



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:00 AM UTC

PDB ID : 1RTK / pdb_00001rtk
Title : Crystal Structure Analysis of the Bb segment of Factor B complexed with 4-guanidinobenzoic acid
Authors : Ponnuraj, K.; Xu, Y.; Macon, K.; Moore, D.; Volanakis, J.E.; Narayana, S.V.
Deposited on : 2003-12-10
Resolution : 2.30 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
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

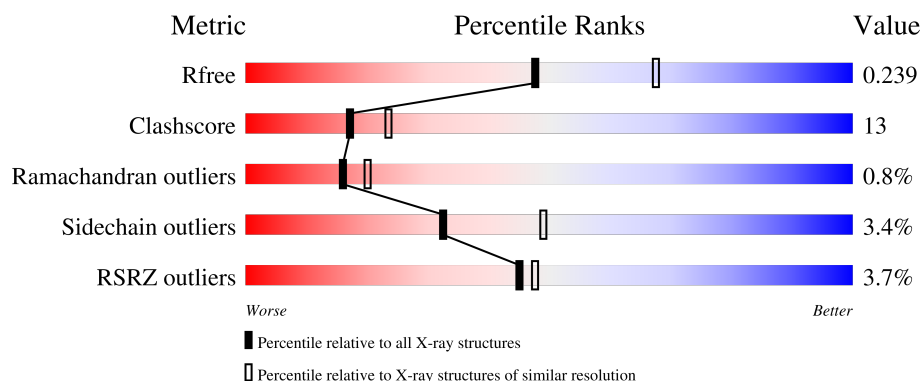
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	497	4% 72% 24% ..

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 4063 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 Complement factor B Bb fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	486	Total	C	N	O	S	0	0	0
			3769	2391	646	711	21			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	267	VAL	CYS	engineered mutation	UNP P00751
A	428	CYS	PHE	engineered mutation	UNP P00751
A	435	CYS	ASN	engineered mutation	UNP P00751

- Molecule 2 is IODIDE ION (CCD ID: IOD) (formula: I).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total I 2 2	0	0

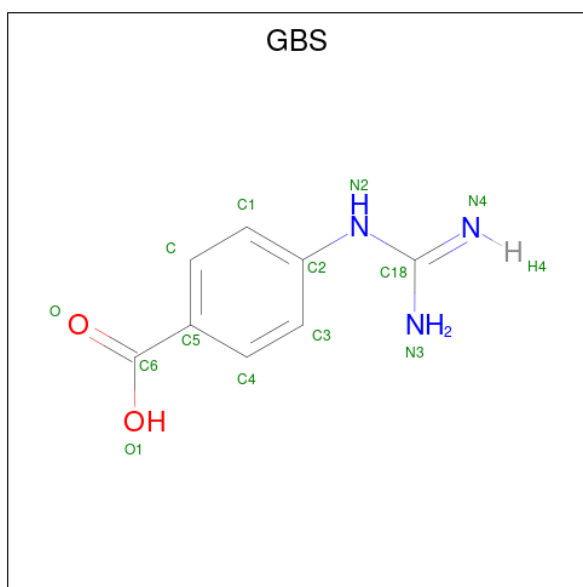
- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	7	Total Na 7 7	0	0

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 5 is 4-carbamimidamidobenzoic acid (CCD ID: GBS) (formula: C₈H₉N₃O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			12	8	3	1		

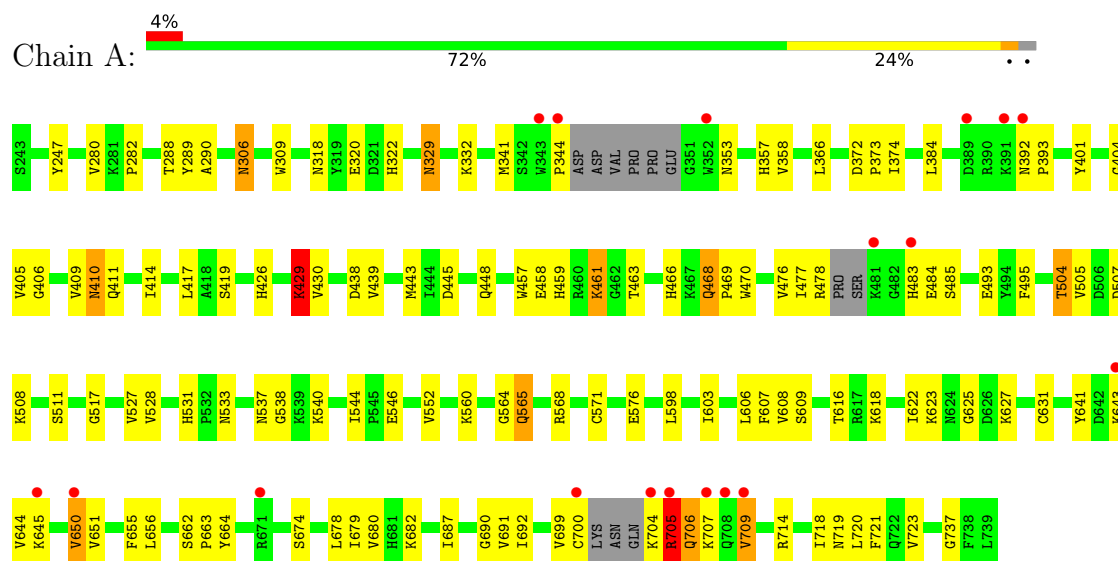
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	272	Total	O	0	0
			272	272		

3 Residue-property plots

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: Complement factor B Bb fragment



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	98.36Å 98.36Å 125.66Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.81 – 2.30 19.81 – 2.30	Depositor EDS
% Data completeness (in resolution range)	95.8 (19.81-2.30) 95.8 (19.81-2.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.72 (at 2.29Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.208 , 0.245 0.201 , 0.239	Depositor DCC
R_{free} test set	861 reflections (2.83%)	wwPDB-VP
Wilson B-factor (Å ²)	27.3	Xtriage
Anisotropy	0.110	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 53.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.049 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4063	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.91% 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: NA, MG, IOD, GBS

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	0.41	0/3843	0.94	17/5204 (0.3%)

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	679	ILE	N-CA-C	7.59	120.42	108.89
1	A	705	ARG	N-CA-C	-7.53	99.88	110.50
1	A	468	GLN	CA-C-N	7.13	127.13	119.28
1	A	468	GLN	C-N-CA	7.13	127.13	119.28
1	A	495	PHE	N-CA-C	7.13	121.13	109.72
1	A	429	LYS	N-CA-C	-7.03	99.10	110.20
1	A	656	LEU	N-CA-C	-6.75	99.22	109.95
1	A	344	PRO	N-CA-CB	6.55	110.21	103.00
1	A	678	LEU	N-CA-C	-6.51	98.12	108.73
1	A	538	GLY	N-CA-C	5.86	121.11	114.67
1	A	655	PHE	N-CA-C	5.81	119.52	110.17
1	A	288	THR	N-CA-C	-5.43	101.27	109.85
1	A	406	GLY	CA-C-N	5.42	125.11	119.64
1	A	406	GLY	C-N-CA	5.42	125.11	119.64
1	A	458	GLU	N-CA-C	5.27	119.69	113.16
1	A	366	LEU	N-CA-C	5.23	117.94	110.68
1	A	517	GLY	N-CA-C	5.18	122.34	114.61

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	3769	0	3690	94	0
2	A	2	0	0	2	0
3	A	7	0	0	0	0
4	A	1	0	0	0	0
5	A	12	0	0	0	0
6	A	272	0	0	3	0
All	All	4063	0	3690	94	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 (94) 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:508:LYS:HA	1:A:508:LYS:HE2	1.40	1.00
1:A:374:ILE:HD13	1:A:417:LEU:HD23	1.58	0.86
1:A:341:MET:HE2	1:A:384:LEU:HD11	1.61	0.80
1:A:459:HIS:HD2	1:A:461:LYS:H	1.30	0.80
1:A:392:ASN:N	1:A:393:PRO:HD3	1.97	0.79
1:A:643:LYS:HG2	6:A:833:HOH:O	1.86	0.74
1:A:306:ASN:C	1:A:306:ASN:HD22	1.97	0.73
1:A:531:HIS:HD2	1:A:533:ASN:H	1.35	0.72
1:A:410:ASN:C	1:A:410:ASN:HD22	1.97	0.72
1:A:623:LYS:HE3	1:A:662:SER:O	1.91	0.71
1:A:463:THR:H	1:A:466:HIS:HD2	1.39	0.69
1:A:576:GLU:HG3	1:A:737:GLY:O	1.94	0.68
1:A:476:VAL:HB	1:A:484:GLU:HB3	1.79	0.65
1:A:417:LEU:HD12	6:A:918:HOH:O	1.98	0.64
1:A:463:THR:H	1:A:466:HIS:CD2	2.15	0.64
1:A:429:LYS:HE3	1:A:429:LYS:HA	1.81	0.62
1:A:373:PRO:HB2	1:A:417:LEU:HD21	1.82	0.61
1:A:247:TYR:OH	1:A:357:HIS:HD2	1.83	0.61
1:A:644:VAL:HG11	1:A:650:VAL:CG1	2.32	0.60
1:A:409:VAL:CG1	1:A:414:ILE:HD12	2.32	0.60
1:A:374:ILE:HD13	1:A:417:LEU:CD2	2.32	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:459:HIS:CD2	1:A:461:LYS:H	2.16	0.59
1:A:329:ASN:ND2	1:A:332:LYS:H	2.01	0.58
1:A:544:ILE:CD1	1:A:650:VAL:HG12	2.36	0.56
1:A:603:ILE:HB	1:A:622:ILE:HB	1.89	0.55
1:A:392:ASN:N	1:A:393:PRO:CD	2.69	0.54
1:A:409:VAL:HG13	1:A:414:ILE:HD12	1.90	0.53
1:A:609:SER:OG	1:A:705:ARG:NH1	2.37	0.53
1:A:663:PRO:HG2	1:A:664:TYR:HD1	1.74	0.53
1:A:508:LYS:HA	1:A:508:LYS:CE	2.25	0.52
1:A:477:ILE:HG23	1:A:511:SER:HB2	1.91	0.52
1:A:616:THR:HG23	1:A:618:LYS:HE3	1.92	0.52
1:A:404:GLY:O	1:A:430:VAL:HG23	2.10	0.52
1:A:306:ASN:ND2	1:A:309:TRP:H	2.08	0.51
1:A:329:ASN:C	1:A:329:ASN:HD22	2.17	0.51
1:A:606:LEU:C	1:A:606:LEU:HD12	2.36	0.51
1:A:405:VAL:HG22	1:A:430:VAL:HG21	1.92	0.50
1:A:445:ASP:OD2	1:A:448:GLN:HG2	2.11	0.50
1:A:700:CYS:HA	1:A:704:LYS:O	2.12	0.50
1:A:358:VAL:HG11	1:A:443:MET:HE3	1.92	0.50
1:A:318:ASN:ND2	1:A:320:GLU:H	2.09	0.49
1:A:438:ASP:OD2	1:A:459:HIS:HE1	1.95	0.49
1:A:289:TYR:HA	1:A:322:HIS:CE1	2.46	0.49
1:A:531:HIS:CD2	1:A:533:ASN:H	2.24	0.49
1:A:691:VAL:HG23	1:A:720:LEU:HD21	1.94	0.49
1:A:699:VAL:O	1:A:706:GLN:HB2	2.12	0.49
1:A:484:GLU:HG3	1:A:485:SER:N	2.26	0.49
1:A:410:ASN:C	1:A:410:ASN:ND2	2.69	0.49
1:A:527:VAL:HG22	1:A:528:VAL:N	2.27	0.49
1:A:571:CYS:HB2	1:A:687:ILE:HD13	1.94	0.49
1:A:341:MET:CE	1:A:384:LEU:HD11	2.40	0.49
1:A:598:LEU:HA	1:A:603:ILE:HD13	1.94	0.49
1:A:564:GLY:O	1:A:568:ARG:HG2	2.13	0.48
1:A:466:HIS:O	1:A:469:PRO:HG3	2.13	0.48
1:A:478:ARG:H	1:A:483:HIS:CD2	2.31	0.48
1:A:537:ASN:OD1	1:A:540:LYS:HE3	2.13	0.48
1:A:372:ASP:OD2	1:A:374:ILE:HG12	2.14	0.47
1:A:318:ASN:HD22	1:A:320:GLU:H	1.62	0.47
1:A:401:TYR:CE1	1:A:439:VAL:HG13	2.49	0.47
1:A:699:VAL:HG11	1:A:709:VAL:HB	1.97	0.47
1:A:477:ILE:HG13	1:A:483:HIS:NE2	2.30	0.47
1:A:609:SER:HG	1:A:705:ARG:HH12	1.61	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:280:VAL:HG12	1:A:282:PRO:HD3	1.97	0.46
1:A:468:GLN:HG2	1:A:608:VAL:HG21	1.97	0.46
1:A:616:THR:CG2	1:A:618:LYS:HE3	2.45	0.46
1:A:680:VAL:HG12	1:A:682:LYS:HG3	1.97	0.46
1:A:466:HIS:HE1	2:A:741:IOD:I	2.68	0.46
1:A:618:LYS:NZ	1:A:705:ARG:CZ	2.79	0.46
1:A:392:ASN:H	1:A:393:PRO:HD3	1.77	0.45
1:A:457:TRP:CH2	1:A:459:HIS:HB3	2.51	0.45
1:A:625:GLY:C	1:A:627:LYS:H	2.24	0.45
1:A:419:SER:HB2	1:A:426:HIS:CD2	2.51	0.45
1:A:650:VAL:HG22	1:A:651:VAL:HG23	1.99	0.45
1:A:457:TRP:CZ3	1:A:459:HIS:HB3	2.52	0.44
1:A:411:GLN:OE1	1:A:429:LYS:NZ	2.48	0.44
1:A:631:CYS:SG	1:A:714:ARG:HD2	2.58	0.44
1:A:493:GLU:CD	1:A:560:LYS:HE2	2.42	0.44
1:A:565:GLN:NE2	1:A:565:GLN:H	2.16	0.44
1:A:663:PRO:HG2	1:A:664:TYR:CD1	2.53	0.44
1:A:531:HIS:HE1	2:A:740:IOD:I	2.71	0.43
1:A:719:ASN:OD1	1:A:721:PHE:HB2	2.18	0.43
1:A:618:LYS:HZ3	1:A:705:ARG:NE	2.16	0.43
1:A:504:THR:HG22	1:A:507:ASP:OD2	2.18	0.43
1:A:607:PHE:HB2	1:A:692:ILE:HD11	2.01	0.43
1:A:531:HIS:HB2	1:A:552:VAL:HG23	2.01	0.42
1:A:720:LEU:HA	1:A:723:VAL:HG22	2.00	0.42
1:A:429:LYS:HA	1:A:429:LYS:CE	2.48	0.42
1:A:544:ILE:HD11	1:A:650:VAL:HG12	2.01	0.42
1:A:641:TYR:CE2	1:A:650:VAL:HG21	2.54	0.42
1:A:318:ASN:HD22	1:A:320:GLU:N	2.18	0.41
1:A:682:LYS:HE3	6:A:912:HOH:O	2.21	0.41
1:A:470:TRP:CD1	1:A:470:TRP:C	2.98	0.41
1:A:690:GLY:HA2	1:A:718:ILE:O	2.21	0.41
1:A:247:TYR:OH	1:A:357:HIS:CD2	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	478/497 (96%)	458 (96%)	16 (3%)	4 (1%)	16	20

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	707	LYS
1	A	645	LYS
1	A	290	ALA
1	A	353	ASN

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	407/439 (93%)	393 (97%)	14 (3%)	32	49

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	306	ASN
1	A	329	ASN
1	A	410	ASN
1	A	429	LYS
1	A	461	LYS
1	A	504	THR
1	A	505	VAL
1	A	546	GLU
1	A	565	GLN
1	A	650	VAL
1	A	674	SER
1	A	705	ARG
1	A	706	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	709	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	270	ASN
1	A	306	ASN
1	A	313	GLN
1	A	315	ASN
1	A	318	ASN
1	A	329	ASN
1	A	357	HIS
1	A	410	ASN
1	A	413	ASN
1	A	415	ASN
1	A	423	ASN
1	A	459	HIS
1	A	466	HIS
1	A	531	HIS
1	A	565	GLN
1	A	591	GLN
1	A	668	ASN
1	A	681	HIS

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 11 ligands modelled in this entry, 10 are monoatomic - leaving 1 for Mogul analysis.

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	$\# Z > 2$
5	GBS	A	750	1	12,12,13	1.99	5 (41%)	14,15,17	0.96	0

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
5	GBS	A	750	1	-	0/6/6/8	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	750	GBS	C18-N2	-2.94	1.34	1.38
5	A	750	GBS	O-C6	2.91	1.31	1.21
5	A	750	GBS	C-C1	2.87	1.43	1.38
5	A	750	GBS	C3-C2	2.78	1.43	1.39
5	A	750	GBS	C-C5	2.04	1.43	1.39

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data [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	486/497 (97%)	0.01	18 (3%) 45 48	15, 30, 46, 63	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	352	TRP	3.9
1	A	650	VAL	3.6
1	A	700	CYS	3.3
1	A	707	LYS	3.3
1	A	389	ASP	3.1
1	A	344	PRO	3.1
1	A	704	LYS	3.1
1	A	391	LYS	3.1
1	A	709	VAL	2.9
1	A	643	LYS	2.8
1	A	645	LYS	2.7
1	A	705	ARG	2.6
1	A	481	LYS	2.5
1	A	708	GLN	2.4
1	A	483	HIS	2.3
1	A	671	ARG	2.3
1	A	343	TRP	2.1
1	A	392	ASN	2.1

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

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

6.3 Carbohydrates [i](#)

There are no 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
4	MG	A	749	1/1	0.92	0.11	32,32,32,32	0
3	NA	A	748	1/1	0.94	0.12	32,32,32,32	0
3	NA	A	747	1/1	0.94	0.18	38,38,38,38	0
5	GBS	A	750	12/13	0.94	0.09	24,26,30,30	0
3	NA	A	744	1/1	0.95	0.16	14,14,14,14	0
3	NA	A	745	1/1	0.95	0.20	7,7,7,7	0
3	NA	A	746	1/1	0.97	0.15	16,16,16,16	0
3	NA	A	743	1/1	0.98	0.15	4,4,4,4	0
3	NA	A	742	1/1	0.98	0.16	4,4,4,4	0
2	IOD	A	741	1/1	0.99	0.04	29,29,29,29	0
2	IOD	A	740	1/1	0.99	0.02	31,31,31,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.