



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:51 PM UTC

PDB ID : 5WJ6 / pdb_00005wj6
Title : Crystal structure of glutaminase C in complex with inhibitor 2-phenyl-N-{5-[4-({5-[(phenylacetyl)amino]-1,3,4-thiadiazol-2-yl}amino)piperidin-1-yl]-1,3,4-thiadiazol-2-yl}acetamide (UPGL-00004)
Authors : Huang, Q.; Cerione, R.A.
Deposited on : 2017-07-21
Resolution : 2.44 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

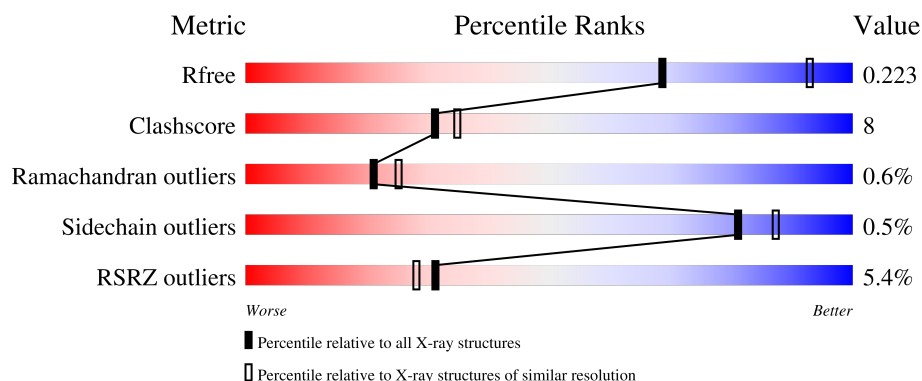
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.44 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	2340 (2.46-2.42)
Clashscore	190562	2400 (2.46-2.42)
Ramachandran outliers	187476	2379 (2.46-2.42)
Sidechain outliers	187428	2379 (2.46-2.42)
RSRZ outliers	180081	2340 (2.46-2.42)

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	539	4% 65% 11% 24%
1	C	539	3% 62% 13% 24%
1	D	539	5% 65% 10% 24%
2	B	539	4% 64% 12% 24%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13503 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 Glutaminase kidney isoform, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	410	Total	C	N	O	S	1	0	0
			3194	2036	540	590	28			
1	C	410	Total	C	N	O	S	1	0	0
			3197	2037	540	592	28			
1	D	410	Total	C	N	O	S	2	0	0
			3195	2035	540	592	28			

There are 39 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	60	MET	-	expression tag	UNP O94925
A	61	ARG	-	expression tag	UNP O94925
A	62	GLY	-	expression tag	UNP O94925
A	63	SER	-	expression tag	UNP O94925
A	64	HIS	-	expression tag	UNP O94925
A	65	HIS	-	expression tag	UNP O94925
A	66	HIS	-	expression tag	UNP O94925
A	67	HIS	-	expression tag	UNP O94925
A	68	HIS	-	expression tag	UNP O94925
A	69	HIS	-	expression tag	UNP O94925
A	70	GLY	-	expression tag	UNP O94925
A	71	SER	-	expression tag	UNP O94925
A	268	ALA	VAL	conflict	UNP O94925
C	60	MET	-	expression tag	UNP O94925
C	61	ARG	-	expression tag	UNP O94925
C	62	GLY	-	expression tag	UNP O94925
C	63	SER	-	expression tag	UNP O94925
C	64	HIS	-	expression tag	UNP O94925
C	65	HIS	-	expression tag	UNP O94925
C	66	HIS	-	expression tag	UNP O94925
C	67	HIS	-	expression tag	UNP O94925
C	68	HIS	-	expression tag	UNP O94925
C	69	HIS	-	expression tag	UNP O94925

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Chain	Residue	Modelled	Actual	Comment	Reference
C	70	GLY	-	expression tag	UNP O94925
C	71	SER	-	expression tag	UNP O94925
C	268	ALA	VAL	conflict	UNP O94925
D	60	MET	-	expression tag	UNP O94925
D	61	ARG	-	expression tag	UNP O94925
D	62	GLY	-	expression tag	UNP O94925
D	63	SER	-	expression tag	UNP O94925
D	64	HIS	-	expression tag	UNP O94925
D	65	HIS	-	expression tag	UNP O94925
D	66	HIS	-	expression tag	UNP O94925
D	67	HIS	-	expression tag	UNP O94925
D	68	HIS	-	expression tag	UNP O94925
D	69	HIS	-	expression tag	UNP O94925
D	70	GLY	-	expression tag	UNP O94925
D	71	SER	-	expression tag	UNP O94925
D	268	ALA	VAL	conflict	UNP O94925

- Molecule 2 is a protein called Glutaminase kidney isoform, mitochondrial.

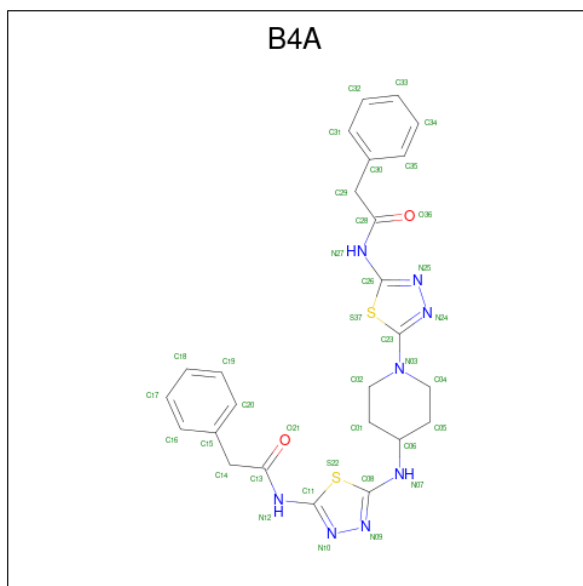
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	410	Total	C	N	O	S	0	0	0
			3193	2035	540	590	28			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	60	MET	-	expression tag	UNP O94925
B	61	ARG	-	expression tag	UNP O94925
B	62	GLY	-	expression tag	UNP O94925
B	63	SER	-	expression tag	UNP O94925
B	64	HIS	-	expression tag	UNP O94925
B	65	HIS	-	expression tag	UNP O94925
B	66	HIS	-	expression tag	UNP O94925
B	67	HIS	-	expression tag	UNP O94925
B	68	HIS	-	expression tag	UNP O94925
B	69	HIS	-	expression tag	UNP O94925
B	70	GLY	-	expression tag	UNP O94925
B	71	SER	-	expression tag	UNP O94925
B	205	ASN	GLN	conflict	UNP O94925
B	268	ALA	VAL	conflict	UNP O94925

- Molecule 3 is 2-phenyl-N-{5-[4-({5-[(phenylacetyl)amino]-1,3,4-thiadiazol-2-yl}amino)piperi

din-1-yl]-1,3,4-thiadiazol-2-yl}acetamide (CCD ID: B4A) (formula: C₂₅H₂₆N₈O₂S₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			37	25	8	2	2		
3	C	1	Total	C	N	O	S	0	0
			37	25	8	2	2		

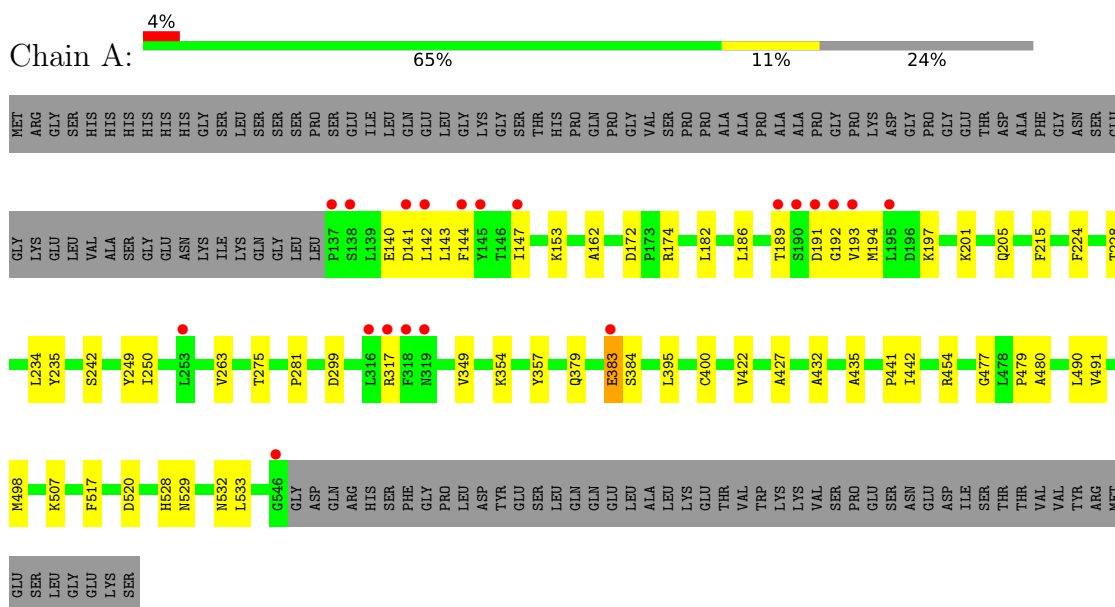
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	138	Total	O	0	0
			138	138		
4	B	161	Total	O	0	0
			161	161		
4	C	171	Total	O	0	0
			171	171		
4	D	180	Total	O	0	0
			180	180		

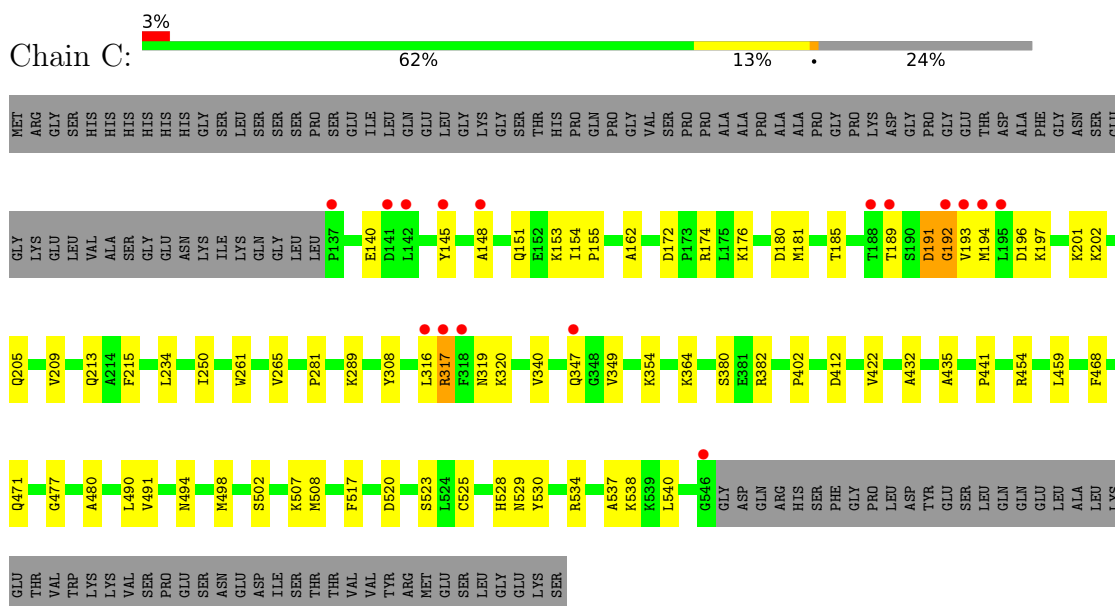
3 Residue-property plots [i](#)

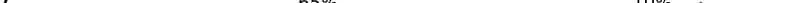
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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.

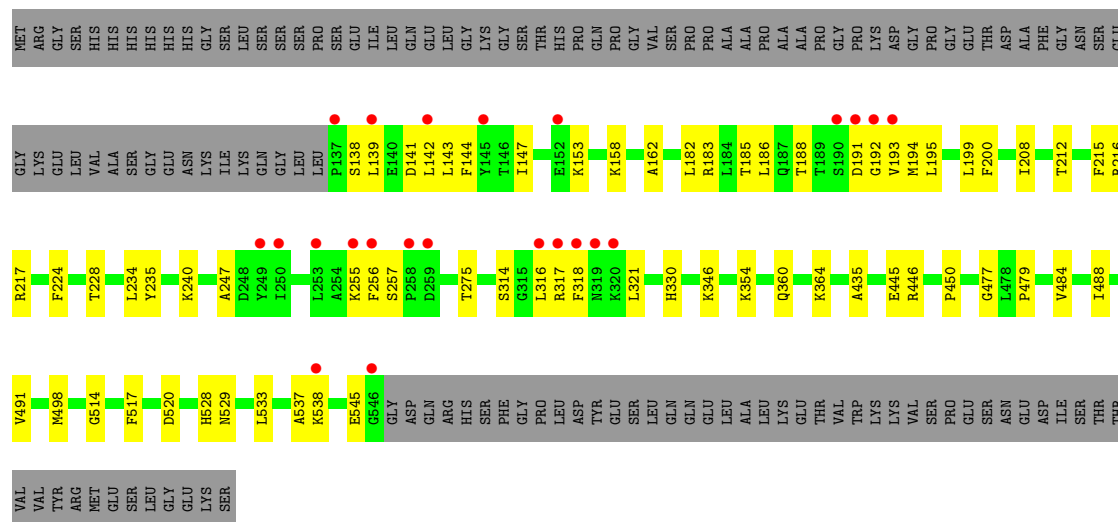
- Molecule 1: Glutaminase kidney isoform, mitochondrial



- Molecule 1: Glutaminase kidney isoform, mitochondrial



- Chain D:  5% 65% 10% 24%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	97.65Å 138.01Å 175.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.05 – 2.44 47.05 – 2.44	Depositor EDS
% Data completeness (in resolution range)	99.3 (47.05-2.44) 99.4 (47.05-2.44)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.36 (at 2.45Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.180 , 0.223 0.182 , 0.223	Depositor DCC
R_{free} test set	2000 reflections (2.28%)	wwPDB-VP
Wilson B-factor (Å ²)	37.7	Xtriage
Anisotropy	0.116	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.9	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	0.96	EDS
Total number of atoms	13503	wwPDB-VP
Average B, all atoms (Å ²)	41.0	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 51.12 % 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 5.8677e-05. 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: B4A

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	0.39	0/3266	0.61	1/4408 (0.0%)
1	C	0.39	0/3269	0.58	1/4412 (0.0%)
1	D	0.39	0/3267	0.57	0/4409
2	B	0.40	0/3265	0.59	0/4407
All	All	0.39	0/13067	0.59	2/17636 (0.0%)

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	C	0	1
All	All	0	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	383	GLU	N-CA-CB	-6.08	100.95	110.44
1	C	537	ALA	CB-CA-C	-5.11	110.28	117.23

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	317	ARG	Peptide
1	C	317	ARG	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	3194	0	3170	42	0
1	C	3197	0	3172	54	0
1	D	3195	0	3165	76	0
2	B	3193	0	3168	70	0
3	A	37	0	0	0	0
3	C	37	0	0	1	0
4	A	138	0	0	2	0
4	B	161	0	0	3	0
4	C	171	0	0	5	0
4	D	180	0	0	6	0
All	All	13503	0	12675	208	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 8.

All (208) 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:529:ASN:ND2	1:C:529:ASN:OD1	1.81	1.13
1:D:153:LYS:HE2	1:D:196:ASP:HB3	1.46	0.97
1:D:152:GLU:CD	1:D:153:LYS:HG3	1.91	0.94
1:D:152:GLU:OE1	1:D:153:LYS:HG3	1.69	0.92
2:B:317:ARG:CB	1:D:318:PHE:HE2	1.86	0.87
2:B:255:LYS:HD3	2:B:256:PHE:H	1.43	0.83
1:D:317:ARG:NH2	1:D:318:PHE:CZ	2.53	0.77
1:C:382:ARG:NH2	1:C:412:ASP:OD1	2.18	0.76
2:B:217:ARG:NH1	2:B:545:GLU:OE1	2.21	0.74
1:D:318:PHE:C	1:D:320:LYS:H	1.94	0.74
1:C:189:THR:OG1	1:C:193:VAL:HG21	1.88	0.74
2:B:147:ILE:O	2:B:158:LYS:NZ	2.21	0.74
2:B:153:LYS:HB2	2:B:194:MET:HE3	1.69	0.73
1:D:375:ASN:ND2	4:D:603:HOH:O	2.22	0.72
1:A:234:LEU:HD22	1:A:520:ASP:HB3	1.72	0.71
2:B:533:LEU:H	1:D:494:ASN:ND2	1.88	0.71
1:D:189:THR:HG22	1:D:191:ASP:H	1.55	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:498:MET:HE1	1:C:517:PHE:HE1	1.54	0.70
1:D:234:LEU:HD22	1:D:520:ASP:HB3	1.73	0.70
1:D:153:LYS:HE2	1:D:196:ASP:CB	2.21	0.69
1:A:201:LYS:O	1:A:205:GLN:HG3	1.93	0.68
1:D:253:LEU:HD12	4:D:665:HOH:O	1.93	0.68
2:B:255:LYS:HA	2:B:255:LYS:HE2	1.77	0.67
1:C:145:TYR:OH	1:C:197:LYS:NZ	2.22	0.67
1:A:235:TYR:CD2	1:A:275:THR:HG23	2.30	0.66
2:B:533:LEU:H	1:D:494:ASN:HD21	1.42	0.66
2:B:318:PHE:CD1	2:B:321:LEU:HD13	2.32	0.65
1:D:318:PHE:O	1:D:320:LYS:N	2.28	0.65
1:D:317:ARG:CZ	1:D:318:PHE:CE2	2.80	0.65
2:B:317:ARG:HB3	1:D:318:PHE:HE2	1.63	0.64
2:B:317:ARG:CB	1:D:318:PHE:CE2	2.77	0.64
1:D:317:ARG:NH2	1:D:318:PHE:CE2	2.66	0.64
2:B:235:TYR:CD2	2:B:275:THR:HG23	2.33	0.63
2:B:234:LEU:HD22	2:B:520:ASP:HB3	1.80	0.63
1:C:234:LEU:HD22	1:C:520:ASP:HB3	1.79	0.63
2:B:318:PHE:CD1	1:D:317:ARG:NH2	2.65	0.63
2:B:143:LEU:HD22	2:B:200:PHE:HZ	1.64	0.62
2:B:183:ARG:HA	2:B:186:LEU:HD12	1.81	0.62
1:C:151:GLN:NE2	4:C:703:HOH:O	2.31	0.62
1:A:533:LEU:H	1:C:494:ASN:ND2	1.98	0.62
1:C:202:LYS:NZ	4:C:704:HOH:O	2.32	0.62
1:D:317:ARG:C	1:D:317:ARG:HD3	2.25	0.61
1:D:313:PRO:HG3	1:D:462:SER:HB2	1.81	0.61
2:B:538:LYS:HD3	2:B:538:LYS:H	1.66	0.61
1:C:181:MET:O	1:C:185:THR:HG23	2.00	0.61
2:B:318:PHE:CE2	1:D:317:ARG:HG3	2.35	0.60
2:B:316:LEU:CD1	1:D:312:GLU:OE1	2.50	0.59
1:D:180:ASP:OD1	1:D:183:ARG:NH1	2.35	0.59
1:A:498:MET:HE1	1:A:517:PHE:CE1	2.37	0.59
1:D:139:LEU:HB3	1:D:140:GLU:OE1	2.03	0.59
2:B:498:MET:HE1	2:B:517:PHE:CE1	2.37	0.58
2:B:318:PHE:CE1	1:D:317:ARG:NH2	2.72	0.58
2:B:317:ARG:HB2	1:D:318:PHE:HE2	1.65	0.57
1:D:188:THR:O	4:D:601:HOH:O	2.17	0.57
2:B:143:LEU:CD2	2:B:200:PHE:HZ	2.16	0.57
1:D:317:ARG:HD3	1:D:318:PHE:N	2.21	0.56
2:B:162:ALA:HB1	2:B:215:PHE:HE1	1.71	0.56
2:B:317:ARG:HG2	1:D:318:PHE:CZ	2.41	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:529:ASN:ND2	4:B:602:HOH:O	2.39	0.56
1:D:317:ARG:CZ	1:D:318:PHE:CZ	2.90	0.55
1:A:153:LYS:HB2	1:A:194:MET:HE2	1.88	0.55
1:D:169:ARG:HG3	1:D:272:ARG:HG3	1.89	0.55
2:B:317:ARG:HB3	1:D:318:PHE:CE2	2.40	0.55
1:A:182:LEU:O	1:A:186:LEU:HD12	2.07	0.54
1:C:162:ALA:HB1	1:C:215:PHE:HE1	1.71	0.54
1:C:364:LYS:NZ	4:C:705:HOH:O	2.32	0.54
1:A:162:ALA:HB1	1:A:215:PHE:HE1	1.72	0.54
1:D:152:GLU:OE2	1:D:153:LYS:CB	2.56	0.54
1:C:201:LYS:O	1:C:205:GLN:HB3	2.08	0.53
1:A:140:GLU:OE1	1:A:140:GLU:N	2.38	0.53
1:A:249:TYR:CD2	1:A:250:ILE:HG23	2.44	0.53
2:B:139:LEU:HD21	2:B:212:THR:OG1	2.09	0.53
2:B:139:LEU:HD22	2:B:208:ILE:HG13	1.90	0.53
1:D:315:GLY:O	1:D:317:ARG:N	2.42	0.53
2:B:346:LYS:HB3	2:B:354:LYS:HG2	1.91	0.53
2:B:138:SER:HB3	2:B:141:ASP:OD2	2.08	0.53
2:B:182:LEU:O	2:B:186:LEU:HD12	2.09	0.53
1:C:480:ALA:HB2	1:C:490:LEU:HD12	1.89	0.53
2:B:217:ARG:HD3	2:B:545:GLU:OE1	2.10	0.52
2:B:316:LEU:HD13	1:D:312:GLU:OE1	2.10	0.52
2:B:316:LEU:HD11	1:D:312:GLU:OE1	2.09	0.52
1:A:172:ASP:OD1	1:A:174:ARG:HD3	2.09	0.51
2:B:318:PHE:CE2	2:B:321:LEU:HD22	2.46	0.51
1:A:479:PRO:HD2	1:A:491:VAL:O	2.11	0.51
1:A:533:LEU:H	1:C:494:ASN:HD21	1.57	0.50
1:C:140:GLU:OE1	1:C:140:GLU:N	2.38	0.50
1:D:152:GLU:OE2	1:D:153:LYS:HG3	2.10	0.50
1:C:507:LYS:N	1:C:507:LYS:HD3	2.26	0.50
1:D:317:ARG:NE	1:D:318:PHE:CE1	2.80	0.49
1:C:265:VAL:HG22	1:C:498:MET:SD	2.52	0.49
1:A:224:PHE:O	1:A:228:THR:HG23	2.13	0.49
1:A:532:ASN:HA	1:C:494:ASN:HD21	1.77	0.49
2:B:317:ARG:HG2	1:D:318:PHE:HZ	1.78	0.49
2:B:538:LYS:H	2:B:538:LYS:CD	2.26	0.49
1:D:152:GLU:OE2	1:D:153:LYS:HB2	2.13	0.49
1:A:477:GLY:O	1:A:529:ASN:HB2	2.13	0.49
2:B:240:LYS:NZ	1:C:523:SER:O	2.46	0.48
2:B:318:PHE:CD2	1:D:317:ARG:NH1	2.80	0.48
2:B:185:THR:HA	2:B:188:THR:OG1	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:318:PHE:C	1:D:320:LYS:N	2.62	0.48
1:A:432:ALA:HB1	1:A:441:PRO:HG2	1.95	0.48
2:B:317:ARG:HB3	1:D:317:ARG:NH2	2.28	0.48
1:C:172:ASP:OD1	1:C:174:ARG:HD3	2.14	0.48
2:B:141:ASP:O	2:B:144:PHE:HB3	2.14	0.47
2:B:537:ALA:HB2	1:D:450:PRO:HG2	1.96	0.47
1:D:152:GLU:OE1	1:D:153:LYS:CG	2.54	0.47
1:A:192:GLY:HA3	1:A:193:VAL:HA	1.57	0.47
1:C:477:GLY:O	1:C:529:ASN:HB2	2.13	0.47
2:B:479:PRO:HD2	2:B:491:VAL:O	2.14	0.47
2:B:255:LYS:HA	2:B:255:LYS:CE	2.44	0.47
1:A:349:VAL:O	1:A:354:LYS:HE3	2.15	0.47
1:C:317:ARG:HA	1:C:319:ASN:CG	2.39	0.46
1:D:148:ALA:O	1:D:150:GLY:N	2.48	0.46
1:D:287:CYS:O	1:D:290:PRO:HD2	2.15	0.46
2:B:224:PHE:O	2:B:228:THR:HG23	2.15	0.46
1:D:152:GLU:OE2	1:D:153:LYS:N	2.48	0.46
2:B:193:VAL:O	2:B:193:VAL:HG23	2.15	0.46
2:B:318:PHE:CZ	2:B:321:LEU:HD22	2.51	0.46
1:A:144:PHE:CE2	1:A:197:LYS:HB2	2.50	0.46
2:B:528:HIS:CD2	1:D:454:ARG:HD2	2.51	0.46
1:D:152:GLU:OE2	1:D:153:LYS:CG	2.64	0.46
2:B:195:LEU:HA	2:B:199:LEU:HD23	1.97	0.46
1:C:432:ALA:HB1	1:C:441:PRO:HG2	1.97	0.46
2:B:317:ARG:CG	1:D:318:PHE:CE2	2.99	0.45
1:C:468:PHE:HD1	1:C:508:MET:HE1	1.82	0.45
1:C:176:LYS:NZ	1:C:180:ASP:OD2	2.49	0.45
1:A:263:VAL:HG13	1:A:498:MET:HE3	1.98	0.45
2:B:435:ALA:HB2	2:B:491:VAL:HG13	1.97	0.45
1:C:525:CYS:HA	1:C:540:LEU:O	2.17	0.45
1:C:347:GLN:NE2	1:C:402:PRO:HG2	2.32	0.45
4:A:735:HOH:O	1:C:471:GLN:HG2	2.17	0.45
1:C:534:ARG:NH2	4:C:715:HOH:O	2.50	0.45
1:D:142:LEU:O	1:D:146:THR:HG23	2.17	0.45
1:D:152:GLU:CD	1:D:153:LYS:CG	2.79	0.44
1:D:259:ASP:OD1	4:D:602:HOH:O	2.21	0.44
1:A:189:THR:O	1:A:189:THR:OG1	2.31	0.44
1:C:250:ILE:HB	1:C:380:SER:HB3	1.99	0.44
1:A:379:GLN:O	1:A:383:GLU:HG3	2.17	0.44
1:A:528:HIS:CG	1:C:454:ARG:HD2	2.53	0.44
2:B:314:SER:HB3	2:B:330:HIS:CG	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:435:ALA:HB2	1:C:491:VAL:HG13	1.98	0.44
1:D:412:ASP:O	1:D:416:GLN:HG3	2.17	0.44
1:A:143:LEU:O	1:A:147:ILE:HD12	2.17	0.44
1:D:176:LYS:NZ	1:D:180:ASP:OD2	2.49	0.44
1:C:234:LEU:CD2	1:C:520:ASP:HB3	2.46	0.44
2:B:317:ARG:CG	1:D:318:PHE:HE2	2.30	0.44
1:C:281:PRO:HA	1:C:422:VAL:O	2.18	0.44
2:B:192:GLY:HA2	2:B:193:VAL:HA	1.83	0.44
1:C:155:PRO:N	1:C:194:MET:HE3	2.33	0.43
1:A:174:ARG:HD2	1:A:442:ILE:C	2.43	0.43
1:C:209:VAL:O	1:C:213:GLN:HG3	2.19	0.43
1:C:349:VAL:O	1:C:354:LYS:HE3	2.17	0.43
1:D:281:PRO:HA	1:D:422:VAL:O	2.18	0.43
2:B:450:PRO:HG2	1:D:537:ALA:HB2	2.00	0.43
1:D:182:LEU:O	1:D:186:LEU:HB2	2.19	0.43
2:B:138:SER:O	2:B:142:LEU:HG	2.19	0.43
1:D:143:LEU:HD11	1:D:147:ILE:HD11	2.00	0.43
1:C:459:LEU:HD23	1:C:459:LEU:HA	1.83	0.43
1:D:498:MET:HE2	1:D:498:MET:HB3	1.68	0.43
1:A:479:PRO:HG3	1:C:530:TYR:CE1	2.54	0.43
1:D:317:ARG:NE	1:D:318:PHE:CZ	2.87	0.43
1:D:261:TRP:CE3	1:D:502:SER:HB2	2.54	0.43
2:B:318:PHE:CD2	1:D:317:ARG:CZ	3.02	0.42
1:A:480:ALA:HB2	1:A:490:LEU:HD12	2.01	0.42
1:C:490:LEU:HD23	1:C:498:MET:CE	2.49	0.42
1:D:250:ILE:HG13	1:D:251:PRO:HD2	2.01	0.42
1:C:316:LEU:HA	1:C:316:LEU:HD23	1.87	0.42
2:B:318:PHE:CG	1:D:317:ARG:NH1	2.88	0.42
2:B:364:LYS:HD3	2:B:445:GLU:OE1	2.19	0.42
1:C:153:LYS:HD3	1:C:196:ASP:HB3	2.01	0.42
1:D:140:GLU:HB3	1:D:208:ILE:HG12	2.01	0.42
1:D:250:ILE:HG23	4:D:609:HOH:O	2.18	0.42
1:D:545:GLU:OE1	1:D:545:GLU:N	2.53	0.42
1:A:299:ASP:OD2	1:A:357:TYR:OH	2.30	0.42
1:A:454:ARG:HD2	1:C:528:HIS:CG	2.55	0.42
1:C:191:ASP:HB3	1:C:192:GLY:H	1.63	0.42
2:B:317:ARG:HB3	1:D:317:ARG:HH22	1.84	0.42
2:B:446:ARG:HD2	4:B:748:HOH:O	2.20	0.42
1:C:289:LYS:NZ	4:C:708:HOH:O	2.41	0.42
1:D:317:ARG:C	1:D:317:ARG:CD	2.92	0.42
1:C:308:TYR:HB3	1:C:340:VAL:CG1	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:247:ALA:HB2	2:B:484:VAL:HG22	2.02	0.41
1:C:261:TRP:CE3	1:C:502:SER:HB2	2.55	0.41
1:D:202:LYS:NZ	4:D:632:HOH:O	2.52	0.41
1:A:142:LEU:HD23	1:A:142:LEU:HA	1.83	0.41
1:A:281:PRO:HA	1:A:422:VAL:O	2.20	0.41
2:B:528:HIS:CG	1:D:454:ARG:HD2	2.55	0.41
2:B:477:GLY:O	2:B:529:ASN:HB2	2.20	0.41
1:A:194:MET:HE2	1:A:194:MET:HB3	1.73	0.41
1:A:422:VAL:HG21	1:A:427:ALA:HB2	2.03	0.41
1:A:249:TYR:CE2	1:A:250:ILE:HG23	2.55	0.41
2:B:216:ARG:NE	2:B:216:ARG:HA	2.35	0.41
2:B:364:LYS:NZ	4:B:612:HOH:O	2.50	0.41
1:C:194:MET:HB3	1:C:194:MET:HE2	1.57	0.41
1:C:347:GLN:HE22	1:C:402:PRO:HG2	1.85	0.41
1:C:148:ALA:HA	1:C:154:ILE:HG12	2.03	0.41
1:C:320:LYS:HD3	3:C:601:B4A:O36	2.20	0.41
1:A:435:ALA:HB2	1:A:491:VAL:HG13	2.03	0.41
1:A:507:LYS:HE2	4:A:723:HOH:O	2.21	0.41
1:A:395:LEU:HB3	1:A:400:CYS:HB2	2.02	0.41
1:A:141:ASP:O	1:A:144:PHE:HB3	2.20	0.40
1:A:153:LYS:CB	1:A:194:MET:HE2	2.51	0.40
1:D:307:ARG:O	1:D:328:LYS:HE2	2.21	0.40
2:B:255:LYS:HD3	2:B:256:PHE:N	2.23	0.40
1:A:498:MET:HE3	1:A:498:MET:HB3	1.87	0.40
2:B:488:ILE:HD12	2:B:514:GLY:HA3	2.03	0.40
1:C:507:LYS:HB2	1:C:507:LYS:HE2	1.86	0.40
1:C:538:LYS:HA	1:C:538:LYS:HD3	1.85	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	408/539 (76%)	399 (98%)	8 (2%)	1 (0%)	43	53
1	C	408/539 (76%)	399 (98%)	7 (2%)	2 (0%)	24	29
1	D	408/539 (76%)	381 (93%)	21 (5%)	6 (2%)	8	7
2	B	408/539 (76%)	400 (98%)	7 (2%)	1 (0%)	43	53
All	All	1632/2156 (76%)	1579 (97%)	43 (3%)	10 (1%)	21	25

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	191	ASP
1	D	191	ASP
1	D	316	LEU
1	D	319	ASN
1	C	191	ASP
1	D	149	GLU
1	D	317	ARG
1	A	191	ASP
1	C	192	GLY
1	D	193	VAL

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	353/461 (77%)	351 (99%)	2 (1%)	78	85
1	C	354/461 (77%)	354 (100%)	0	100	100
1	D	353/461 (77%)	350 (99%)	3 (1%)	73	81
2	B	353/461 (77%)	351 (99%)	2 (1%)	78	85
All	All	1413/1844 (77%)	1406 (100%)	7 (0%)	81	87

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	242	SER

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Mol	Chain	Res	Type
1	A	384	SER
2	B	257	SER
2	B	360	GLN
1	D	257	SER
1	D	259	ASP
1	D	529	ASN

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	187	GLN
1	A	271	GLN
1	A	335	ASN
1	A	388	ASN
1	A	516	HIS
2	B	187	GLN
2	B	205	ASN
2	B	347	GLN
2	B	388	ASN
2	B	516	HIS
1	C	241	GLN
1	C	330	HIS
1	C	347	GLN
1	C	368	ASN
1	C	494	ASN
1	C	516	HIS
1	D	494	ASN
1	D	516	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

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
3	B4A	C	601	-	41,41,41	2.20	14 (34%)	51,55,55	3.51	25 (49%)
3	B4A	A	601	-	41,41,41	2.29	13 (31%)	51,55,55	3.50	25 (49%)

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
3	B4A	C	601	-	-	5/24/34/34	0/5/5/5
3	B4A	A	601	-	-	4/24/34/34	0/5/5/5

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	601	B4A	C08-N07	8.53	1.47	1.34
3	A	601	B4A	C08-N07	8.42	1.47	1.34
3	A	601	B4A	C13-N12	4.44	1.45	1.37
3	A	601	B4A	C28-N27	4.17	1.44	1.37
3	C	601	B4A	C13-N12	3.90	1.44	1.37
3	C	601	B4A	C28-N27	3.81	1.44	1.37
3	A	601	B4A	C26-N27	3.32	1.43	1.38
3	A	601	B4A	C23-N03	3.19	1.41	1.34
3	A	601	B4A	C11-N12	3.14	1.42	1.38
3	C	601	B4A	C11-N12	2.84	1.42	1.38
3	C	601	B4A	N25-N24	-2.80	1.33	1.39
3	C	601	B4A	C23-N03	2.79	1.40	1.34
3	A	601	B4A	C06-N07	-2.74	1.41	1.46
3	A	601	B4A	C08-N09	2.70	1.35	1.31
3	C	601	B4A	C26-N27	2.69	1.42	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	601	B4A	C14-C13	2.59	1.57	1.52
3	A	601	B4A	N25-N24	-2.58	1.34	1.39
3	A	601	B4A	C14-C13	2.40	1.56	1.52
3	C	601	B4A	C06-N07	-2.34	1.42	1.46
3	A	601	B4A	C26-N25	2.34	1.34	1.31
3	C	601	B4A	N10-N09	-2.30	1.34	1.39
3	A	601	B4A	C11-N10	2.29	1.34	1.31
3	C	601	B4A	C01-C06	-2.25	1.46	1.52
3	C	601	B4A	C11-N10	2.19	1.34	1.31
3	C	601	B4A	C34-C33	2.16	1.43	1.38
3	C	601	B4A	C08-N09	2.15	1.34	1.31
3	A	601	B4A	C01-C06	-2.12	1.47	1.52

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	601	B4A	C06-N07-C08	-10.91	105.34	124.48
3	A	601	B4A	C06-N07-C08	-10.23	106.53	124.48
3	A	601	B4A	S22-C11-N10	-8.33	105.22	114.52
3	C	601	B4A	S22-C11-N10	-8.19	105.38	114.52
3	A	601	B4A	S22-C08-N09	-7.28	106.39	114.52
3	C	601	B4A	S22-C08-N09	-7.24	106.44	114.52
3	A	601	B4A	S37-C26-N25	-7.17	106.52	114.52
3	C	601	B4A	S37-C26-N25	-6.66	107.09	114.52
3	A	601	B4A	N03-C23-N24	6.35	129.19	116.91
3	A	601	B4A	C14-C13-N12	5.46	121.06	114.40
3	C	601	B4A	N03-C23-N24	5.45	127.45	116.91
3	A	601	B4A	N12-C11-N10	5.17	125.43	120.81
3	C	601	B4A	C04-C05-C06	-5.14	102.55	110.67
3	A	601	B4A	C23-N24-N25	5.09	117.12	107.08
3	C	601	B4A	C23-N24-N25	5.07	117.07	107.08
3	C	601	B4A	C14-C13-N12	4.83	120.29	114.40
3	A	601	B4A	S37-C23-N24	-4.82	106.10	114.44
3	C	601	B4A	S22-C11-N12	4.61	129.90	123.72
3	C	601	B4A	S37-C23-N24	-4.45	106.75	114.44
3	C	601	B4A	N12-C11-N10	4.37	124.72	120.81
3	C	601	B4A	C29-C28-N27	4.35	119.70	114.40
3	C	601	B4A	S37-C26-N27	4.34	129.53	123.72
3	A	601	B4A	S22-C11-N12	4.20	129.35	123.72
3	A	601	B4A	C04-C05-C06	-4.10	104.19	110.67
3	C	601	B4A	S37-C23-N03	4.08	125.77	120.31
3	A	601	B4A	N27-C26-N25	3.97	124.36	120.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	601	B4A	C04-N03-C23	-3.96	112.05	122.09
3	A	601	B4A	S37-C26-N27	3.94	128.99	123.72
3	A	601	B4A	C11-N10-N09	3.80	116.86	112.02
3	A	601	B4A	C29-C30-C35	-3.73	115.41	120.89
3	A	601	B4A	C05-C06-N07	-3.71	103.17	110.57
3	C	601	B4A	C11-N10-N09	3.68	116.71	112.02
3	C	601	B4A	C08-N09-N10	3.45	116.41	112.02
3	A	601	B4A	C29-C28-N27	3.32	118.45	114.40
3	A	601	B4A	S37-C23-N03	3.28	124.69	120.31
3	C	601	B4A	C01-C06-N07	3.23	117.03	110.57
3	C	601	B4A	C05-C06-N07	-3.22	104.14	110.57
3	A	601	B4A	C08-N09-N10	3.18	116.06	112.02
3	A	601	B4A	C04-N03-C23	-3.15	114.11	122.09
3	C	601	B4A	C05-C04-N03	-2.87	105.04	110.66
3	C	601	B4A	N27-C26-N25	2.70	123.23	120.81
3	A	601	B4A	C26-N25-N24	2.66	115.40	112.02
3	A	601	B4A	C30-C29-C28	2.63	120.18	112.33
3	C	601	B4A	C26-N25-N24	2.58	115.30	112.02
3	A	601	B4A	C02-C01-C06	-2.56	106.63	110.67
3	C	601	B4A	C35-C30-C31	2.56	122.03	118.23
3	C	601	B4A	C02-C01-C06	-2.50	106.73	110.67
3	A	601	B4A	C29-C30-C31	2.49	124.55	120.89
3	A	601	B4A	C01-C06-N07	2.38	115.32	110.57
3	C	601	B4A	C02-N03-C23	-2.07	116.85	122.09

There are no chirality outliers.

All (9) torsion outliers are listed below:

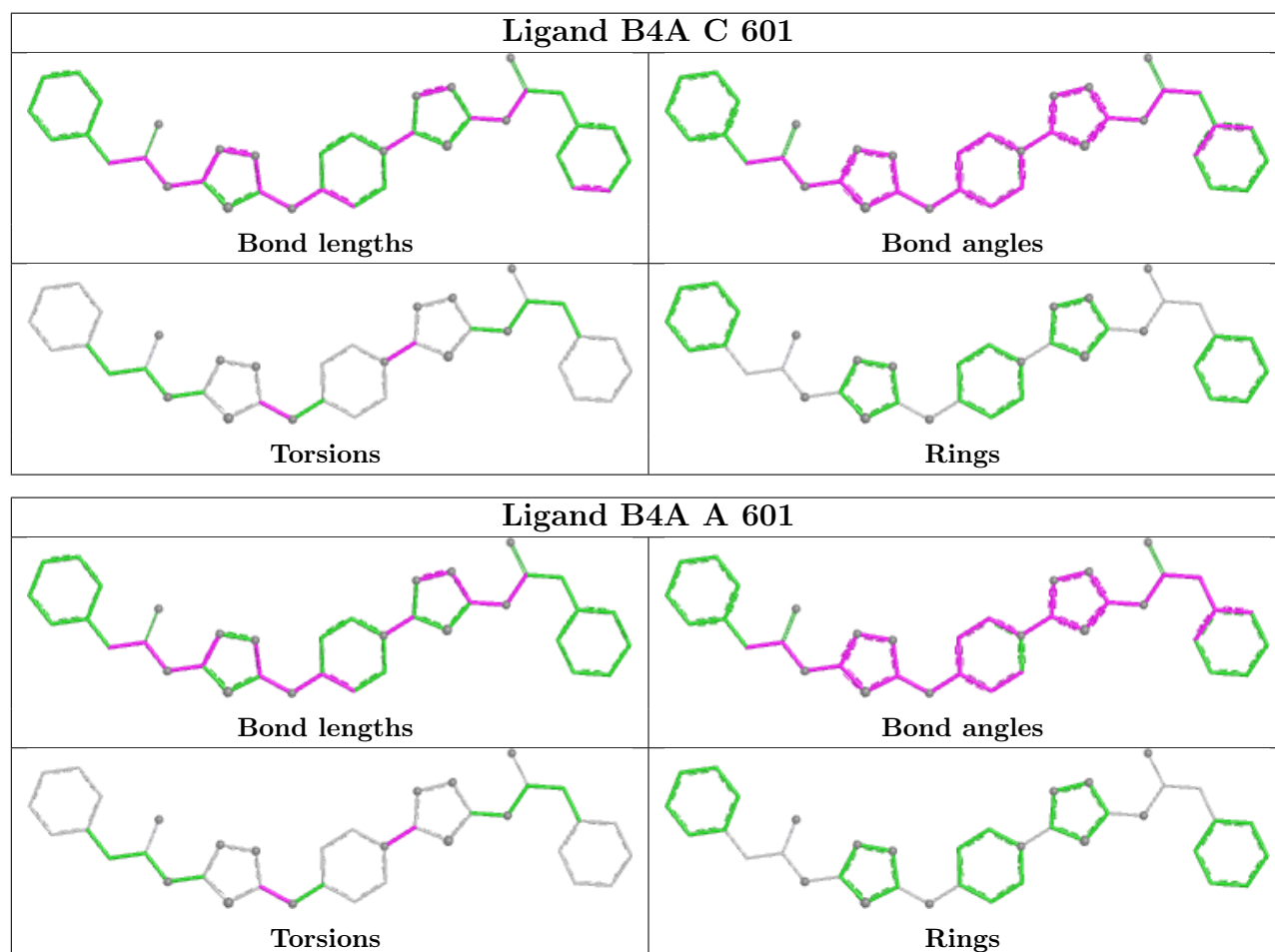
Mol	Chain	Res	Type	Atoms
3	A	601	B4A	N24-C23-N03-C04
3	A	601	B4A	S37-C23-N03-C04
3	A	601	B4A	S22-C08-N07-C06
3	C	601	B4A	N24-C23-N03-C04
3	C	601	B4A	S37-C23-N03-C04
3	C	601	B4A	N09-C08-N07-C06
3	C	601	B4A	S22-C08-N07-C06
3	A	601	B4A	N09-C08-N07-C06
3	C	601	B4A	N24-C23-N03-C02

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	601	B4A	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	410/539 (76%)	-0.14	20 (4%)	35 32	24, 34, 77, 134	0
1	C	410/539 (76%)	-0.13	16 (3%)	43 40	23, 33, 78, 134	1 (0%)
1	D	410/539 (76%)	0.01	29 (7%)	22 18	23, 35, 90, 170	2 (0%)
2	B	410/539 (76%)	-0.05	23 (5%)	30 27	22, 34, 81, 153	0
All	All	1640/2156 (76%)	-0.08	88 (5%)	31 28	22, 34, 81, 170	3 (0%)

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	318	PHE	7.4
1	D	316	LEU	7.3
1	D	546	GLY	6.5
2	B	193	VAL	5.8
1	D	318	PHE	5.8
1	C	546	GLY	5.5
1	C	137	PRO	5.3
1	C	317	ARG	5.3
1	C	318	PHE	5.2
1	A	191	ASP	5.1
1	A	316	LEU	5.1
1	D	139	LEU	5.1
1	A	318	PHE	5.0
1	D	192	GLY	4.9
1	A	546	GLY	4.4
1	A	137	PRO	4.4
2	B	250	ILE	4.3
1	D	250	ILE	4.1
2	B	546	GLY	4.1
1	D	317	ARG	4.0
1	D	138	SER	3.9

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Mol	Chain	Res	Type	RSRZ
1	D	137	PRO	3.8
1	D	188	THR	3.8
1	A	193	VAL	3.8
1	D	315	GLY	3.8
2	B	191	ASP	3.7
2	B	316	LEU	3.7
1	C	193	VAL	3.5
1	D	253	LEU	3.5
1	D	193	VAL	3.5
1	C	192	GLY	3.5
2	B	317	ARG	3.4
2	B	253	LEU	3.3
2	B	190	SER	3.2
1	C	316	LEU	3.2
2	B	320	LYS	3.2
1	C	189	THR	3.2
2	B	137	PRO	3.1
1	A	142	LEU	3.0
1	D	258	PRO	2.9
1	C	347	GLN	2.9
1	D	251	PRO	2.9
2	B	258	PRO	2.8
1	D	186	LEU	2.8
1	A	192	GLY	2.8
2	B	538	LYS	2.8
1	A	317	ARG	2.8
2	B	255	LYS	2.7
1	C	141	ASP	2.7
2	B	259	ASP	2.7
2	B	249	TYR	2.7
1	A	190	SER	2.7
2	B	142	LEU	2.7
1	D	181	MET	2.6
1	A	253	LEU	2.6
1	A	383	GLU	2.6
1	D	185	THR	2.6
1	A	145	TYR	2.6
1	D	152	GLU	2.5
1	A	138	SER	2.5
1	D	146	THR	2.5
1	C	145	TYR	2.5
1	A	147	ILE	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	141	ASP	2.4
1	D	184	LEU	2.4
2	B	192	GLY	2.4
1	D	145	TYR	2.4
1	A	144	PHE	2.4
2	B	319	ASN	2.4
2	B	139	LEU	2.4
1	C	188	THR	2.3
1	C	195	LEU	2.3
1	A	319	ASN	2.3
1	D	319	ASN	2.3
1	D	187	GLN	2.3
1	C	148	ALA	2.2
2	B	145	TYR	2.2
1	D	256	PHE	2.2
1	C	142	LEU	2.1
2	B	152	GLU	2.1
2	B	256	PHE	2.1
1	A	189	THR	2.1
1	A	195	LEU	2.1
1	D	275	THR	2.1
1	C	194	MET	2.1
1	D	156	VAL	2.0
1	D	314	SER	2.0
1	D	148	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

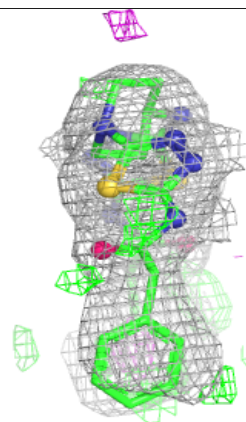
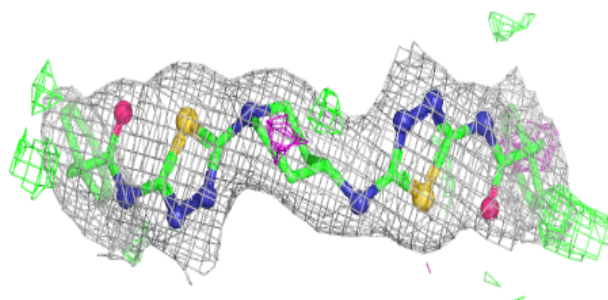
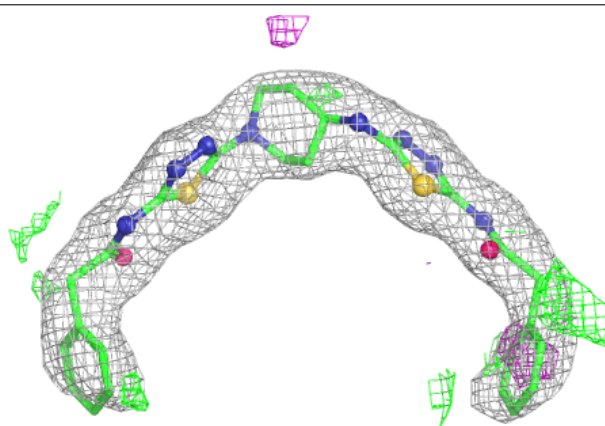
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	B4A	C	601	37/37	0.93	0.11	38,47,60,65	0
3	B4A	A	601	37/37	0.94	0.11	40,46,89,91	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

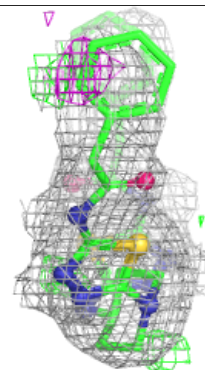
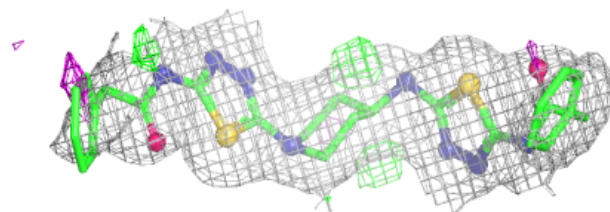
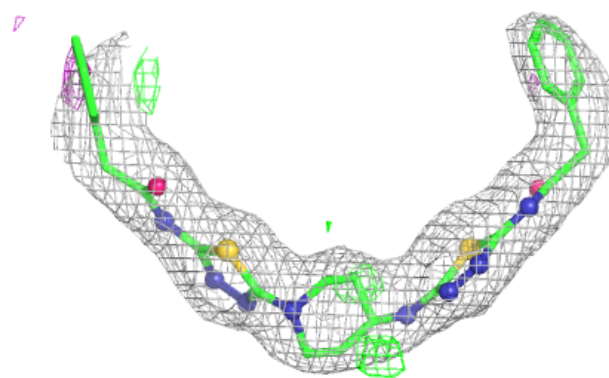
Electron density around B4A C 601:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
 and green (positive)



Electron density around B4A A 601:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.