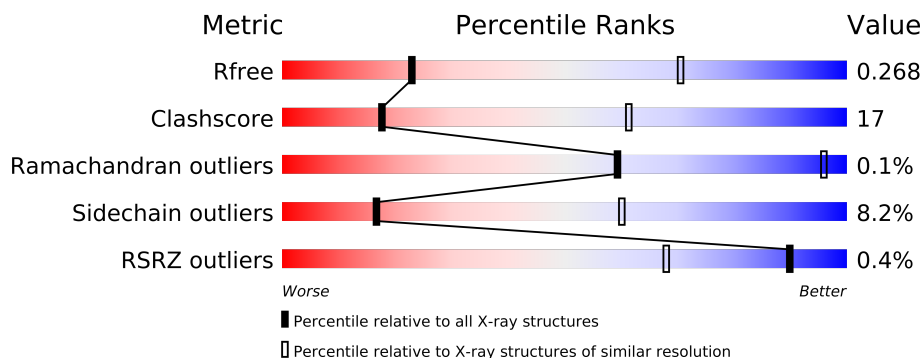
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 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 3.40 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	1017 (3.52-3.28)
Clashscore	79885	1214 (3.50-3.30)
Ramachandran outliers	78287	1177 (3.50-3.30)
Sidechain outliers	78261	1177 (3.50-3.30)
RSRZ outliers	66119	1017 (3.52-3.28)

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	203	
1	B	203	
1	C	203	
1	D	203	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 5748 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 Interferon-activable protein 202.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	180	Total	C	N	O	S	0	0	0
			1414	905	229	271	9			
1	B	185	Total	C	N	O	S	0	0	0
			1463	935	240	277	11			
1	C	182	Total	C	N	O	S	0	0	0
			1438	919	236	272	11			
1	D	183	Total	C	N	O	S	0	0	0
			1433	913	237	272	11			

There are 20 discrepancies between the modelled and reference sequences:

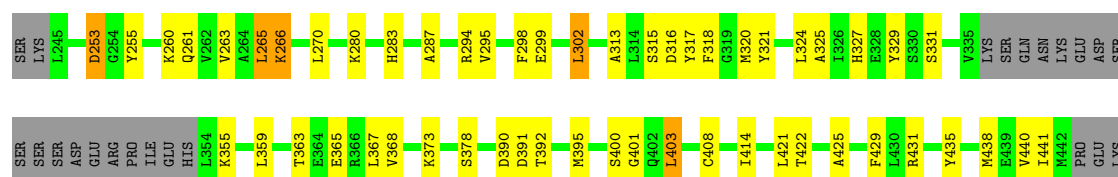
Chain	Residue	Modelled	Actual	Comment	Reference
A	243	SER	-	EXPRESSION TAG	UNP Q9R002
A	350	PRO	LEU	VARIANT	UNP Q9R002
A	364	GLU	LYS	VARIANT	UNP Q9R002
A	379	SER	THR	VARIANT	UNP Q9R002
A	432	ALA	SER	VARIANT	UNP Q9R002
B	243	SER	-	EXPRESSION TAG	UNP Q9R002
B	350	PRO	LEU	VARIANT	UNP Q9R002
B	364	GLU	LYS	VARIANT	UNP Q9R002
B	379	SER	THR	VARIANT	UNP Q9R002
B	432	ALA	SER	VARIANT	UNP Q9R002
C	243	SER	-	EXPRESSION TAG	UNP Q9R002
C	350	PRO	LEU	VARIANT	UNP Q9R002
C	364	GLU	LYS	VARIANT	UNP Q9R002
C	379	SER	THR	VARIANT	UNP Q9R002
C	432	ALA	SER	VARIANT	UNP Q9R002
D	243	SER	-	EXPRESSION TAG	UNP Q9R002
D	350	PRO	LEU	VARIANT	UNP Q9R002
D	364	GLU	LYS	VARIANT	UNP Q9R002
D	379	SER	THR	VARIANT	UNP Q9R002
D	432	ALA	SER	VARIANT	UNP Q9R002

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.

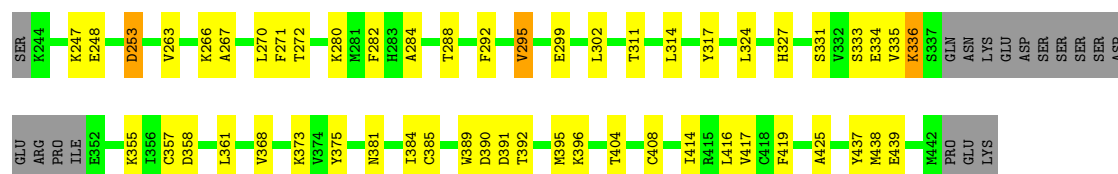
• Molecule 1: Interferon-activable protein 202

Chain A:



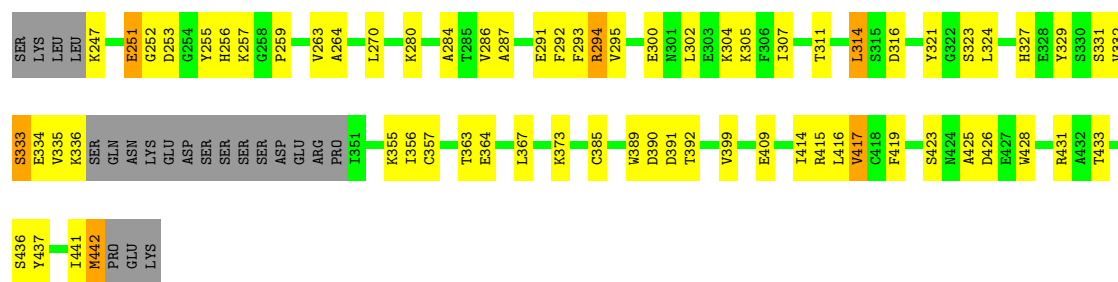
• Molecule 1: Interferon-activable protein 202

Chain B:



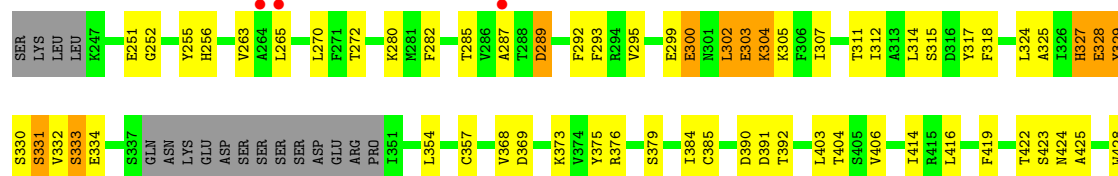
• Molecule 1: Interferon-activable protein 202

Chain C:



• Molecule 1: Interferon-activable protein 202

Chain D:



M438	M441	M442	PRO	GLU	LYS
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	61.89Å 71.86Å 95.47Å 90.00° 102.38° 90.00°	Depositor
Resolution (Å)	46.39 – 3.40 46.39 – 3.41	Depositor EDS
% Data completeness (in resolution range)	97.9 (46.39-3.40) 94.0 (46.39-3.41)	Depositor EDS
R_{merge}	0.21	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.47 (at 3.40Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.206 , 0.268 0.210 , 0.268	Depositor DCC
R_{free} test set	1085 reflections (9.98%)	DCC
Wilson B-factor (Å ²)	80.5	Xtriage
Anisotropy	0.600	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 51.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 11110 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	5748	wwPDB-VP
Average B, all atoms (Å ²)	102.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.87% 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	0.39	0/1441	0.58	0/1943
1	B	0.41	0/1491	0.60	0/2007
1	C	0.40	0/1466	0.60	0/1975
1	D	0.38	0/1460	0.58	1/1966 (0.1%)
All	All	0.40	0/5858	0.59	1/7891 (0.0%)

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	D	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	D	304	LYS	C-N-CA	6.26	137.34	121.70

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	329	TYR	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	1414	0	1375	47	0
1	B	1463	0	1436	32	0
1	C	1438	0	1402	41	0
1	D	1433	0	1392	91	0
All	All	5748	0	5605	196	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 17.

All (196) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:315:SER:N	1:D:331:SER:OG	1.58	1.35
1:D:302:LEU:HD11	1:D:328:GLU:O	1.33	1.26
1:D:282:PHE:CE1	1:D:303:GLU:CB	2.23	1.21
1:D:302:LEU:CD1	1:D:328:GLU:O	1.91	1.18
1:A:255:TYR:HB2	1:D:327:HIS:ND1	1.68	1.08
1:D:282:PHE:CE1	1:D:303:GLU:HB2	1.87	1.07
1:A:265:LEU:O	1:A:266:LYS:NZ	1.88	1.06
1:A:365:GLU:N	1:A:365:GLU:OE1	1.91	1.03
1:D:282:PHE:CZ	1:D:303:GLU:HA	1.94	1.00
1:D:302:LEU:HD22	1:D:332:VAL:HG21	1.45	0.99
1:D:270:LEU:HD11	1:D:300:GLU:HB3	1.46	0.97
1:D:299:GLU:OE1	1:D:300:GLU:N	1.99	0.96
1:D:282:PHE:CE1	1:D:303:GLU:HB3	1.99	0.95
1:D:315:SER:CA	1:D:331:SER:OG	2.16	0.94
1:B:357:CYS:HG	1:B:392:THR:HG1	1.10	0.92
1:D:302:LEU:CD2	1:D:332:VAL:CG2	2.48	0.91
1:D:302:LEU:HD22	1:D:332:VAL:CG2	2.01	0.89
1:D:302:LEU:CD2	1:D:332:VAL:HB	2.03	0.89
1:A:255:TYR:HB2	1:D:327:HIS:CE1	2.09	0.88
1:D:328:GLU:HB2	1:D:330:SER:HB2	1.55	0.86
1:A:255:TYR:HD2	1:D:327:HIS:HE1	1.24	0.84
1:A:255:TYR:CD2	1:D:327:HIS:HE1	1.98	0.81
1:D:315:SER:OG	1:D:331:SER:OG	2.00	0.80
1:A:359:LEU:HD23	1:A:421:LEU:HD22	1.64	0.79
1:D:302:LEU:HD21	1:D:332:VAL:HB	1.65	0.79
1:B:295:VAL:HB	1:B:324:LEU:HB2	1.64	0.78
1:D:424:ASN:OD1	1:D:425:ALA:N	2.16	0.78
1:D:302:LEU:CD2	1:D:332:VAL:HG21	2.11	0.77

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:327:HIS:C	1:D:328:GLU:HG2	2.06	0.76
1:C:357:CYS:H	1:C:392:THR:HG23	1.48	0.76
1:D:282:PHE:CZ	1:D:303:GLU:CA	2.67	0.76
1:B:327:HIS:HD2	1:C:255:TYR:HB2	1.48	0.76
1:D:302:LEU:CD2	1:D:332:VAL:CB	2.64	0.75
1:B:355:LYS:HB3	1:B:392:THR:HG21	1.68	0.75
1:D:302:LEU:HD11	1:D:328:GLU:C	2.08	0.74
1:A:255:TYR:CD2	1:D:327:HIS:CE1	2.75	0.74
1:D:282:PHE:CE1	1:D:303:GLU:CA	2.71	0.73
1:D:282:PHE:HZ	1:D:303:GLU:HA	1.46	0.73
1:A:265:LEU:HD23	1:A:265:LEU:N	2.03	0.73
1:C:334:GLU:HG3	1:C:335:VAL:N	2.03	0.72
1:D:315:SER:N	1:D:331:SER:CB	2.53	0.71
1:A:265:LEU:HD23	1:A:265:LEU:H	1.57	0.70
1:D:282:PHE:CZ	1:D:303:GLU:HB2	2.27	0.69
1:A:265:LEU:HB2	1:A:266:LYS:HD2	1.75	0.69
1:B:414:ILE:HD12	1:B:416:LEU:HD21	1.76	0.68
1:D:289:ASP:N	1:D:289:ASP:OD1	2.21	0.68
1:D:282:PHE:HE1	1:D:303:GLU:HB3	1.56	0.67
1:D:282:PHE:CD1	1:D:303:GLU:HB2	2.29	0.67
1:D:304:LYS:O	1:D:307:ILE:HG12	1.94	0.67
1:B:253:ASP:OD1	1:B:253:ASP:N	2.27	0.67
1:A:298:PHE:HB2	1:A:327:HIS:HA	1.76	0.66
1:D:302:LEU:CD1	1:D:328:GLU:C	2.64	0.66
1:D:282:PHE:CZ	1:D:303:GLU:CB	2.78	0.65
1:D:327:HIS:O	1:D:328:GLU:HG2	1.96	0.65
1:C:355:LYS:HB3	1:C:392:THR:HG21	1.79	0.64
1:D:299:GLU:OE1	1:D:300:GLU:HG3	1.98	0.64
1:A:265:LEU:C	1:A:266:LYS:HD2	2.18	0.63
1:D:252:GLY:O	1:D:256:HIS:NE2	2.26	0.63
1:D:315:SER:CB	1:D:331:SER:OG	2.46	0.63
1:D:314:LEU:HA	1:D:331:SER:HB3	1.80	0.62
1:D:265:LEU:HD11	1:D:287:ALA:HB2	1.81	0.62
1:A:255:TYR:HD2	1:D:327:HIS:CE1	2.11	0.62
1:B:425:ALA:HB2	1:C:385:CYS:SG	2.40	0.62
1:D:327:HIS:H	1:D:327:HIS:CD2	2.16	0.61
1:D:270:LEU:HD11	1:D:300:GLU:CB	2.27	0.61
1:B:417:VAL:HG22	1:B:437:TYR:HB3	1.82	0.61
1:A:403:LEU:HD13	1:A:438:MET:HB2	1.81	0.61
1:C:255:TYR:HE2	1:C:257:LYS:HB2	1.65	0.61
1:D:272:THR:HG22	1:D:280:LYS:HG2	1.82	0.61
1:A:355:LYS:HB3	1:A:392:THR:HG21	1.82	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:315:SER:O	1:A:331:SER:HB2	2.01	0.60
1:D:302:LEU:HD13	1:D:328:GLU:O	1.94	0.60
1:A:390:ASP:OD1	1:A:391:ASP:N	2.35	0.60
1:B:385:CYS:SG	1:C:425:ALA:HB2	2.42	0.60
1:D:295:VAL:HG12	1:D:324:LEU:HB2	1.84	0.60
1:D:318:PHE:HB3	1:D:325:ALA:HB3	1.84	0.59
1:D:302:LEU:HD23	1:D:332:VAL:HB	1.80	0.59
1:D:270:LEU:HA	1:D:282:PHE:HB3	1.83	0.59
1:A:320:MET:HG2	1:A:321:TYR:HD2	1.68	0.58
1:B:358:ASP:HA	1:B:361:LEU:HD12	1.86	0.58
1:B:270:LEU:HD21	1:B:280:LYS:HB3	1.85	0.58
1:C:251:GLU:HG2	1:C:293:PHE:HE2	1.69	0.58
1:C:270:LEU:HD21	1:C:280:LYS:HB3	1.85	0.57
1:A:403:LEU:HB3	1:A:438:MET:HE3	1.87	0.56
1:C:334:GLU:HG3	1:C:335:VAL:H	1.71	0.56
1:C:294:ARG:HH22	1:C:367:LEU:HB2	1.71	0.56
1:D:414:ILE:HD12	1:D:416:LEU:HD21	1.86	0.55
1:D:282:PHE:CD1	1:D:303:GLU:CB	2.87	0.55
1:A:425:ALA:HB2	1:D:385:CYS:SG	2.46	0.55
1:A:408:CYS:SG	1:A:414:ILE:HD11	2.47	0.54
1:A:265:LEU:HD21	1:A:287:ALA:HB3	1.89	0.54
1:D:251:GLU:HG2	1:D:293:PHE:CE2	2.42	0.54
1:B:390:ASP:OD1	1:B:391:ASP:N	2.41	0.54
1:B:267:ALA:HA	1:B:284:ALA:HA	1.89	0.54
1:D:328:GLU:CB	1:D:330:SER:HB2	2.31	0.54
1:D:270:LEU:HD21	1:D:280:LYS:HB3	1.90	0.54
1:D:384:ILE:HG12	1:D:404:THR:HG22	1.90	0.54
1:C:441:ILE:HG22	1:C:442:MET:HG3	1.90	0.53
1:D:403:LEU:HB3	1:D:438:MET:HE3	1.90	0.53
1:B:263:VAL:HG22	1:B:311:THR:HG22	1.90	0.53
1:D:328:GLU:C	1:D:330:SER:HB2	2.29	0.53
1:C:414:ILE:HD12	1:C:416:LEU:HD21	1.90	0.53
1:C:252:GLY:O	1:C:256:HIS:NE2	2.32	0.53
1:C:294:ARG:NH2	1:C:367:LEU:HB2	2.23	0.53
1:A:263:VAL:HG23	1:A:287:ALA:O	2.09	0.53
1:D:357:CYS:H	1:D:392:THR:HG23	1.72	0.53
1:C:390:ASP:OD1	1:C:391:ASP:N	2.42	0.52
1:B:248:GLU:HG3	1:B:292:PHE:CZ	2.45	0.52
1:A:395:MET:HE2	1:A:429:PHE:HA	1.92	0.52
1:B:368:VAL:N	1:B:419:PHE:O	2.41	0.52
1:A:318:PHE:HB3	1:A:325:ALA:HB3	1.92	0.51
1:A:270:LEU:HD21	1:A:280:LYS:HB3	1.92	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:302:LEU:HD11	1:D:329:TYR:CA	2.41	0.50
1:A:299:GLU:HB3	1:A:302:LEU:HD12	1.94	0.49
1:D:270:LEU:HD12	1:D:303:GLU:HG2	1.95	0.49
1:C:264:ALA:HA	1:C:286:VAL:HG12	1.95	0.49
1:D:406:VAL:HG21	1:D:438:MET:HB3	1.94	0.48
1:C:270:LEU:CD2	1:C:280:LYS:HB3	2.43	0.48
1:C:416:LEU:HD22	1:C:419:PHE:CZ	2.49	0.48
1:C:259:PRO:HA	1:C:314:LEU:O	2.14	0.48
1:A:265:LEU:CD2	1:A:265:LEU:H	2.19	0.48
1:D:265:LEU:HD12	1:D:285:THR:HG22	1.96	0.48
1:A:261:GLN:HA	1:A:313:ALA:HA	1.95	0.47
1:D:263:VAL:HG22	1:D:311:THR:HG22	1.95	0.47
1:D:373:LYS:NZ	1:D:373:LYS:HB3	2.29	0.47
1:A:294:ARG:HH22	1:A:367:LEU:HB3	1.79	0.47
1:B:419:PHE:HZ	1:B:438:MET:HE2	1.79	0.47
1:C:332:VAL:HG12	1:C:333:SER:O	2.14	0.47
1:A:400:SER:OG	1:A:401:GLY:N	2.47	0.47
1:A:295:VAL:HG12	1:A:324:LEU:HD12	1.96	0.47
1:C:305:LYS:NZ	1:C:329:TYR:OH	2.35	0.47
1:C:433:THR:N	1:C:436:SER:OG	2.45	0.47
1:D:368:VAL:N	1:D:419:PHE:O	2.48	0.47
1:D:328:GLU:HB2	1:D:330:SER:CB	2.35	0.47
1:B:271:PHE:CZ	1:B:437:TYR:HB2	2.50	0.47
1:D:251:GLU:OE2	1:D:317:TYR:OH	2.30	0.46
1:D:302:LEU:HD21	1:D:332:VAL:CG2	2.41	0.46
1:A:431:ARG:HD2	1:D:422:THR:HG23	1.97	0.46
1:C:423:SER:HA	1:C:428:TRP:CE2	2.50	0.46
1:A:378:SER:HB2	1:B:381:ASN:OD1	2.15	0.46
1:D:390:ASP:OD2	1:D:392:THR:HG22	2.15	0.46
1:D:369:ASP:HA	1:D:416:LEU:O	2.15	0.46
1:A:283:HIS:NE2	1:A:435:TYR:O	2.44	0.46
1:C:263:VAL:HG22	1:C:311:THR:HG22	1.96	0.46
1:A:327:HIS:CD2	1:D:255:TYR:HB2	2.50	0.46
1:C:295:VAL:HG12	1:C:324:LEU:HB2	1.98	0.46
1:B:416:LEU:HD22	1:B:419:PHE:CZ	2.51	0.46
1:A:265:LEU:HD21	1:A:287:ALA:CB	2.46	0.46
1:B:299:GLU:HB3	1:B:302:LEU:HD12	1.99	0.45
1:D:302:LEU:HD21	1:D:332:VAL:CB	2.34	0.45
1:A:320:MET:HG2	1:A:321:TYR:CD2	2.50	0.45
1:B:384:ILE:HG12	1:B:404:THR:HG22	1.97	0.45
1:A:253:ASP:OD1	1:A:253:ASP:N	2.41	0.45
1:D:302:LEU:O	1:D:305:LYS:HB2	2.17	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:284:ALA:O	1:C:295:VAL:HG22	2.16	0.45
1:C:426:ASP:HB3	1:D:425:ALA:HB3	1.99	0.45
1:D:302:LEU:HD11	1:D:329:TYR:HA	1.98	0.44
1:B:373:LYS:HB3	1:B:373:LYS:HE2	1.79	0.44
1:B:335:VAL:HG12	1:B:336:LYS:N	2.31	0.44
1:D:312:ILE:HD12	1:D:332:VAL:HG13	1.98	0.44
1:A:316:ASP:HB2	1:A:331:SER:HB3	2.00	0.43
1:A:421:LEU:HD12	1:A:422:THR:N	2.33	0.43
1:C:287:ALA:HA	1:C:292:PHE:HA	2.01	0.43
1:C:247:LYS:HA	1:C:291:GLU:HA	1.99	0.43
1:D:331:SER:O	1:D:333:SER:OG	2.33	0.43
1:B:375:TYR:HD2	1:B:389:TRP:HB2	1.83	0.43
1:A:422:THR:OG1	1:A:431:ARG:HG2	2.18	0.43
1:B:271:PHE:CE2	1:B:437:TYR:HB2	2.54	0.43
1:B:395:MET:HG2	1:B:396:LYS:H	1.82	0.43
1:B:416:LEU:HD22	1:B:419:PHE:CE1	2.54	0.43
1:B:270:LEU:HA	1:B:282:PHE:HB3	2.01	0.43
1:C:442:MET:HB2	1:C:442:MET:HE2	1.90	0.43
1:D:287:ALA:HA	1:D:292:PHE:HA	2.01	0.42
1:A:373:LYS:HB3	1:A:373:LYS:HE2	1.70	0.42
1:A:317:TYR:CD2	1:A:324:LEU:HD22	2.54	0.42
1:C:295:VAL:HG12	1:C:324:LEU:HD12	2.02	0.42
1:C:356:ILE:HG13	1:C:390:ASP:OD2	2.19	0.42
1:A:359:LEU:HD21	1:A:368:VAL:HG11	2.02	0.42
1:B:247:LYS:HE3	1:B:247:LYS:HB2	1.84	0.42
1:D:357:CYS:H	1:D:392:THR:CG2	2.32	0.42
1:C:257:LYS:HG3	1:C:316:ASP:OD1	2.20	0.42
1:B:302:LEU:HD23	1:B:302:LEU:HA	1.90	0.41
1:D:375:TYR:CE1	1:D:376:ARG:HD2	2.55	0.41
1:D:423:SER:HA	1:D:428:TRP:CE2	2.55	0.41
1:C:251:GLU:HG2	1:C:293:PHE:CE2	2.51	0.41
1:C:270:LEU:HD11	1:C:300:GLU:HA	2.02	0.41
1:D:441:ILE:HG22	1:D:442:MET:N	2.36	0.41
1:C:417:VAL:HG22	1:C:437:TYR:HB3	2.01	0.41
1:C:373:LYS:HE2	1:C:373:LYS:HB3	1.78	0.41
1:B:317:TYR:CE2	1:B:324:LEU:HD22	2.56	0.41
1:C:415:ARG:HB2	1:C:441:ILE:HD11	2.03	0.41
1:A:255:TYR:CB	1:D:327:HIS:CE1	2.93	0.41
1:D:390:ASP:OD1	1:D:391:ASP:N	2.54	0.40
1:D:302:LEU:HD23	1:D:332:VAL:CB	2.43	0.40
1:C:357:CYS:HB3	1:C:392:THR:O	2.21	0.40
1:C:304:LYS:O	1:C:307:ILE:HG12	2.21	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	176/203 (87%)	164 (93%)	12 (7%)	0	100	100
1	B	181/203 (89%)	170 (94%)	11 (6%)	0	100	100
1	C	178/203 (88%)	165 (93%)	12 (7%)	1 (1%)	33	85
1	D	179/203 (88%)	166 (93%)	13 (7%)	0	100	100
All	All	714/812 (88%)	665 (93%)	48 (7%)	1 (0%)	59	96

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	321	TYR

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	153/182 (84%)	143 (94%)	10 (6%)	24	71
1	B	160/182 (88%)	148 (92%)	12 (8%)	19	64
1	C	156/182 (86%)	138 (88%)	18 (12%)	8	38
1	D	155/182 (85%)	144 (93%)	11 (7%)	21	67
All	All	624/728 (86%)	573 (92%)	51 (8%)	17	60

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	253	ASP
1	A	260	LYS
1	A	265	LEU
1	A	266	LYS
1	A	302	LEU
1	A	329	TYR
1	A	363	THR
1	A	403	LEU
1	A	440	VAL
1	A	441	ILE
1	B	253	ASP
1	B	266	LYS
1	B	272	THR
1	B	288	THR
1	B	295	VAL
1	B	314	LEU
1	B	331	SER
1	B	333	SER
1	B	334	GLU
1	B	336	LYS
1	B	408	CYS
1	B	439	GLU
1	C	251	GLU
1	C	253	ASP
1	C	294	ARG
1	C	302	LEU
1	C	314	LEU
1	C	323	SER
1	C	327	HIS
1	C	331	SER
1	C	333	SER
1	C	336	LYS
1	C	363	THR
1	C	364	GLU
1	C	389	TRP
1	C	399	VAL
1	C	409	GLU
1	C	417	VAL
1	C	431	ARG
1	C	442	MET
1	D	289	ASP
1	D	300	GLU
1	D	302	LEU

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Mol	Chain	Res	Type
1	D	303	GLU
1	D	327	HIS
1	D	328	GLU
1	D	331	SER
1	D	333	SER
1	D	334	GLU
1	D	354	LEU
1	D	379	SER

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

Mol	Chain	Res	Type
1	D	327	HIS
1	D	360	HIS

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	180/203 (88%)	0.04	0	100 100	73, 114, 145, 153	0
1	B	185/203 (91%)	-0.11	0	100 100	60, 88, 116, 135	0
1	C	182/203 (89%)	-0.10	0	100 100	66, 89, 119, 133	0
1	D	183/203 (90%)	0.11	3 (1%)	68 32	84, 116, 142, 158	0
All	All	730/812 (89%)	-0.02	3 (0%)	90 68	60, 99, 141, 158	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	287	ALA	2.6
1	D	264	ALA	2.4
1	D	265	LEU	2.3

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.