



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 07:49 PM UTC

PDB ID : 1NLT / pdb_00001nlt
Title : The crystal structure of Hsp40 Ydj1
Authors : Li, J.; Sha, B.
Deposited on : 2003-01-07
Resolution : 2.70 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

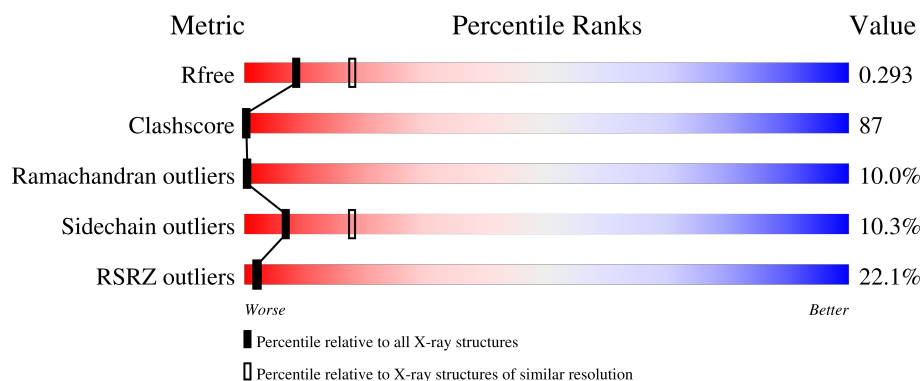
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	248	21% 29% 47% 14% 8%
2	B	7	29% 57% 14%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 1962 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 Mitochondrial protein import protein MAS5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	228	Total	C	N	O	S	0	0	0
			1752	1101	311	327	13			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	335	ASP	PHE	engineered mutation	UNP P25491

- Molecule 2 is a protein called Seven residue peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	7	Total	C	N	O	0	0	0
			61	42	8	11			

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Zn	0	0
			2	2		

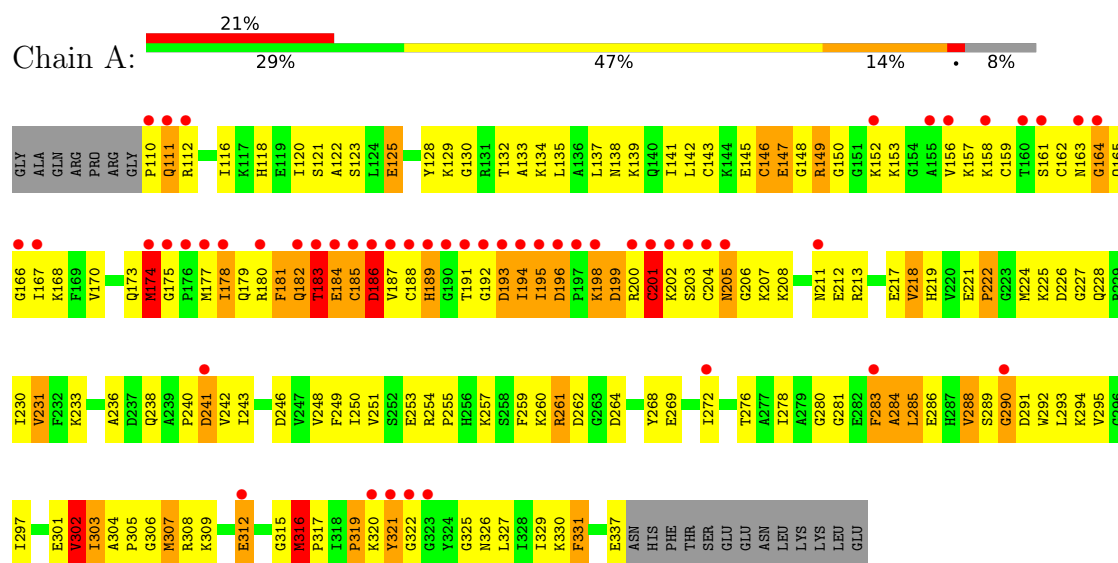
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	142	Total	O	0	0
			142	142		
4	B	5	Total	O	0	0
			5	5		

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: Mitochondrial protein import protein MAS5



- Molecule 2: Seven residue peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	55.02Å 55.02Å 161.87Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	27.50 – 2.70 27.50 – 2.70	Depositor EDS
% Data completeness (in resolution range)	93.6 (27.50-2.70) 93.6 (27.50-2.70)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.50 (at 2.72Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.269 , 0.296 0.267 , 0.293	Depositor DCC
R_{free} test set	796 reflections (10.03%)	wwPDB-VP
Wilson B-factor (Å ²)	39.0	Xtriage
Anisotropy	0.399	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 47.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.035 for -h,-k,l	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	1962	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.96% 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: ZN

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	0.49	0/1781	1.12	14/2385 (0.6%)
2	B	0.39	0/63	0.92	1/85 (1.2%)
All	All	0.49	0/1844	1.11	15/2470 (0.6%)

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	199	ASP	N-CA-C	-7.58	102.34	114.09
1	A	316	MET	CA-C-N	7.46	127.39	119.85
1	A	316	MET	C-N-CA	7.46	127.39	119.85
1	A	243	ILE	N-CA-C	-6.93	101.25	107.56
1	A	231	VAL	N-CA-C	6.93	118.56	108.58
1	A	147	GLU	N-CA-C	-6.89	104.17	112.58
1	A	175	GLY	N-CA-C	-6.79	98.48	112.34
1	A	303	ILE	N-CA-C	6.26	120.80	112.04
1	A	302	VAL	N-CA-C	6.20	122.39	113.16
1	A	149	ARG	N-CA-C	5.75	117.22	111.07
1	A	331	PHE	N-CA-C	5.64	118.09	108.90
1	A	284	ALA	N-CA-C	5.54	117.83	109.23
1	A	198	LYS	N-CA-C	5.38	117.94	107.71
1	A	174	MET	N-CA-C	-5.26	99.60	110.80
2	B	3	LEU	N-CA-C	5.01	116.61	109.14

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	1752	0	1783	296	0
2	B	61	0	57	10	0
3	A	2	0	0	1	0
4	A	142	0	0	152	0
4	B	5	0	0	3	0
All	All	1962	0	1840	301	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 87.

All (301) 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:177:MET:HB2	4:A:477:HOH:O	1.55	1.03
1:A:184:GLU:HA	1:A:189:HIS:HB2	1.43	0.99
1:A:142:LEU:H	1:A:238:GLN:HE21	1.05	0.99
1:A:251:VAL:HB	4:A:378:HOH:O	1.64	0.98
1:A:194:ILE:HG23	1:A:195:ILE:H	1.29	0.97
1:A:195:ILE:HG22	1:A:196:ASP:H	1.31	0.96
1:A:231:VAL:HA	4:A:421:HOH:O	1.68	0.93
1:A:157:LYS:HB2	4:A:426:HOH:O	1.67	0.93
1:A:134:LYS:HB2	4:B:101:HOH:O	1.70	0.92
1:A:191:THR:HG22	4:A:394:HOH:O	1.70	0.92
1:A:142:LEU:H	1:A:238:GLN:NE2	1.69	0.89
1:A:316:MET:HB3	4:A:425:HOH:O	1.70	0.89
1:A:146:CYS:HB3	4:A:355:HOH:O	1.72	0.88
1:A:218:VAL:HG13	4:A:484:HOH:O	1.73	0.88
1:A:178:ILE:HD12	1:A:180:ARG:HE	1.39	0.87
1:A:329:ILE:HG23	4:A:398:HOH:O	1.74	0.87
1:A:183:THR:HG22	4:A:391:HOH:O	1.75	0.87
1:A:148:GLY:HA2	4:A:468:HOH:O	1.73	0.86
1:A:173:GLN:HG3	1:A:177:MET:HG2	1.58	0.85
1:A:142:LEU:N	1:A:238:GLN:HE21	1.72	0.85
1:A:194:ILE:HD11	1:A:200:ARG:HG2	1.56	0.84
1:A:251:VAL:HA	4:A:461:HOH:O	1.76	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:ARG:HB2	4:A:412:HOH:O	1.76	0.84
1:A:200:ARG:HB3	4:A:474:HOH:O	1.79	0.83
1:A:198:LYS:HB3	4:A:424:HOH:O	1.78	0.82
1:A:238:GLN:HG3	4:A:409:HOH:O	1.80	0.81
1:A:289:SER:O	1:A:291:ASP:N	2.15	0.80
1:A:120:ILE:HB	4:A:461:HOH:O	1.83	0.79
1:A:159:CYS:CB	1:A:162:CYS:SG	2.71	0.79
1:A:228:GLN:N	4:A:378:HOH:O	2.13	0.79
1:A:288:VAL:N	4:A:492:HOH:O	2.15	0.79
1:A:143:CYS:SG	1:A:146:CYS:HB2	2.22	0.78
1:A:178:ILE:HD12	1:A:180:ARG:NE	1.99	0.78
1:A:248:VAL:HA	4:A:421:HOH:O	1.83	0.78
1:A:337:GLU:HB2	4:A:360:HOH:O	1.84	0.77
1:A:143:CYS:SG	1:A:204:CYS:HB3	2.25	0.77
1:A:286:GLU:HB2	4:A:466:HOH:O	1.84	0.77
1:A:330:LYS:HG3	4:A:481:HOH:O	1.86	0.76
1:A:133:ALA:HB3	4:A:484:HOH:O	1.84	0.76
1:A:143:CYS:HG	3:A:351:ZN:ZN	0.42	0.75
1:A:195:ILE:HB	4:A:426:HOH:O	1.86	0.75
1:A:236:ALA:HB2	4:A:488:HOH:O	1.87	0.75
1:A:217:GLU:HB2	4:A:404:HOH:O	1.85	0.75
1:A:315:GLY:O	1:A:316:MET:HB2	1.85	0.75
1:A:158:LYS:HD2	4:A:480:HOH:O	1.86	0.74
1:A:194:ILE:HG23	1:A:195:ILE:N	2.02	0.74
1:A:221:GLU:HG3	4:A:463:HOH:O	1.87	0.74
1:A:134:LYS:HB3	4:A:444:HOH:O	1.88	0.73
1:A:268:TYR:HD2	4:A:398:HOH:O	1.72	0.72
1:A:268:TYR:HB3	4:A:398:HOH:O	1.89	0.72
1:A:224:MET:HE1	1:A:230:ILE:HD11	1.72	0.72
1:A:121:SER:HB2	4:A:452:HOH:O	1.89	0.72
1:A:159:CYS:HB3	1:A:162:CYS:SG	2.29	0.72
1:A:143:CYS:SG	1:A:146:CYS:CB	2.77	0.71
4:A:444:HOH:O	2:B:6:ILE:HG13	1.90	0.71
1:A:224:MET:HA	4:A:459:HOH:O	1.90	0.71
1:A:218:VAL:O	4:A:484:HOH:O	2.08	0.71
1:A:224:MET:HE2	1:A:228:GLN:HG2	1.73	0.70
1:A:304:ALA:HA	4:A:407:HOH:O	1.90	0.70
1:A:208:LYS:HB3	4:A:468:HOH:O	1.92	0.70
1:A:286:GLU:N	4:A:466:HOH:O	2.25	0.70
1:A:132:THR:HA	4:A:399:HOH:O	1.91	0.70
1:A:128:TYR:CE2	1:A:289:SER:HB3	2.27	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:CYS:SG	1:A:204:CYS:CB	2.81	0.69
1:A:162:CYS:CB	1:A:185:CYS:SG	2.80	0.69
2:B:6:ILE:HB	4:B:101:HOH:O	1.91	0.69
1:A:249:PHE:C	1:A:250:ILE:HD12	2.18	0.68
1:A:222:PRO:HA	4:A:449:HOH:O	1.91	0.68
1:A:276:THR:HG23	4:A:447:HOH:O	1.95	0.66
1:A:118:HIS:O	1:A:249:PHE:HA	1.95	0.66
1:A:128:TYR:C	4:A:449:HOH:O	2.38	0.66
1:A:204:CYS:SG	4:A:379:HOH:O	2.52	0.66
1:A:141:ILE:HG13	1:A:238:GLN:NE2	2.10	0.66
1:A:181:PHE:C	1:A:181:PHE:CD2	2.73	0.66
1:A:150:GLY:N	4:A:355:HOH:O	2.20	0.66
1:A:264:ASP:HB2	4:A:415:HOH:O	1.94	0.66
1:A:178:ILE:HG21	1:A:180:ARG:NH2	2.10	0.66
1:A:179:GLN:HG2	4:A:455:HOH:O	1.95	0.65
1:A:168:LYS:HE2	4:A:456:HOH:O	1.95	0.65
1:A:194:ILE:HG22	4:A:405:HOH:O	1.96	0.65
1:A:194:ILE:HB	4:A:383:HOH:O	1.96	0.65
1:A:138:ASN:HA	4:A:482:HOH:O	1.95	0.65
1:A:248:VAL:HG13	4:A:421:HOH:O	1.97	0.64
1:A:305:PRO:HD3	4:A:407:HOH:O	1.97	0.64
1:A:195:ILE:HG22	1:A:196:ASP:N	2.08	0.64
1:A:139:LYS:HE3	1:A:141:ILE:HD13	1.80	0.64
1:A:181:PHE:C	1:A:181:PHE:HD2	2.04	0.64
1:A:189:HIS:C	1:A:191:THR:H	2.04	0.64
2:B:7:SER:HA	4:B:111:HOH:O	1.98	0.63
1:A:283:PHE:CD2	1:A:295:VAL:HB	2.34	0.63
1:A:240:PRO:HG3	4:A:468:HOH:O	1.98	0.62
1:A:315:GLY:HA3	1:A:326:ASN:O	1.98	0.62
1:A:268:TYR:CD2	4:A:398:HOH:O	2.49	0.62
2:B:6:ILE:O	2:B:7:SER:HB3	1.99	0.62
1:A:110:PRO:O	1:A:111:GLN:HB2	1.98	0.62
1:A:132:THR:HG23	4:A:399:HOH:O	2.00	0.62
1:A:142:LEU:N	4:A:409:HOH:O	2.33	0.61
1:A:170:VAL:HG11	4:A:431:HOH:O	2.00	0.61
1:A:272:ILE:HD12	1:A:331:PHE:HD2	1.65	0.61
1:A:240:PRO:O	1:A:241:ASP:HB3	2.00	0.61
1:A:206:GLY:N	4:A:474:HOH:O	2.33	0.61
1:A:302:VAL:HA	1:A:307:MET:HE3	1.82	0.61
1:A:233:LYS:HD2	1:A:246:ASP:OD1	2.01	0.61
1:A:240:PRO:O	1:A:241:ASP:CB	2.49	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:ALA:N	4:A:484:HOH:O	2.32	0.60
1:A:162:CYS:HB3	1:A:185:CYS:SG	2.40	0.60
1:A:305:PRO:CD	4:A:407:HOH:O	2.49	0.60
1:A:211:ASN:OD1	4:A:482:HOH:O	2.16	0.60
1:A:134:LYS:NZ	4:A:444:HOH:O	2.34	0.60
1:A:207:LYS:N	4:A:361:HOH:O	2.35	0.60
1:A:135:LEU:HD21	2:B:5:GLU:HG2	1.84	0.60
1:A:264:ASP:O	1:A:325:GLY:HA3	2.01	0.60
1:A:181:PHE:HD2	1:A:182:GLN:N	2.00	0.59
1:A:161:SER:HB3	4:A:456:HOH:O	2.01	0.59
1:A:233:LYS:HA	1:A:246:ASP:OD1	2.02	0.59
1:A:166:GLY:HA3	1:A:189:HIS:HA	1.83	0.59
1:A:130:GLY:N	4:A:449:HOH:O	2.35	0.59
1:A:116:ILE:N	4:A:488:HOH:O	2.36	0.58
1:A:188:CYS:HB2	1:A:192:GLY:HA2	1.85	0.58
1:A:156:VAL:HG12	1:A:194:ILE:HG13	1.86	0.58
1:A:205:ASN:C	4:A:361:HOH:O	2.46	0.58
1:A:264:ASP:O	1:A:315:GLY:O	2.20	0.58
1:A:195:ILE:HD12	4:A:426:HOH:O	2.02	0.58
1:A:201:CYS:C	4:A:357:HOH:O	2.47	0.57
1:A:320:LYS:O	1:A:321:TYR:CB	2.53	0.56
1:A:185:CYS:O	1:A:188:CYS:N	2.39	0.56
1:A:228:GLN:CD	4:A:381:HOH:O	2.48	0.56
1:A:143:CYS:CB	1:A:146:CYS:HB2	2.34	0.56
1:A:183:THR:N	4:A:451:HOH:O	2.39	0.56
1:A:145:GLU:O	1:A:201:CYS:SG	2.64	0.56
1:A:201:CYS:N	4:A:458:HOH:O	2.38	0.56
1:A:194:ILE:HG12	1:A:195:ILE:N	2.19	0.56
1:A:188:CYS:O	1:A:189:HIS:HB3	2.05	0.56
1:A:293:LEU:HD12	4:A:397:HOH:O	2.06	0.55
1:A:303:ILE:O	4:A:407:HOH:O	2.18	0.55
1:A:272:ILE:HD12	1:A:331:PHE:CD2	2.40	0.55
1:A:135:LEU:HD12	1:A:218:VAL:CG1	2.36	0.55
1:A:191:THR:N	4:A:445:HOH:O	2.39	0.55
1:A:240:PRO:O	1:A:241:ASP:OD1	2.24	0.55
1:A:219:HIS:HA	4:A:399:HOH:O	2.05	0.55
1:A:221:GLU:O	1:A:224:MET:HG3	2.07	0.55
1:A:188:CYS:O	1:A:189:HIS:CB	2.55	0.54
1:A:276:THR:CG2	4:A:447:HOH:O	2.54	0.54
1:A:196:ASP:CG	4:A:460:HOH:O	2.50	0.54
1:A:284:ALA:HB2	4:A:396:HOH:O	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:ARG:NH1	4:A:479:HOH:O	2.42	0.53
1:A:194:ILE:HD11	1:A:200:ARG:CG	2.36	0.53
1:A:116:ILE:HG13	4:A:488:HOH:O	2.08	0.53
1:A:133:ALA:CB	4:A:484:HOH:O	2.52	0.52
1:A:195:ILE:CG2	1:A:196:ASP:H	2.14	0.52
1:A:135:LEU:CD2	2:B:5:GLU:HG2	2.39	0.52
1:A:112:ARG:HB3	4:A:469:HOH:O	2.09	0.52
1:A:152:LYS:HE3	4:A:418:HOH:O	2.10	0.52
1:A:162:CYS:O	1:A:164:GLY:N	2.38	0.51
1:A:145:GLU:HB2	1:A:203:SER:OG	2.10	0.51
1:A:272:ILE:HG23	1:A:276:THR:HB	1.93	0.51
1:A:241:ASP:OD1	1:A:242:VAL:HG23	2.10	0.51
1:A:149:ARG:N	4:A:355:HOH:O	2.42	0.51
1:A:135:LEU:HD13	1:A:249:PHE:CE2	2.45	0.51
1:A:226:ASP:HB2	1:A:253:GLU:HB2	1.93	0.51
1:A:189:HIS:CD2	4:A:391:HOH:O	2.64	0.51
1:A:148:GLY:C	4:A:355:HOH:O	2.53	0.51
1:A:142:LEU:HD22	4:A:468:HOH:O	2.11	0.50
1:A:188:CYS:C	4:A:367:HOH:O	2.54	0.50
1:A:183:THR:O	1:A:184:GLU:C	2.55	0.50
1:A:329:ILE:HG12	4:A:398:HOH:O	2.11	0.50
1:A:228:GLN:NE2	4:A:459:HOH:O	2.44	0.50
1:A:292:TRP:CZ3	4:A:466:HOH:O	2.64	0.50
1:A:233:LYS:NZ	4:A:434:HOH:O	2.45	0.49
1:A:281:GLY:C	4:A:447:HOH:O	2.54	0.49
1:A:302:VAL:HG22	1:A:331:PHE:CD1	2.47	0.49
1:A:137:LEU:HD13	2:B:3:LEU:HD21	1.94	0.49
1:A:167:ILE:HG12	1:A:181:PHE:CZ	2.47	0.49
1:A:146:CYS:CA	4:A:412:HOH:O	2.61	0.49
1:A:248:VAL:HG12	1:A:250:ILE:HD11	1.95	0.49
1:A:173:GLN:C	1:A:174:MET:HG3	2.37	0.49
1:A:308:ARG:HG3	4:A:481:HOH:O	2.13	0.49
1:A:152:LYS:N	1:A:199:ASP:O	2.46	0.48
1:A:200:ARG:O	1:A:201:CYS:O	2.30	0.48
1:A:312:GLU:O	1:A:326:ASN:HA	2.14	0.48
1:A:330:LYS:NZ	4:A:382:HOH:O	2.46	0.48
1:A:207:LYS:HG3	4:A:361:HOH:O	2.12	0.48
1:A:138:ASN:CA	4:A:482:HOH:O	2.56	0.48
1:A:138:ASN:HB3	4:A:482:HOH:O	2.13	0.48
1:A:138:ASN:ND2	1:A:213:ARG:HG3	2.29	0.48
1:A:135:LEU:HD12	1:A:218:VAL:HG11	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:CYS:O	1:A:147:GLU:HB2	2.13	0.48
1:A:122:ALA:HA	1:A:254:ARG:HH21	1.78	0.48
1:A:143:CYS:HB3	1:A:146:CYS:HB2	1.96	0.48
1:A:143:CYS:HB3	1:A:146:CYS:CB	2.44	0.47
1:A:146:CYS:HA	4:A:412:HOH:O	2.14	0.47
1:A:185:CYS:O	1:A:186:ASP:C	2.57	0.47
1:A:227:GLY:N	4:A:378:HOH:O	2.47	0.47
1:A:257:LYS:NZ	4:A:387:HOH:O	2.31	0.47
1:A:141:ILE:C	4:A:409:HOH:O	2.58	0.47
1:A:189:HIS:C	1:A:191:THR:N	2.71	0.47
1:A:162:CYS:C	1:A:164:GLY:H	2.21	0.47
1:A:198:LYS:HA	1:A:198:LYS:HD3	1.65	0.47
1:A:250:ILE:HD12	1:A:250:ILE:N	2.29	0.47
1:A:285:LEU:HD13	1:A:285:LEU:N	2.30	0.47
1:A:285:LEU:HD13	1:A:285:LEU:H	1.79	0.47
1:A:222:PRO:HD2	4:A:411:HOH:O	2.14	0.47
1:A:228:GLN:NE2	4:A:486:HOH:O	2.49	0.46
1:A:259:PHE:HB3	4:A:478:HOH:O	2.14	0.46
1:A:276:THR:O	1:A:280:GLY:N	2.42	0.46
1:A:308:ARG:CD	4:A:481:HOH:O	2.62	0.46
1:A:248:VAL:HG12	1:A:250:ILE:CD1	2.45	0.46
1:A:222:PRO:N	4:A:411:HOH:O	2.49	0.46
1:A:135:LEU:HD13	1:A:249:PHE:HE2	1.79	0.46
1:A:168:LYS:HE3	1:A:185:CYS:HA	1.97	0.46
1:A:178:ILE:HD13	1:A:180:ARG:HH21	1.81	0.46
1:A:194:ILE:C	1:A:195:ILE:HG13	2.41	0.46
1:A:315:GLY:O	1:A:316:MET:CB	2.58	0.46
1:A:179:GLN:C	4:A:455:HOH:O	2.58	0.46
1:A:182:GLN:C	4:A:451:HOH:O	2.59	0.46
1:A:261:ARG:HG3	1:A:262:ASP:N	2.30	0.46
1:A:286:GLU:CA	4:A:466:HOH:O	2.61	0.45
1:A:168:LYS:NZ	1:A:186:ASP:OD1	2.49	0.45
1:A:260:LYS:C	4:A:478:HOH:O	2.58	0.45
1:A:110:PRO:O	1:A:111:GLN:CB	2.64	0.45
1:A:288:VAL:HG21	1:A:316:MET:CE	2.47	0.45
1:A:255:PRO:HA	4:A:385:HOH:O	2.16	0.45
1:A:260:LYS:N	4:A:478:HOH:O	2.49	0.45
1:A:290:GLY:N	4:A:437:HOH:O	2.49	0.45
1:A:135:LEU:CD1	1:A:218:VAL:HG11	2.47	0.45
1:A:278:ILE:CA	4:A:440:HOH:O	2.63	0.45
1:A:123:SER:OG	1:A:125:GLU:HG2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:192:GLY:O	1:A:193:ASP:HB3	2.16	0.45
1:A:285:LEU:C	1:A:285:LEU:HD22	2.42	0.45
1:A:289:SER:C	1:A:291:ASP:H	2.25	0.45
1:A:145:GLU:CG	4:A:423:HOH:O	2.64	0.45
1:A:194:ILE:CG2	1:A:195:ILE:N	2.71	0.45
1:A:307:MET:O	1:A:308:ARG:HD2	2.16	0.45
1:A:292:TRP:CZ2	4:A:466:HOH:O	2.68	0.45
1:A:162:CYS:HB2	1:A:165:GLN:O	2.17	0.45
1:A:116:ILE:CB	4:A:488:HOH:O	2.65	0.44
1:A:192:GLY:N	4:A:445:HOH:O	2.50	0.44
1:A:150:GLY:HA3	4:A:476:HOH:O	2.17	0.44
1:A:193:ASP:O	1:A:195:ILE:HG13	2.18	0.44
1:A:222:PRO:CD	4:A:411:HOH:O	2.66	0.44
1:A:236:ALA:CB	4:A:488:HOH:O	2.55	0.44
1:A:289:SER:C	1:A:291:ASP:N	2.74	0.44
1:A:116:ILE:CG1	4:A:488:HOH:O	2.65	0.44
1:A:183:THR:HB	1:A:184:GLU:H	1.48	0.44
1:A:278:ILE:N	4:A:440:HOH:O	2.50	0.44
1:A:294:LYS:HB3	4:A:396:HOH:O	2.18	0.44
1:A:308:ARG:NE	4:A:481:HOH:O	2.50	0.44
1:A:135:LEU:HD12	1:A:218:VAL:HG12	1.99	0.44
1:A:178:ILE:CD1	1:A:180:ARG:HH21	2.30	0.44
1:A:297:ILE:HG22	1:A:309:LYS:CE	2.46	0.44
1:A:173:GLN:O	1:A:174:MET:HG3	2.17	0.44
1:A:194:ILE:O	1:A:195:ILE:HB	2.17	0.44
1:A:305:PRO:HD2	4:A:429:HOH:O	2.17	0.44
4:A:444:HOH:O	2:B:6:ILE:CG1	2.57	0.43
1:A:116:ILE:CD1	2:B:1:GLY:HA3	2.49	0.43
1:A:178:ILE:HD12	1:A:180:ARG:CZ	2.48	0.43
1:A:129:LYS:HB3	1:A:129:LYS:HE2	1.82	0.43
1:A:302:VAL:HG21	1:A:331:PHE:CE2	2.54	0.43
1:A:288:VAL:HG13	4:A:492:HOH:O	2.18	0.43
1:A:167:ILE:HG23	1:A:181:PHE:CE2	2.53	0.43
1:A:281:GLY:HA3	4:A:447:HOH:O	2.19	0.43
1:A:289:SER:O	1:A:290:GLY:C	2.62	0.43
1:A:306:GLY:O	1:A:308:ARG:HD3	2.19	0.43
1:A:218:VAL:C	4:A:399:HOH:O	2.62	0.43
1:A:259:PHE:C	4:A:478:HOH:O	2.61	0.43
1:A:259:PHE:CB	4:A:478:HOH:O	2.67	0.43
1:A:304:ALA:CA	4:A:407:HOH:O	2.59	0.43
1:A:225:LYS:HG3	1:A:226:ASP:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:255:PRO:CA	4:A:385:HOH:O	2.67	0.42
1:A:174:MET:SD	1:A:178:ILE:HD11	2.60	0.42
1:A:224:MET:HE1	1:A:230:ILE:CD1	2.46	0.42
1:A:316:MET:HA	1:A:317:PRO:HD3	1.70	0.42
1:A:189:HIS:C	4:A:445:HOH:O	2.62	0.42
1:A:261:ARG:NH1	4:A:425:HOH:O	2.52	0.42
1:A:307:MET:SD	1:A:308:ARG:N	2.92	0.42
1:A:294:LYS:HA	4:A:396:HOH:O	2.20	0.42
1:A:187:VAL:HG12	1:A:187:VAL:O	2.19	0.42
1:A:149:ARG:NH1	1:A:153:LYS:HE2	2.35	0.42
1:A:321:TYR:CG	1:A:322:GLY:N	2.86	0.42
1:A:204:CYS:O	1:A:206:GLY:N	2.53	0.42
1:A:221:GLU:CG	4:A:463:HOH:O	2.59	0.42
1:A:293:LEU:CD1	4:A:397:HOH:O	2.64	0.42
1:A:116:ILE:HB	4:A:488:HOH:O	2.19	0.41
1:A:173:GLN:HG3	1:A:177:MET:CG	2.37	0.41
1:A:278:ILE:CD1	4:A:440:HOH:O	2.68	0.41
1:A:194:ILE:O	1:A:195:ILE:CB	2.69	0.41
1:A:255:PRO:N	4:A:385:HOH:O	2.53	0.41
1:A:135:LEU:HD23	2:B:5:GLU:HA	2.02	0.41
1:A:228:GLN:CG	4:A:381:HOH:O	2.69	0.41
1:A:141:ILE:HA	1:A:238:GLN:HG3	2.03	0.41
1:A:150:GLY:CA	4:A:476:HOH:O	2.69	0.41
1:A:201:CYS:HB3	1:A:204:CYS:H	1.85	0.41
1:A:224:MET:CA	4:A:459:HOH:O	2.60	0.41
1:A:121:SER:HB3	1:A:254:ARG:HG3	2.01	0.41
1:A:178:ILE:CD1	1:A:180:ARG:NH2	2.84	0.41
1:A:200:ARG:CD	4:A:487:HOH:O	2.69	0.41
1:A:211:ASN:OD1	1:A:212:GLU:N	2.54	0.41
1:A:134:LYS:HB3	1:A:134:LYS:NZ	2.36	0.41
1:A:221:GLU:N	1:A:224:MET:SD	2.94	0.41
1:A:228:GLN:HG3	4:A:381:HOH:O	2.21	0.40
1:A:182:GLN:CA	4:A:451:HOH:O	2.70	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	226/248 (91%)	181 (80%)	22 (10%)	23 (10%)	0	0
2	B	5/7 (71%)	5 (100%)	0	0	100	100
All	All	231/255 (91%)	186 (80%)	22 (10%)	23 (10%)	0	0

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	163	ASN
1	A	189	HIS
1	A	193	ASP
1	A	195	ILE
1	A	201	CYS
1	A	290	GLY
1	A	312	GLU
1	A	321	TYR
1	A	111	GLN
1	A	174	MET
1	A	183	THR
1	A	184	GLU
1	A	186	ASP
1	A	202	LYS
1	A	205	ASN
1	A	241	ASP
1	A	283	PHE
1	A	146	CYS
1	A	194	ILE
1	A	196	ASP
1	A	316	MET
1	A	319	PRO
1	A	164	GLY

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	189/206 (92%)	170 (90%)	19 (10%)	7	18
2	B	6/6 (100%)	5 (83%)	1 (17%)	2	6
All	All	195/212 (92%)	175 (90%)	20 (10%)	7	18

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	125	GLU
1	A	178	ILE
1	A	181	PHE
1	A	182	GLN
1	A	183	THR
1	A	185	CYS
1	A	186	ASP
1	A	201	CYS
1	A	218	VAL
1	A	222	PRO
1	A	261	ARG
1	A	269	GLU
1	A	285	LEU
1	A	288	VAL
1	A	301	GLU
1	A	302	VAL
1	A	307	MET
1	A	319	PRO
1	A	327	LEU
2	B	3	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	138	ASN
1	A	238	GLN

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Mol	Chain	Res	Type
1	A	326	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 2 ligands modelled in this entry, 2 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	228/248 (91%)	0.84	52 (22%) 2 2	15, 34, 115, 135	0
2	B	7/7 (100%)	0.33	0 100 100	20, 24, 37, 47	0
All	All	235/255 (92%)	0.83	52 (22%) 2 2	15, 34, 115, 135	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	187	VAL	6.3
1	A	194	ILE	5.7
1	A	195	ILE	4.3
1	A	177	MET	4.0
1	A	203	SER	3.9
1	A	160	THR	3.8
1	A	202	LYS	3.8
1	A	211	ASN	3.7
1	A	155	ALA	3.6
1	A	111	GLN	3.4
1	A	183	THR	3.4
1	A	166	GLY	3.4
1	A	175	GLY	3.3
1	A	192	GLY	3.3
1	A	322	GLY	3.3
1	A	158	LYS	3.3
1	A	189	HIS	3.2
1	A	190	GLY	3.2
1	A	164	GLY	3.2
1	A	323	GLY	3.1
1	A	178	ILE	3.1
1	A	204	CYS	3.1
1	A	180	ARG	3.0
1	A	241	ASP	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	197	PRO	2.9
1	A	193	ASP	2.9
1	A	321	TYR	2.7
1	A	191	THR	2.7
1	A	174	MET	2.7
1	A	196	ASP	2.7
1	A	312	GLU	2.6
1	A	163	ASN	2.6
1	A	272	ILE	2.6
1	A	156	VAL	2.6
1	A	283	PHE	2.6
1	A	184	GLU	2.5
1	A	290	GLY	2.5
1	A	176	PRO	2.5
1	A	320	LYS	2.5
1	A	110	PRO	2.5
1	A	188	CYS	2.4
1	A	201	CYS	2.4
1	A	112	ARG	2.4
1	A	182	GLN	2.4
1	A	152	LYS	2.3
1	A	200	ARG	2.3
1	A	205	ASN	2.2
1	A	186	ASP	2.2
1	A	167	ILE	2.2
1	A	161	SER	2.1
1	A	185	CYS	2.1
1	A	198	LYS	2.0

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	ZN	A	352	1/1	0.96	0.04	124,124,124,124	0
3	ZN	A	351	1/1	0.98	0.06	76,76,76,76	0

6.5 Other polymers [i](#)

There are no such residues in this entry.