



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:42 PM UTC

PDB ID : 4FCP / pdb_00004fcp
Title : Targetting conserved water molecules: Design of 4-aryl-5-cyanopyrrolo [2,3-d] pyrimidine Hsp90 inhibitors using fragment-based screening and structure-based optimization
Authors : Davies, N.G.M.; Browne, H.; Davies, B.; Foloppe, N.; Geoffrey, S.; Gibbons, B.; Hart, T.; Drysdale, M.; Mansell, H.; Massey, A.; Matassova, N.; Moore, J.D.; Murray, J.; Pratt, R.; Ray, S.; Roughley, S.D.; Jensen, M.R.; Schoepfer, J.; Scriven, K.; Simmonite, H.; Stokes, S.; Surgenor, A.; Webb, P.; Wright, L.; Brough, P.
Deposited on : 2012-05-25
Resolution : 2.00 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

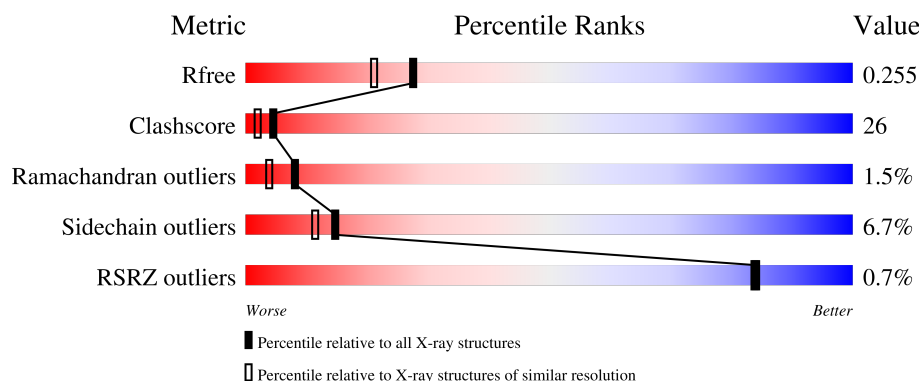
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	236	
1	B	236	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

Ideal geometry (proteins)	: Engh & Huber (2001)
Ideal geometry (DNA, RNA)	: Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	: 2.49

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	42C	A	301	-	X	-	-
2	42C	B	301	-	X	-	-

2 Entry composition [i](#)

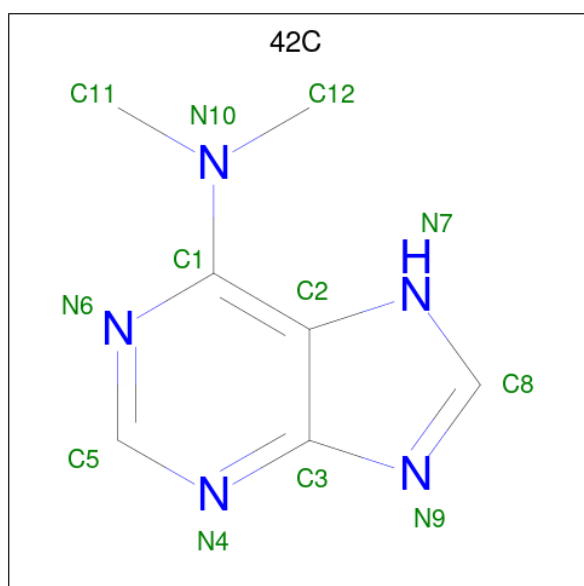
There are 3 unique types of molecules in this entry. The entry contains 3662 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 HSP 90-alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	208	Total	C	N	O	S	0	0	0
			1631	1036	269	321	5			
1	B	208	Total	C	N	O	S	0	0	0
			1631	1036	269	321	5			

- Molecule 2 is N,N-dimethyl-7H-purin-6-amine (CCD ID: 42C) (formula: C₇H₉N₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	0	0
			12	7	5		
2	B	1	Total	C	N	0	0
			12	7	5		

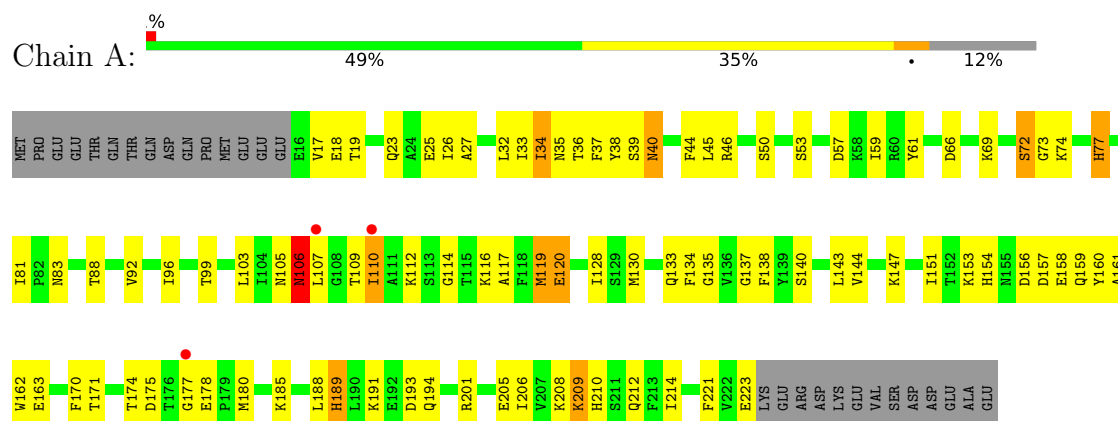
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	179	Total 179	O 179	0	0
3	B	197	Total 197	O 197	0	0

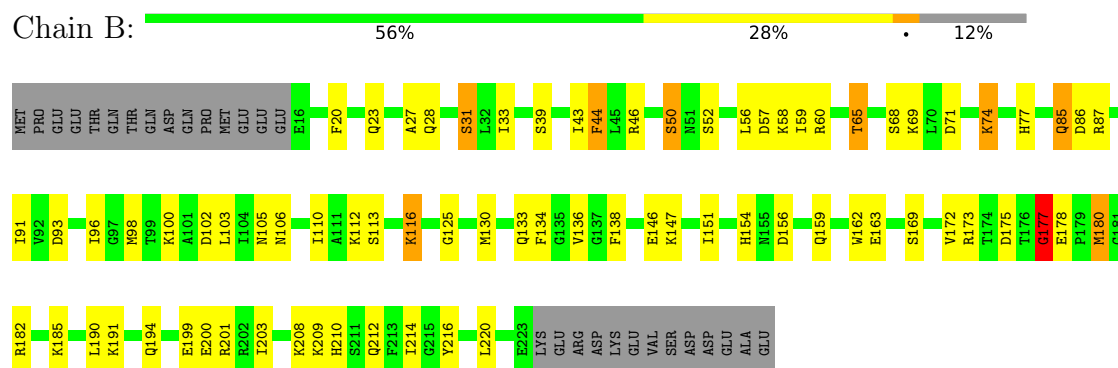
3 Residue-property plots [i](#)

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 HSP 90-alpha



• Molecule 1: Heat shock protein HSP 90-alpha



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	64.17Å 88.69Å 98.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00 20.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	90.9 (20.00-2.00) 90.8 (20.00-2.00)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.11 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.239 , 0.292 0.243 , 0.255	Depositor DCC
R_{free} test set	1779 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	18.7	Xtriage
Anisotropy	0.119	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 45.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3662	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 39.46 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.1633e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: 42C

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	1.25	3/1657 (0.2%)	1.35	7/2235 (0.3%)
1	B	1.32	5/1657 (0.3%)	1.40	10/2235 (0.4%)
All	All	1.29	8/3314 (0.2%)	1.37	17/4470 (0.4%)

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	A	0	1
1	B	0	1
All	All	0	2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	172	VAL	C-O	6.47	1.30	1.23
1	B	154	HIS	CG-CD2	6.31	1.42	1.35
1	B	59	ILE	N-CA	5.67	1.53	1.46
1	A	189	HIS	CG-CD2	5.67	1.42	1.35
1	B	100	LYS	C-O	5.50	1.30	1.24
1	A	77	HIS	CG-CD2	5.47	1.41	1.35
1	A	34	ILE	CA-C	5.16	1.59	1.52
1	B	77	HIS	CG-ND1	-5.07	1.32	1.38

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	106	ASN	CA-C-N	-8.42	108.90	122.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	106	ASN	C-N-CA	-8.42	108.90	122.62
1	B	216	TYR	CA-C-N	8.05	128.54	119.93
1	B	216	TYR	C-N-CA	8.05	128.54	119.93
1	B	31	SER	N-CA-C	8.05	120.77	111.11
1	B	209	LYS	N-CA-C	7.51	119.47	111.28
1	A	35	ASN	N-CA-C	7.04	122.53	113.88
1	A	206	ILE	N-CA-C	7.02	117.11	110.30
1	B	156	ASP	N-CA-C	6.26	120.77	113.20
1	B	44	PHE	N-CA-C	-6.17	104.55	111.28
1	B	125	GLY	N-CA-C	6.10	123.14	115.21
1	B	105	ASN	N-CA-C	5.99	118.30	111.11
1	A	106	ASN	CB-CA-C	5.56	121.49	110.42
1	B	151	ILE	N-CA-C	-5.56	100.41	108.36
1	B	60	ARG	N-CA-C	-5.35	105.35	111.07
1	A	39	SER	N-CA-C	5.14	117.63	111.71
1	A	81	ILE	N-CA-C	5.14	113.24	108.15

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	106	ASN	Peptide
1	B	177	GLY	Peptide

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	1631	0	1630	108	0
1	B	1631	0	1630	64	0
2	A	12	0	9	3	0
2	B	12	0	9	0	0
3	A	179	0	0	62	1
3	B	197	0	0	34	1
All	All	3662	0	3278	171	2

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 26.

All (171) 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:B:27:ALA:HA	3:B:590:HOH:O	1.34	1.24
1:B:116:LYS:HD3	3:B:459:HOH:O	1.35	1.24
1:A:53:SER:HB2	3:A:571:HOH:O	1.36	1.21
1:A:194:GLN:HG3	3:A:570:HOH:O	1.43	1.18
1:A:212:GLN:HG3	3:A:453:HOH:O	1.48	1.12
1:A:105:ASN:HB3	3:A:558:HOH:O	1.48	1.10
1:A:110:ILE:HD11	3:A:573:HOH:O	1.47	1.10
1:A:106:ASN:O	1:A:107:LEU:HD12	1.51	1.10
1:B:212:GLN:HG3	3:B:529:HOH:O	1.50	1.10
1:B:180:MET:HE2	3:B:585:HOH:O	1.60	1.01
1:A:185:LYS:HA	3:A:567:HOH:O	1.66	0.96
1:A:151:ILE:HD12	3:A:563:HOH:O	1.70	0.90
1:A:99:THR:CG2	1:A:154:HIS:HD2	1.85	0.89
1:A:112:LYS:HA	3:A:572:HOH:O	1.72	0.88
1:A:116:LYS:HD3	3:A:503:HOH:O	1.73	0.88
1:B:134:PHE:HE2	3:B:582:HOH:O	1.55	0.87
1:A:46:ARG:HG2	3:A:507:HOH:O	1.78	0.84
1:A:32:LEU:HD12	1:A:119:MET:HE3	1.57	0.84
1:A:99:THR:HG23	1:A:154:HIS:HD2	1.42	0.83
1:A:96:ILE:HG13	3:A:562:HOH:O	1.79	0.83
1:A:50:SER:O	3:A:566:HOH:O	1.96	0.81
1:A:154:HIS:CD2	3:A:550:HOH:O	2.34	0.80
1:A:106:ASN:O	1:A:107:LEU:CD1	2.30	0.79
1:A:161:ALA:HB2	3:A:563:HOH:O	1.83	0.78
1:B:147:LYS:HG2	3:B:586:HOH:O	1.84	0.77
1:A:109:THR:HG21	3:A:527:HOH:O	1.85	0.76
1:A:57:ASP:HB2	3:A:502:HOH:O	1.86	0.76
1:B:147:LYS:HE2	3:B:586:HOH:O	1.85	0.75
1:A:50:SER:C	3:A:566:HOH:O	2.29	0.75
1:B:199:GLU:HG3	3:B:538:HOH:O	1.87	0.75
1:A:151:ILE:HG23	3:A:563:HOH:O	1.87	0.75
1:A:25:GLU:HB2	3:A:572:HOH:O	1.86	0.74
1:A:151:ILE:CD1	3:A:563:HOH:O	2.32	0.74
1:A:99:THR:HG23	1:A:154:HIS:CD2	2.23	0.73
1:A:26:ILE:HD11	3:A:548:HOH:O	1.88	0.72
1:A:134:PHE:HE2	3:A:560:HOH:O	1.72	0.72
1:A:61:TYR:OH	3:A:483:HOH:O	2.08	0.72
1:B:23:GLN:NE2	1:B:110:ILE:H	1.88	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:LYS:HD2	3:A:570:HOH:O	1.91	0.69
1:A:33:ILE:O	3:A:578:HOH:O	2.11	0.68
1:B:74:LYS:HB2	3:B:548:HOH:O	1.96	0.65
1:A:46:ARG:HD3	1:A:210:HIS:HB3	1.78	0.65
1:A:223:GLU:HG2	3:A:403:HOH:O	1.96	0.65
1:A:154:HIS:CG	3:A:550:HOH:O	2.49	0.63
1:B:52:SER:HB3	1:B:93:ASP:HB3	1.81	0.62
1:A:72:SER:HA	3:A:574:HOH:O	1.99	0.62
1:B:96:ILE:C	3:B:583:HOH:O	2.42	0.62
1:B:147:LYS:CE	3:B:586:HOH:O	2.46	0.61
1:A:140:SER:HB2	1:A:143:LEU:HD12	1.80	0.61
1:A:191:LYS:HB2	3:A:570:HOH:O	2.01	0.61
1:B:180:MET:HB2	3:B:585:HOH:O	2.01	0.60
1:A:106:ASN:ND2	3:A:548:HOH:O	2.34	0.60
1:A:153:LYS:HB3	3:A:556:HOH:O	2.02	0.59
1:B:182:ARG:HG2	3:B:583:HOH:O	2.02	0.59
1:B:178:GLU:O	1:B:178:GLU:HG3	2.03	0.59
1:B:182:ARG:CG	3:B:583:HOH:O	2.51	0.59
1:A:154:HIS:CE1	3:A:550:HOH:O	2.55	0.59
1:B:177:GLY:O	3:B:579:HOH:O	2.16	0.58
1:A:209:LYS:HE2	1:A:210:HIS:CE1	2.38	0.58
1:B:130:MET:HB3	3:B:582:HOH:O	2.02	0.58
1:A:66:ASP:HB2	3:A:564:HOH:O	2.02	0.58
1:A:180:MET:HE2	3:A:556:HOH:O	2.03	0.58
1:A:59:ILE:HD13	3:A:562:HOH:O	2.02	0.57
1:B:102:ASP:O	1:B:106:ASN:HB2	2.04	0.57
1:B:113:SER:HB2	1:B:136:VAL:HG13	1.87	0.56
1:B:203:ILE:HG21	1:B:220:LEU:HD13	1.88	0.56
1:A:99:THR:HG22	1:A:154:HIS:HD2	1.68	0.56
1:A:110:ILE:CD1	3:A:445:HOH:O	2.55	0.55
1:A:44:PHE:C	1:A:44:PHE:CD2	2.84	0.55
1:A:185:LYS:CA	3:A:567:HOH:O	2.39	0.55
1:B:212:GLN:NE2	3:B:529:HOH:O	2.38	0.55
1:A:36:THR:N	3:A:578:HOH:O	2.39	0.55
1:B:65:THR:HG23	3:B:533:HOH:O	2.05	0.55
1:B:69:LYS:HE3	3:B:541:HOH:O	2.07	0.55
1:A:120:GLU:HG3	3:A:511:HOH:O	2.08	0.54
1:A:114:GLY:HA2	3:A:465:HOH:O	2.07	0.54
1:A:92:VAL:HG22	3:A:567:HOH:O	2.07	0.54
1:A:107:LEU:O	2:A:301:42C:C11	2.56	0.54
1:A:205:GLU:OE1	1:B:39:SER:OG	2.24	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:ILE:HD13	3:A:445:HOH:O	2.08	0.54
1:A:92:VAL:HG13	1:A:185:LYS:HG2	1.91	0.53
1:B:85:GLN:HE21	1:B:85:GLN:H	1.57	0.53
1:B:68:SER:O	1:B:71:ASP:HB2	2.07	0.53
1:A:103:LEU:HD13	1:A:162:TRP:HB2	1.90	0.53
1:A:159:GLN:NE2	1:A:175:ASP:OD2	2.42	0.52
1:B:46:ARG:HD3	1:B:210:HIS:HB3	1.90	0.52
1:B:180:MET:HE3	1:B:185:LYS:HE3	1.91	0.52
1:A:99:THR:CG2	1:A:154:HIS:CD2	2.78	0.52
1:A:83:ASN:OD1	1:A:83:ASN:C	2.54	0.51
1:B:159:GLN:HG2	1:B:175:ASP:HB3	1.93	0.51
1:B:134:PHE:CE2	3:B:582:HOH:O	2.43	0.50
1:A:106:ASN:O	1:A:107:LEU:CG	2.59	0.50
1:B:98:MET:HB3	1:B:102:ASP:HB2	1.94	0.50
1:A:37:PHE:HB2	1:B:201:ARG:NH1	2.27	0.49
1:B:96:ILE:HA	3:B:583:HOH:O	2.12	0.49
1:A:130:MET:O	1:A:133:GLN:HB2	2.13	0.49
1:A:128:ILE:HD12	3:A:414:HOH:O	2.12	0.49
1:B:130:MET:CB	3:B:582:HOH:O	2.61	0.49
1:A:53:SER:N	3:A:566:HOH:O	2.45	0.49
1:B:163:GLU:HB3	3:B:460:HOH:O	2.11	0.49
1:A:116:LYS:HB2	3:A:543:HOH:O	2.13	0.49
1:B:43:ILE:O	1:B:44:PHE:C	2.55	0.48
1:A:221:PHE:HD1	1:A:221:PHE:N	2.11	0.48
1:A:130:MET:HB3	3:A:560:HOH:O	2.13	0.48
2:A:301:42C:H8	3:A:417:HOH:O	2.12	0.48
1:B:178:GLU:CD	3:B:543:HOH:O	2.57	0.48
1:B:50:SER:HB2	3:B:490:HOH:O	2.13	0.48
1:A:88:THR:HA	1:A:188:LEU:O	2.14	0.48
1:A:147:LYS:HE3	1:A:189:HIS:CD2	2.48	0.48
1:A:73:GLY:HA2	3:A:494:HOH:O	2.13	0.47
1:A:88:THR:HG22	1:A:189:HIS:HA	1.95	0.47
1:B:85:GLN:H	1:B:85:GLN:NE2	2.12	0.47
1:A:221:PHE:N	1:A:221:PHE:CD1	2.81	0.47
1:A:208:LYS:O	1:A:212:GLN:HB3	2.14	0.47
1:B:20:PHE:O	1:B:169:SER:HB2	2.15	0.47
1:A:106:ASN:C	1:A:107:LEU:HD12	2.34	0.47
1:A:26:ILE:HD13	1:A:106:ASN:OD1	2.15	0.46
1:B:146:GLU:HB2	3:B:588:HOH:O	2.14	0.46
1:A:18:GLU:O	1:A:171:THR:HA	2.15	0.46
1:B:178:GLU:HA	3:B:579:HOH:O	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:103:LEU:HD13	1:B:162:TRP:HB2	1.97	0.46
1:A:69:LYS:HE2	3:A:531:HOH:O	2.15	0.46
1:B:43:ILE:C	1:B:43:ILE:HD12	2.40	0.46
1:A:107:LEU:C	1:A:109:THR:H	2.23	0.46
1:A:156:ASP:O	1:A:157:ASP:HB2	2.14	0.46
1:A:133:GLN:HG2	3:A:530:HOH:O	2.15	0.45
1:B:182:ARG:N	3:B:587:HOH:O	2.48	0.45
1:A:194:GLN:CG	3:A:570:HOH:O	2.27	0.45
1:A:160:TYR:CE1	1:A:174:THR:HG22	2.52	0.45
1:A:106:ASN:ND2	3:A:565:HOH:O	2.49	0.45
1:A:210:HIS:HB2	3:A:438:HOH:O	2.17	0.45
1:A:107:LEU:O	2:A:301:42C:H11B	2.17	0.44
1:A:112:LYS:N	3:A:548:HOH:O	2.51	0.44
1:B:28:GLN:O	1:B:31:SER:HB2	2.16	0.44
1:A:44:PHE:CG	1:A:45:LEU:N	2.85	0.44
1:A:77:HIS:CD2	1:A:77:HIS:C	2.95	0.44
1:B:214:ILE:HD13	1:B:214:ILE:HA	1.88	0.44
1:A:92:VAL:HA	3:A:567:HOH:O	2.17	0.44
1:A:135:GLY:HA3	3:A:445:HOH:O	2.18	0.44
1:B:113:SER:HB2	1:B:136:VAL:CG1	2.48	0.44
1:A:44:PHE:CD2	1:A:45:LEU:N	2.86	0.44
1:A:133:GLN:NE2	3:A:499:HOH:O	2.50	0.44
1:B:74:LYS:HE3	3:B:589:HOH:O	2.16	0.44
1:B:159:GLN:OE1	1:B:180:MET:HG3	2.19	0.43
1:B:96:ILE:CA	3:B:583:HOH:O	2.66	0.43
1:A:147:LYS:HD2	1:A:163:GLU:OE2	2.19	0.43
1:B:180:MET:CE	1:B:185:LYS:HG3	2.48	0.43
1:A:140:SER:HB2	1:A:143:LEU:CD1	2.46	0.43
1:A:26:ILE:O	1:A:27:ALA:C	2.61	0.42
1:A:19:THR:HA	1:A:170:PHE:O	2.19	0.42
1:B:190:LEU:HD13	1:B:194:GLN:O	2.19	0.42
1:A:23:GLN:NE2	1:A:106:ASN:HD21	2.16	0.42
1:A:144:VAL:HA	3:A:570:HOH:O	2.19	0.42
1:A:174:THR:HG23	3:A:491:HOH:O	2.20	0.42
1:A:38:TYR:CZ	1:A:40:ASN:HB3	2.55	0.42
1:B:91:ILE:HD12	1:B:91:ILE:N	2.35	0.42
1:B:191:LYS:HE2	3:B:559:HOH:O	2.19	0.42
1:B:52:SER:HB3	1:B:93:ASP:CB	2.50	0.41
1:A:36:THR:HG21	1:A:128:ILE:HD11	2.02	0.41
1:A:25:GLU:HG2	3:A:543:HOH:O	2.20	0.41
1:A:53:SER:HB3	3:A:566:HOH:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:116:LYS:O	1:A:117:ALA:C	2.64	0.41
1:B:212:GLN:CG	3:B:529:HOH:O	2.33	0.41
1:B:182:ARG:CD	3:B:583:HOH:O	2.69	0.41
1:A:107:LEU:C	1:A:109:THR:N	2.79	0.41
1:B:173:ARG:HD3	3:B:537:HOH:O	2.21	0.41
1:B:56:LEU:O	1:B:57:ASP:C	2.65	0.41
1:B:86:ASP:N	1:B:86:ASP:OD1	2.53	0.41
1:B:138:PHE:CD2	1:B:138:PHE:C	3.00	0.40
1:A:137:GLY:O	1:A:138:PHE:C	2.64	0.40
1:A:185:LYS:CB	3:A:567:HOH:O	2.66	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:431:HOH:O	3:A:431:HOH:O[2_555]	2.07	0.13
3:B:423:HOH:O	3:B:578:HOH:O[4_455]	2.09	0.11

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	206/236 (87%)	184 (89%)	18 (9%)	4 (2%)	6	3
1	B	206/236 (87%)	188 (91%)	16 (8%)	2 (1%)	12	8
All	All	412/472 (87%)	372 (90%)	34 (8%)	6 (2%)	8	4

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	40	ASN
1	A	72	SER

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Mol	Chain	Res	Type
1	A	106	ASN
1	A	177	GLY
1	B	133	GLN
1	B	177	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	178/206 (86%)	166 (93%)	12 (7%)	15	11
1	B	178/206 (86%)	166 (93%)	12 (7%)	15	11
All	All	356/412 (86%)	332 (93%)	24 (7%)	15	11

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	VAL
1	A	34	ILE
1	A	74	LYS
1	A	110	ILE
1	A	119	MET
1	A	120	GLU
1	A	158	GLU
1	A	178	GLU
1	A	193	ASP
1	A	201	ARG
1	A	209	LYS
1	A	214	ILE
1	B	33	ILE
1	B	50	SER
1	B	58	LYS
1	B	65	THR
1	B	74	LYS
1	B	85	GLN
1	B	87	ARG

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Mol	Chain	Res	Type
1	B	112	LYS
1	B	116	LYS
1	B	180	MET
1	B	200	GLU
1	B	208	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	ASN
1	A	77	HIS
1	A	79	ASN
1	A	105	ASN
1	A	106	ASN
1	A	133	GLN
1	A	210	HIS
1	B	23	GLN
1	B	79	ASN
1	B	85	GLN
1	B	105	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](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	42C	A	301	-	13,13,13	1.76	3 (23%)	17,18,18	3.71	13 (76%)
2	42C	B	301	-	13,13,13	2.12	4 (30%)	17,18,18	3.47	11 (64%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	42C	A	301	-	-	0/4/4/4	0/2/2/2
2	42C	B	301	-	-	0/4/4/4	0/2/2/2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	301	42C	C3-N9	-5.07	1.34	1.38
2	A	301	42C	C3-N9	-4.50	1.34	1.38
2	B	301	42C	C1-N10	2.93	1.44	1.36
2	A	301	42C	C2-N7	-2.91	1.32	1.37
2	A	301	42C	C1-N6	2.71	1.39	1.34
2	B	301	42C	C5-N6	2.70	1.38	1.33
2	B	301	42C	C2-C1	2.51	1.45	1.40

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	42C	C3-N9-C8	8.85	109.47	103.84
2	A	301	42C	C3-N9-C8	7.28	108.47	103.84
2	A	301	42C	N4-C5-N6	-5.19	120.72	128.58
2	A	301	42C	C2-C3-N9	-4.91	106.55	110.46
2	A	301	42C	C5-N6-C1	4.78	123.50	111.83
2	B	301	42C	C5-N6-C1	4.65	123.18	111.83
2	B	301	42C	C2-N7-C8	4.33	109.89	106.27
2	A	301	42C	N4-C3-N9	4.14	132.63	125.78
2	A	301	42C	N6-C1-N10	4.01	121.74	116.86
2	B	301	42C	C2-C1-N6	-3.95	112.42	121.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	42C	C2-C3-N4	-3.91	120.82	124.21
2	B	301	42C	N7-C8-N9	-3.47	108.59	113.87
2	A	301	42C	C5-N4-C3	3.46	119.62	113.19
2	B	301	42C	C2-C3-N9	-3.30	107.84	110.46
2	B	301	42C	N4-C5-N6	-3.15	123.82	128.58
2	A	301	42C	N7-C8-N9	-3.10	109.15	113.87
2	A	301	42C	C3-C2-N7	3.05	107.33	105.28
2	B	301	42C	C1-C2-C3	3.05	120.67	118.14
2	A	301	42C	C12-N10-C1	2.70	127.39	120.52
2	A	301	42C	C2-N7-C8	2.65	108.49	106.27
2	B	301	42C	N4-C3-N9	2.65	130.17	125.78
2	B	301	42C	C2-C3-N4	-2.60	121.95	124.21
2	A	301	42C	C2-C1-N6	-2.54	115.72	121.68
2	B	301	42C	C5-N4-C3	2.51	117.85	113.19

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	42C	3	0

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	208/236 (88%)	-0.27	3 (1%) 73 73	16, 28, 51, 74	0
1	B	208/236 (88%)	-0.60	0 100 100	14, 22, 40, 54	0
All	All	416/472 (88%)	-0.43	3 (0%) 84 84	14, 25, 48, 74	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	177	GLY	2.2
1	A	110	ILE	2.1
1	A	107	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(Å ²)	Q<0.9
2	42C	A	301	12/12	0.95	0.06	30,33,35,37	0
2	42C	B	301	12/12	0.95	0.06	18,23,25,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.