



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:14 AM GMT

PDB ID : 2C6E
Title : AURORA A KINASE ACTIVATED MUTANT (T287D) IN COMPLEX WITH A 5-AMINOPYRIMIDINYL QUINAZOLINE INHIBITOR
Authors : Pauptit, R.A.; Pannifer, A.D.; Breed, J.; McMiken, H.H.J.; Rowsell, S.; Anderson, M.
Deposited on : 2005-11-09
Resolution : 2.10 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

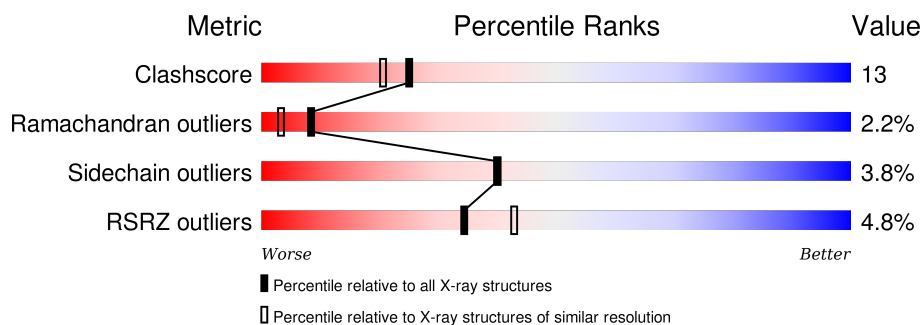
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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))
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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	283	<div> <div>4%</div> <div>72%</div> <div>15%</div> <div>•</div> <div>11%</div> </div>
1	B	283	<div> <div>5%</div> <div>63%</div> <div>19%</div> <div>•</div> <div>13%</div> </div>

2 Entry composition [i](#)

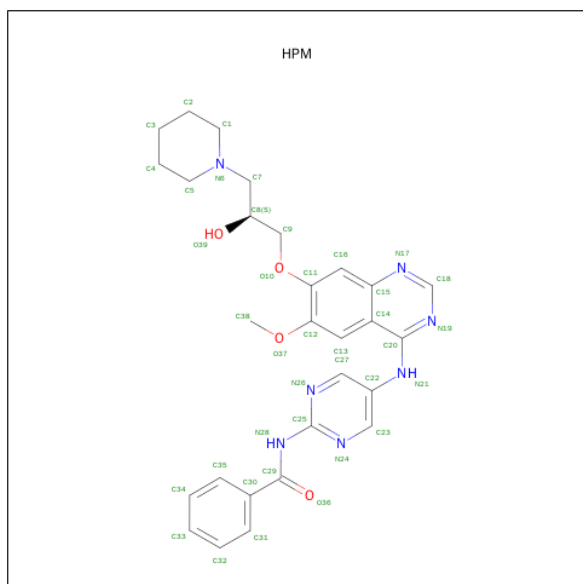
There are 3 unique types of molecules in this entry. The entry contains 4197 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 SERINE/THREONINE-PROTEIN KINASE 6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	251	Total	C	N	O	S	0	0	0
			2040	1317	352	366	5			
1	B	246	Total	C	N	O	S	0	0	0
			1977	1275	339	358	5			

- Molecule 2 is N-{5-[(7-[(2S)-2-HYDROXY-3-PIPERIDIN-1-YLPROPYL]OXY}-6-METHOXYQUINAZOLIN-4-YL)AMINO]PYRIMIDIN-2-YL}BENZAMIDE (three-letter code: HPM) (formula: C₂₈H₃₁N₇O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			39	28	7	4		
2	B	1	Total	C	N	O	0	0
			39	28	7	4		

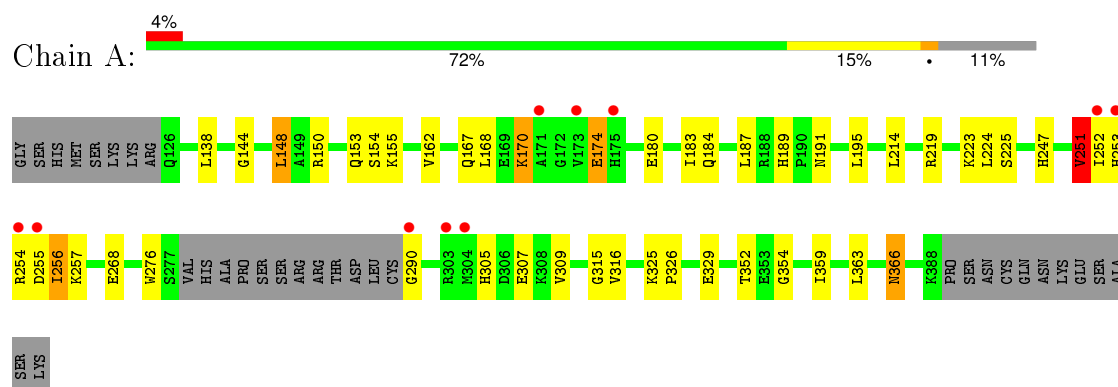
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	56	Total 56	O 56	0	0
3	B	46	Total 46	O 46	0	0

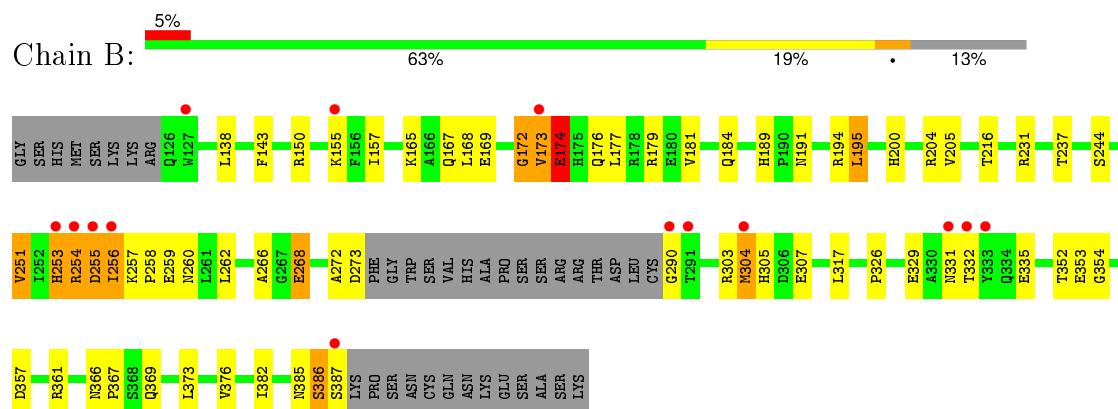
3 Residue-property plots [i](#)

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: SERINE/THREONINE-PROTEIN KINASE 6



• Molecule 1: SERINE/THREONINE-PROTEIN KINASE 6



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	52.60Å 88.42Å 67.83Å 90.00° 90.01° 90.00°	Depositor
Resolution (Å)	53.00 – 2.10 52.60 – 2.10	Depositor EDS
% Data completeness (in resolution range)	(Not available) (53.00-2.10) 72.5 (52.60-2.10)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.01 (at 2.10Å)	Xtriage
Refinement program	CNX2000	Depositor
R, R_{free}	0.200 , 0.270 0.223 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	27.5	Xtriage
Anisotropy	0.245	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 56.9	EDS
Estimated twinning fraction	0.038 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 26287 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4197	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.65% 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

5.1 Standard geometry

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

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	0.38	0/2091	0.69	3/2829 (0.1%)
1	B	0.38	0/2025	0.64	2/2743 (0.1%)
All	All	0.38	0/4116	0.67	5/5572 (0.1%)

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

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	251	VAL	O-C-N	-14.07	100.18	122.70
1	B	251	VAL	O-C-N	-9.01	108.29	122.70
1	A	251	VAL	CA-C-N	7.60	133.92	117.20
1	A	251	VAL	C-N-CA	5.80	136.19	121.70
1	B	251	VAL	CA-C-N	5.35	128.97	117.20

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	251	VAL	Mainchain

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Mol	Chain	Res	Type	Group
1	B	251	VAL	Mainchain

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	2040	0	2017	41	0
1	B	1977	0	1942	61	0
2	A	39	0	31	2	0
2	B	39	0	31	1	0
3	A	56	0	0	2	0
3	B	46	0	0	2	0
All	All	4197	0	4021	103	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 13.

All (103) 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:363:LEU:HD22	3:A:2035:HOH:O	1.57	1.02
1:B:169:GLU:HA	1:B:173:VAL:HA	1.51	0.92
1:B:272:ALA:O	1:B:273:ASP:HB2	1.79	0.82
1:B:189:HIS:HE1	1:B:191:ASN:HD22	1.33	0.77
1:B:184:GLN:HG2	2:B:1388:HPM:H32	1.71	0.71
1:B:165:LYS:O	1:B:169:GLU:HG2	1.90	0.70
1:B:150:ARG:HH11	1:B:155:LYS:NZ	1.90	0.69
1:A:252:ILE:N	1:A:252:ILE:HD12	2.08	0.69
1:B:258:PRO:HD2	1:B:259:GLU:OE2	1.94	0.68
1:B:257:LYS:HB3	1:B:259:GLU:OE2	1.95	0.67
1:A:168:LEU:O	1:A:174:GLU:HG2	1.96	0.66
1:B:143:PHE:CD2	1:B:168:LEU:HD23	2.30	0.66
1:B:150:ARG:HH11	1:B:155:LYS:HZ3	1.42	0.66
1:A:219:ARG:O	1:A:223:LYS:HD3	1.96	0.66
1:B:200:HIS:HB2	1:B:204:ARG:O	1.95	0.65
1:B:257:LYS:HD3	1:B:290:GLY:HA3	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:315:GLY:HA3	1:A:363:LEU:HD21	1.78	0.65
1:B:366:ASN:HD22	1:B:369:GLN:HG3	1.63	0.64
1:B:253:HIS:O	1:B:254:ARG:HB2	1.95	0.64
1:A:155:LYS:O	1:A:155:LYS:HG2	1.98	0.63
1:B:366:ASN:HB3	1:B:369:GLN:HE21	1.63	0.63
1:A:180:GLU:HG2	1:A:276:TRP:HB2	1.79	0.63
1:B:366:ASN:ND2	1:B:369:GLN:HG3	2.16	0.61
1:A:183:ILE:O	1:A:187:LEU:HD13	2.01	0.61
1:B:357:ASP:O	1:B:361:ARG:HG3	2.02	0.60
1:A:325:LYS:HB2	1:A:326:PRO:HD2	1.83	0.60
1:B:326:PRO:HG2	1:B:329:GLU:HB2	1.84	0.60
1:B:143:PHE:HA	1:B:167:GLN:HG2	1.84	0.59
1:A:247:HIS:CE1	1:A:307:GLU:HB2	2.38	0.59
1:B:143:PHE:CD2	1:B:167:GLN:HG3	2.40	0.57
1:A:254:ARG:HG2	1:A:309:VAL:HG21	1.85	0.57
1:B:352:THR:HG22	1:B:354:GLY:H	1.71	0.55
1:A:167:GLN:O	1:A:170:LYS:HG3	2.06	0.55
1:B:382:ILE:HG22	1:B:387:SER:HB2	1.88	0.55
1:A:326:PRO:HG2	1:A:329:GLU:HB2	1.88	0.54
1:B:335:GLU:HA	1:B:335:GLU:OE2	2.08	0.54
1:A:253:HIS:O	1:A:254:ARG:HB2	2.07	0.54
1:A:256:ILE:O	1:A:256:ILE:HG23	2.06	0.54
1:B:366:ASN:HD22	1:B:369:GLN:HE21	1.56	0.53
1:B:169:GLU:HA	1:B:173:VAL:CA	2.34	0.53
1:B:155:LYS:O	1:B:155:LYS:HD3	2.09	0.53
1:B:256:ILE:O	1:B:256:ILE:HG23	2.09	0.52
1:B:179:ARG:HH11	1:B:179:ARG:HG2	1.74	0.52
1:A:138:LEU:HD11	1:A:148:LEU:HB2	1.92	0.52
1:A:153:GLN:HA	1:A:153:GLN:NE2	2.25	0.52
1:A:153:GLN:HA	1:A:153:GLN:HE21	1.74	0.52
1:A:352:THR:HG22	1:A:354:GLY:H	1.75	0.51
1:B:174:GLU:HG3	1:B:176:GLN:HG3	1.93	0.51
1:B:326:PRO:HB2	3:B:2031:HOH:O	2.11	0.51
1:A:150:ARG:HD2	1:A:155:LYS:HG3	1.92	0.50
1:A:254:ARG:CZ	1:A:305:HIS:NE2	2.74	0.50
1:B:259:GLU:OE2	1:B:259:GLU:N	2.35	0.50
1:B:165:LYS:NZ	1:B:200:HIS:HD2	2.10	0.50
1:B:366:ASN:HB3	1:B:369:GLN:NE2	2.27	0.49
1:B:385:ASN:O	1:B:386:SER:HB3	2.11	0.49
1:B:177:LEU:HD13	1:B:177:LEU:C	2.33	0.49
1:B:352:THR:HG23	3:B:2038:HOH:O	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:GLY:HA3	1:A:162:VAL:O	2.13	0.49
1:B:172:GLY:O	1:B:173:VAL:C	2.51	0.48
1:B:255:ASP:OD1	1:B:290:GLY:N	2.45	0.48
1:A:154:SER:O	1:A:155:LYS:HB3	2.13	0.48
1:B:150:ARG:HD3	1:B:155:LYS:NZ	2.29	0.47
1:B:155:LYS:HZ1	1:B:157:ILE:HD11	1.78	0.47
1:B:237:THR:HG21	1:B:387:SER:C	2.35	0.47
1:A:254:ARG:CG	1:A:309:VAL:HG21	2.44	0.47
1:A:189:HIS:CD2	1:A:191:ASN:H	2.32	0.47
1:B:304:MET:SD	1:B:305:HIS:N	2.81	0.47
1:B:189:HIS:CE1	1:B:191:ASN:HD22	2.23	0.46
1:A:359:ILE:O	1:A:363:LEU:HG	2.15	0.46
1:A:366:ASN:C	1:A:366:ASN:HD22	2.19	0.46
1:A:257:LYS:NZ	1:A:290:GLY:HA2	2.31	0.46
1:B:200:HIS:HB3	1:B:205:VAL:HA	1.98	0.45
1:A:315:GLY:HA3	1:A:363:LEU:CD2	2.46	0.45
1:B:257:LYS:NZ	1:B:260:ASN:HD21	2.14	0.45
1:B:332:THR:HG23	1:B:335:GLU:HB2	1.98	0.45
1:B:177:LEU:O	1:B:181:VAL:HG23	2.17	0.45
1:A:316:VAL:HG23	3:A:2035:HOH:O	2.15	0.45
1:A:174:GLU:HA	1:A:174:GLU:OE1	2.16	0.45
1:B:253:HIS:O	1:B:254:ARG:CB	2.65	0.45
1:A:189:HIS:HD2	1:A:191:ASN:H	1.64	0.45
1:A:252:ILE:N	1:A:252:ILE:CD1	2.77	0.45
1:A:184:GLN:HG2	2:A:1389:HPM:H32	2.00	0.44
1:B:216:THR:HG22	1:B:262:LEU:CD2	2.48	0.43
1:B:184:GLN:HB3	1:B:195:LEU:HD12	2.00	0.43
1:B:150:ARG:CD	1:B:155:LYS:HZ3	2.32	0.43
1:A:155:LYS:NZ	1:A:155:LYS:HB3	2.33	0.43
1:B:385:ASN:O	1:B:386:SER:CB	2.66	0.43
1:A:150:ARG:HD2	1:A:155:LYS:CG	2.49	0.43
1:B:257:LYS:HB2	1:B:260:ASN:ND2	2.34	0.43
1:B:216:THR:HG22	1:B:262:LEU:HD23	2.01	0.43
1:B:376:VAL:O	1:B:382:ILE:HD12	2.19	0.42
1:B:352:THR:HG22	1:B:353:GLU:N	2.34	0.42
1:B:155:LYS:O	1:B:155:LYS:CD	2.67	0.42
1:B:244:SER:HA	1:B:373:LEU:HD12	2.02	0.42
1:A:214:LEU:N	1:A:214:LEU:HD12	2.35	0.41
1:B:237:THR:OG1	1:B:387:SER:HA	2.20	0.41
1:A:251:VAL:C	1:A:252:ILE:HD12	2.40	0.41
1:B:266:ALA:N	1:B:268:GLU:OE2	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:363:LEU:HD23	1:A:363:LEU:N	2.36	0.41
1:A:153:GLN:CA	1:A:153:GLN:HE21	2.33	0.41
1:A:224:LEU:O	1:A:225:SER:CB	2.69	0.41
2:A:1389:HPM:H5C2	2:A:1389:HPM:H8	1.95	0.40
1:B:366:ASN:HA	1:B:367:PRO:HD2	1.89	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	247/283 (87%)	235 (95%)	10 (4%)	2 (1%)	24	17
1	B	242/283 (86%)	222 (92%)	11 (4%)	9 (4%)	4	1
All	All	489/566 (86%)	457 (94%)	21 (4%)	11 (2%)	8	3

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	256	ILE
1	B	173	VAL
1	A	174	GLU
1	B	255	ASP
1	B	256	ILE
1	B	303	ARG
1	B	253	HIS
1	B	386	SER
1	B	174	GLU
1	B	254	ARG
1	B	172	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	215/250 (86%)	209 (97%)	6 (3%)	51	55
1	B	207/250 (83%)	197 (95%)	10 (5%)	31	29
All	All	422/500 (84%)	406 (96%)	16 (4%)	40	40

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	148	LEU
1	A	170	LYS
1	A	195	LEU
1	A	255	ASP
1	A	268	GLU
1	A	366	ASN
1	B	138	LEU
1	B	174	GLU
1	B	194	ARG
1	B	195	LEU
1	B	231	ARG
1	B	268	GLU
1	B	304	MET
1	B	307	GLU
1	B	317	LEU
1	B	331	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	153	GLN
1	A	175	HIS
1	A	176	GLN
1	A	189	HIS
1	A	191	ASN
1	A	222	GLN

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Mol	Chain	Res	Type
1	A	334	GLN
1	A	366	ASN
1	A	369	GLN
1	B	126	GLN
1	B	167	GLN
1	B	176	GLN
1	B	191	ASN
1	B	200	HIS
1	B	260	ASN
1	B	331	ASN
1	B	365	HIS
1	B	366	ASN
1	B	369	GLN

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 carbohydrates 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	HPM	A	1389	-	43,43,43	1.98	9 (20%)	57,58,58	2.02	14 (24%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	HPM	B	1388	-	43,43,43	2.07	13 (30%)	57,58,58	2.06	15 (26%)

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	HPM	A	1389	-	-	0/23/31/31	0/5/5/5
2	HPM	B	1388	-	-	0/23/31/31	0/5/5/5

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1388	HPM	C1-N6	2.32	1.53	1.46
2	B	1388	HPM	C23-C22	2.38	1.43	1.39
2	B	1388	HPM	O37-C12	2.40	1.41	1.37
2	B	1388	HPM	C7-C8	2.47	1.57	1.52
2	A	1389	HPM	C7-N6	2.48	1.52	1.47
2	A	1389	HPM	C35-C30	2.59	1.43	1.39
2	B	1388	HPM	C5-N6	2.65	1.54	1.46
2	B	1388	HPM	C35-C30	2.65	1.43	1.39
2	A	1389	HPM	C27-C22	2.82	1.44	1.39
2	B	1388	HPM	C31-C30	2.85	1.44	1.39
2	A	1389	HPM	C31-C30	2.96	1.44	1.39
2	B	1388	HPM	C16-C11	2.99	1.42	1.36
2	B	1388	HPM	C7-N6	3.04	1.53	1.47
2	B	1388	HPM	C13-C12	3.11	1.42	1.36
2	A	1389	HPM	C13-C12	3.13	1.42	1.36
2	A	1389	HPM	C16-C11	3.53	1.42	1.36
2	A	1389	HPM	C14-C15	3.85	1.48	1.42
2	B	1388	HPM	C25-N28	4.12	1.43	1.39
2	B	1388	HPM	C14-C15	4.36	1.49	1.42
2	A	1389	HPM	C25-N28	4.37	1.43	1.39
2	A	1389	HPM	C20-N21	6.06	1.47	1.36
2	B	1388	HPM	C20-N21	6.49	1.47	1.36

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1389	HPM	C22-C23-N24	-5.11	120.56	124.57
2	A	1389	HPM	O37-C12-C13	-5.03	118.51	125.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1388	HPM	C22-C23-N24	-4.57	120.98	124.57
2	B	1388	HPM	O37-C12-C13	-4.49	119.23	125.25
2	A	1389	HPM	N17-C18-N19	-3.96	125.86	128.89
2	A	1389	HPM	C22-C27-N26	-3.50	121.82	124.57
2	A	1389	HPM	O10-C11-C16	-3.36	119.02	125.21
2	B	1388	HPM	N17-C18-N19	-3.36	126.33	128.89
2	B	1388	HPM	C22-C27-N26	-3.09	122.15	124.57
2	B	1388	HPM	O10-C11-C16	-2.70	120.23	125.21
2	A	1389	HPM	C8-C7-N6	2.02	115.82	112.10
2	B	1388	HPM	C4-C5-N6	2.03	114.86	111.47
2	A	1389	HPM	C13-C14-C20	2.27	126.10	124.89
2	B	1388	HPM	C5-N6-C1	2.33	113.94	108.90
2	A	1389	HPM	O10-C11-C12	2.55	121.15	115.78
2	A	1389	HPM	C14-C20-N21	2.61	121.90	119.63
2	B	1388	HPM	C7-C8-C9	2.64	116.60	111.22
2	A	1389	HPM	C9-O10-C11	2.85	123.84	118.03
2	B	1388	HPM	C2-C1-N6	3.05	116.56	111.47
2	B	1388	HPM	C18-N17-C15	3.21	118.15	115.19
2	A	1389	HPM	C18-N17-C15	3.30	118.23	115.19
2	B	1388	HPM	O37-C12-C11	3.60	120.53	115.40
2	B	1388	HPM	C18-N19-C20	4.05	119.40	116.48
2	A	1389	HPM	O37-C12-C11	4.08	121.21	115.40
2	B	1388	HPM	C8-C7-N6	4.38	120.16	112.10
2	A	1389	HPM	C18-N19-C20	4.59	119.78	116.48
2	B	1388	HPM	C25-N28-C29	4.85	137.28	128.18
2	B	1388	HPM	C9-O10-C11	4.90	128.00	118.03
2	A	1389	HPM	C25-N28-C29	5.54	138.59	128.18

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1389	HPM	2	0
2	B	1388	HPM	1	0

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	251/283 (88%)	0.32	10 (3%) 42 51	13, 28, 51, 69	0
1	B	246/283 (86%)	0.36	14 (5%) 27 35	15, 28, 53, 72	0
All	All	497/566 (87%)	0.34	24 (4%) 34 43	13, 28, 53, 72	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	173	VAL	8.6
1	A	173	VAL	7.7
1	A	254	ARG	6.8
1	B	254	ARG	6.2
1	B	387	SER	5.7
1	B	290	GLY	5.4
1	B	255	ASP	5.4
1	A	255	ASP	4.7
1	B	333	TYR	4.6
1	A	290	GLY	4.4
1	A	252	ILE	4.3
1	A	253	HIS	3.4
1	A	171	ALA	3.3
1	B	256	ILE	3.0
1	B	291	THR	2.7
1	B	304	MET	2.4
1	A	303	ARG	2.3
1	B	127	TRP	2.2
1	B	155	LYS	2.1
1	B	331	ASN	2.1
1	A	304	MET	2.1
1	B	332	THR	2.0
1	A	175	HIS	2.0
1	B	253	HIS	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 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(\AA^2)	Q<0.9
2	HPM	B	1388	39/39	0.95	0.14	1.11	16,19,37,37	0
2	HPM	A	1389	39/39	0.96	0.13	-0.03	15,19,28,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.