



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 10:42 PM UTC

PDB ID : 8YI6 / pdb_00008yi6
Title : BesA wild-type from Streptomyces cattleyicolor
Authors : Fujishiro, T.
Deposited on : 2024-02-29
Resolution : 3.60 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

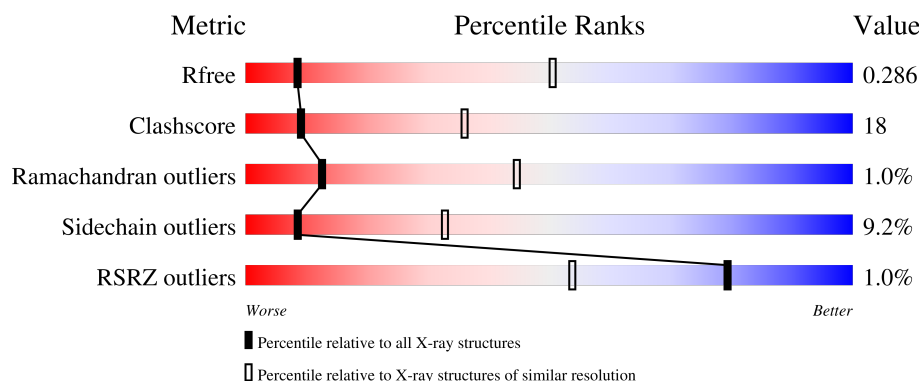
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.60 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	1747 (3.70-3.50)
Clashscore	190562	1827 (3.70-3.50)
Ramachandran outliers	187476	1773 (3.70-3.50)
Sidechain outliers	187428	1772 (3.70-3.50)
RSRZ outliers	180081	1745 (3.70-3.50)

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	475	 59% 32% 5% . .
1	B	475	 % 60% 32% . . .

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ADE	B	501	-	-	X	-

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7010 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.

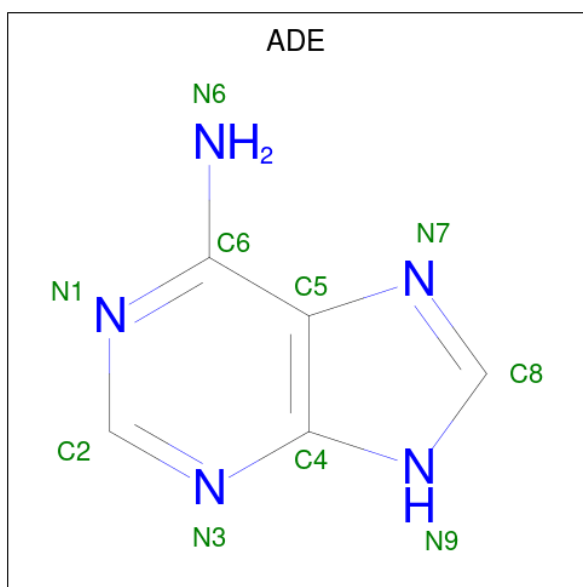
- Molecule 1 is a protein called L-propargylglycine--L-glutamate ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	458	Total	C	N	O	S	0	0	0
			3495	2179	647	660	9			
1	B	458	Total	C	N	O	S	0	0	0
			3495	2179	647	660	9			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	483	VAL	-	expression tag	UNP G8XHD8
A	484	ASP	-	expression tag	UNP G8XHD8
A	485	LEU	-	expression tag	UNP G8XHD8
A	486	GLU	-	expression tag	UNP G8XHD8
A	487	HIS	-	expression tag	UNP G8XHD8
A	488	HIS	-	expression tag	UNP G8XHD8
A	489	HIS	-	expression tag	UNP G8XHD8
A	490	HIS	-	expression tag	UNP G8XHD8
A	491	HIS	-	expression tag	UNP G8XHD8
A	492	HIS	-	expression tag	UNP G8XHD8
B	483	VAL	-	expression tag	UNP G8XHD8
B	484	ASP	-	expression tag	UNP G8XHD8
B	485	LEU	-	expression tag	UNP G8XHD8
B	486	GLU	-	expression tag	UNP G8XHD8
B	487	HIS	-	expression tag	UNP G8XHD8
B	488	HIS	-	expression tag	UNP G8XHD8
B	489	HIS	-	expression tag	UNP G8XHD8
B	490	HIS	-	expression tag	UNP G8XHD8
B	491	HIS	-	expression tag	UNP G8XHD8
B	492	HIS	-	expression tag	UNP G8XHD8

- Molecule 2 is ADENINE (CCD ID: ADE) (formula: C₅H₅N₅) (labeled as "Ligand of Interest" by depositor).

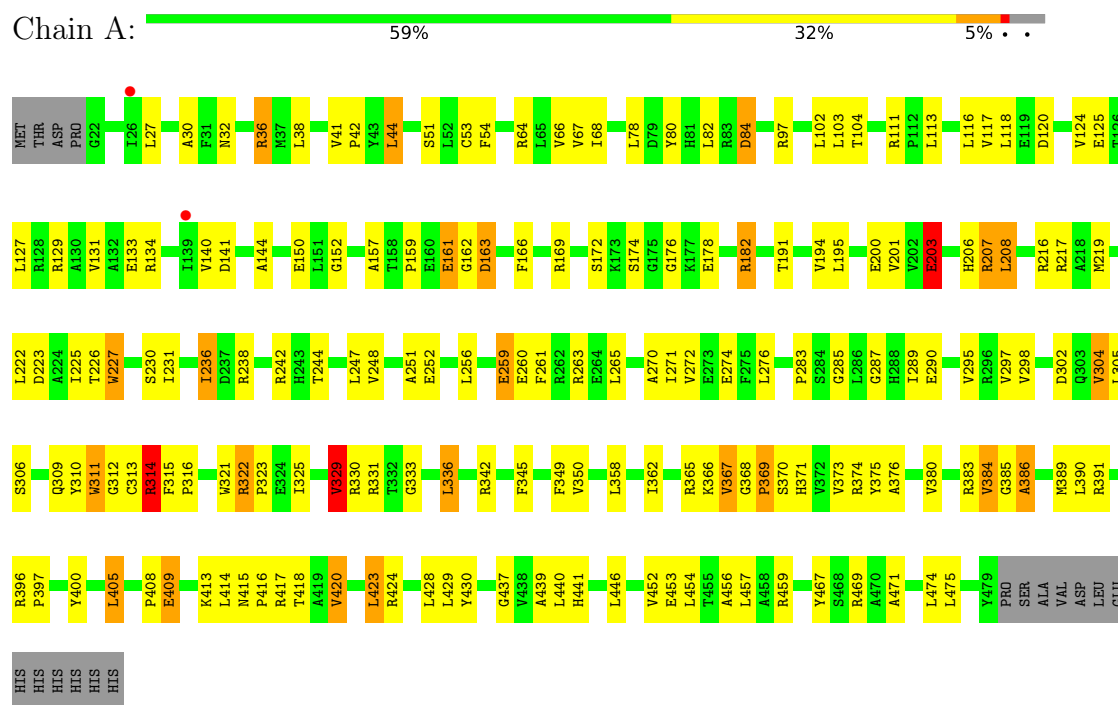


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	0	0
			10	5	5		
2	B	1	Total	C	N	0	0
			10	5	5		

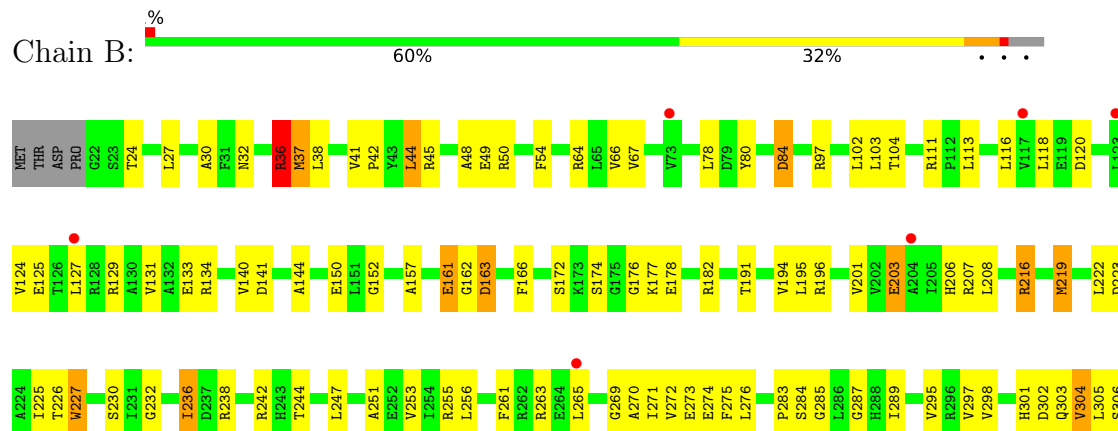
3 Residue-property plots

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: L-propargylglycine--L-glutamate ligase



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4 Data and refinement statistics

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	138.68Å 138.68Å 144.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.45 – 3.60 46.45 – 3.60	Depositor EDS
% Data completeness (in resolution range)	99.9 (46.45-3.60) 99.9 (46.45-3.60)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 3.57Å)	Xtriage
Refinement program	REFMAC 5.8.0419	Depositor
R, R_{free}	0.239 , 0.290 0.231 , 0.286	Depositor DCC
R_{free} test set	848 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	143.6	Xtriage
Anisotropy	0.457	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 160.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.017 for -h,l,k 0.015 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7010	wwPDB-VP
Average B, all atoms (Å ²)	170.0	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 29.97 % 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 1.4166e-03. 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: ADE

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	0.56	0/3557	1.50	27/4825 (0.6%)
1	B	0.56	0/3557	1.49	24/4825 (0.5%)
All	All	0.56	0/7114	1.49	51/9650 (0.5%)

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	4
1	B	0	3
All	All	0	7

There are no bond length outliers.

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	97	ARG	N-CA-CB	10.78	125.97	110.12
1	B	97	ARG	N-CA-CB	10.06	124.90	110.12
1	B	420	VAL	N-CA-CB	9.46	123.41	110.54
1	A	182	ARG	CB-CA-C	9.26	126.16	110.79
1	A	420	VAL	N-CA-CB	9.12	122.95	110.54
1	A	375	TYR	N-CA-CB	8.94	122.97	110.01
1	B	375	TYR	N-CA-CB	7.58	120.99	110.01
1	B	206	HIS	CB-CA-C	7.46	123.52	110.85
1	B	166	PHE	CA-CB-CG	7.22	121.03	113.80
1	A	469	ARG	CB-CA-C	6.90	121.72	110.88
1	B	322	ARG	O-C-N	6.87	125.01	120.27
1	B	104	THR	CA-CB-OG1	-6.76	99.45	109.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	329	VAL	N-CA-CB	6.72	118.41	110.55
1	B	469	ARG	CB-CA-C	6.59	121.22	110.88
1	A	423	LEU	N-CA-CB	6.48	119.64	110.12
1	A	97	ARG	CD-NE-CZ	6.27	133.18	124.40
1	A	231	ILE	N-CA-C	-6.22	106.89	112.12
1	A	124	VAL	N-CA-CB	6.14	117.73	110.55
1	A	342	ARG	CB-CA-C	6.11	120.37	109.38
1	A	104	THR	CA-CB-OG1	-6.10	100.45	109.60
1	B	329	VAL	N-CA-CB	6.08	117.66	110.55
1	B	97	ARG	CD-NE-CZ	6.06	132.89	124.40
1	B	424	ARG	CB-CA-C	6.04	120.81	110.79
1	B	124	VAL	N-CA-CB	5.95	117.51	110.55
1	A	84	ASP	CB-CA-C	5.91	119.93	109.65
1	A	369	PRO	N-CD-CG	-5.91	94.34	103.20
1	B	461	PRO	N-CA-CB	5.86	110.00	103.44
1	B	423	LEU	N-CA-CB	5.81	118.67	110.12
1	A	169	ARG	CG-CD-NE	-5.75	99.36	112.00
1	A	120	ASP	CA-CB-CG	5.73	118.33	112.60
1	A	166	PHE	CA-CB-CG	5.70	119.50	113.80
1	A	182	ARG	CA-CB-CG	-5.65	102.80	114.10
1	B	322	ARG	N-CA-CB	5.56	115.67	110.39
1	B	406	LEU	N-CA-CB	-5.51	102.11	110.77
1	A	203	GLU	CB-CA-C	5.51	119.53	110.88
1	B	413	LYS	CG-CD-CE	5.50	123.96	111.30
1	B	120	ASP	CA-CB-CG	5.50	118.09	112.60
1	B	84	ASP	CB-CA-C	5.43	119.10	109.65
1	A	342	ARG	CB-CG-CD	5.35	123.61	111.30
1	A	413	LYS	CG-CD-CE	5.34	123.58	111.30
1	B	36	ARG	CA-CB-CG	5.29	124.68	114.10
1	B	37	MET	CG-SD-CE	5.28	112.51	100.90
1	B	64	ARG	CB-CA-C	5.26	119.22	109.70
1	A	322	ARG	O-C-N	5.25	124.53	120.38
1	A	408	PRO	CA-C-N	5.20	127.56	120.54
1	A	408	PRO	C-N-CA	5.20	127.56	120.54
1	B	45	ARG	CB-CA-C	5.13	120.13	110.63
1	A	336	LEU	N-CA-CB	5.12	117.64	110.12
1	A	206	HIS	CB-CA-C	-5.09	104.09	112.03
1	A	290	GLU	N-CA-C	-5.06	104.18	110.41
1	B	423	LEU	CB-CA-C	5.03	119.14	110.79

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	207	ARG	Sidechain
1	A	314	ARG	Sidechain
1	A	36	ARG	Sidechain
1	A	424	ARG	Sidechain
1	B	314	ARG	Sidechain
1	B	36	ARG	Sidechain
1	B	424	ARG	Sidechain

5.2 Too-close contacts [i](#)

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	3495	0	3481	132	0
1	B	3495	0	3481	129	0
2	A	10	0	4	3	0
2	B	10	0	4	6	0
All	All	7010	0	6970	256	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 18.

All (256) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:30:ALA:HB2	1:B:113:LEU:HB3	1.49	0.95
1:A:30:ALA:HB2	1:A:113:LEU:HB3	1.48	0.94
1:B:423:LEU:HD12	1:B:440:LEU:HD11	1.60	0.82
1:A:27:LEU:HD12	1:A:67:VAL:HG22	1.64	0.80
1:B:322:ARG:HB3	1:B:323:PRO:HD3	1.63	0.79
1:A:365:ARG:HH12	1:A:367:VAL:HG13	1.49	0.77
1:B:389:MET:HE2	1:B:397:PRO:HB3	1.64	0.77
1:A:423:LEU:HD12	1:A:440:LEU:HD11	1.66	0.77
1:A:440:LEU:N	1:A:440:LEU:HD12	2.00	0.76
1:A:322:ARG:HB3	1:A:323:PRO:HD3	1.68	0.75
1:A:162:GLY:O	1:A:163:ASP:HB2	1.84	0.75
1:B:36:ARG:HB2	1:B:36:ARG:CZ	2.18	0.72
1:A:314:ARG:NH2	1:A:314:ARG:HG3	2.05	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:440:LEU:HD12	1:B:440:LEU:N	2.03	0.71
1:A:310:TYR:O	1:A:312:GLY:N	2.24	0.71
1:A:310:TYR:C	1:A:312:GLY:H	2.00	0.70
1:A:244:THR:HG21	1:B:311:TRP:NE1	2.07	0.69
1:B:275:PHE:HA	2:B:501:ADE:N1	2.07	0.69
1:B:310:TYR:O	1:B:312:GLY:N	2.25	0.69
1:B:314:ARG:NH2	1:B:314:ARG:HG3	2.07	0.68
1:B:310:TYR:C	1:B:312:GLY:H	2.01	0.68
1:B:321:TRP:O	1:B:325:ILE:HG13	1.94	0.68
1:A:310:TYR:C	1:A:312:GLY:N	2.50	0.67
1:A:314:ARG:HG3	1:A:314:ARG:HH21	1.58	0.67
1:B:421:GLU:OE1	1:B:421:GLU:HA	1.92	0.67
1:B:118:LEU:HD21	1:B:150:GLU:HG2	1.78	0.66
1:B:310:TYR:C	1:B:312:GLY:N	2.53	0.66
1:B:423:LEU:HD23	1:B:474:LEU:HD23	1.77	0.66
1:A:195:LEU:CD1	1:A:272:VAL:HG23	2.27	0.65
1:B:314:ARG:HG3	1:B:314:ARG:HH21	1.61	0.65
1:B:125:GLU:O	1:B:129:ARG:HG3	1.96	0.65
1:A:219:MET:HE1	2:A:501:ADE:C4	2.33	0.64
1:A:195:LEU:HD13	1:A:201:VAL:HA	1.79	0.64
1:A:227:TRP:CE3	1:A:227:TRP:HA	2.33	0.64
1:B:80:TYR:HD1	1:B:84:ASP:HB2	1.64	0.63
1:B:195:LEU:HD13	1:B:201:VAL:HA	1.79	0.63
1:A:118:LEU:HD21	1:A:150:GLU:HG2	1.81	0.63
1:B:30:ALA:CB	1:B:113:LEU:HB3	2.27	0.62
1:A:144:ALA:CB	1:A:162:GLY:HA2	2.28	0.62
1:A:287:GLY:HA3	1:A:345:PHE:CZ	2.35	0.62
1:A:200:GLU:O	1:A:203:GLU:HG3	2.00	0.61
1:A:176:GLY:HA2	1:A:362:ILE:HG21	1.82	0.61
1:B:276:LEU:HD12	1:B:350:VAL:HG21	1.82	0.61
1:A:41:VAL:HG23	1:A:44:LEU:HG	1.81	0.61
1:A:321:TRP:O	1:A:325:ILE:HG13	2.01	0.61
1:A:276:LEU:HD13	1:A:358:LEU:HD12	1.81	0.61
1:A:222:LEU:HD21	1:A:261:PHE:CE1	2.36	0.61
1:B:176:GLY:HA2	1:B:362:ILE:HG21	1.83	0.61
1:B:287:GLY:HA3	1:B:345:PHE:CZ	2.36	0.60
1:A:423:LEU:HD23	1:A:474:LEU:HD23	1.83	0.60
1:B:227:TRP:HA	1:B:227:TRP:CE3	2.35	0.60
1:A:80:TYR:HD1	1:A:84:ASP:HB2	1.65	0.60
1:A:125:GLU:O	1:A:129:ARG:HG3	2.02	0.60
1:B:274:GLU:O	2:B:501:ADE:N6	2.25	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:400:TYR:HA	1:B:456:ALA:O	2.02	0.60
1:A:376:ALA:O	1:A:380:VAL:HG23	2.02	0.60
1:A:276:LEU:HD12	1:A:350:VAL:HG21	1.84	0.59
1:A:314:ARG:HH21	1:A:314:ARG:CG	2.16	0.59
1:A:400:TYR:HA	1:A:456:ALA:O	2.02	0.59
1:B:285:GLY:HA3	1:B:329:VAL:HG21	1.85	0.58
1:A:54:PHE:HA	1:A:374:ARG:NH1	2.17	0.58
1:B:440:LEU:HD23	1:B:452:VAL:HG21	1.86	0.58
1:A:415:ASN:C	1:A:415:ASN:OD1	2.45	0.58
1:A:440:LEU:HD23	1:A:452:VAL:HG21	1.86	0.58
1:B:314:ARG:HH21	1:B:314:ARG:CG	2.15	0.58
1:A:315:PHE:CD2	1:A:316:PRO:HA	2.39	0.57
1:A:430:TYR:HA	1:A:437:GLY:HA3	1.86	0.57
1:B:430:TYR:HA	1:B:437:GLY:HA3	1.85	0.57
1:B:80:TYR:CD1	1:B:84:ASP:HB2	2.38	0.57
1:B:428:LEU:HD12	1:B:474:LEU:CD2	2.35	0.57
1:A:252:GLU:O	1:B:253:VAL:HG13	2.04	0.57
1:B:152:GLY:HA2	1:B:157:ALA:O	2.03	0.57
1:B:376:ALA:O	1:B:380:VAL:HG23	2.04	0.57
1:A:428:LEU:HD12	1:A:474:LEU:CD2	2.35	0.57
1:B:315:PHE:CD2	1:B:316:PRO:HA	2.40	0.57
1:A:295:VAL:HG11	1:A:333:GLY:HA3	1.87	0.57
1:A:414:LEU:HD23	1:A:418:THR:CG2	2.35	0.57
1:B:222:LEU:HD21	1:B:261:PHE:CE1	2.40	0.57
1:B:283:PRO:HB2	1:B:325:ILE:CD1	2.35	0.57
1:A:80:TYR:CD1	1:A:84:ASP:HB2	2.40	0.56
1:B:144:ALA:CB	1:B:162:GLY:HA2	2.36	0.56
1:B:276:LEU:HD13	1:B:358:LEU:HD12	1.86	0.56
1:A:80:TYR:OH	1:A:439:ALA:HA	2.05	0.56
1:A:440:LEU:HD12	1:A:440:LEU:H	1.71	0.56
1:A:144:ALA:HB2	1:A:162:GLY:HA2	1.87	0.56
1:B:417:ARG:HG3	1:B:417:ARG:HH11	1.71	0.56
1:A:30:ALA:CB	1:A:113:LEU:HB3	2.28	0.55
1:A:227:TRP:HA	1:A:227:TRP:HE3	1.71	0.55
1:A:310:TYR:O	1:A:310:TYR:CD1	2.60	0.55
1:B:80:TYR:OH	1:B:439:ALA:HA	2.07	0.55
1:B:428:LEU:HD12	1:B:474:LEU:HD23	1.89	0.55
1:B:367:VAL:O	1:B:368:GLY:C	2.50	0.55
1:B:66:VAL:HG13	1:B:103:LEU:HD12	1.87	0.55
1:B:414:LEU:HD23	1:B:418:THR:CG2	2.37	0.55
1:B:415:ASN:OD1	1:B:415:ASN:C	2.47	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:244:THR:HG21	1:B:311:TRP:HE1	1.73	0.54
1:A:219:MET:HE1	2:A:501:ADE:C5	2.43	0.54
1:B:297:VAL:O	1:B:297:VAL:HG13	2.07	0.54
1:A:428:LEU:HD12	1:A:474:LEU:HD23	1.89	0.53
1:B:41:VAL:HG23	1:B:44:LEU:HG	1.90	0.53
1:B:54:PHE:HA	1:B:374:ARG:NH1	2.23	0.53
1:B:227:TRP:HA	1:B:227:TRP:HE3	1.73	0.53
1:B:301:HIS:CE1	1:B:369:PRO:HB2	2.44	0.53
1:B:219:MET:HE1	2:B:501:ADE:C4	2.44	0.53
1:B:178:GLU:HB3	1:B:182:ARG:NH2	2.23	0.53
1:B:310:TYR:O	1:B:310:TYR:CD1	2.62	0.53
1:A:152:GLY:HA2	1:A:157:ALA:O	2.08	0.53
1:A:414:LEU:HD23	1:A:418:THR:HG22	1.90	0.53
1:B:48:ALA:C	1:B:50:ARG:N	2.66	0.53
1:A:417:ARG:HH11	1:A:417:ARG:HG3	1.72	0.53
1:B:141:ASP:O	1:B:366:LYS:HE2	2.09	0.53
1:B:298:VAL:O	1:B:384:VAL:HG11	2.08	0.53
1:A:298:VAL:O	1:A:384:VAL:HG11	2.10	0.52
1:B:275:PHE:CD1	2:B:501:ADE:C2	2.97	0.52
1:A:141:ASP:O	1:A:366:LYS:HE2	2.09	0.52
1:A:385:GLY:O	1:A:386:ALA:C	2.53	0.52
1:B:238:ARG:O	1:B:242:ARG:HG3	2.09	0.51
1:B:314:ARG:NH2	1:B:314:ARG:CG	2.73	0.51
1:A:297:VAL:O	1:A:297:VAL:HG13	2.10	0.51
1:B:304:VAL:HG22	1:B:306:SER:HB3	1.92	0.51
1:A:236:ILE:HG13	1:A:251:ALA:HA	1.92	0.51
1:B:414:LEU:HD23	1:B:418:THR:HG22	1.93	0.51
1:B:27:LEU:HD23	1:B:140:VAL:HB	1.93	0.51
1:B:295:VAL:HG11	1:B:333:GLY:HA3	1.93	0.50
1:A:441:HIS:HB2	1:A:453:GLU:HB2	1.92	0.50
1:B:346:GLY:HA3	1:B:367:VAL:HG21	1.92	0.50
1:A:195:LEU:HD11	1:A:272:VAL:HG23	1.93	0.50
1:B:42:PRO:C	1:B:44:LEU:H	2.19	0.50
1:B:48:ALA:O	1:B:49:GLU:C	2.52	0.50
1:B:400:TYR:HB3	1:B:457:LEU:HG	1.93	0.50
1:A:127:LEU:O	1:A:131:VAL:HG23	2.12	0.49
1:A:42:PRO:C	1:A:44:LEU:H	2.18	0.49
1:B:144:ALA:HB2	1:B:162:GLY:HA2	1.94	0.49
1:B:195:LEU:CD1	1:B:272:VAL:HG23	2.42	0.49
1:A:283:PRO:HB2	1:A:325:ILE:CD1	2.42	0.48
1:A:322:ARG:CB	1:A:323:PRO:HD3	2.41	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:284:SER:HB3	1:B:303:GLN:NE2	2.28	0.48
1:A:322:ARG:N	1:A:323:PRO:CD	2.75	0.48
1:A:362:ILE:O	1:A:362:ILE:CG2	2.61	0.48
1:B:219:MET:HE1	2:B:501:ADE:C5	2.48	0.48
1:B:373:VAL:HG22	1:B:400:TYR:OH	2.13	0.48
1:B:322:ARG:N	1:B:323:PRO:CD	2.76	0.48
1:A:386:ALA:CB	1:A:391:ARG:HH21	2.27	0.48
1:B:78:LEU:HD22	1:B:102:LEU:HD11	1.96	0.48
1:A:309:GLN:HB2	1:A:311:TRP:CE2	2.49	0.48
1:B:384:VAL:O	1:B:390:LEU:HA	2.14	0.48
1:A:172:SER:HB2	1:A:223:ASP:O	2.14	0.48
1:B:111:ARG:HB2	1:B:116:LEU:HD21	1.96	0.48
1:B:236:ILE:HG13	1:B:251:ALA:HA	1.96	0.48
1:A:174:SER:O	1:A:178:GLU:HG3	2.14	0.47
1:A:144:ALA:HB2	1:A:161:GLU:O	2.14	0.47
1:A:222:LEU:HD21	1:A:261:PHE:HE1	1.79	0.47
1:A:380:VAL:HG21	1:A:390:LEU:HD21	1.97	0.47
1:A:111:ARG:HB2	1:A:116:LEU:HD21	1.96	0.47
1:A:384:VAL:O	1:A:390:LEU:HA	2.14	0.47
1:B:48:ALA:O	1:B:50:ARG:N	2.48	0.47
1:A:405:LEU:HD21	1:A:454:LEU:HD23	1.96	0.47
1:A:310:TYR:HH	1:A:313:CYS:HG	1.62	0.47
1:A:176:GLY:CA	1:A:362:ILE:HG21	2.45	0.47
1:A:201:VAL:HG21	1:A:265:LEU:HD21	1.95	0.46
1:B:172:SER:HB2	1:B:223:ASP:O	2.15	0.46
1:B:284:SER:CB	1:B:303:GLN:NE2	2.78	0.46
1:A:27:LEU:HD23	1:A:140:VAL:HB	1.97	0.46
1:B:309:GLN:HB2	1:B:311:TRP:CE2	2.51	0.46
1:A:178:GLU:HB3	1:A:182:ARG:NH2	2.31	0.46
1:A:386:ALA:HB2	1:A:391:ARG:HH21	1.80	0.46
1:B:177:LYS:NZ	1:B:273:GLU:HG2	2.30	0.46
1:B:208:LEU:HB3	1:B:238:ARG:NH1	2.31	0.46
1:B:405:LEU:HD21	1:B:454:LEU:HD23	1.96	0.46
1:A:159:PRO:HB2	1:A:162:GLY:HA3	1.98	0.46
1:B:222:LEU:HD23	1:B:270:ALA:HB2	1.97	0.46
1:B:27:LEU:HD12	1:B:67:VAL:HG22	1.96	0.46
1:A:66:VAL:HG13	1:A:103:LEU:HD12	1.98	0.45
1:B:144:ALA:HB2	1:B:161:GLU:O	2.16	0.45
1:B:362:ILE:O	1:B:362:ILE:CG2	2.63	0.45
1:B:417:ARG:HG3	1:B:417:ARG:NH1	2.30	0.45
1:B:66:VAL:CG1	1:B:103:LEU:HD12	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:365:ARG:HG3	1:A:365:ARG:HH11	1.81	0.45
1:B:325:ILE:HG23	1:B:349:PHE:CE2	2.52	0.45
1:A:315:PHE:CG	1:A:316:PRO:HA	2.52	0.45
1:A:310:TYR:CE2	1:A:313:CYS:HB3	2.52	0.45
1:A:274:GLU:O	2:A:501:ADE:N6	2.47	0.45
1:A:244:THR:CG2	1:B:311:TRP:NE1	2.78	0.44
1:B:201:VAL:HG21	1:B:265:LEU:HD21	1.99	0.44
1:B:256:LEU:HD23	1:B:256:LEU:HA	1.80	0.44
1:B:441:HIS:HB2	1:B:453:GLU:HB2	1.99	0.44
1:A:222:LEU:HD23	1:A:270:ALA:HB2	1.99	0.44
1:A:396:ARG:HA	1:A:397:PRO:HD3	1.80	0.44
1:A:259:GLU:HG3	1:A:260:GLU:N	2.33	0.44
1:B:230:SER:OG	1:B:255:ARG:NH1	2.47	0.44
1:B:283:PRO:CG	1:B:325:ILE:HD11	2.48	0.44
1:B:304:VAL:O	1:B:305:LEU:C	2.61	0.44
1:A:306:SER:HA	1:B:244:THR:HG21	2.00	0.43
1:A:400:TYR:HB3	1:A:457:LEU:HG	2.00	0.43
1:B:152:GLY:CA	1:B:157:ALA:O	2.66	0.43
1:B:216:ARG:HH11	1:B:238:ARG:HB2	1.83	0.43
1:B:208:LEU:HD12	1:B:208:LEU:HA	1.85	0.43
1:A:42:PRO:C	1:A:44:LEU:N	2.76	0.43
1:B:127:LEU:O	1:B:131:VAL:HG23	2.18	0.43
1:B:176:GLY:CA	1:B:362:ILE:HG21	2.47	0.43
1:B:315:PHE:CG	1:B:316:PRO:HA	2.53	0.43
1:A:285:GLY:HA3	1:A:329:VAL:HG21	2.00	0.43
1:A:336:LEU:HD23	1:A:336:LEU:HA	1.81	0.43
1:A:417:ARG:HG3	1:A:417:ARG:NH1	2.33	0.43
1:A:238:ARG:O	1:A:242:ARG:HG3	2.19	0.42
1:A:265:LEU:HD12	1:A:265:LEU:HA	1.87	0.42
1:A:304:VAL:O	1:A:305:LEU:C	2.62	0.42
1:A:330:ARG:HH21	1:A:331:ARG:HG3	1.84	0.42
1:B:174:SER:O	1:B:178:GLU:HG3	2.19	0.42
1:A:370:SER:OG	1:A:371:HIS:N	2.52	0.42
1:B:336:LEU:HD23	1:B:336:LEU:HA	1.80	0.42
1:B:423:LEU:HD11	1:B:475:LEU:HD11	2.01	0.42
1:A:78:LEU:HD22	1:A:102:LEU:HD11	2.00	0.42
1:A:416:PRO:O	1:A:420:VAL:HG23	2.19	0.42
1:A:446:LEU:HD23	1:A:446:LEU:HA	1.84	0.42
1:A:66:VAL:CG1	1:A:103:LEU:HD12	2.49	0.42
1:A:405:LEU:HD21	1:A:454:LEU:CD2	2.50	0.42
1:B:222:LEU:HD12	1:B:232:GLY:HA2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:ILE:HA	1:A:103:LEU:O	2.19	0.42
1:A:256:LEU:HA	1:A:256:LEU:HD23	1.78	0.42
1:A:152:GLY:CA	1:A:157:ALA:O	2.68	0.42
1:A:208:LEU:HD12	1:A:208:LEU:HA	1.86	0.42
1:A:248:VAL:HG12	1:A:248:VAL:O	2.20	0.42
1:A:452:VAL:HG22	1:A:453:GLU:O	2.20	0.42
1:A:222:LEU:HD21	1:A:261:PHE:CD1	2.55	0.41
1:B:42:PRO:C	1:B:44:LEU:N	2.76	0.41
1:B:203:GLU:CG	1:B:207:ARG:HH12	2.33	0.41
1:B:330:ARG:HH21	1:B:331:ARG:HG3	1.85	0.41
1:B:428:LEU:HD12	1:B:474:LEU:HD22	2.02	0.41
1:A:113:LEU:O	1:A:117:VAL:HG23	2.20	0.41
1:A:369:PRO:O	1:A:373:VAL:HG23	2.20	0.41
1:B:196:ARG:HA	1:B:265:LEU:O	2.19	0.41
1:B:429:LEU:HD23	1:B:430:TYR:N	2.35	0.41
1:A:51:SER:O	1:A:54:PHE:HD2	2.02	0.41
1:B:388:GLY:O	1:B:389:MET:CB	2.69	0.41
1:B:423:LEU:CD1	1:B:440:LEU:HD11	2.42	0.41
1:B:416:PRO:O	1:B:420:VAL:HG23	2.20	0.41
1:A:409:GLU:H	1:A:409:GLU:HG2	1.53	0.41
1:B:265:LEU:HA	1:B:269:GLY:HA2	2.02	0.41
1:B:360:VAL:HB	2:B:501:ADE:C8	2.56	0.41
1:B:429:LEU:HD23	1:B:430:TYR:H	1.85	0.41
1:A:64:ARG:HH11	1:A:134:ARG:HG2	1.85	0.41
1:A:423:LEU:HD11	1:A:475:LEU:HD11	2.01	0.41
1:B:316:PRO:HD2	1:B:399:TYR:HD1	1.86	0.41
1:A:389:MET:HE2	1:A:391:ARG:HG2	2.01	0.41
1:A:428:LEU:HD13	1:A:471:ALA:HA	2.03	0.40
1:B:162:GLY:O	1:B:163:ASP:CB	2.68	0.40
1:A:78:LEU:O	1:A:82:LEU:HG	2.22	0.40
1:A:428:LEU:HD12	1:A:474:LEU:HD22	2.02	0.40
1:A:41:VAL:HA	1:A:42:PRO:HD3	1.96	0.40
1:A:217:ARG:HD3	1:A:252:GLU:OE1	2.22	0.40
1:B:111:ARG:HE	1:B:111:ARG:HB3	1.55	0.40
1:A:325:ILE:HG23	1:A:349:PHE:CE2	2.56	0.40
1:A:367:VAL:HG12	1:A:368:GLY:H	1.87	0.40
1:A:414:LEU:CD2	1:A:418:THR:HG22	2.50	0.40
1:B:24:THR:OG1	1:B:134:ARG:HG3	2.21	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	456/475 (96%)	421 (92%)	31 (7%)	4 (1%)	14	46
1	B	456/475 (96%)	418 (92%)	33 (7%)	5 (1%)	11	43
All	All	912/950 (96%)	839 (92%)	64 (7%)	9 (1%)	12	45

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	163	ASP
1	A	207	ARG
1	B	163	ASP
1	B	389	MET
1	A	311	TRP
1	A	386	ALA
1	B	311	TRP
1	B	386	ALA
1	B	390	LEU

5.3.2 Protein sidechains [i](#)

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	355/371 (96%)	321 (90%)	34 (10%)	8	32
1	B	355/371 (96%)	324 (91%)	31 (9%)	9	34
All	All	710/742 (96%)	645 (91%)	65 (9%)	8	33

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	ASN
1	A	36	ARG
1	A	38	LEU
1	A	44	LEU
1	A	53	CYS
1	A	133	GLU
1	A	161	GLU
1	A	191	THR
1	A	194	VAL
1	A	203	GLU
1	A	208	LEU
1	A	216	ARG
1	A	225	ILE
1	A	226	THR
1	A	227	TRP
1	A	230	SER
1	A	236	ILE
1	A	247	LEU
1	A	259	GLU
1	A	263	ARG
1	A	271	ILE
1	A	289	ILE
1	A	302	ASP
1	A	304	VAL
1	A	314	ARG
1	A	329	VAL
1	A	367	VAL
1	A	383	ARG
1	A	384	VAL
1	A	405	LEU
1	A	409	GLU
1	A	429	LEU
1	A	459	ARG
1	A	467	TYR
1	B	32	ASN
1	B	36	ARG
1	B	37	MET
1	B	38	LEU
1	B	44	LEU
1	B	133	GLU
1	B	161	GLU
1	B	191	THR

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Mol	Chain	Res	Type
1	B	194	VAL
1	B	203	GLU
1	B	216	ARG
1	B	219	MET
1	B	225	ILE
1	B	226	THR
1	B	227	TRP
1	B	236	ILE
1	B	247	LEU
1	B	263	ARG
1	B	271	ILE
1	B	289	ILE
1	B	302	ASP
1	B	304	VAL
1	B	314	ARG
1	B	329	VAL
1	B	384	VAL
1	B	405	LEU
1	B	417	ARG
1	B	420	VAL
1	B	424	ARG
1	B	459	ARG
1	B	467	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	288	HIS
1	A	303	GLN
1	A	441	HIS
1	A	472	GLN
1	B	303	GLN
1	B	441	HIS
1	B	472	GLN

5.3.3 RNA ⓘ

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	ADE	A	501	-	11,11,11	0.30	0	15,15,15	0.38	0
2	ADE	B	501	-	11,11,11	0.30	0	15,15,15	0.38	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
2	ADE	A	501	-	-	-	0/2/2/2
2	ADE	B	501	-	-	-	0/2/2/2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

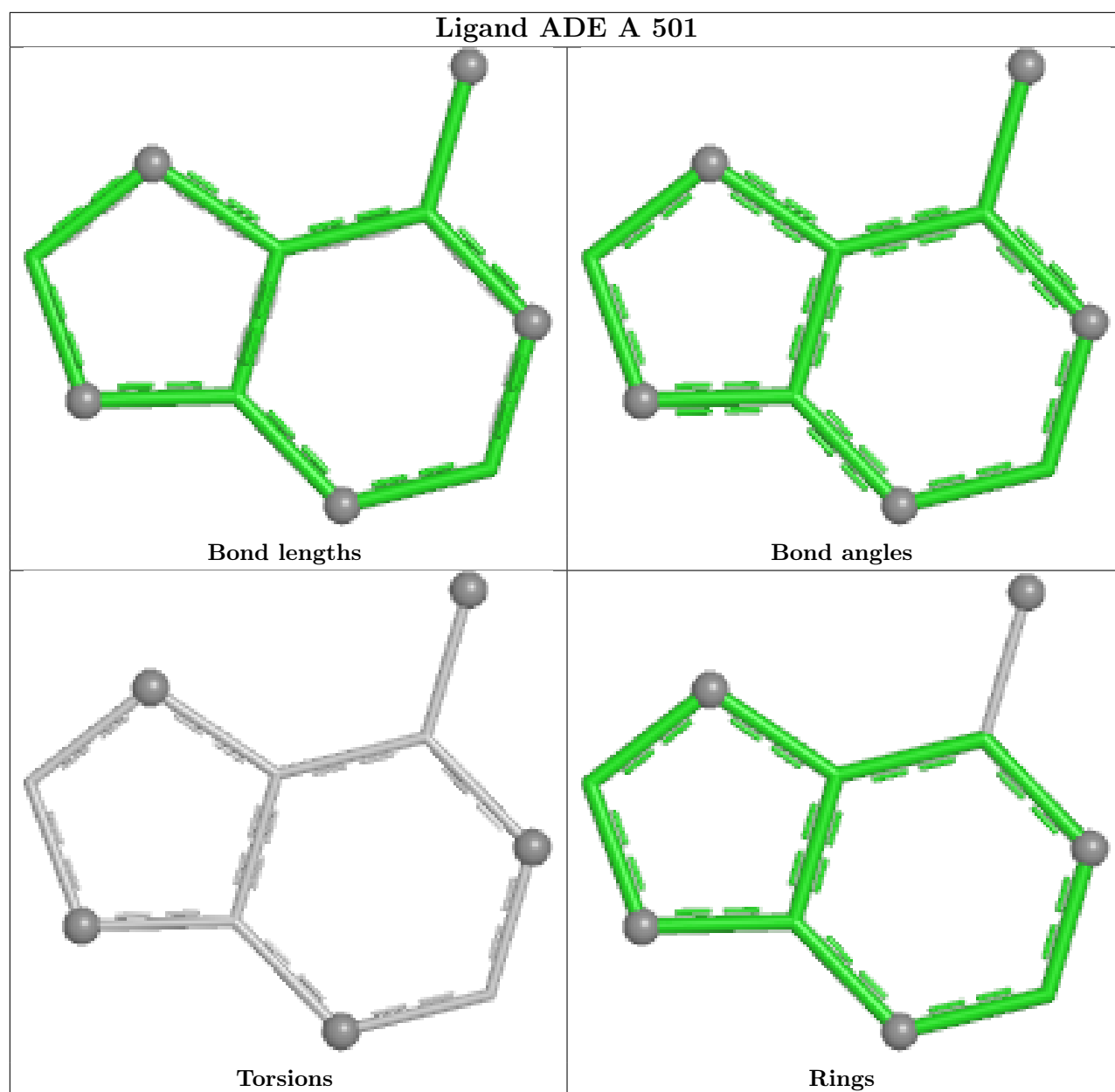
There are no torsion outliers.

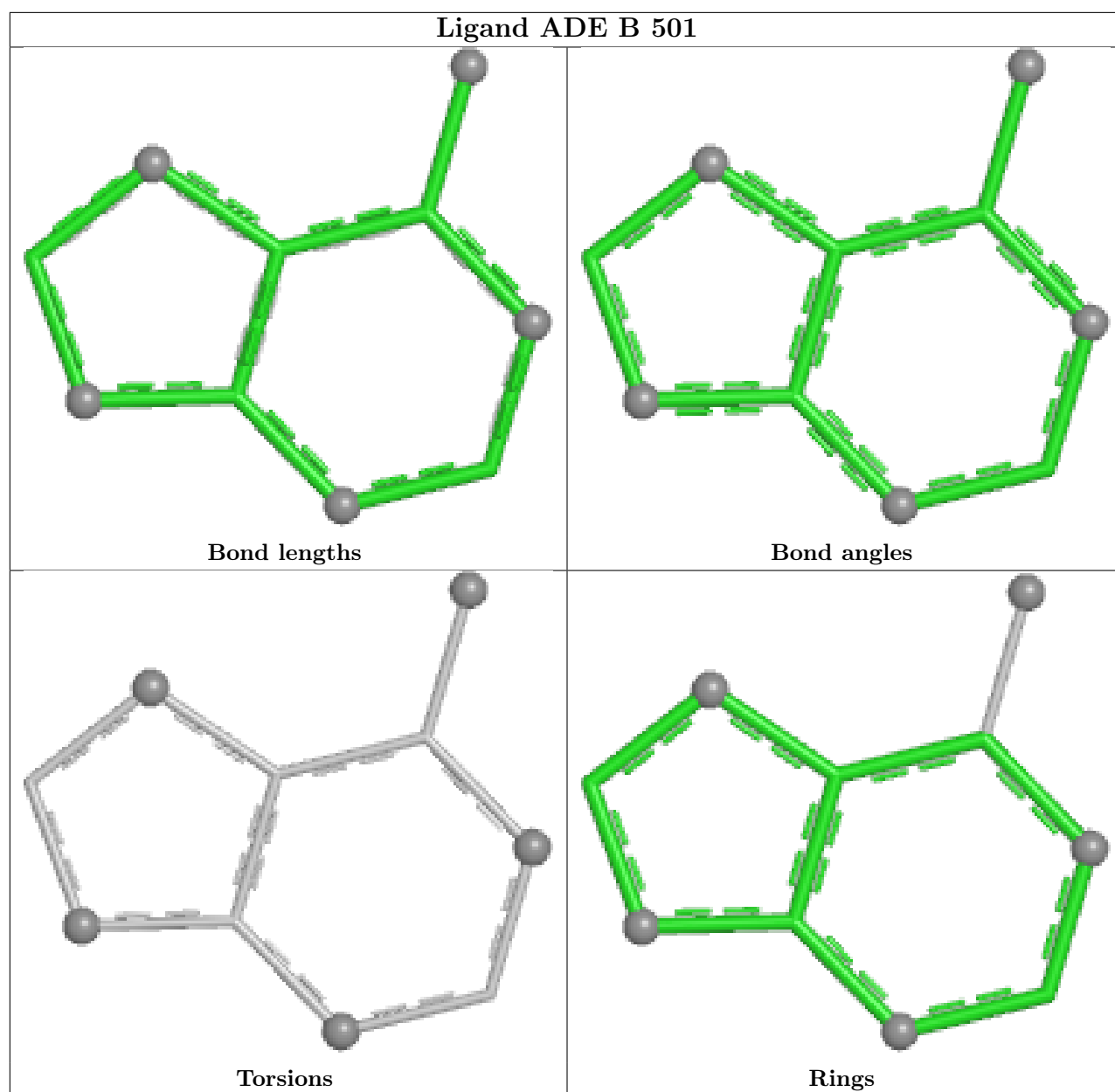
There are no ring outliers.

2 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	ADE	3	0
2	B	501	ADE	6	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 [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	458/475 (96%)	-0.47	2 (0%)	88 70	81, 161, 233, 321	0
1	B	458/475 (96%)	-0.36	7 (1%)	72 44	91, 172, 240, 380	0
All	All	916/950 (96%)	-0.41	9 (0%)	79 54	81, 167, 238, 380	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	123	LEU	3.6
1	B	127	LEU	2.7
1	B	470	ALA	2.5
1	B	265	LEU	2.2
1	B	73	VAL	2.2
1	A	139	ILE	2.2
1	B	204	ALA	2.2
1	A	26	ILE	2.2
1	B	117	VAL	2.1

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

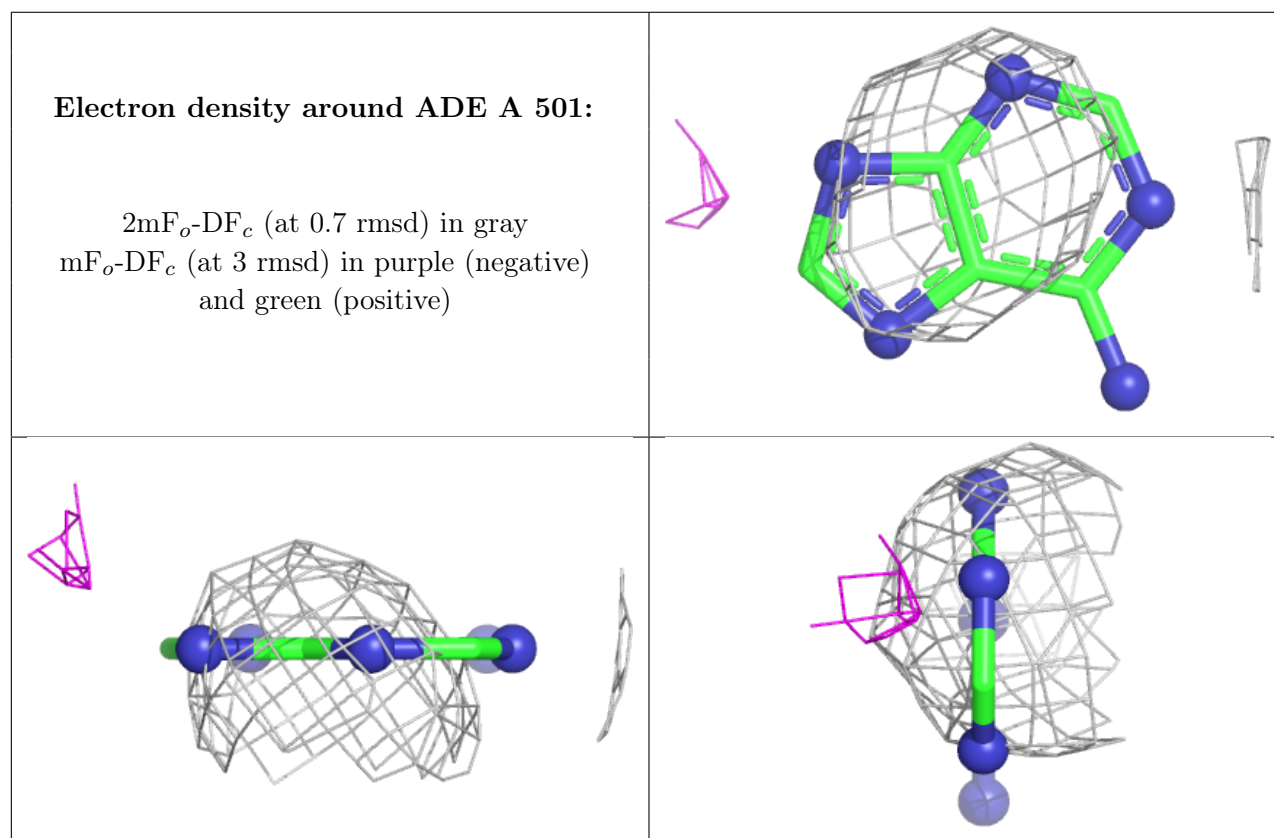
6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

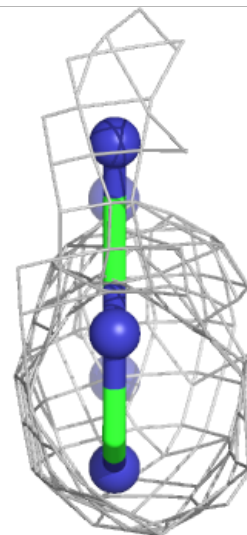
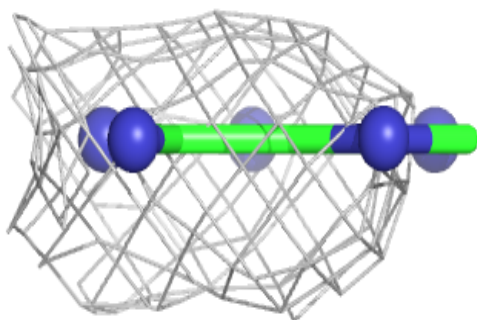
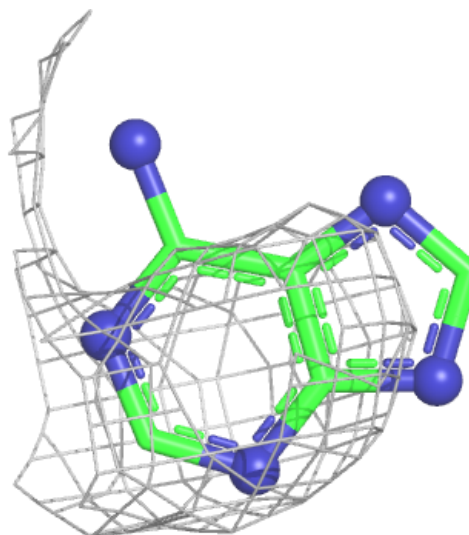
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ADE	A	501	10/10	0.93	0.10	151,169,240,291	0
2	ADE	B	501	10/10	0.94	0.09	133,168,209,243	0

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 ADE B 501:

$2mF_o-DF_c$ (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.