



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:51 PM GMT

PDB ID : 1HK7
Title : MIDDLE DOMAIN OF HSP90
Authors : Meyer, P.; Prodromou, C.; Roe, S.M.; Pearl, L.H.
Deposited on : 2003-03-06
Resolution : 2.50 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

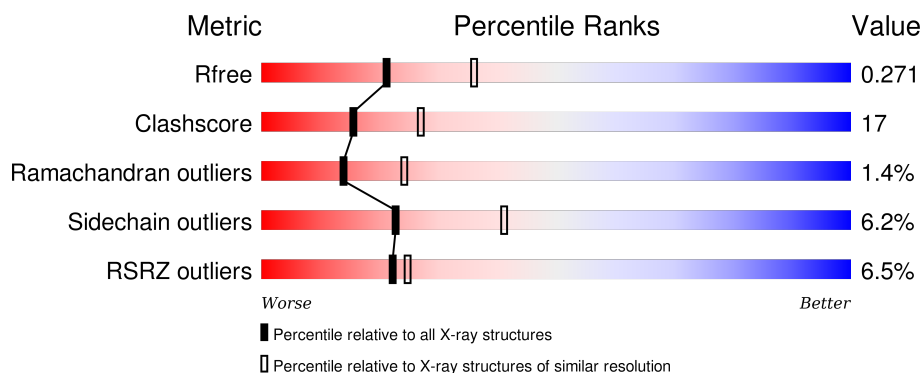
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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	288	<div> <div>6%</div> <div> <div></div> <div>51%</div> <div>31%</div> <div>•</div> <div>14%</div> </div> </div>
1	B	288	<div> <div>5%</div> <div> <div></div> <div>59%</div> <div>23%</div> <div>•</div> <div>14%</div> </div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4299 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 HEAT SHOCK PROTEIN HSP82.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	249	Total	C	N	O	S	0	0	1
			2053	1328	334	389	2			
1	B	247	Total	C	N	O	S	0	0	0
			2040	1321	329	387	3			

- Molecule 2 is CADMIUM ION (three-letter code: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Cd	0	0
			1	1		
2	A	5	Total	Cd	0	0
			5	5		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		

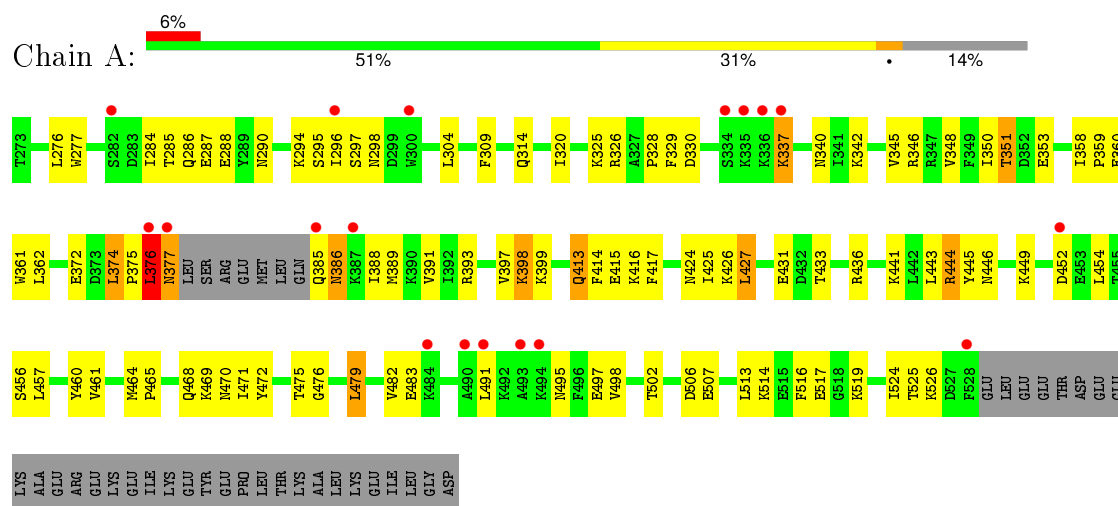
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	104	Total	O	0	0
			104	104		
4	B	95	Total	O	0	0
			95	95		

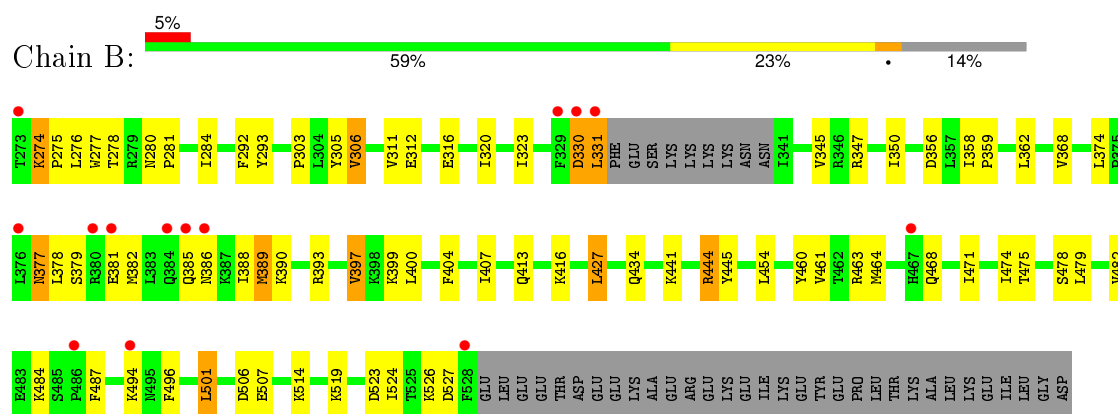
3 Residue-property plots

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: HEAT SHOCK PROTEIN HSP82



• Molecule 1: HEAT SHOCK PROTEIN HSP82



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	112.88Å 112.88Å 112.06Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.68 – 2.50 29.68 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.6 (29.68-2.50) 99.7 (29.68-2.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 2.51Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.240 , 0.277 0.235 , 0.271	Depositor DCC
R_{free} test set	1344 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	68.3	Xtriage
Anisotropy	0.210	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 52.5	EDS
Estimated twinning fraction	0.064 for h,-h-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 55207 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4299	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.01% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.40	0/2097	0.62	1/2830 (0.0%)
1	B	0.43	0/2084	0.61	0/2816
All	All	0.42	0/4181	0.61	1/5646 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	376	LEU	N-CA-C	8.03	132.67	111.00

There are no chirality outliers.

There are no planarity outliers.

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	2053	0	2052	85	0
1	B	2040	0	2029	56	0
2	A	5	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
4	A	104	0	0	6	0
4	B	95	0	0	5	0
All	All	4299	0	4081	141	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 17.

All (141) 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:469:LYS:HD3	1:A:469:LYS:O	1.65	0.96
1:A:446:ASN:ND2	1:A:454:LEU:HD23	1.88	0.89
1:B:501:LEU:HD13	1:B:506:ASP:HB3	1.58	0.84
1:B:275:PRO:O	1:B:278:THR:HG22	1.80	0.82
1:A:374:LEU:HD12	1:A:376:LEU:HA	1.60	0.81
1:A:359:PRO:HG2	1:A:362:LEU:HD12	1.62	0.79
1:A:491:LEU:HD11	1:A:524:ILE:HG21	1.66	0.77
1:B:359:PRO:HD3	1:B:393:ARG:HH11	1.56	0.71
1:B:460:TYR:CE2	1:B:471:ILE:HG23	2.26	0.70
1:A:519:LYS:HG2	4:A:2096:HOH:O	1.90	0.70
1:A:449:LYS:HD2	1:A:497:GLU:HB2	1.75	0.68
1:A:374:LEU:CD1	1:A:376:LEU:HA	2.22	0.68
1:A:374:LEU:C	1:A:376:LEU:H	1.96	0.67
1:B:377:ASN:HD22	1:B:377:ASN:H	1.40	0.67
1:A:491:LEU:HD13	1:A:498:VAL:HG21	1.77	0.67
1:B:501:LEU:CD1	1:B:506:ASP:HB3	2.26	0.66
1:A:350:ILE:HG22	1:A:351:THR:HG22	1.78	0.65
1:B:345:VAL:CG2	1:B:350:ILE:HD12	2.29	0.63
1:B:312:GLU:HA	1:B:316:GLU:HG2	1.82	0.62
1:A:337:LYS:HG3	1:A:360:GLU:HG3	1.83	0.61
1:B:460:TYR:HA	1:B:463:ARG:NH1	2.15	0.61
1:A:296:ILE:HG23	1:A:342:LYS:HE2	1.81	0.61
1:A:479:LEU:O	1:A:483:GLU:HG3	1.99	0.60
1:A:443:LEU:HD13	1:A:445:TYR:OH	2.01	0.60
1:B:474:ILE:HB	1:B:524:ILE:HD11	1.85	0.59
1:A:296:ILE:O	1:A:296:ILE:HG22	2.02	0.59
1:B:461:VAL:HG13	1:B:464:MET:HE1	1.83	0.59
1:A:358:ILE:HG13	1:A:359:PRO:HD2	1.83	0.59
1:A:470:ASN:HB2	1:A:472:TYR:CZ	2.38	0.59
1:A:376:LEU:HD23	1:A:377:ASN:H	1.68	0.58
1:B:388:ILE:O	1:B:390:LYS:N	2.30	0.58
1:B:311:VAL:O	1:B:316:GLU:HA	2.04	0.58
1:A:359:PRO:CG	1:A:362:LEU:HD12	2.32	0.58
1:B:444:ARG:HG2	1:B:454:LEU:HB3	1.86	0.58
1:A:287:GLU:HA	1:A:290:ASN:HB2	1.85	0.58
1:A:386:ASN:HD21	1:A:388:ILE:HB	1.67	0.58
1:B:305:TYR:CB	1:B:407:ILE:HD11	2.33	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:345:VAL:O	1:A:345:VAL:HG13	2.05	0.57
1:A:372:GLU:HB2	4:A:2034:HOH:O	2.04	0.56
1:A:393:ARG:O	1:A:397:VAL:HG23	2.05	0.56
1:A:374:LEU:C	1:A:376:LEU:N	2.59	0.56
1:B:389:MET:HA	4:B:2037:HOH:O	2.06	0.56
1:A:446:ASN:HD21	1:A:454:LEU:HD23	1.69	0.55
1:B:514:LYS:HE3	4:B:2093:HOH:O	2.07	0.55
1:A:491:LEU:CD1	1:A:524:ILE:HG21	2.35	0.55
1:B:362:LEU:HD22	1:B:400:LEU:HD13	1.88	0.55
1:B:523:ASP:HB3	1:B:526:LYS:HD3	1.87	0.55
1:A:290:ASN:O	1:A:294:LYS:HG2	2.07	0.54
1:A:417:PHE:CE2	1:A:425:ILE:HD11	2.42	0.54
1:B:306:VAL:HG13	4:B:2009:HOH:O	2.07	0.54
1:A:475:THR:CG2	1:A:507:GLU:HG2	2.38	0.53
1:A:398:LYS:HE3	4:A:2042:HOH:O	2.07	0.53
1:A:469:LYS:HD3	1:A:469:LYS:C	2.27	0.53
1:B:293:TYR:CD2	1:B:303:PRO:HD3	2.44	0.53
1:A:277:TRP:HA	1:A:320:ILE:HD11	1.89	0.53
1:A:491:LEU:HD11	1:A:524:ILE:CG2	2.39	0.53
1:A:441:LYS:NZ	1:A:441:LYS:HB2	2.23	0.53
1:A:286:GLN:C	1:A:288:GLU:H	2.12	0.52
1:A:413:GLN:HG2	4:A:2012:HOH:O	2.09	0.52
1:B:347:ARG:HG3	4:B:2024:HOH:O	2.10	0.52
1:A:361:TRP:O	1:A:424:ASN:HB3	2.11	0.51
1:A:460:TYR:CE2	1:A:471:ILE:HG23	2.45	0.51
1:A:457:LEU:O	1:A:461:VAL:HG23	2.11	0.51
1:A:345:VAL:O	1:A:348:VAL:HG22	2.11	0.51
1:A:416:LYS:NZ	1:A:416:LYS:HB2	2.25	0.51
1:B:445:TYR:CD2	1:B:501:LEU:HD22	2.47	0.50
1:A:461:VAL:HA	1:A:464:MET:CE	2.42	0.49
1:A:465:PRO:HD2	1:A:468:GLN:OE1	2.12	0.49
1:A:524:ILE:HG13	1:A:525:THR:N	2.27	0.49
1:A:348:VAL:HG23	1:A:348:VAL:O	2.12	0.49
1:A:426:LYS:HE2	1:A:506:ASP:OD1	2.12	0.49
1:A:456:SER:HB2	4:A:2080:HOH:O	2.12	0.49
1:A:491:LEU:HD13	1:A:498:VAL:CG2	2.42	0.48
1:A:374:LEU:HB3	1:A:376:LEU:HG	1.94	0.48
1:A:296:ILE:CG2	1:A:342:LYS:HE2	2.43	0.48
1:B:358:ILE:HD11	1:B:362:LEU:HB3	1.95	0.48
1:B:379:SER:OG	1:B:381:GLU:HG3	2.14	0.48
1:A:375:PRO:HB3	1:A:389:MET:CE	2.43	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:377:ASN:ND2	1:B:377:ASN:H	2.10	0.47
1:B:464:MET:HE1	1:B:519:LYS:HD2	1.95	0.47
1:A:417:PHE:HE2	1:A:425:ILE:HD11	1.77	0.47
1:B:345:VAL:HG23	1:B:350:ILE:HD12	1.96	0.47
1:B:478:SER:O	1:B:482:VAL:HG23	2.15	0.47
1:B:441:LYS:O	1:B:444:ARG:NH2	2.48	0.47
1:B:280:ASN:OD1	1:B:281:PRO:HD2	2.15	0.47
1:A:516:PHE:CE2	1:A:517:GLU:HG3	2.49	0.47
1:A:465:PRO:HG2	1:A:468:GLN:HB2	1.95	0.47
1:B:374:LEU:HD13	1:B:378:LEU:HD11	1.97	0.47
1:A:346:ARG:HB3	1:A:346:ARG:HH11	1.79	0.47
1:B:474:ILE:HD12	1:B:524:ILE:HD11	1.97	0.47
1:A:375:PRO:HB3	1:A:389:MET:HE2	1.97	0.46
1:B:399:LYS:NZ	4:B:2041:HOH:O	2.48	0.46
1:A:285:THR:O	1:A:288:GLU:HB2	2.15	0.46
1:A:414:PHE:HE1	1:A:444:ARG:HD2	1.80	0.46
1:A:433:THR:HG22	1:A:436:ARG:NH2	2.30	0.46
1:A:475:THR:HG21	1:A:507:GLU:HG2	1.98	0.46
1:B:330:ASP:O	1:B:331:LEU:HG	2.14	0.46
1:B:381:GLU:O	1:B:385:GLN:HG3	2.15	0.46
1:A:479:LEU:O	1:A:482:VAL:HG12	2.16	0.46
1:A:304:LEU:HB2	1:A:325:LYS:HA	1.98	0.45
1:A:427:LEU:O	1:A:431:GLU:HG2	2.17	0.45
1:A:385:GLN:HG2	1:A:386:ASN:N	2.32	0.45
1:B:474:ILE:HG23	1:B:482:VAL:HG11	1.99	0.44
1:B:468:GLN:HG3	1:B:496:PHE:CE1	2.53	0.44
1:B:393:ARG:O	1:B:397:VAL:HG13	2.17	0.44
1:B:356:ASP:N	1:B:356:ASP:OD2	2.50	0.44
1:A:513:LEU:O	1:A:514:LYS:HB2	2.17	0.44
1:B:358:ILE:HG13	1:B:359:PRO:HD2	2.00	0.44
1:B:292:PHE:HZ	1:B:368:VAL:HG21	1.83	0.44
1:B:464:MET:HE1	1:B:519:LYS:CD	2.49	0.43
1:A:309:PHE:CE1	1:A:399:LYS:HG3	2.54	0.43
1:A:337:LYS:HB2	1:A:360:GLU:HG3	2.01	0.43
1:B:305:TYR:HB2	1:B:407:ILE:HD11	2.00	0.43
1:A:326:ARG:HH11	1:A:326:ARG:HG3	1.83	0.43
1:A:471:ILE:HD11	1:A:519:LYS:HB2	2.01	0.42
1:A:276:LEU:HD11	1:A:284:ILE:CG2	2.49	0.42
1:B:427:LEU:HD23	1:B:427:LEU:HA	1.92	0.42
1:B:413:GLN:NE2	1:B:416:LYS:HE3	2.34	0.42
1:A:297:SER:O	1:A:298:ASN:HB2	2.20	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:304:LEU:O	1:A:304:LEU:HD23	2.20	0.42
1:A:328:PRO:HG3	4:A:2030:HOH:O	2.19	0.42
1:B:362:LEU:HD13	1:B:400:LEU:HD12	2.02	0.42
1:B:382:MET:HG2	1:B:386:ASN:HD21	1.85	0.42
1:A:376:LEU:O	1:A:377:ASN:HB2	2.20	0.41
1:A:353:GLU:O	1:A:353:GLU:HG2	2.20	0.41
1:A:286:GLN:C	1:A:288:GLU:N	2.74	0.41
1:A:388:ILE:O	1:A:391:VAL:HG22	2.20	0.41
1:B:277:TRP:HA	1:B:320:ILE:HD11	2.02	0.41
1:B:276:LEU:HD11	1:B:284:ILE:CG2	2.51	0.41
1:A:471:ILE:CD1	1:A:519:LYS:HB2	2.51	0.41
1:A:286:GLN:O	1:A:287:GLU:HB3	2.20	0.41
1:A:461:VAL:HA	1:A:464:MET:HE1	2.03	0.41
1:B:475:THR:CG2	1:B:507:GLU:HG2	2.51	0.41
1:A:476:GLY:O	1:A:502:THR:HA	2.21	0.41
1:B:460:TYR:HA	1:B:463:ARG:HH12	1.82	0.40
1:B:274:LYS:HG2	1:B:278:THR:CG2	2.51	0.40
1:A:345:VAL:HG12	1:A:350:ILE:CG1	2.51	0.40
1:A:346:ARG:HB3	1:A:346:ARG:NH1	2.36	0.40
1:B:323:ILE:HD12	1:B:404:PHE:HE1	1.86	0.40
1:B:482:VAL:C	1:B:484:LYS:H	2.24	0.40
1:B:305:TYR:HB3	1:B:407:ILE:HD11	2.04	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	245/288 (85%)	222 (91%)	20 (8%)	3 (1%)	16	29
1	B	243/288 (84%)	222 (91%)	17 (7%)	4 (2%)	12	21
All	All	488/576 (85%)	444 (91%)	37 (8%)	7 (1%)	14	24

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	337	LYS
1	B	389	MET
1	A	376	LEU
1	A	386	ASN
1	B	494	LYS
1	B	274	LYS
1	B	527	ASP

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	228/266 (86%)	211 (92%)	17 (8%)	17	31
1	B	226/266 (85%)	215 (95%)	11 (5%)	31	55
All	All	454/532 (85%)	426 (94%)	28 (6%)	23	41

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

Mol	Chain	Res	Type
1	A	295	SER
1	A	314	GLN
1	A	329	PHE
1	A	330	ASP
1	A	340	ASN
1	A	351	THR
1	A	374	LEU
1	A	377	ASN
1	A	398	LYS
1	A	413	GLN
1	A	415	GLU
1	A	427	LEU
1	A	444	ARG
1	A	452	ASP
1	A	479	LEU
1	A	495	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	526	LYS
1	B	306	VAL
1	B	330	ASP
1	B	331	LEU
1	B	377	ASN
1	B	397	VAL
1	B	427	LEU
1	B	434	GLN
1	B	444	ARG
1	B	479	LEU
1	B	487	PHE
1	B	501	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	286	GLN
1	A	385	GLN
1	A	386	ASN
1	A	413	GLN
1	A	434	GLN
1	A	446	ASN
1	A	495	ASN
1	B	286	GLN
1	B	377	ASN
1	B	385	GLN
1	B	386	ASN
1	B	413	GLN
1	B	434	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

Of 7 ligands modelled in this entry, 7 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	249/288 (86%)	0.24	18 (7%)	18 20	45, 72, 114, 131	0
1	B	247/288 (85%)	0.18	14 (5%)	27 31	40, 65, 110, 135	0
All	All	496/576 (86%)	0.21	32 (6%)	22 25	40, 69, 113, 135	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	528	PHE	7.3
1	B	273	THR	5.5
1	A	300	TRP	4.7
1	A	337	LYS	4.4
1	A	452	ASP	4.3
1	B	384	GLN	4.3
1	A	377	ASN	4.1
1	A	385	GLN	3.9
1	A	493	ALA	3.7
1	B	331	LEU	3.6
1	B	528	PHE	3.5
1	A	335	LYS	3.5
1	B	486	PRO	3.3
1	B	330	ASP	2.9
1	B	380	ARG	2.9
1	B	385	GLN	2.8
1	A	334	SER	2.7
1	B	494	LYS	2.7
1	B	329	PHE	2.6
1	A	376	LEU	2.6
1	A	490	ALA	2.6
1	B	376	LEU	2.5
1	A	336	LYS	2.5
1	B	467	HIS	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	494	LYS	2.3
1	A	484	LYS	2.2
1	A	387	LYS	2.2
1	A	296	ILE	2.2
1	A	282	SER	2.2
1	A	491	LEU	2.1
1	B	381	GLU	2.1
1	B	386	ASN	2.1

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	CD	A	1530	1/1	0.99	0.13	-0.53	68,68,68,68	0
2	CD	A	1531	1/1	0.98	0.09	-1.60	74,74,74,74	0
2	CD	B	1529	1/1	0.94	0.04	-	117,117,117,117	0
2	CD	A	1528	1/1	1.00	0.07	-	44,44,44,44	0
2	CD	A	1529	1/1	0.98	0.02	-	145,145,145,145	0
3	MG	A	1533	1/1	0.90	0.20	-	90,90,90,90	0
2	CD	A	1532	1/1	0.95	0.09	-	114,114,114,114	0

6.5 Other polymers [i](#)

There are no such residues in this entry.