



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 12:43 PM UTC

PDB ID : 1HK7 / pdb_00001hk7
Title : Middle Domain of HSP90
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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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

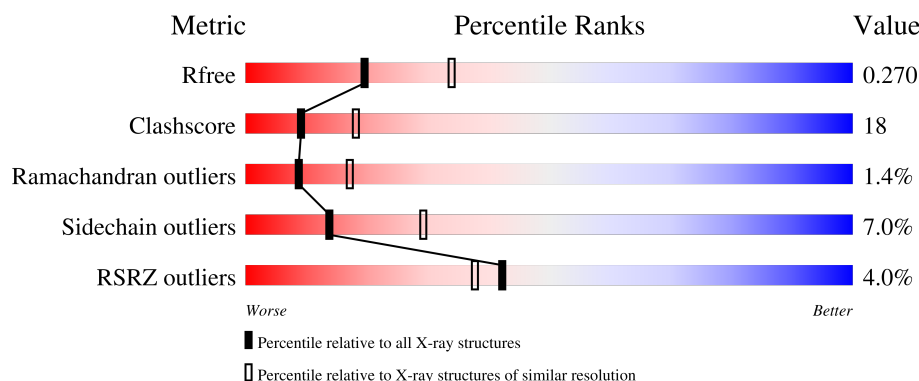
1 Overall quality at a glance

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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	5% 49% 32% 5% 14%
1	B	288	2% 56% 26% • • 14%

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 (CCD ID: CD) (formula: Cd).

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

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

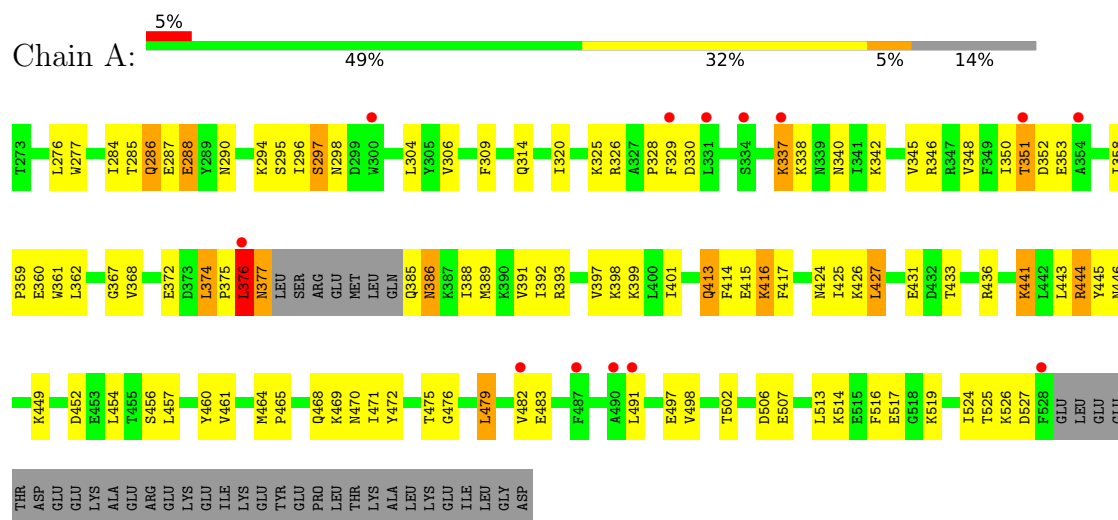
- Molecule 4 is water.

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

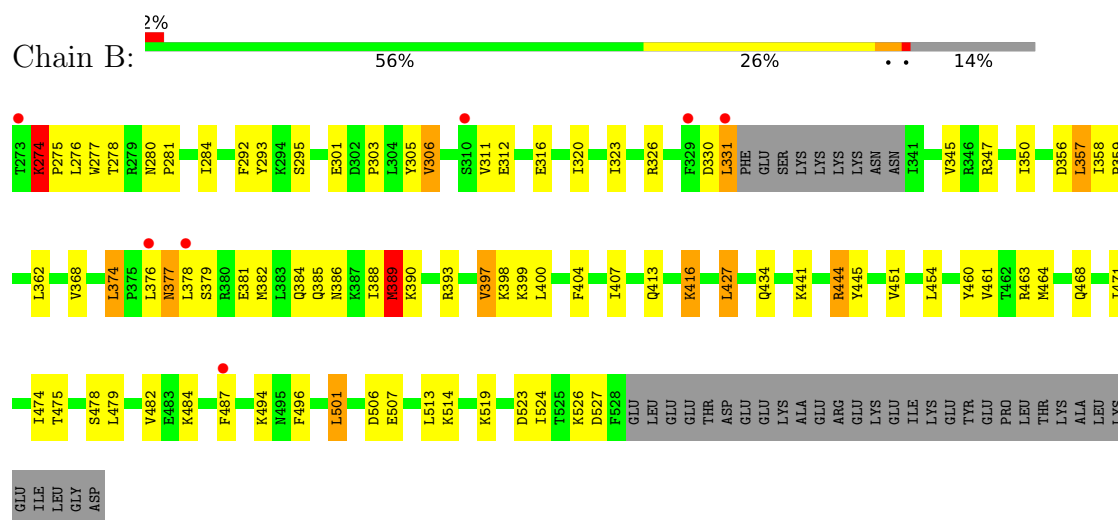
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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) 99.6 (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 0.233 , 0.270	Depositor DCC
R_{free} test set	2656 reflections (4.81%)	wwPDB-VP
Wilson B-factor (Å ²)	68.3	Xtriage
Anisotropy	0.210	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 42.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.064 for h,-h-k,-l	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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0.46	1/2097 (0.0%)	0.96	10/2830 (0.4%)
1	B	0.49	0/2084	0.93	8/2816 (0.3%)
All	All	0.48	1/4181 (0.0%)	0.94	18/5646 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	527	ASP	C-N	-5.40	1.25	1.33

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	376	LEU	N-CA-C	10.27	132.67	110.80
1	A	288	GLU	N-CA-C	-7.00	103.09	111.69
1	A	401	ILE	N-CA-C	-6.68	103.80	110.62
1	A	367	GLY	N-CA-C	6.43	119.11	110.43
1	B	295	SER	N-CA-C	6.39	117.93	110.97
1	B	389	MET	N-CA-C	6.39	124.40	110.80
1	A	352	ASP	N-CA-C	-5.99	105.28	112.59
1	A	286	GLN	N-CA-C	-5.98	102.78	110.19
1	A	392	ILE	N-CA-C	-5.84	105.15	110.82
1	A	297	SER	N-CA-C	5.80	120.51	113.38
1	B	451	VAL	N-CA-C	5.68	117.06	111.67
1	A	368	VAL	N-CA-C	5.54	116.65	108.45
1	B	274	LYS	CA-C-N	5.41	125.71	119.93
1	B	274	LYS	C-N-CA	5.41	125.71	119.93
1	B	374	LEU	CA-C-N	5.15	124.87	119.82
1	B	374	LEU	C-N-CA	5.15	124.87	119.82
1	B	357	LEU	N-CA-C	-5.11	106.21	112.90
1	A	338	LYS	N-CA-C	-5.10	104.32	110.44

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-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 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	87	0
1	B	2040	0	2029	63	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	150	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 18.

All (150) 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:O	1:A:469:LYS:HD3	1.65	0.96
1:A:461:VAL:HA	1:A:464:MET:HE2	1.52	0.91
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:A:374:LEU:C	1:A:376:LEU:H	1.96	0.73
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:464:MET:HE1	1:A:519:LYS:HG3	1.73	0.69
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:B:377:ASN:HD22	1:B:377:ASN:H	1.40	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:461:VAL:HG13	1:B:464:MET:HE2	1.81	0.62
1:A:337:LYS:HG3	1:A:360:GLU:HG3	1.83	0.61
1:A:374:LEU:C	1:A:376:LEU:N	2.59	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:B:464:MET:HE1	1:B:519:LYS:CD	2.31	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:HG22	1:A:296:ILE:O	2.02	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:A:469:LYS:HD3	1:A:469:LYS:C	2.27	0.58
1:B:311:VAL:O	1:B:316:GLU:HA	2.04	0.58
1:A:461:VAL:CA	1:A:464:MET:HE2	2.31	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
1:A:286:GLN:C	1:A:288:GLU:H	2.12	0.57
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: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:B:464:MET:HE1	1:B:519:LYS:HE3	1.88	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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: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:B:357:LEU:HD13	1:B:389:MET:HE3	1.92	0.52
1:A:361:TRP:O	1:A:424:ASN:HB3	2.11	0.51
1:A:457:LEU:O	1:A:461:VAL:HG23	2.11	0.51
1:A:460:TYR:CE2	1:A:471:ILE:HG23	2.45	0.51
1:A:345:VAL:O	1:A:348:VAL:HG22	2.11	0.51
1:A:416:LYS:HB2	1:A:416:LYS:NZ	2.25	0.51
1:B:445:TYR:CD2	1:B:501:LEU:HD22	2.47	0.50
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:O	1:A:348:VAL:HG23	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
1:B:377:ASN:H	1:B:377:ASN:ND2	2.10	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:280:ASN:OD1	1:B:281:PRO:HD2	2.15	0.47
1:B:441:LYS:O	1:B:444:ARG:NH2	2.48	0.47
1:A:516:PHE:CE2	1:A:517:GLU:HG3	2.49	0.47
1:A:346:ARG:HH11	1:A:346:ARG:HB3	1.79	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:B:464:MET:HE1	1:B:519:LYS:CE	2.44	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:286:GLN:C	1:A:288:GLU:N	2.74	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:B:482:VAL:C	1:B:484:LYS:H	2.24	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:356:ASP:OD2	1:B:356:ASP:N	2.50	0.44
1:B:393:ARG:O	1:B:397:VAL:HG13	2.17	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:376:LEU:C	1:B:378:LEU:H	2.25	0.44
1:A:441:LYS:HB2	1:A:441:LYS:HZ3	1.83	0.44
1:B:357:LEU:CD1	1:B:389:MET:HE3	2.47	0.43
1:A:309:PHE:CE1	1:A:399:LYS:HG3	2.54	0.43
1:A:326:ARG:HH11	1:A:326:ARG:HG3	1.83	0.43
1:A:337:LYS:HB2	1:A:360:GLU:HG3	2.01	0.43
1:B:301:GLU:CD	1:B:326:ARG:HH21	2.25	0.43
1:B:305:TYR:HB2	1:B:407:ILE:HD11	2.00	0.43
1:A:276:LEU:HD11	1:A:284:ILE:CG2	2.49	0.42
1:A:471:ILE:HD11	1:A:519:LYS:HB2	2.01	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
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:B:376:LEU:C	1:B:376:LEU:HD23	2.45	0.41
1:B:464:MET:HE1	1:B:519:LYS:HD2	2.02	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:276:LEU:HD11	1:B:284:ILE:CG2	2.51	0.41
1:A:286:GLN:O	1:A:287:GLU:HB3	2.20	0.41
1:A:471:ILE:CD1	1:A:519:LYS:HB2	2.51	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:A:345:VAL:HG12	1:A:350:ILE:CG1	2.51	0.40
1:B:274:LYS:HG2	1:B:278:THR:CG2	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: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%)	10	20
1	B	243/288 (84%)	222 (91%)	17 (7%)	4 (2%)	7	14
All	All	488/576 (85%)	444 (91%)	37 (8%)	7 (1%)	9	17

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%)	210 (92%)	18 (8%)	11	24
1	B	226/266 (85%)	212 (94%)	14 (6%)	16	34
All	All	454/532 (85%)	422 (93%)	32 (7%)	14	29

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	295	SER
1	A	306	VAL
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	413	GLN
1	A	415	GLU
1	A	416	LYS
1	A	427	LEU
1	A	441	LYS
1	A	444	ARG
1	A	452	ASP
1	A	479	LEU
1	A	526	LYS
1	B	306	VAL
1	B	331	LEU
1	B	377	ASN
1	B	384	GLN
1	B	397	VAL
1	B	398	LYS
1	B	416	LYS
1	B	427	LEU
1	B	434	GLN

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Mol	Chain	Res	Type
1	B	444	ARG
1	B	479	LEU
1	B	487	PHE
1	B	501	LEU
1	B	513	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) 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
1	B	446	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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 [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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%)	0.47	13 (5%) 33 29	45, 72, 114, 131	0
1	B	247/288 (85%)	0.34	7 (2%) 55 50	40, 65, 110, 135	0
All	All	496/576 (86%)	0.40	20 (4%) 42 38	40, 69, 113, 135	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	528	PHE	4.5
1	B	376	LEU	4.2
1	A	337	LYS	4.1
1	A	490	ALA	3.4
1	A	491	LEU	3.3
1	A	300	TRP	3.1
1	B	487	PHE	3.1
1	A	331	LEU	2.7
1	A	354	ALA	2.5
1	B	378	LEU	2.5
1	B	329	PHE	2.5
1	A	487	PHE	2.4
1	A	329	PHE	2.3
1	B	310	SER	2.2
1	B	273	THR	2.2
1	A	482	VAL	2.2
1	A	376	LEU	2.2
1	A	334	SER	2.1
1	A	351	THR	2.1
1	B	331	LEU	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 oligosaccharides 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. 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	B-factors($\text{\AA}^2$)	Q<0.9
3	MG	A	1533	1/1	0.85	0.26	90,90,90,90	0
2	CD	B	1529	1/1	0.97	0.05	117,117,117,117	0
2	CD	A	1529	1/1	0.97	0.04	145,145,145,145	0
2	CD	A	1531	1/1	0.98	0.04	74,74,74,74	0
2	CD	A	1532	1/1	0.98	0.07	114,114,114,114	0
2	CD	A	1530	1/1	0.99	0.04	68,68,68,68	0
2	CD	A	1528	1/1	1.00	0.02	44,44,44,44	0

6.5 Other polymers [i](#)

There are no such residues in this entry.