



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2026 – 11:25 PM UTC

PDB ID : 7REN / pdb_00007ren
Title : Room temperature serial crystal structure of Glutaminase C in complex with inhibitor UPGL-00004
Authors : Milano, S.K.; Finke, A.; Cerione, R.A.
Deposited on : 2021-07-13
Resolution : 2.80 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

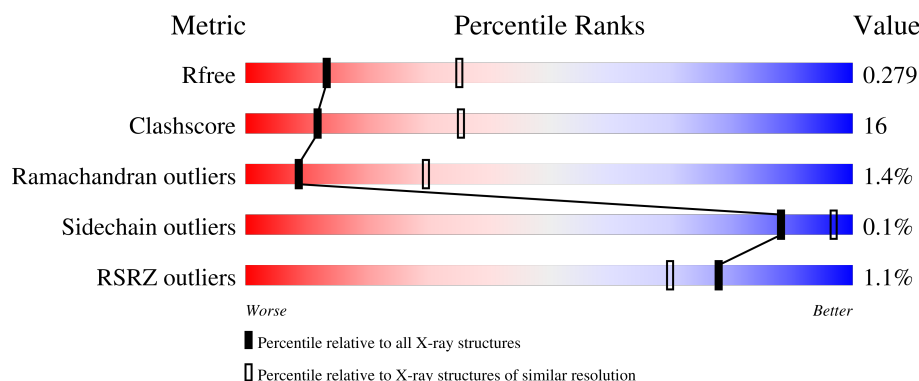
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	52% 22% • 25%
1	B	539	% 50% 25% • 25%
1	C	539	% 54% 21% • 25%
1	D	539	% 47% 27% • 25%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12675 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	404	Total	C	N	O	S	1	0	0
			3148	2003	533	584	28			
1	B	404	Total	C	N	O	S	1	0	0
			3148	2003	533	584	28			
1	C	404	Total	C	N	O	S	1	0	0
			3148	2003	533	584	28			
1	D	404	Total	C	N	O	S	1	0	0
			3148	2003	533	584	28			

There are 48 discrepancies between the modelled and reference sequences:

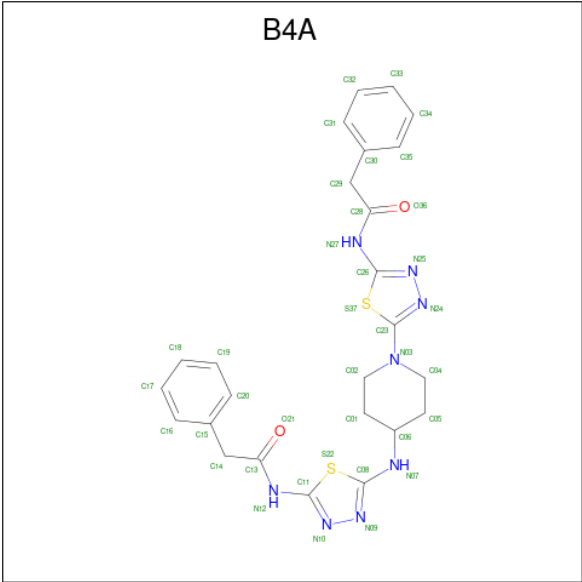
Chain	Residue	Modelled	Actual	Comment	Reference
A	60	MET	-	initiating methionine	UNP O94925
A	61	ARG	-	expression tag	UNP O94925
A	62	GLY	-	expression tag	UNP O94925
A	63	SER	-	expression tag	UNP O94925
A	64	HIS	-	expression tag	UNP O94925
A	65	HIS	-	expression tag	UNP O94925
A	66	HIS	-	expression tag	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
B	60	MET	-	initiating methionine	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

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Chain	Residue	Modelled	Actual	Comment	Reference
B	69	HIS	-	expression tag	UNP O94925
B	70	GLY	-	expression tag	UNP O94925
B	71	SER	-	expression tag	UNP O94925
C	60	MET	-	initiating methionine	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
C	70	GLY	-	expression tag	UNP O94925
C	71	SER	-	expression tag	UNP O94925
D	60	MET	-	initiating methionine	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

- Molecule 2 is 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 (CCD ID: B4A) (formula: C₂₅H₂₆N₈O₂S₂) (labeled as "Ligand of Interest" by depositor).



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

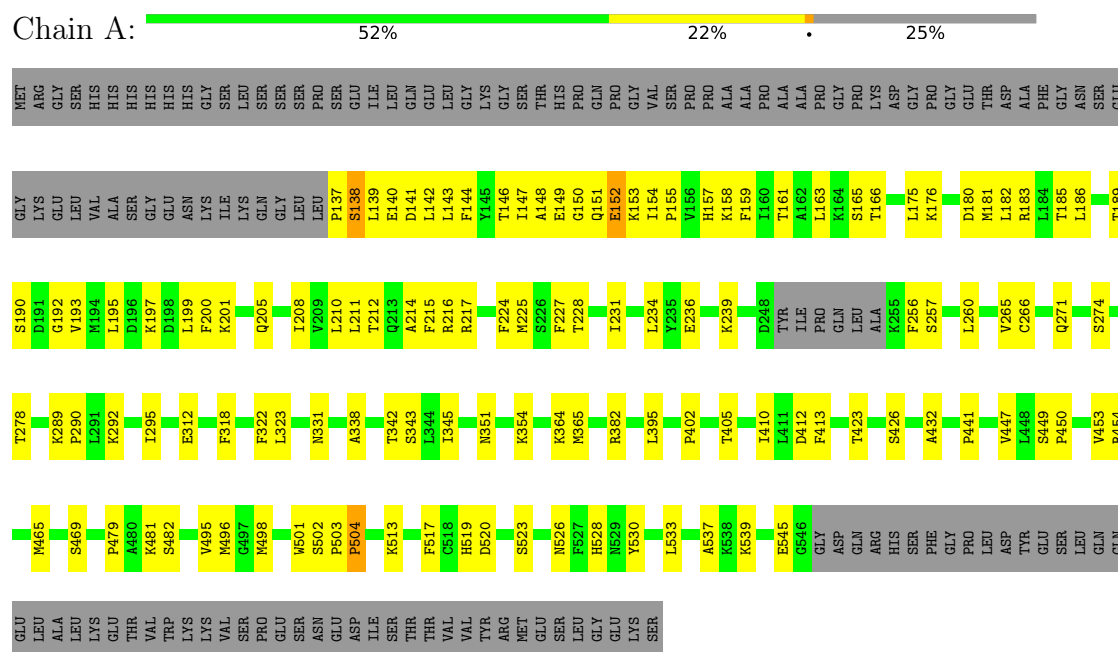
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total	O	0	0
			3	3		
3	C	3	Total	O	0	0
			3	3		
3	D	3	Total	O	0	0
			3	3		

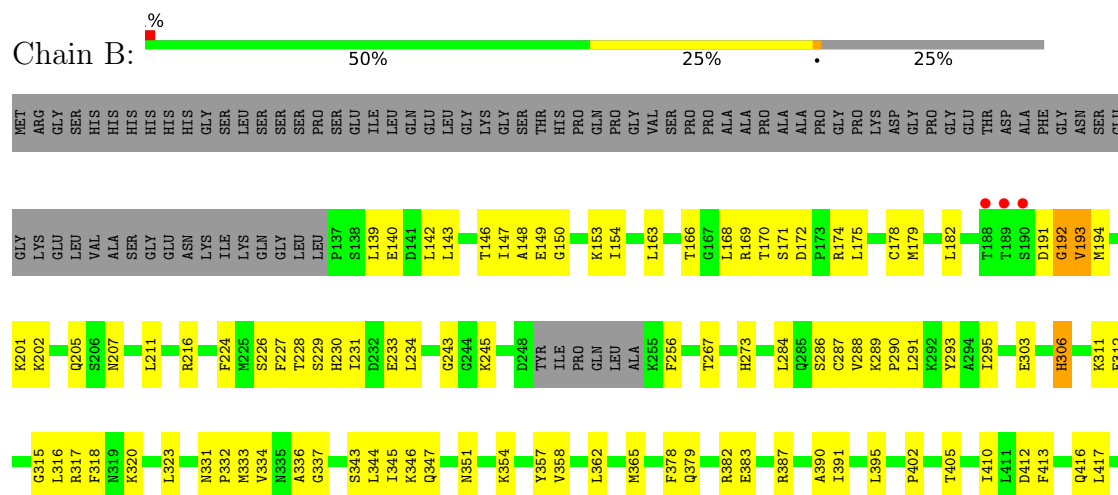
3 Residue-property plots

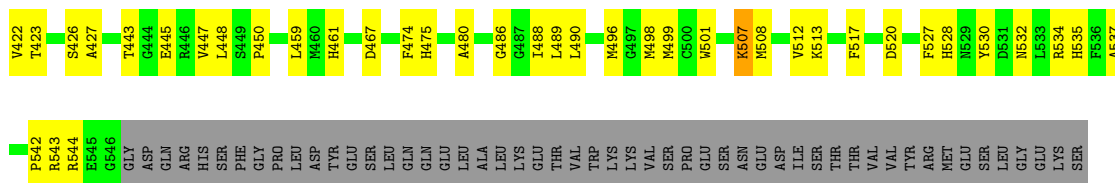
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

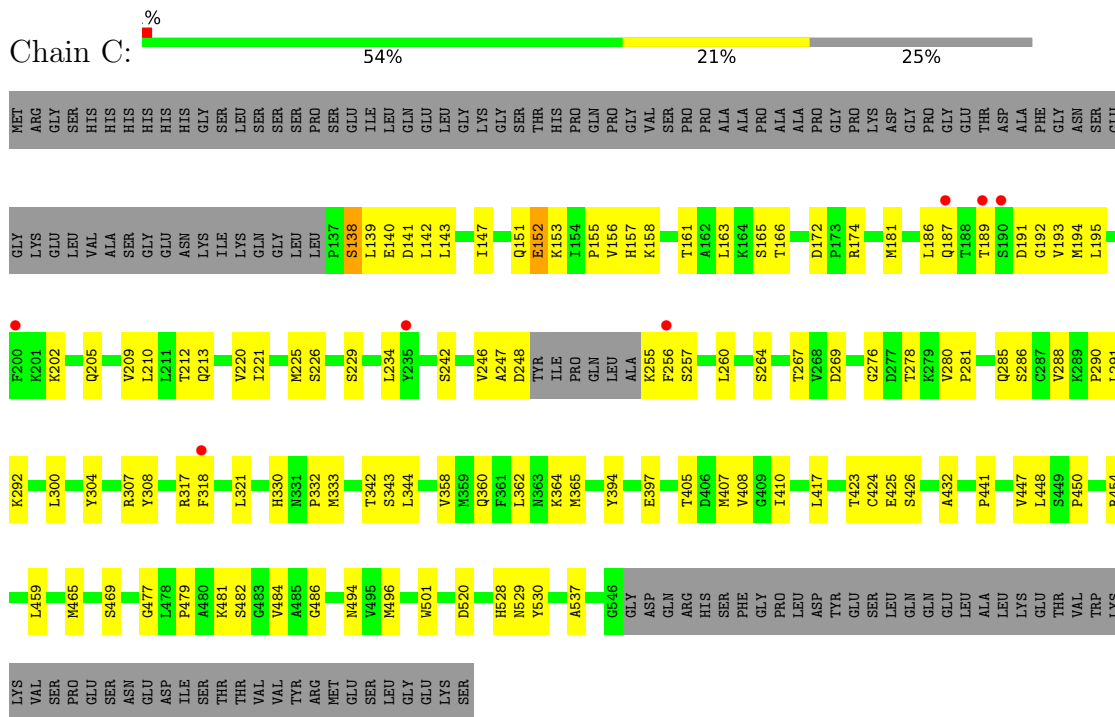


- Molecule 1: Glutaminase kidney isoform, mitochondrial

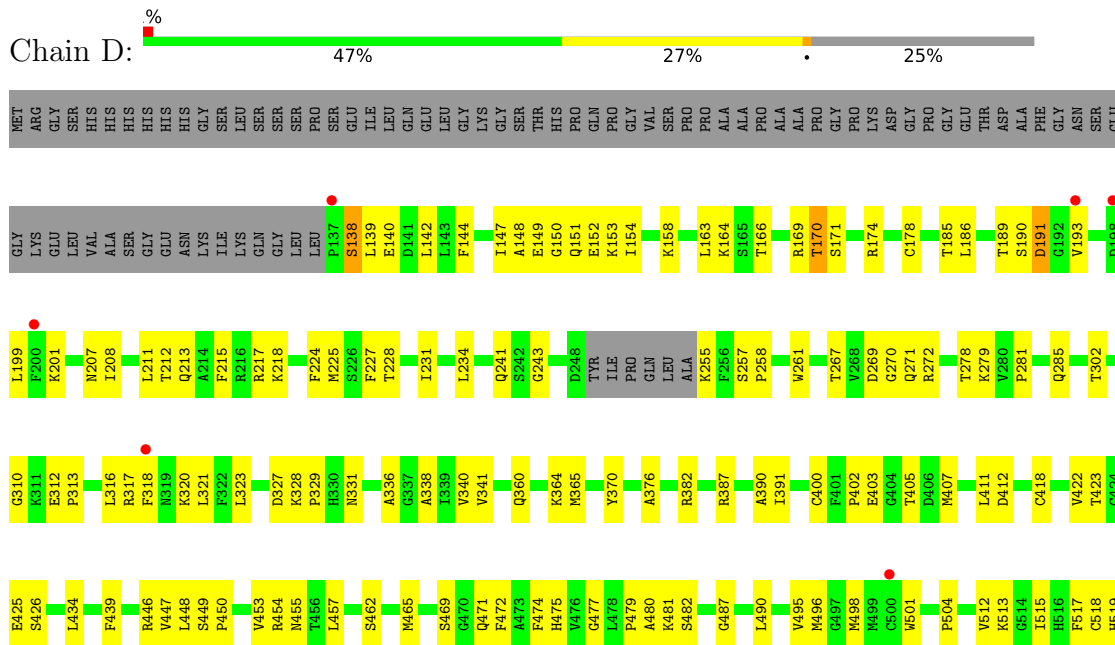




- Molecule 1: Glutaminase kidney isoform, mitochondrial



- Molecule 1: Glutaminase kidney isoform, mitochondrial



THR THR VAL VAL TYR ARG MET GLU SER LEU GLY GLU LYS SER

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	53.20Å 138.44Å 177.49Å 90.00° 93.44° 90.00°	Depositor
Resolution (Å)	20.44 – 2.80 20.44 – 2.80	Depositor EDS
% Data completeness (in resolution range)	89.1 (20.44-2.80) 78.4 (20.44-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.37 (at 2.79Å)	Xtriage
Refinement program	PHENIX 1.16_3549	Depositor
R, R_{free}	0.210 , 0.278 0.212 , 0.279	Depositor DCC
R_{free} test set	2013 reflections (3.19%)	wwPDB-VP
Wilson B-factor (Å ²)	54.4	Xtriage
Anisotropy	0.484	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 41.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12675	wwPDB-VP
Average B, all atoms (Å ²)	86.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.38% 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: 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	0.55	1/3217 (0.0%)	0.68	0/4338
1	B	0.43	0/3217	0.67	0/4338
1	C	0.40	0/3217	0.66	0/4338
1	D	0.41	0/3217	0.67	0/4338
All	All	0.45	1/12868 (0.0%)	0.67	0/17352

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	2
All	All	0	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	504	PRO	N-CD	-20.02	1.19	1.47

There are no bond angle outliers.

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	151	GLN	Peptide
1	B	306	HIS	Peptide
1	B	320	LYS	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	3148	0	3117	100	0
1	B	3148	0	3117	111	0
1	C	3148	0	3116	91	0
1	D	3148	0	3117	117	0
2	B	37	0	0	2	0
2	C	37	0	0	2	0
3	A	3	0	0	2	0
3	C	3	0	0	2	0
3	D	3	0	0	7	0
All	All	12675	0	12467	397	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 16.

All (397) 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:C:255:LYS:HA	3:C:703:HOH:O	1.26	1.32
2:C:601:B4A:N09	2:C:601:B4A:N10	1.60	1.19
2:B:601:B4A:N24	2:B:601:B4A:N25	1.60	1.16
1:D:255:LYS:HA	3:D:601:HOH:O	1.47	1.11
1:A:195:LEU:HD23	1:A:199:LEU:HG	1.08	1.08
1:A:165:SER:HA	1:A:225:MET:HE3	1.47	0.96
1:A:195:LEU:CD2	1:A:199:LEU:HG	1.96	0.93
1:A:195:LEU:HD23	1:A:199:LEU:CG	1.98	0.92
1:D:255:LYS:CA	3:D:601:HOH:O	2.08	0.88
1:D:525:CYS:HB2	3:D:602:HOH:O	1.72	0.88
1:D:525:CYS:SG	3:D:602:HOH:O	2.32	0.85
1:D:525:CYS:CB	3:D:602:HOH:O	2.25	0.84
1:D:481:LYS:NZ	1:D:482:SER:O	2.11	0.83
1:B:234:LEU:HD22	1:B:520:ASP:HB3	1.61	0.82
1:B:532:ASN:HD22	1:B:535:HIS:H	1.23	0.82
1:B:174:ARG:O	1:B:207:ASN:ND2	2.12	0.82
1:B:163:LEU:O	1:B:166:THR:HG22	1.80	0.81
1:D:189:THR:HG22	1:D:190:SER:H	1.46	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:351:ASN:HB3	3:A:602:HOH:O	1.81	0.80
1:B:316:LEU:HD21	1:D:312:GLU:OE1	1.81	0.80
1:A:537:ALA:HB2	1:C:450:PRO:HG2	1.63	0.79
1:C:394:TYR:CZ	2:C:601:B4A:N09	2.52	0.77
1:A:165:SER:HA	1:A:225:MET:CE	2.17	0.75
1:A:152:GLU:OE1	1:A:153:LYS:HB2	1.87	0.74
1:B:148:ALA:HB2	1:B:154:ILE:HG12	1.67	0.74
1:B:543:ARG:C	1:B:544:ARG:HD2	2.12	0.74
1:D:387:ARG:H	1:D:387:ARG:HD2	1.53	0.74
1:A:265:VAL:HG22	1:A:498:MET:HG2	1.70	0.73
1:A:163:LEU:O	1:A:166:THR:HG22	1.87	0.73
1:B:532:ASN:ND2	1:B:535:HIS:H	1.87	0.73
1:A:143:LEU:HD12	1:A:200:PHE:HZ	1.54	0.72
1:C:481:LYS:NZ	1:C:482:SER:O	2.22	0.72
1:B:346:LYS:NZ	1:B:357:TYR:CD1	2.57	0.72
1:A:142:LEU:O	1:A:146:THR:OG1	2.08	0.71
1:D:423:THR:HG23	1:D:426:SER:H	1.55	0.70
1:B:143:LEU:O	1:B:147:ILE:HD12	1.93	0.69
1:A:189:THR:HB	1:A:193:VAL:HG22	1.76	0.68
1:A:217:ARG:NH2	1:A:545:GLU:OE2	2.27	0.68
1:C:163:LEU:O	1:C:166:THR:HG22	1.93	0.68
1:B:346:LYS:NZ	1:B:357:TYR:CE1	2.61	0.68
1:C:138:SER:HB3	1:C:141:ASP:HB2	1.75	0.68
1:C:343:SER:HA	1:C:410:ILE:HD12	1.76	0.68
1:D:144:PHE:HE1	1:D:154:ILE:HG12	1.59	0.68
1:B:153:LYS:HB3	1:B:194:MET:HB3	1.75	0.67
1:D:213:GLN:HB2	1:D:218:LYS:HB2	1.75	0.67
1:D:279:LYS:HE2	1:D:425:GLU:OE2	1.94	0.67
1:D:147:ILE:O	1:D:158:LYS:NZ	2.26	0.67
1:D:178:CYS:HB3	1:D:207:ASN:ND2	2.09	0.66
1:A:503:PRO:N	1:A:504:PRO:HD2	2.10	0.66
1:C:191:ASP:HB3	1:C:193:VAL:HB	1.78	0.66
1:A:139:LEU:HD11	1:A:212:THR:HG21	1.78	0.65
1:A:351:ASN:CB	3:A:602:HOH:O	2.38	0.65
1:B:387:ARG:HD3	1:B:387:ARG:H	1.61	0.65
1:A:481:LYS:NZ	1:A:482:SER:O	2.30	0.65
1:A:519:HIS:O	1:A:523:SER:OG	2.10	0.65
1:A:323:LEU:HD21	1:A:395:LEU:HD21	1.79	0.64
1:D:255:LYS:C	3:D:601:HOH:O	2.36	0.64
1:B:382:ARG:NH1	1:B:412:ASP:OD1	2.29	0.64
1:A:143:LEU:HD12	1:A:200:PHE:CZ	2.32	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:423:THR:HG23	1:B:426:SER:H	1.62	0.63
1:C:202:LYS:HA	1:C:205:GLN:HG2	1.81	0.63
1:C:174:ARG:HB3	3:C:701:HOH:O	1.98	0.63
1:B:316:LEU:HD22	1:B:316:LEU:H	1.64	0.63
1:D:434:LEU:HB3	1:D:457:LEU:HD11	1.80	0.63
1:A:175:LEU:HD21	1:A:210:LEU:HD23	1.80	0.63
1:A:450:PRO:HG2	1:C:537:ALA:HB2	1.81	0.63
1:B:267:THR:HA	1:B:496:MET:HA	1.79	0.63
1:D:234:LEU:HD22	1:D:520:ASP:HB3	1.80	0.62
1:C:278:THR:O	1:C:425:GLU:HG3	2.00	0.61
1:A:274:SER:HB3	1:A:278:THR:HG21	1.82	0.61
1:B:530:TYR:OH	1:D:479:PRO:HD3	2.00	0.61
1:C:304:TYR:HE2	1:C:344:LEU:HD22	1.66	0.61
1:B:422:VAL:HG21	1:B:427:ALA:HB2	1.82	0.61
1:A:498:MET:HE1	1:A:517:PHE:HE1	1.66	0.61
1:B:332:PRO:HG2	1:B:333:MET:HE2	1.83	0.61
1:C:285:GLN:HG3	1:C:484:VAL:HG23	1.83	0.61
1:D:270:GLY:O	1:D:272:ARG:HG2	2.00	0.61
1:D:403:GLU:CD	1:D:403:GLU:H	2.09	0.61
1:D:360:GLN:O	1:D:364:LYS:HG3	2.00	0.60
1:A:533:LEU:H	1:C:494:ASN:HD21	1.50	0.60
1:D:148:ALA:O	1:D:150:GLY:N	2.35	0.60
1:B:148:ALA:O	1:B:150:GLY:N	2.34	0.60
1:C:156:VAL:HG23	1:C:195:LEU:HD11	1.83	0.60
1:C:290:PRO:HA	1:C:333:MET:HE1	1.81	0.60
1:D:336:ALA:HA	1:D:391:ILE:HG21	1.84	0.60
1:D:465:MET:H	1:D:469:SER:HB3	1.66	0.60
1:A:502:SER:OG	1:A:513:LYS:NZ	2.35	0.60
1:B:379:GLN:O	1:B:383:GLU:HG2	2.02	0.60
1:C:397:GLU:HG3	1:D:387:ARG:HB3	1.83	0.60
1:D:382:ARG:NH2	1:D:412:ASP:OD1	2.35	0.60
1:B:226:SER:O	1:B:229:SER:OG	2.18	0.60
1:B:286:SER:HB3	1:B:289:LYS:HD2	1.84	0.59
1:D:423:THR:HG22	1:D:426:SER:HB3	1.84	0.59
1:B:346:LYS:NZ	1:B:357:TYR:CG	2.70	0.59
1:B:317:ARG:HG2	1:D:318:PHE:CE1	2.38	0.59
1:B:334:VAL:O	1:B:337:GLY:N	2.35	0.59
1:D:138:SER:O	1:D:142:LEU:HG	2.03	0.59
1:D:169:ARG:HD3	1:D:170:THR:H	1.68	0.59
1:C:365:MET:HG3	1:C:447:VAL:HG11	1.85	0.59
1:D:224:PHE:O	1:D:228:THR:HG23	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:498:MET:HE1	1:B:517:PHE:HE1	1.67	0.58
1:C:234:LEU:HD22	1:C:520:ASP:HB3	1.85	0.58
1:D:498:MET:HE1	1:D:517:PHE:HE1	1.67	0.58
1:A:176:LYS:O	1:A:180:ASP:N	2.34	0.58
1:B:507:LYS:H	1:B:507:LYS:HD2	1.68	0.58
1:D:474:PHE:HD2	1:D:475:HIS:CD2	2.20	0.58
1:B:293:TYR:OH	1:B:306:HIS:NE2	2.26	0.58
1:D:365:MET:HG3	1:D:447:VAL:HG11	1.84	0.58
1:B:450:PRO:HG2	1:D:537:ALA:HB2	1.84	0.58
1:D:498:MET:HE1	1:D:517:PHE:CE1	2.38	0.58
1:D:522:VAL:HG21	1:D:528:HIS:HB2	1.85	0.58
1:D:189:THR:HG22	1:D:190:SER:N	2.17	0.57
1:B:139:LEU:HA	1:B:142:LEU:HG	1.84	0.57
1:B:528:HIS:CG	1:D:454:ARG:HD2	2.39	0.57
1:D:495:VAL:HG12	1:D:496:MET:HG2	1.86	0.57
1:B:486:GLY:O	1:B:501:TRP:HA	2.04	0.56
1:B:498:MET:HE1	1:B:517:PHE:CE1	2.39	0.56
1:A:234:LEU:HD22	1:A:520:ASP:HB3	1.85	0.56
1:A:465:MET:H	1:A:469:SER:HB3	1.69	0.56
1:B:343:SER:HA	1:B:410:ILE:HD12	1.88	0.56
1:C:477:GLY:O	1:C:529:ASN:HB2	2.05	0.56
1:D:139:LEU:HA	1:D:142:LEU:HD12	1.87	0.56
1:B:201:LYS:O	1:B:205:GLN:HB2	2.05	0.56
1:A:338:ALA:O	1:A:342:THR:HG23	2.06	0.56
1:A:382:ARG:NH2	1:A:412:ASP:OD2	2.38	0.56
1:C:423:THR:H	1:C:426:SER:HB3	1.71	0.56
1:D:211:LEU:O	1:D:215:PHE:HD2	1.89	0.56
1:A:454:ARG:HD2	1:C:528:HIS:CD2	2.41	0.56
1:A:154:ILE:HG22	1:A:155:PRO:O	2.06	0.56
1:B:387:ARG:H	1:B:387:ARG:CD	2.17	0.56
1:D:512:VAL:HG13	1:D:513:LYS:H	1.71	0.56
1:B:480:ALA:HB2	1:B:490:LEU:HD12	1.87	0.56
1:C:143:LEU:HD21	1:C:212:THR:HG22	1.88	0.55
1:C:165:SER:HA	1:C:225:MET:HE3	1.87	0.55
1:D:163:LEU:O	1:D:166:THR:HG22	2.07	0.55
1:C:288:VAL:HG21	1:C:417:LEU:O	2.06	0.55
1:A:530:TYR:OH	1:C:479:PRO:HD3	2.07	0.55
1:A:343:SER:HA	1:A:410:ILE:HD12	1.88	0.55
1:D:370:TYR:O	1:D:426:SER:HB2	2.06	0.55
1:A:147:ILE:HG12	1:A:158:LYS:NZ	2.22	0.54
1:B:346:LYS:NZ	1:B:357:TYR:CZ	2.75	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:140:GLU:HG2	1:A:205:GLN:HG3	1.87	0.54
1:A:141:ASP:OD2	1:A:201:LYS:NZ	2.40	0.54
1:B:512:VAL:HG13	1:B:513:LYS:H	1.73	0.54
1:C:172:ASP:OD1	1:C:174:ARG:HD3	2.07	0.54
1:C:248:ASP:OD1	1:C:248:ASP:N	2.38	0.54
1:C:358:VAL:HG21	1:C:417:LEU:HD21	1.89	0.54
1:D:407:MET:SD	1:D:411:LEU:HD11	2.47	0.54
1:B:427:ALA:HB3	1:B:499:MET:HG2	1.90	0.54
1:B:139:LEU:H	1:B:139:LEU:HD23	1.73	0.54
2:B:601:B4A:C19	1:C:317:ARG:HD2	2.37	0.54
1:D:261:TRP:HA	1:D:501:TRP:O	2.08	0.54
1:B:224:PHE:O	1:B:228:THR:HG23	2.08	0.54
1:B:230:HIS:HA	1:B:233:GLU:HB2	1.89	0.54
1:C:174:ARG:HB2	1:C:210:LEU:HD21	1.89	0.53
1:A:312:GLU:O	1:A:331:ASN:HB3	2.08	0.53
1:C:140:GLU:N	1:C:140:GLU:OE1	2.40	0.53
1:D:329:PRO:HG2	1:D:340:VAL:HG21	1.89	0.53
1:C:191:ASP:CB	1:C:193:VAL:HB	2.39	0.53
1:B:534:ARG:HD2	1:B:535:HIS:NE2	2.24	0.53
1:A:147:ILE:HD11	1:A:159:PHE:HA	1.90	0.53
1:B:193:VAL:HG22	1:B:194:MET:H	1.74	0.53
1:C:143:LEU:N	1:C:143:LEU:HD22	2.24	0.53
1:D:302:THR:HG22	1:D:455:ASN:ND2	2.24	0.53
1:B:312:GLU:OE1	1:D:316:LEU:CD2	2.57	0.52
1:B:312:GLU:OE1	1:D:316:LEU:HD21	2.09	0.52
1:C:226:SER:O	1:C:229:SER:OG	2.26	0.52
1:B:537:ALA:HB2	1:D:450:PRO:HG2	1.90	0.52
1:C:155:PRO:HG3	1:C:194:MET:HE3	1.90	0.52
1:A:533:LEU:H	1:C:494:ASN:ND2	2.07	0.52
1:A:318:PHE:HE2	1:C:317:ARG:HB2	1.74	0.52
1:A:227:PHE:O	1:A:231:ILE:HG12	2.09	0.52
1:C:407:MET:HE3	1:C:408:VAL:HG22	1.91	0.52
1:D:447:VAL:HG12	1:D:448:LEU:HG	1.92	0.51
1:C:151:GLN:HG2	1:C:152:GLU:H	1.75	0.51
1:B:143:LEU:HD12	1:B:146:THR:HB	1.93	0.51
1:B:405:THR:HG22	1:B:410:ILE:HG13	1.93	0.51
1:A:152:GLU:CD	1:A:153:LYS:HB2	2.36	0.51
1:A:528:HIS:CG	1:C:454:ARG:HD2	2.45	0.51
1:B:544:ARG:HD2	1:B:544:ARG:N	2.25	0.51
1:C:151:GLN:HG2	1:C:152:GLU:N	2.26	0.51
1:B:243:GLY:O	1:B:512:VAL:HG11	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:387:ARG:O	1:B:390:ALA:HB3	2.11	0.51
1:B:354:LYS:HB3	1:B:413:PHE:CZ	2.45	0.50
1:A:224:PHE:O	1:A:228:THR:HG23	2.11	0.50
1:B:143:LEU:HG	1:B:147:ILE:HD11	1.92	0.50
1:B:284:LEU:HB3	1:B:287:CYS:HB2	1.94	0.50
1:A:292:LYS:HA	1:A:295:ILE:HD12	1.92	0.50
1:B:228:THR:HB	1:B:273:HIS:CE1	2.47	0.50
1:D:140:GLU:HG3	1:D:208:ILE:HG12	1.93	0.50
1:A:182:LEU:O	1:A:186:LEU:HG	2.12	0.50
1:C:486:GLY:O	1:C:501:TRP:HA	2.12	0.50
1:D:189:THR:HB	1:D:191:ASP:H	1.76	0.50
1:B:402:PRO:O	1:B:405:THR:OG1	2.31	0.49
1:A:181:MET:O	1:A:185:THR:HG23	2.13	0.49
1:D:313:PRO:HG3	1:D:462:SER:HB2	1.94	0.49
1:D:320:LYS:HG3	1:D:321:LEU:HD12	1.94	0.49
1:D:465:MET:HG3	1:D:472:PHE:CG	2.48	0.49
1:A:479:PRO:HG3	1:C:530:TYR:CE1	2.47	0.49
1:C:365:MET:HE2	1:C:448:LEU:HD11	1.93	0.49
1:C:304:TYR:CE2	1:C:344:LEU:HD22	2.47	0.49
1:D:518:CYS:O	1:D:522:VAL:HG12	2.13	0.49
1:B:496:MET:HE3	1:B:498:MET:HG3	1.94	0.49
1:D:217:ARG:HH22	1:D:545:GLU:CG	2.26	0.49
1:B:474:PHE:HD2	1:B:475:HIS:CD2	2.31	0.49
1:D:519:HIS:O	1:D:523:SER:OG	2.19	0.49
1:B:336:ALA:HA	1:B:391:ILE:HG21	1.95	0.48
1:C:209:VAL:O	1:C:213:GLN:HG3	2.12	0.48
1:D:482:SER:HA	1:D:487:GLY:O	2.13	0.48
1:B:143:LEU:HG	1:B:147:ILE:CD1	2.43	0.48
1:A:148:ALA:C	1:A:150:GLY:H	2.21	0.48
1:A:158:LYS:O	1:A:158:LYS:HG2	2.11	0.48
1:A:186:LEU:HD22	1:A:193:VAL:HB	1.95	0.48
1:B:532:ASN:HD22	1:B:535:HIS:N	2.02	0.48
1:B:291:LEU:O	1:B:295:ILE:HG13	2.13	0.48
1:D:225:MET:HE2	1:D:225:MET:HA	1.96	0.48
1:A:260:LEU:HD13	1:A:501:TRP:CH2	2.49	0.48
1:C:153:LYS:HD2	1:C:194:MET:HB2	1.95	0.48
1:C:358:VAL:O	1:C:362:LEU:HG	2.14	0.48
1:D:449:SER:O	1:D:453:VAL:HG23	2.13	0.48
1:A:365:MET:HG3	1:A:447:VAL:HG11	1.96	0.47
1:A:256:PHE:HB3	1:A:257:SER:H	1.56	0.47
1:B:346:LYS:NZ	1:B:357:TYR:CD2	2.82	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:521:LEU:HD12	3:D:602:HOH:O	2.14	0.47
1:C:276:GLY:O	1:C:278:THR:HG23	2.15	0.47
1:B:489:LEU:HD12	1:B:499:MET:HE2	1.97	0.47
1:C:220:VAL:HG13	1:C:221:ILE:HG13	1.97	0.47
1:B:166:THR:HG23	1:B:168:LEU:H	1.80	0.47
1:D:169:ARG:C	1:D:171:SER:H	2.23	0.47
1:A:289:LYS:HA	1:A:292:LYS:HE2	1.96	0.47
1:A:354:LYS:HB3	1:A:413:PHE:CZ	2.50	0.47
1:A:449:SER:O	1:A:453:VAL:HG23	2.14	0.47
1:D:185:THR:HG21	1:D:199:LEU:HD11	1.97	0.47
1:C:405:THR:HG22	1:C:410:ILE:HG13	1.97	0.47
1:A:498:MET:HE1	1:A:517:PHE:CE1	2.48	0.46
1:B:288:VAL:HG21	1:B:417:LEU:O	2.15	0.46
1:C:307:ARG:HB2	1:C:308:TYR:CE1	2.51	0.46
1:C:143:LEU:CD2	1:C:212:THR:HG22	2.45	0.46
1:D:174:ARG:H	1:D:174:ARG:HG3	1.50	0.46
1:D:541:ASP:OD1	1:D:543:ARG:HD3	2.15	0.46
1:B:228:THR:HB	1:B:273:HIS:ND1	2.31	0.46
1:D:217:ARG:HH22	1:D:545:GLU:HG3	1.79	0.46
1:D:387:ARG:O	1:D:390:ALA:HB3	2.15	0.46
1:A:423:THR:HG23	1:A:426:SER:H	1.79	0.46
1:B:169:ARG:C	1:B:171:SER:H	2.23	0.46
1:B:172:ASP:HB3	1:B:175:LEU:HG	1.96	0.46
1:B:311:LYS:HE3	1:D:471:GLN:HG3	1.98	0.46
1:C:189:THR:HB	1:C:191:ASP:OD1	2.15	0.46
1:C:332:PRO:HD2	1:C:459:LEU:HD13	1.96	0.46
1:D:178:CYS:HB3	1:D:207:ASN:HD22	1.80	0.46
1:B:207:ASN:O	1:B:211:LEU:HB2	2.16	0.46
1:A:152:GLU:OE1	1:A:153:LYS:HE2	2.15	0.46
1:A:147:ILE:HG12	1:A:158:LYS:HZ3	1.80	0.46
1:A:260:LEU:HD13	1:A:501:TRP:HH2	1.81	0.46
1:B:315:GLY:O	1:B:318:PHE:N	2.44	0.46
1:C:300:LEU:HD22	1:C:304:TYR:CE2	2.50	0.46
1:B:346:LYS:O	1:B:354:LYS:HE3	2.15	0.45
1:D:227:PHE:O	1:D:231:ILE:HG12	2.15	0.45
1:A:318:PHE:CE2	1:C:317:ARG:HB2	2.51	0.45
1:D:302:THR:HG22	1:D:455:ASN:HD21	1.81	0.45
1:A:140:GLU:HG3	1:A:208:ILE:HG12	1.98	0.45
1:D:164:LYS:HE2	1:D:169:ARG:NH1	2.32	0.45
1:D:257:SER:HA	1:D:258:PRO:HD3	1.78	0.45
1:A:138:SER:OG	1:A:141:ASP:HB2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:454:ARG:HD2	1:C:528:HIS:CG	2.52	0.45
1:B:543:ARG:O	1:B:544:ARG:HD2	2.17	0.45
1:A:354:LYS:HD3	1:A:413:PHE:CE2	2.51	0.45
1:C:187:GLN:C	1:C:189:THR:H	2.24	0.45
1:D:139:LEU:HD21	1:D:212:THR:HG21	1.99	0.45
1:D:504:PRO:O	1:D:512:VAL:HG12	2.17	0.45
1:C:181:MET:HG2	1:C:202:LYS:O	2.16	0.45
1:B:227:PHE:CZ	1:B:231:ILE:HD11	2.51	0.45
1:B:245:LYS:HA	1:B:245:LYS:HD2	1.80	0.45
1:C:192:GLY:N	1:C:193:VAL:HB	2.32	0.45
1:D:243:GLY:O	1:D:512:VAL:HG11	2.17	0.45
1:B:527:PHE:CZ	1:B:542:PRO:HG2	2.52	0.44
1:A:154:ILE:HG23	1:A:158:LYS:NZ	2.32	0.44
1:D:387:ARG:O	1:D:391:ILE:HG13	2.17	0.44
1:D:422:VAL:HB	1:D:426:SER:OG	2.17	0.44
1:A:351:ASN:HA	1:A:354:LYS:HB2	2.00	0.44
1:A:143:LEU:HD13	1:A:143:LEU:O	2.17	0.44
1:A:197:LYS:O	1:A:201:LYS:HE2	2.18	0.44
1:B:140:GLU:OE1	1:B:140:GLU:N	2.47	0.44
1:C:432:ALA:HB1	1:C:441:PRO:HG3	1.98	0.44
1:A:266:CYS:HA	1:A:271:GLN:O	2.17	0.44
1:B:378:PHE:CE1	1:B:416:GLN:HG3	2.52	0.44
1:C:153:LYS:HB2	1:C:194:MET:HE2	1.99	0.44
1:B:512:VAL:HG13	1:B:513:LYS:N	2.32	0.44
1:C:220:VAL:N	1:C:269:ASP:OD2	2.36	0.44
1:D:317:ARG:NE	1:D:317:ARG:HA	2.32	0.44
1:C:264:SER:HB2	1:C:424:CYS:HB3	1.99	0.44
1:A:152:GLU:C	1:A:153:LYS:HD3	2.43	0.44
1:D:512:VAL:HG13	1:D:513:LYS:N	2.33	0.44
1:A:190:SER:O	1:A:190:SER:OG	2.28	0.43
1:B:443:THR:OG1	1:B:445:GLU:HG2	2.18	0.43
1:D:302:THR:CG2	1:D:455:ASN:HD21	2.31	0.43
1:B:202:LYS:HA	1:B:205:GLN:HB2	2.00	0.43
1:C:157:HIS:O	1:C:161:THR:HG23	2.18	0.43
1:C:417:LEU:HD23	1:C:417:LEU:HA	1.69	0.43
1:D:281:PRO:HA	1:D:422:VAL:O	2.18	0.43
1:A:364:LYS:HB3	1:A:447:VAL:HG22	2.00	0.43
1:B:365:MET:HG3	1:B:447:VAL:HG11	2.01	0.43
1:A:323:LEU:HD21	1:A:395:LEU:CD2	2.48	0.43
1:B:507:LYS:H	1:B:507:LYS:CD	2.29	0.43
1:B:293:TYR:CE2	1:B:459:LEU:HD12	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:292:LYS:HD3	1:C:342:THR:OG1	2.18	0.43
1:D:323:LEU:HD22	1:D:328:LYS:C	2.43	0.43
1:A:503:PRO:N	1:A:504:PRO:CD	2.78	0.43
1:C:143:LEU:O	1:C:147:ILE:HG13	2.18	0.43
1:A:166:THR:HG21	1:A:214:ALA:HB1	2.01	0.43
1:D:186:LEU:HA	1:D:193:VAL:HG21	1.99	0.43
1:D:323:LEU:HD11	1:D:400:CYS:SG	2.58	0.43
1:D:423:THR:CG2	1:D:426:SER:H	2.25	0.43
1:A:180:ASP:OD1	1:A:183:ARG:NH2	2.50	0.43
1:D:189:THR:HB	1:D:191:ASP:HB2	2.00	0.43
1:A:495:VAL:HG12	1:A:496:MET:HB3	2.01	0.43
1:B:317:ARG:HA	1:B:317:ARG:HD2	1.85	0.43
1:C:360:GLN:O	1:C:364:LYS:HG3	2.19	0.43
1:D:338:ALA:O	1:D:341:VAL:HG12	2.18	0.43
1:B:312:GLU:O	1:B:331:ASN:HB3	2.19	0.43
1:A:530:TYR:CE1	1:C:479:PRO:HG3	2.53	0.42
1:B:178:CYS:SG	1:B:179:MET:HE2	2.59	0.42
1:B:354:LYS:HD3	1:B:413:PHE:CE2	2.54	0.42
1:C:246:VAL:HG12	1:C:247:ALA:H	1.84	0.42
1:A:137:PRO:O	1:A:138:SER:HB3	2.18	0.42
1:B:182:LEU:HD23	1:B:182:LEU:HA	1.86	0.42
1:C:260:LEU:HB3	1:C:501:TRP:CZ3	2.54	0.42
1:C:280:VAL:HA	1:C:281:PRO:HD3	1.89	0.42
1:D:531:ASP:OD2	1:D:539:LYS:NZ	2.51	0.42
1:B:345:ILE:O	1:B:354:LYS:HE2	2.19	0.42
1:D:423:THR:N	1:D:426:SER:OG	2.48	0.42
1:A:322:PHE:O	1:A:323:LEU:HD23	2.18	0.42
1:B:289:LYS:N	1:B:290:PRO:HD2	2.34	0.42
1:C:139:LEU:HB2	1:C:140:GLU:OE1	2.19	0.42
1:D:323:LEU:HD13	1:D:327:ASP:C	2.45	0.42
1:B:467:ASP:HB2	1:B:508:MET:HE3	2.01	0.42
1:C:290:PRO:HD3	1:C:481:LYS:HD3	2.01	0.42
1:D:139:LEU:O	1:D:139:LEU:HD23	2.20	0.42
1:A:236:GLU:O	1:A:239:LYS:HB3	2.20	0.42
1:B:169:ARG:O	1:B:171:SER:N	2.52	0.42
1:B:216:ARG:HD3	1:B:216:ARG:HA	1.89	0.42
1:A:197:LYS:HE3	1:A:201:LYS:HZ1	1.84	0.42
1:D:525:CYS:HB3	1:D:540:LEU:O	2.20	0.42
1:A:144:PHE:HD1	1:A:154:ILE:HD11	1.85	0.42
1:C:291:LEU:H	1:C:291:LEU:HG	1.66	0.42
1:A:402:PRO:O	1:A:405:THR:OG1	2.22	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:201:LYS:O	1:B:201:LYS:HG2	2.19	0.41
1:B:346:LYS:HD3	1:B:346:LYS:HA	1.93	0.41
1:C:158:LYS:HA	1:C:158:LYS:HD2	1.66	0.41
1:D:278:THR:O	1:D:423:THR:OG1	2.34	0.41
1:D:320:LYS:H	1:D:320:LYS:HG2	1.63	0.41
1:A:289:LYS:N	1:A:290:PRO:HD2	2.34	0.41
1:B:303:GLU:H	1:B:303:GLU:CD	2.28	0.41
1:B:488:ILE:O	1:B:499:MET:HA	2.20	0.41
1:C:256:PHE:HB3	1:C:257:SER:H	1.76	0.41
1:D:545:GLU:OE1	1:D:545:GLU:N	2.52	0.41
1:A:432:ALA:HB1	1:A:441:PRO:HG3	2.02	0.41
1:B:508:MET:HE2	1:B:508:MET:HB3	1.88	0.41
1:D:140:GLU:HB2	1:D:201:LYS:HG3	2.01	0.41
1:D:522:VAL:CG2	1:D:528:HIS:HB2	2.50	0.41
1:A:157:HIS:O	1:A:161:THR:HG23	2.21	0.41
1:A:292:LYS:HD2	1:A:342:THR:HG22	2.02	0.41
1:B:317:ARG:HG2	1:D:318:PHE:CZ	2.55	0.41
1:C:465:MET:H	1:C:469:SER:CB	2.34	0.41
1:D:267:THR:C	1:D:269:ASP:H	2.28	0.41
1:D:271:GLN:H	1:D:271:GLN:HG3	1.67	0.41
1:B:358:VAL:O	1:B:362:LEU:HG	2.21	0.41
1:C:465:MET:H	1:C:469:SER:HB2	1.85	0.41
1:D:310:GLY:O	1:D:331:ASN:HA	2.20	0.41
1:B:323:LEU:HD21	1:B:395:LEU:HD21	2.01	0.41
1:D:164:LYS:HE2	1:D:169:ARG:CZ	2.51	0.41
1:B:344:LEU:HA	1:B:347:GLN:HE21	1.85	0.41
1:C:286:SER:C	1:C:288:VAL:N	2.79	0.41
1:A:526:ASN:HA	1:A:539:LYS:HE3	2.02	0.41
1:C:138:SER:O	1:C:142:LEU:HG	2.21	0.41
1:C:267:THR:HA	1:C:496:MET:HA	2.03	0.41
1:C:318:PHE:CD1	1:C:321:LEU:HD13	2.56	0.41
1:D:439:PHE:CE2	1:D:446:ARG:HB3	2.56	0.41
1:D:477:GLY:O	1:D:529:ASN:HB2	2.20	0.41
1:D:285:GLN:OE1	1:D:418:CYS:HB3	2.20	0.41
1:A:140:GLU:HG2	1:A:205:GLN:CG	2.49	0.40
1:C:152:GLU:HB2	1:C:153:LYS:H	1.60	0.40
1:C:257:SER:O	1:C:260:LEU:HD12	2.21	0.40
1:D:480:ALA:HB2	1:D:490:LEU:HD12	2.03	0.40
1:C:186:LEU:HD12	1:C:186:LEU:O	2.22	0.40
1:D:151:GLN:OE1	1:D:153:LYS:O	2.39	0.40
1:A:216:ARG:HD2	1:A:216:ARG:HA	1.79	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:413:PHE:O	1:B:417:LEU:HD13	2.20	0.40
1:B:448:LEU:HA	1:B:448:LEU:HD23	1.79	0.40
1:A:211:LEU:O	1:A:215:PHE:HD2	2.04	0.40
1:A:479:PRO:HD3	1:C:530:TYR:OH	2.21	0.40
1:B:192:GLY:HA3	1:B:193:VAL:HA	1.92	0.40
1:D:490:LEU:HB3	1:D:498:MET:HB2	2.02	0.40
1:D:528:HIS:HB3	1:D:531:ASP:OD2	2.22	0.40
1:A:139:LEU:HA	1:A:142:LEU:HG	2.04	0.40
1:D:402:PRO:O	1:D:405:THR:OG1	2.24	0.40
1:D:515:ILE:HD13	1:D:515:ILE:HA	1.94	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	400/539 (74%)	361 (90%)	34 (8%)	5 (1%)	9	31
1	B	400/539 (74%)	358 (90%)	34 (8%)	8 (2%)	6	21
1	C	400/539 (74%)	358 (90%)	39 (10%)	3 (1%)	16	44
1	D	400/539 (74%)	358 (90%)	35 (9%)	7 (2%)	6	23
All	All	1600/2156 (74%)	1435 (90%)	142 (9%)	23 (1%)	9	30

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	138	SER
1	B	149	GLU
1	B	191	ASP
1	C	152	GLU
1	C	330	HIS

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Mol	Chain	Res	Type
1	D	138	SER
1	D	149	GLU
1	A	152	GLU
1	A	345	ILE
1	D	152	GLU
1	A	192	GLY
1	B	256	PHE
1	C	138	SER
1	D	170	THR
1	D	191	ASP
1	D	241	GLN
1	A	149	GLU
1	B	170	THR
1	B	351	ASN
1	B	461	HIS
1	D	376	ALA
1	B	192	GLY
1	B	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	349/462 (76%)	349 (100%)	0	100	100
1	B	349/462 (76%)	348 (100%)	1 (0%)	86	95
1	C	349/462 (76%)	348 (100%)	1 (0%)	86	95
1	D	349/462 (76%)	349 (100%)	0	100	100
All	All	1396/1848 (76%)	1394 (100%)	2 (0%)	88	97

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	507	LYS
1	C	242	SER

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	510	ASN
1	A	516	HIS
1	B	375	ASN
1	B	510	ASN
1	B	532	ASN
1	C	213	GLN
1	C	347	GLN
1	C	351	ASN
1	C	494	ASN
1	C	519	HIS
1	D	207	ASN
1	D	319	ASN
1	D	347	GLN
1	D	388	ASN
1	D	416	GLN
1	D	455	ASN
1	D	471	GLN
1	D	475	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](#)

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	B4A	C	601	-	41,41,41	5.04	17 (41%)	51,55,55	5.47	22 (43%)
2	B4A	B	601	-	41,41,41	4.98	18 (43%)	51,55,55	5.42	19 (37%)

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	B4A	C	601	-	-	12/24/34/34	1/5/5/5
2	B4A	B	601	-	-	12/24/34/34	1/5/5/5

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	601	B4A	C26-S37	-13.09	1.52	1.73
2	B	601	B4A	C26-S37	-12.53	1.52	1.73
2	B	601	B4A	C11-S22	-12.47	1.53	1.73
2	C	601	B4A	C08-S22	-11.89	1.54	1.73
2	C	601	B4A	C11-S22	-11.38	1.54	1.73
2	B	601	B4A	C08-S22	-11.02	1.55	1.73
2	B	601	B4A	C23-S37	-10.59	1.51	1.75
2	B	601	B4A	N25-N24	9.96	1.60	1.39
2	C	601	B4A	C23-S37	-9.90	1.52	1.75
2	C	601	B4A	N10-N09	9.85	1.60	1.39
2	C	601	B4A	N25-N24	9.69	1.60	1.39
2	B	601	B4A	N10-N09	9.38	1.59	1.39
2	B	601	B4A	C08-N07	7.37	1.45	1.34
2	C	601	B4A	C11-N10	7.30	1.42	1.31
2	C	601	B4A	C26-N25	6.12	1.40	1.31
2	C	601	B4A	C08-N07	5.83	1.43	1.34
2	B	601	B4A	C26-N25	5.65	1.39	1.31
2	C	601	B4A	C08-N09	5.61	1.39	1.31
2	B	601	B4A	C11-N10	5.53	1.39	1.31
2	B	601	B4A	C08-N09	5.49	1.39	1.31
2	C	601	B4A	C11-N12	5.00	1.45	1.38
2	C	601	B4A	C23-N03	4.98	1.45	1.34
2	C	601	B4A	C23-N24	4.81	1.40	1.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	601	B4A	C28-N27	4.61	1.45	1.37
2	B	601	B4A	C23-N24	4.55	1.39	1.31
2	B	601	B4A	C28-N27	4.55	1.45	1.37
2	B	601	B4A	C23-N03	4.39	1.44	1.34
2	C	601	B4A	C26-N27	4.36	1.44	1.38
2	B	601	B4A	C26-N27	4.20	1.44	1.38
2	B	601	B4A	C11-N12	4.14	1.44	1.38
2	C	601	B4A	C13-N12	4.11	1.44	1.37
2	B	601	B4A	C13-N12	3.89	1.44	1.37
2	B	601	B4A	O21-C13	-2.55	1.18	1.23
2	C	601	B4A	O21-C13	-2.14	1.19	1.23
2	B	601	B4A	C29-C30	2.01	1.54	1.51

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	B4A	N27-C26-N25	-25.09	98.41	120.81
2	C	601	B4A	N27-C26-N25	-25.06	98.44	120.81
2	C	601	B4A	S37-C26-N27	22.51	153.87	123.72
2	B	601	B4A	S37-C26-N27	22.24	153.50	123.72
2	C	601	B4A	C04-N03-C23	-7.75	102.45	122.09
2	B	601	B4A	N03-C23-N24	6.71	129.88	116.91
2	C	601	B4A	N03-C23-N24	6.51	129.49	116.91
2	C	601	B4A	S37-C26-N25	-6.12	107.69	114.52
2	B	601	B4A	S37-C26-N25	-5.77	108.08	114.52
2	B	601	B4A	S22-C11-N10	-5.72	108.13	114.52
2	B	601	B4A	C13-N12-C11	-5.61	114.09	123.69
2	C	601	B4A	C28-N27-C26	-5.32	114.59	123.69
2	B	601	B4A	C14-C13-N12	5.27	120.83	114.40
2	B	601	B4A	S22-C08-N09	-5.21	108.71	114.52
2	B	601	B4A	C06-N07-C08	-5.11	115.52	124.48
2	B	601	B4A	N12-C11-N10	5.09	125.36	120.81
2	B	601	B4A	C04-N03-C23	-4.95	109.54	122.09
2	C	601	B4A	C29-C28-N27	4.85	120.31	114.40
2	C	601	B4A	C02-N03-C23	-4.54	110.59	122.09
2	C	601	B4A	S22-C11-N10	-4.46	109.55	114.52
2	C	601	B4A	C06-N07-C08	-4.22	117.08	124.48
2	B	601	B4A	S37-C23-N24	-4.18	107.20	114.44
2	C	601	B4A	C11-N10-N09	-3.98	106.95	112.02
2	B	601	B4A	C26-N25-N24	-3.93	107.02	112.02
2	C	601	B4A	S22-C08-N09	-3.91	110.16	114.52
2	C	601	B4A	S37-C23-N24	-3.90	107.70	114.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	B4A	C08-N09-N10	-3.66	107.36	112.02
2	C	601	B4A	N12-C11-N10	3.59	124.02	120.81
2	C	601	B4A	C05-C06-N07	-3.30	103.99	110.57
2	B	601	B4A	C02-C01-C06	3.15	115.64	110.67
2	C	601	B4A	C08-N09-N10	-3.09	108.09	112.02
2	C	601	B4A	C01-C06-N07	-3.08	104.41	110.57
2	C	601	B4A	C14-C13-N12	3.03	118.10	114.40
2	C	601	B4A	C26-N25-N24	-3.00	108.20	112.02
2	B	601	B4A	O21-C13-N12	-2.83	117.08	122.44
2	C	601	B4A	C04-C05-C06	-2.32	107.01	110.67
2	B	601	B4A	C01-C02-N03	-2.20	106.36	110.66
2	B	601	B4A	C11-N10-N09	-2.18	109.24	112.02
2	B	601	B4A	S22-C11-N12	2.07	126.50	123.72
2	C	601	B4A	S22-C11-N12	2.03	126.44	123.72
2	C	601	B4A	O36-C28-N27	-2.02	118.60	122.44

There are no chirality outliers.

All (24) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	601	B4A	N10-C11-N12-C13
2	B	601	B4A	S22-C11-N12-C13
2	B	601	B4A	N25-C26-N27-C28
2	B	601	B4A	S37-C26-N27-C28
2	B	601	B4A	N24-C23-N03-C02
2	B	601	B4A	S37-C23-N03-C02
2	B	601	B4A	N24-C23-N03-C04
2	B	601	B4A	S37-C23-N03-C04
2	B	601	B4A	N09-C08-N07-C06
2	B	601	B4A	S22-C08-N07-C06
2	C	601	B4A	N10-C11-N12-C13
2	C	601	B4A	S22-C11-N12-C13
2	C	601	B4A	N25-C26-N27-C28
2	C	601	B4A	N24-C23-N03-C04
2	C	601	B4A	S37-C23-N03-C04
2	C	601	B4A	N09-C08-N07-C06
2	C	601	B4A	S22-C08-N07-C06
2	C	601	B4A	S37-C26-N27-C28
2	B	601	B4A	O36-C28-C29-C30
2	C	601	B4A	O36-C28-C29-C30
2	C	601	B4A	C13-C14-C15-C20
2	C	601	B4A	C13-C14-C15-C16

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Mol	Chain	Res	Type	Atoms
2	B	601	B4A	N27-C28-C29-C30
2	C	601	B4A	N27-C28-C29-C30

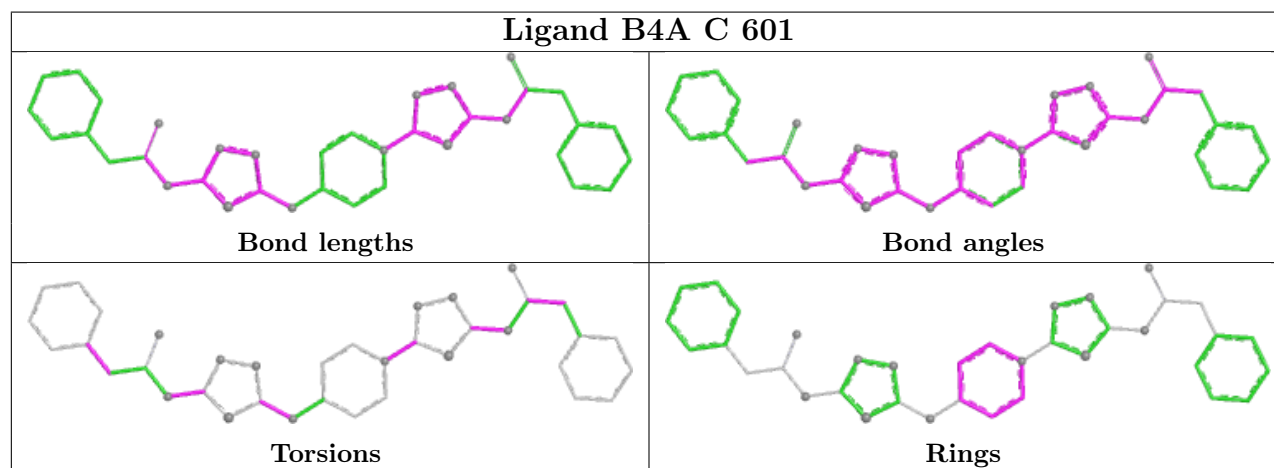
All (2) ring outliers are listed below:

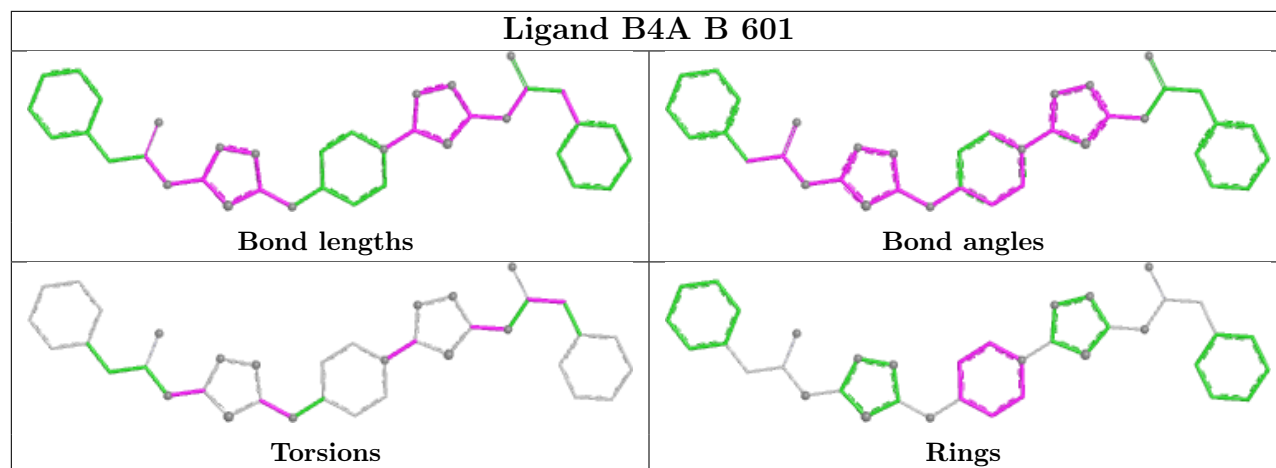
Mol	Chain	Res	Type	Atoms
2	C	601	B4A	C01-C02-C04-C05-C06-N03
2	B	601	B4A	C01-C02-C04-C05-C06-N03

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	601	B4A	2	0
2	B	601	B4A	2	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 [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	404/539 (74%)	-0.01	0 100 100	48, 78, 139, 245	0
1	B	404/539 (74%)	-0.01	3 (0%) 84 77	45, 77, 141, 245	0
1	C	404/539 (74%)	-0.01	7 (1%) 69 60	45, 78, 144, 243	0
1	D	404/539 (74%)	0.01	7 (1%) 69 60	45, 78, 149, 228	0
All	All	1616/2156 (74%)	-0.00	17 (1%) 78 70	45, 78, 147, 245	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	189	THR	4.7
1	B	189	THR	4.0
1	D	318	PHE	3.4
1	D	137	PRO	3.1
1	B	190	SER	2.9
1	B	188	THR	2.9
1	C	190	SER	2.8
1	C	256	PHE	2.7
1	C	187	GLN	2.6
1	C	318	PHE	2.6
1	D	193	VAL	2.4
1	D	200	PHE	2.4
1	C	200	PHE	2.3
1	D	500	CYS	2.3
1	C	235	TYR	2.2
1	D	537	ALA	2.2
1	D	198	ASP	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 [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

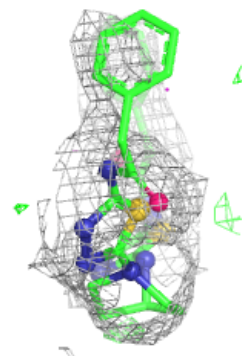
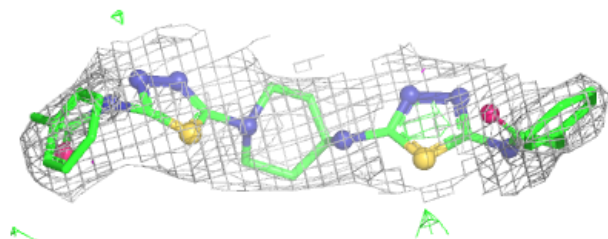
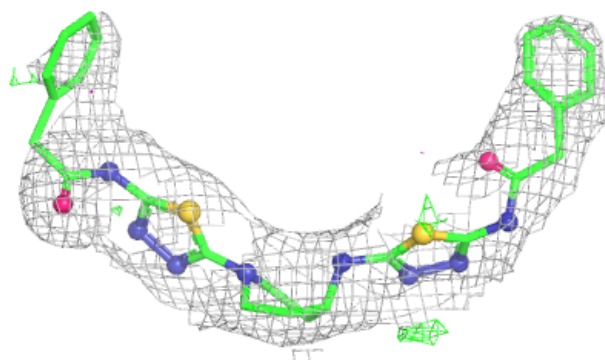
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	B4A	B	601	37/37	0.91	0.12	78,99,117,121	0
2	B4A	C	601	37/37	0.91	0.12	59,103,118,120	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.

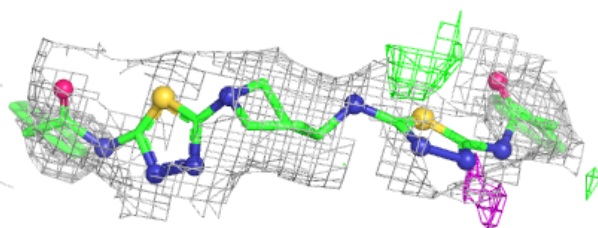
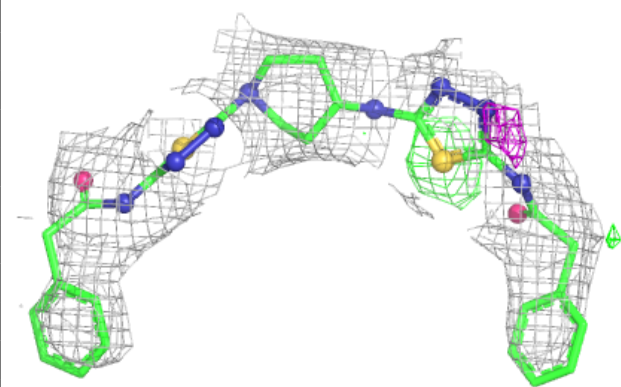
Electron density around B4A B 601:

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 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.