



Full wwPDB X-ray Structure Validation Report ⓘ

May 7, 2026 – 12:01 PM EDT

PDB ID : 1FCN / pdb_00001fcn
Title : Crystal Structure of the E. Coli AMPC Beta-Lactamase Mutant Q120L/Y150E Covalently Acylated with the Substrate Beta-Lactam LO-RACARBEF
Authors : Patera, A.; Blaszcak, L.C.; Shoichet, B.K.
Deposited on : 2000-07-18
Resolution : 2.35 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)
Xtrriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

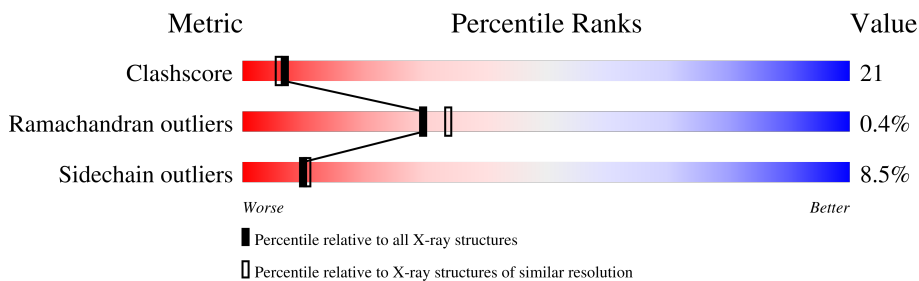
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	358	 61% 31% 6% ..
1	B	358	 63% 31% 6%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5715 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.

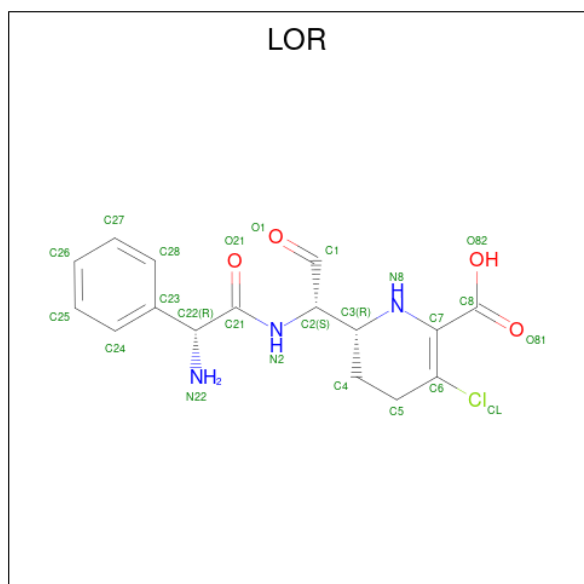
- Molecule 1 is a protein called BETA-LACTAMASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	352	Total	C	N	O	S	0	0	0
			2749	1776	465	502	6			
1	B	358	Total	C	N	O	S	0	0	0
			2792	1799	474	513	6			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	117	LEU	GLN	engineered mutation	UNP P00811
A	147	GLU	TYR	engineered mutation	UNP P00811
B	117	LEU	GLN	engineered mutation	UNP P00811
B	147	GLU	TYR	engineered mutation	UNP P00811

- Molecule 2 is LORACABEF (Open form) (CCD ID: LOR) (formula: C₁₆H₁₈ClN₃O₄).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Cl	N	O	0	0
			24	16	1	3	4		
2	B	1	Total	C	Cl	N	O	0	0
			24	16	1	3	4		

- Molecule 3 is water.

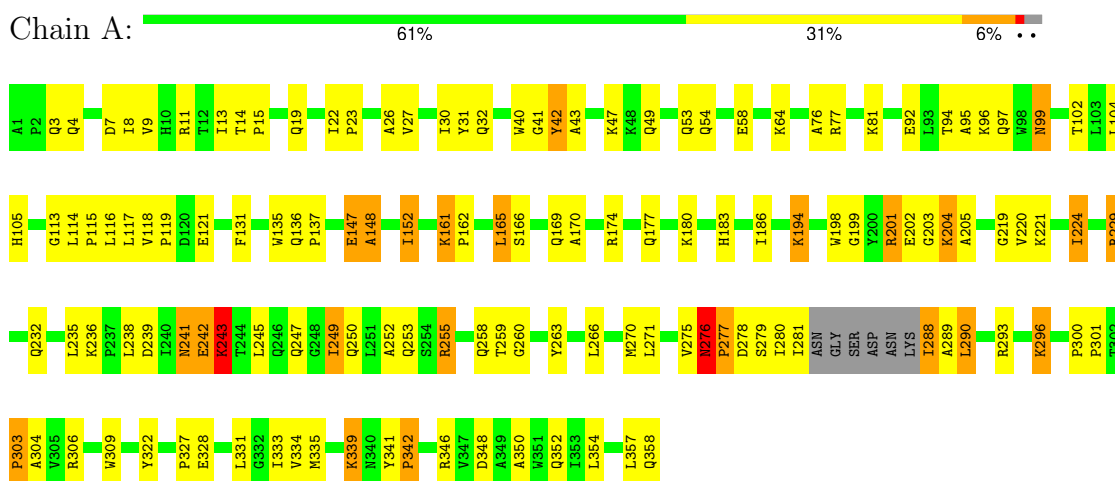
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	57	Total	O	0	0
			57	57		
3	B	69	Total	O	0	0
			69	69		

3 Residue-property plots

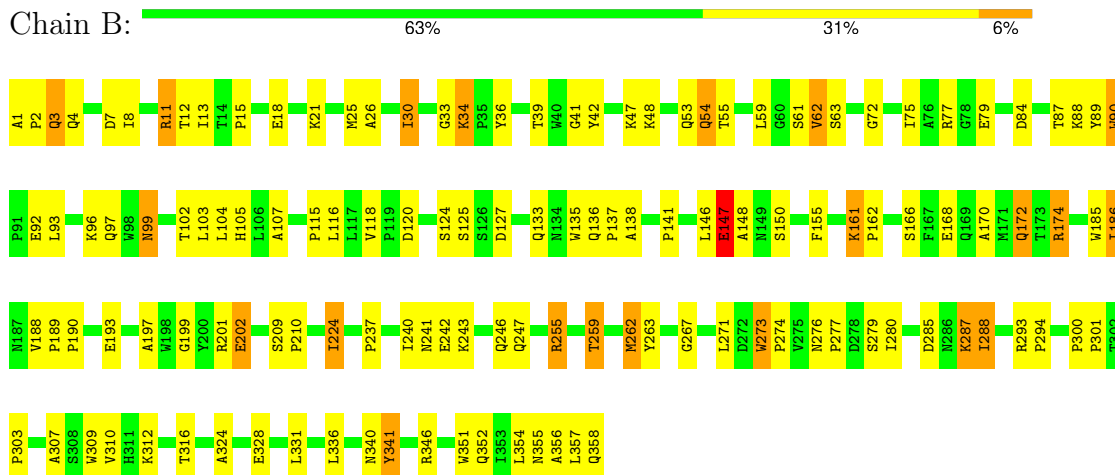
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

Note EDS was not executed.

• Molecule 1: BETA-LACTAMASE



• Molecule 1: BETA-LACTAMASE



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	119.36Å 76.26Å 98.36Å 90.00° 116.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.35	Depositor
% Data completeness (in resolution range)	95.1 (8.00-2.35)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.208 , 0.257	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5715	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	0.81	0/2827	1.23	23/3861 (0.6%)
1	B	0.82	0/2871	1.22	26/3921 (0.7%)
All	All	0.81	0/5698	1.23	49/7782 (0.6%)

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

There are no bond length outliers.

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	241	ASN	N-CA-C	12.56	125.05	111.36
1	A	219	GLY	N-CA-C	12.47	129.25	114.16
1	A	118	VAL	N-CA-C	-10.83	99.31	108.63
1	A	279	SER	N-CA-C	-10.65	99.46	111.71
1	B	118	VAL	N-CA-C	-9.76	99.39	108.95
1	B	186	ILE	N-CA-C	-9.49	101.52	111.58
1	A	148	ALA	N-CA-C	9.14	124.87	108.69
1	A	41	GLY	N-CA-C	8.47	126.09	111.78
1	B	125	SER	N-CA-C	-8.09	102.85	112.89
1	B	118	VAL	CA-C-N	-7.97	111.57	119.78
1	B	118	VAL	C-N-CA	-7.97	111.57	119.78
1	B	288	ILE	N-CA-C	-7.87	105.41	111.62
1	A	186	ILE	N-CA-C	-7.86	104.14	111.45
1	A	42	TYR	N-CA-C	7.74	122.05	109.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	220	VAL	N-CA-C	7.51	119.26	109.58
1	B	42	TYR	N-CA-C	7.46	121.60	109.59
1	A	342	PRO	N-CA-C	7.39	122.41	111.03
1	B	99	ASN	CB-CA-C	-7.21	99.77	110.16
1	A	276	ASN	CA-C-N	6.88	128.44	119.84
1	A	276	ASN	C-N-CA	6.88	128.44	119.84
1	B	41	GLY	N-CA-C	6.74	123.39	111.73
1	B	147	GLU	CA-C-O	-6.41	114.15	121.07
1	B	34	LYS	CA-C-N	6.20	126.71	120.14
1	B	34	LYS	C-N-CA	6.20	126.71	120.14
1	B	199	GLY	N-CA-C	-5.99	103.49	113.02
1	A	147	GLU	CA-C-N	5.94	132.31	122.33
1	A	147	GLU	C-N-CA	5.94	132.31	122.33
1	A	243	LYS	N-CA-C	5.87	117.36	110.97
1	B	316	THR	N-CA-C	-5.86	100.59	108.86
1	A	309	TRP	N-CA-C	-5.80	96.18	107.62
1	B	13	ILE	N-CA-C	5.67	115.86	110.42
1	B	33	GLY	N-CA-C	5.67	123.07	115.59
1	B	166	SER	N-CA-C	-5.53	103.61	110.41
1	A	43	ALA	N-CA-C	-5.52	104.90	111.69
1	A	22	ILE	CA-C-N	5.43	124.62	118.97
1	A	22	ILE	C-N-CA	5.43	124.62	118.97
1	B	240	ILE	N-CA-C	5.42	116.05	108.89
1	B	185	TRP	N-CA-C	5.38	117.57	109.23
1	A	241	ASN	N-CA-C	5.34	118.26	111.69
1	B	273	TRP	N-CA-C	-5.32	102.97	110.31
1	A	161	LYS	CA-C-N	-5.26	113.21	119.05
1	A	161	LYS	C-N-CA	-5.26	113.21	119.05
1	B	138	ALA	N-CA-C	-5.26	106.70	113.23
1	A	22	ILE	CA-C-O	5.22	122.23	119.15
1	B	309	TRP	N-CA-C	-5.20	99.31	108.20
1	A	99	ASN	N-CA-C	-5.20	102.22	109.96
1	B	341	TYR	CA-C-N	5.09	125.37	119.93
1	B	341	TYR	C-N-CA	5.09	125.37	119.93
1	B	90	TRP	N-CA-C	-5.02	100.22	108.75

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	42	TYR	Sidechain

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	2749	0	2736	136	0
1	B	2792	0	2774	107	0
2	A	24	0	16	7	0
2	B	24	0	16	1	0
3	A	57	0	0	4	0
3	B	69	0	0	3	0
All	All	5715	0	5542	235	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 21.

All (235) 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:243:LYS:H	1:A:243:LYS:CE	1.67	1.05
1:A:243:LYS:H	1:A:243:LYS:HE3	0.89	1.05
1:A:335:MET:HE1	1:A:350:ALA:HB2	1.41	1.02
1:A:194:LYS:HZ2	1:A:194:LYS:HB3	1.20	1.01
1:A:243:LYS:HE3	1:A:243:LYS:N	1.74	1.01
1:B:26:ALA:HB1	1:B:224:ILE:HG12	1.44	0.99
1:A:165:LEU:HB3	1:A:169:GLN:HG3	1.45	0.97
1:B:102:THR:H	1:B:105:HIS:HD2	1.05	0.95
1:A:116:LEU:HD21	2:A:961:LOR:H5A	1.54	0.87
1:A:259:THR:OG1	1:A:293:ARG:HD3	1.76	0.86
1:B:53:GLN:HA	1:B:224:ILE:HD11	1.58	0.85
1:B:102:THR:H	1:B:105:HIS:CD2	1.96	0.82
1:A:245:LEU:O	1:A:249:ILE:HD12	1.80	0.82
1:A:26:ALA:HB1	1:A:224:ILE:HG12	1.62	0.81
1:A:277:PRO:O	1:A:281:ILE:N	2.14	0.81
1:A:194:LYS:HZ2	1:A:194:LYS:CB	1.94	0.81
1:A:229:ARG:HA	1:A:232:GLN:HE21	1.45	0.80
1:A:288:ILE:HG22	1:A:288:ILE:O	1.80	0.80
1:B:96:LYS:HD2	1:B:99:ASN:HD21	1.46	0.80
1:A:341:TYR:HB2	1:A:342:PRO:HD2	1.63	0.79
1:A:166:SER:OG	1:A:169:GLN:HG2	1.81	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:277:PRO:O	1:A:281:ILE:HG13	1.83	0.79
1:A:102:THR:H	1:A:105:HIS:HD2	1.31	0.79
1:A:116:LEU:HD11	2:A:961:LOR:H4	1.63	0.79
1:B:96:LYS:HD2	1:B:99:ASN:ND2	1.98	0.78
1:B:11:ARG:HH11	1:B:11:ARG:HB2	1.50	0.77
1:B:190:PRO:O	1:B:193:GLU:HG3	1.85	0.76
1:A:194:LYS:HB3	1:A:194:LYS:NZ	1.93	0.75
1:B:287:LYS:HD2	1:B:288:ILE:HG13	1.67	0.75
1:B:328:GLU:O	1:B:328:GLU:HG2	1.87	0.75
1:B:34:LYS:HD3	1:B:36:TYR:OH	1.88	0.74
1:A:53:GLN:HA	1:A:224:ILE:HD11	1.70	0.73
1:A:183:HIS:HD2	3:A:998:HOH:O	1.70	0.73
1:A:331:LEU:HG	1:A:354:LEU:HD22	1.68	0.73
1:A:236:LYS:HE2	1:A:239:ASP:OD1	1.88	0.73
1:A:14:THR:HB	1:A:15:PRO:HD3	1.72	0.72
1:B:84:ASP:OD1	1:B:89:TYR:OH	2.09	0.71
1:B:259:THR:HG23	1:B:294:PRO:O	1.89	0.71
1:A:180:LYS:HB2	1:A:229:ARG:NH2	2.06	0.70
1:B:11:ARG:HB2	1:B:11:ARG:NH1	2.06	0.70
1:B:34:LYS:HD3	1:B:36:TYR:CZ	2.28	0.68
1:A:253:GLN:HG3	3:A:974:HOH:O	1.92	0.68
1:A:271:LEU:HD22	1:A:280:ILE:CG2	2.22	0.68
1:B:79:GLU:HB3	1:B:162:PRO:HB2	1.77	0.67
1:A:333:ILE:HD13	1:A:335:MET:HE2	1.76	0.66
1:A:252:ALA:HA	1:A:266:LEU:HB2	1.76	0.66
1:A:102:THR:H	1:A:105:HIS:CD2	2.15	0.65
1:A:277:PRO:HB3	1:A:281:ILE:HD12	1.79	0.65
1:A:335:MET:CE	1:A:350:ALA:HB2	2.20	0.65
1:A:341:TYR:CE2	1:A:346:ARG:HG2	2.33	0.64
1:A:247:GLN:HA	1:A:250:GLN:HE21	1.60	0.64
1:B:351:TRP:CH2	1:B:355:ASN:ND2	2.66	0.64
1:B:2:PRO:HD2	1:B:36:TYR:CE1	2.34	0.63
1:B:87:THR:O	1:B:87:THR:HG22	1.96	0.63
1:A:64:LYS:HE2	1:A:152:ILE:HG21	1.79	0.63
1:A:276:ASN:N	1:A:276:ASN:HD22	1.95	0.62
1:A:277:PRO:C	1:A:281:ILE:HG13	2.23	0.62
1:A:180:LYS:HB2	1:A:229:ARG:HH21	1.64	0.62
1:B:190:PRO:HA	1:B:193:GLU:HG3	1.80	0.62
1:A:96:LYS:HE3	1:A:99:ASN:HD21	1.65	0.62
1:A:243:LYS:N	1:A:243:LYS:CD	2.62	0.62
1:A:275:VAL:HG22	1:A:276:ASN:N	2.14	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1:ALA:CB	1:B:36:TYR:CD1	2.84	0.61
1:A:116:LEU:HD21	2:A:961:LOR:C5	2.28	0.61
1:A:4:GLN:O	1:A:8:ILE:HG22	2.00	0.61
1:A:136:GLN:HA	1:A:136:GLN:NE2	2.15	0.61
1:B:255:ARG:HG2	1:B:263:TYR:CB	2.32	0.60
1:A:115:PRO:HG2	1:A:131:PHE:HZ	1.65	0.60
1:A:194:LYS:CB	1:A:194:LYS:NZ	2.57	0.60
1:B:124:SER:OG	1:B:127:ASP:N	2.33	0.60
1:B:202:GLU:HA	1:B:202:GLU:OE1	2.00	0.60
1:A:276:ASN:N	1:A:276:ASN:ND2	2.49	0.59
1:B:59:LEU:HB3	1:B:62:VAL:HG12	1.85	0.59
1:A:276:ASN:ND2	1:A:276:ASN:H	2.01	0.59
1:B:53:GLN:HA	1:B:224:ILE:CD1	2.31	0.59
1:B:237:PRO:O	1:B:246:GLN:HG3	2.03	0.59
1:A:116:LEU:HD13	1:A:147:GLU:O	2.04	0.58
1:A:243:LYS:CE	1:A:243:LYS:N	2.47	0.58
1:A:229:ARG:HA	1:A:232:GLN:NE2	2.18	0.58
1:B:89:TYR:CE2	1:B:103:LEU:HD11	2.38	0.58
1:A:255:ARG:HD2	3:A:1001:HOH:O	2.02	0.58
1:A:255:ARG:HG2	1:A:263:TYR:HB3	1.84	0.58
1:A:7:ASP:O	1:A:11:ARG:HD3	2.04	0.58
1:A:255:ARG:HG2	1:A:263:TYR:CB	2.34	0.58
1:A:348:ASP:O	1:A:352:GLN:HB2	2.03	0.58
1:A:259:THR:HG1	1:A:293:ARG:HD3	1.69	0.58
1:B:303:PRO:HD3	3:B:980:HOH:O	2.02	0.58
1:B:11:ARG:NH1	1:B:11:ARG:CB	2.66	0.57
1:B:161:LYS:NZ	1:B:161:LYS:HB3	2.19	0.57
1:A:96:LYS:HE3	1:A:99:ASN:ND2	2.19	0.57
1:A:242:GLU:CD	1:A:242:GLU:H	2.11	0.57
1:A:357:LEU:O	1:A:358:GLN:CB	2.52	0.57
1:B:30:ILE:HD12	1:B:30:ILE:N	2.20	0.57
1:B:307:ALA:HB2	1:B:328:GLU:OE1	2.05	0.56
1:B:190:PRO:HA	1:B:193:GLU:CG	2.35	0.56
1:A:8:ILE:HG23	1:A:9:VAL:N	2.20	0.56
1:A:115:PRO:O	1:A:148:ALA:HB1	2.06	0.55
1:A:290:LEU:C	1:A:290:LEU:HD23	2.31	0.55
1:B:293:ARG:HG3	1:B:293:ARG:HH11	1.70	0.55
1:B:168:GLU:O	1:B:172:GLN:HB2	2.06	0.55
1:B:26:ALA:CB	1:B:224:ILE:HG12	2.29	0.55
1:B:170:ALA:O	1:B:174:ARG:HB2	2.08	0.54
1:A:301:PRO:HG2	1:B:300:PRO:CB	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:SER:H	1:A:169:GLN:CG	2.21	0.54
1:B:102:THR:N	1:B:105:HIS:HD2	1.89	0.54
1:B:116:LEU:HD11	2:B:961:LOR:H5A	1.90	0.54
1:A:8:ILE:HG23	1:A:9:VAL:H	1.73	0.54
1:B:1:ALA:HB2	1:B:36:TYR:CD1	2.43	0.53
1:B:26:ALA:HB1	1:B:224:ILE:CG1	2.30	0.53
1:B:161:LYS:HB2	1:B:162:PRO:HD3	1.90	0.53
1:A:288:ILE:O	1:A:290:LEU:N	2.41	0.53
1:A:64:LYS:HE2	1:A:152:ILE:CG2	2.39	0.53
1:B:168:GLU:HA	1:B:186:ILE:HD12	1.90	0.53
1:A:116:LEU:CD2	2:A:961:LOR:H5A	2.33	0.53
1:B:93:LEU:HD12	1:B:133:GLN:OE1	2.09	0.53
1:A:271:LEU:HD22	1:A:280:ILE:HG23	1.90	0.52
1:B:328:GLU:O	1:B:328:GLU:CG	2.56	0.52
1:B:357:LEU:O	1:B:358:GLN:CB	2.58	0.52
1:A:322:TYR:HD2	1:A:335:MET:HE3	1.75	0.52
1:A:170:ALA:O	1:A:174:ARG:HB2	2.10	0.51
1:A:306:ARG:HB3	1:A:328:GLU:HB3	1.91	0.51
1:B:188:VAL:HG23	1:B:188:VAL:O	2.10	0.51
1:B:12:THR:O	1:B:15:PRO:HD2	2.11	0.51
1:A:242:GLU:CD	1:A:242:GLU:N	2.69	0.50
1:B:147:GLU:O	1:B:148:ALA:HB2	2.12	0.50
1:A:115:PRO:CG	1:A:131:PHE:HZ	2.24	0.50
1:A:260:GLY:HA3	1:A:293:ARG:HH12	1.76	0.50
1:A:47:LYS:O	1:A:49:GLN:NE2	2.36	0.50
1:A:116:LEU:HD12	1:A:148:ALA:HA	1.94	0.50
1:A:247:GLN:HA	1:A:250:GLN:NE2	2.27	0.50
1:B:1:ALA:O	1:B:2:PRO:C	2.55	0.50
1:B:97:GLN:NE2	1:B:135:TRP:O	2.45	0.50
1:B:1:ALA:HB2	1:B:36:TYR:HD1	1.77	0.49
1:A:116:LEU:HD11	2:A:961:LOR:C4	2.40	0.49
1:B:147:GLU:HG3	1:B:148:ALA:N	2.28	0.49
1:A:104:LEU:HD22	1:B:300:PRO:HD3	1.95	0.49
1:A:58:GLU:HA	1:A:221:LYS:HG3	1.95	0.49
1:B:30:ILE:O	1:B:331:LEU:HD12	2.13	0.49
1:B:271:LEU:HD22	1:B:280:ILE:HG22	1.94	0.49
1:A:277:PRO:HB3	1:A:281:ILE:CD1	2.41	0.49
1:A:116:LEU:HD12	1:A:116:LEU:O	2.13	0.48
1:B:262:MET:HE3	1:B:262:MET:HB3	1.71	0.48
1:A:77:ARG:HH21	1:A:174:ARG:HG2	1.78	0.48
1:B:287:LYS:CD	1:B:288:ILE:HG13	2.42	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:90:TRP:CE3	1:B:93:LEU:HD22	2.48	0.48
1:B:55:THR:CG2	1:B:197:ALA:HB2	2.43	0.48
1:A:258:GLN:HB3	1:A:296:LYS:HG2	1.94	0.48
1:A:54:GLN:HB3	3:A:1006:HOH:O	2.12	0.48
1:A:117:LEU:HD22	2:A:961:LOR:C25	2.43	0.48
1:B:202:GLU:OE1	1:B:202:GLU:CA	2.62	0.48
1:A:31:TYR:CE1	1:A:32:GLN:NE2	2.81	0.48
1:A:301:PRO:HG2	1:B:300:PRO:HB2	1.94	0.48
1:B:1:ALA:HB1	1:B:36:TYR:CD1	2.49	0.48
1:A:199:GLY:O	1:A:205:ALA:HA	2.14	0.48
1:A:235:LEU:CD2	1:A:327:PRO:HA	2.44	0.47
1:A:275:VAL:CG2	1:A:276:ASN:N	2.77	0.47
1:B:259:THR:HG21	1:B:293:ARG:HB3	1.96	0.47
1:B:285:ASP:OD1	1:B:287:LYS:HG3	2.14	0.47
1:A:27:VAL:HG23	1:A:40:TRP:HZ3	1.80	0.47
1:A:238:LEU:HD12	1:B:243:LYS:HD3	1.95	0.47
1:B:271:LEU:HD22	1:B:280:ILE:CG2	2.44	0.47
1:B:331:LEU:HG	1:B:354:LEU:HD22	1.96	0.47
1:A:135:TRP:CZ3	1:A:137:PRO:HG3	2.50	0.47
1:B:161:LYS:HB3	1:B:161:LYS:HZ3	1.78	0.47
1:A:166:SER:H	1:A:169:GLN:CD	2.23	0.47
1:B:255:ARG:HG2	1:B:263:TYR:HB3	1.96	0.47
1:B:141:PRO:HD3	3:B:974:HOH:O	2.14	0.46
1:A:335:MET:HE1	1:A:350:ALA:CB	2.30	0.46
1:A:306:ARG:O	1:A:327:PRO:HG2	2.15	0.46
1:A:23:PRO:HD2	1:A:339:LYS:HB2	1.98	0.46
1:B:331:LEU:HD13	3:B:994:HOH:O	2.15	0.46
1:A:243:LYS:H	1:A:243:LYS:CD	2.18	0.46
1:B:115:PRO:O	1:B:150:SER:OG	2.28	0.45
1:B:340:ASN:C	1:B:340:ASN:OD1	2.59	0.45
1:A:339:LYS:C	1:A:339:LYS:HD2	2.41	0.45
1:A:30:ILE:N	1:A:30:ILE:HD12	2.32	0.45
1:B:276:ASN:O	1:B:279:SER:HB3	2.17	0.45
1:A:161:LYS:HB2	1:A:162:PRO:HD3	1.99	0.45
1:B:87:THR:O	1:B:87:THR:CG2	2.63	0.45
1:B:277:PRO:HG3	1:B:351:TRP:CZ2	2.52	0.45
1:A:9:VAL:O	1:A:13:ILE:HG12	2.17	0.44
1:A:31:TYR:CE1	1:A:32:GLN:CD	2.95	0.44
1:B:341:TYR:CZ	1:B:346:ARG:HG2	2.52	0.44
1:A:77:ARG:NH2	1:A:174:ARG:HG2	2.33	0.44
1:A:135:TRP:CH2	1:A:137:PRO:HG3	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:61:SER:C	1:B:63:SER:N	2.75	0.44
1:A:270:MET:C	1:A:271:LEU:HD12	2.43	0.44
1:A:301:PRO:HD2	1:B:301:PRO:HD2	1.99	0.44
1:B:54:GLN:HE21	1:B:54:GLN:HB2	1.58	0.44
1:A:243:LYS:N	1:A:243:LYS:HD3	2.31	0.44
1:B:107:ALA:HB2	1:B:155:PHE:CD1	2.53	0.44
1:B:77:ARG:HH21	1:B:174:ARG:HG2	1.83	0.43
1:A:303:PRO:O	1:A:304:ALA:C	2.61	0.43
1:B:90:TRP:C	1:B:92:GLU:H	2.26	0.43
1:A:7:ASP:HB3	1:A:11:ARG:NH1	2.33	0.43
1:A:76:ALA:HB1	1:A:247:GLN:OE1	2.19	0.43
1:A:136:GLN:NE2	1:A:136:GLN:CA	2.78	0.43
1:A:14:THR:CB	1:A:15:PRO:HD3	2.46	0.43
1:A:238:LEU:CD1	1:B:243:LYS:HD3	2.48	0.43
1:A:201:ARG:NH1	1:A:202:GLU:OE2	2.51	0.43
1:A:341:TYR:CZ	1:A:346:ARG:HG2	2.53	0.43
1:A:116:LEU:HD21	2:A:961:LOR:C4	2.48	0.43
1:B:7:ASP:OD2	1:B:11:ARG:NH2	2.45	0.43
1:B:59:LEU:HB3	1:B:62:VAL:CG1	2.48	0.43
1:A:281:ILE:HG22	1:A:281:ILE:O	2.20	0.42
1:B:8:ILE:HG12	1:B:356:ALA:HB3	2.01	0.42
1:A:203:GLY:C	1:A:204:LYS:HD2	2.44	0.42
1:B:293:ARG:HG3	1:B:293:ARG:NH1	2.33	0.42
1:A:300:PRO:HD3	1:B:104:LEU:HD22	2.01	0.42
1:B:90:TRP:C	1:B:92:GLU:N	2.77	0.42
1:A:301:PRO:HG2	1:B:300:PRO:HB3	2.02	0.42
1:B:72:GLY:HA2	1:B:75:ILE:HD12	2.02	0.42
1:B:267:GLY:O	1:B:312:LYS:HD3	2.20	0.41
1:B:273:TRP:CD2	1:B:274:PRO:HA	2.55	0.41
1:A:94:THR:O	1:A:95:ALA:C	2.63	0.41
1:A:113:GLY:O	1:A:114:LEU:C	2.61	0.41
1:A:333:ILE:HG12	1:A:334:VAL:N	2.36	0.41
1:B:136:GLN:HA	1:B:137:PRO:HD3	1.91	0.41
1:B:25:MET:HG3	1:B:336:LEU:O	2.21	0.41
1:B:84:ASP:CG	1:B:89:TYR:HH	2.23	0.41
1:A:31:TYR:O	1:A:32:GLN:C	2.64	0.41
1:A:92:GLU:N	1:A:92:GLU:CD	2.78	0.41
1:B:209:SER:O	1:B:210:PRO:C	2.64	0.41
1:B:146:LEU:O	1:B:147:GLU:C	2.63	0.41
1:B:310:VAL:O	1:B:324:ALA:HA	2.20	0.41
1:B:341:TYR:HD1	1:B:341:TYR:H	1.67	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:95:ALA:HB1	1:A:97:GLN:OE1	2.20	0.41
1:A:119:PRO:HD2	1:A:131:PHE:CE1	2.56	0.41
1:A:198:TRP:HB3	1:A:205:ALA:HB1	2.03	0.40
1:A:241:ASN:O	1:A:243:LYS:HD3	2.20	0.40
1:B:96:LYS:O	1:B:99:ASN:ND2	2.54	0.40
1:B:201:ARG:HG2	1:B:202:GLU:HG2	2.03	0.40
1:B:188:VAL:HA	1:B:189:PRO:HD3	1.95	0.40
1:A:271:LEU:CD2	1:A:280:ILE:CG2	2.97	0.40
1:A:15:PRO:O	1:A:19:GLN:HB2	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	348/358 (97%)	339 (97%)	8 (2%)	1 (0%)	36	43
1	B	356/358 (99%)	339 (95%)	15 (4%)	2 (1%)	21	24
All	All	704/716 (98%)	678 (96%)	23 (3%)	3 (0%)	30	34

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	289	ALA
1	B	3	GLN
1	B	147	GLU

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	286/292 (98%)	263 (92%)	23 (8%)	11	12
1	B	291/292 (100%)	265 (91%)	26 (9%)	9	10
All	All	577/584 (99%)	528 (92%)	49 (8%)	10	11

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	81	LYS
1	A	121	GLU
1	A	152	ILE
1	A	165	LEU
1	A	177	GLN
1	A	194	LYS
1	A	201	ARG
1	A	204	LYS
1	A	224	ILE
1	A	229	ARG
1	A	242	GLU
1	A	243	LYS
1	A	249	ILE
1	A	255	ARG
1	A	276	ASN
1	A	277	PRO
1	A	278	ASP
1	A	288	ILE
1	A	290	LEU
1	A	296	LYS
1	A	303	PRO
1	A	339	LYS
1	B	3	GLN
1	B	4	GLN
1	B	11	ARG
1	B	18	GLU
1	B	21	LYS
1	B	30	ILE
1	B	39	THR
1	B	47	LYS
1	B	48	LYS
1	B	54	GLN

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Mol	Chain	Res	Type
1	B	62	VAL
1	B	88	LYS
1	B	120	ASP
1	B	147	GLU
1	B	161	LYS
1	B	172	GLN
1	B	174	ARG
1	B	202	GLU
1	B	224	ILE
1	B	242	GLU
1	B	247	GLN
1	B	255	ARG
1	B	259	THR
1	B	262	MET
1	B	287	LYS
1	B	352	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	20	GLN
1	A	32	GLN
1	A	99	ASN
1	A	105	HIS
1	A	134	ASN
1	A	136	GLN
1	A	183	HIS
1	A	232	GLN
1	A	250	GLN
1	A	253	GLN
1	A	276	ASN
1	A	352	GLN
1	B	10	HIS
1	B	19	GLN
1	B	20	GLN
1	B	32	GLN
1	B	54	GLN
1	B	99	ASN
1	B	105	HIS
1	B	134	ASN
1	B	187	ASN

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Mol	Chain	Res	Type
1	B	232	GLN
1	B	250	GLN
1	B	286	ASN
1	B	352	GLN
1	B	355	ASN

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	LOR	A	961	1	21,25,25	4.53	15 (71%)	21,34,34	2.00	6 (28%)
2	LOR	B	961	1	21,25,25	3.78	9 (42%)	21,34,34	1.93	5 (23%)

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	LOR	A	961	1	-	6/19/35/35	0/2/2/2
2	LOR	B	961	1	-	6/19/35/35	0/2/2/2

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	961	LOR	C5-C6	12.67	1.74	1.50
2	B	961	LOR	C5-C6	10.72	1.70	1.50
2	A	961	LOR	O82-C8	10.02	1.57	1.30
2	B	961	LOR	O82-C8	9.19	1.55	1.30
2	A	961	LOR	C2-N2	5.67	1.54	1.46
2	A	961	LOR	C7-N8	5.59	1.50	1.37
2	B	961	LOR	C24-C23	4.28	1.45	1.39
2	B	961	LOR	C26-C25	4.04	1.47	1.38
2	A	961	LOR	C24-C23	4.00	1.45	1.39
2	A	961	LOR	C27-C28	3.71	1.45	1.38
2	A	961	LOR	C27-C26	3.65	1.46	1.38
2	A	961	LOR	C26-C25	3.39	1.45	1.38
2	B	961	LOR	C27-C26	3.29	1.45	1.38
2	A	961	LOR	C28-C23	3.28	1.44	1.39
2	A	961	LOR	C23-C22	-3.22	1.49	1.52
2	B	961	LOR	C2-N2	3.19	1.50	1.46
2	B	961	LOR	C7-N8	3.02	1.44	1.37
2	B	961	LOR	C27-C28	2.79	1.43	1.38
2	B	961	LOR	C28-C23	2.74	1.43	1.39
2	A	961	LOR	C3-N8	2.64	1.52	1.47
2	A	961	LOR	O81-C8	2.52	1.28	1.22
2	A	961	LOR	C7-C8	2.27	1.51	1.48
2	A	961	LOR	C2-C3	2.24	1.57	1.54
2	A	961	LOR	C25-C24	2.07	1.42	1.38

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	961	LOR	C22-C21-N2	4.70	120.44	114.68
2	B	961	LOR	C4-C3-C2	-4.67	107.73	112.61
2	B	961	LOR	C22-C21-N2	3.61	119.10	114.68
2	B	961	LOR	C4-C3-N8	3.52	115.81	108.52
2	A	961	LOR	C4-C3-C2	-3.48	108.98	112.61
2	B	961	LOR	O81-C8-C7	-3.30	114.45	120.20
2	A	961	LOR	C4-C3-N8	3.18	115.10	108.52
2	A	961	LOR	O82-C8-C7	2.95	121.42	116.73
2	A	961	LOR	O81-C8-C7	-2.90	115.15	120.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	961	LOR	O82-C8-C7	2.33	120.42	116.73
2	A	961	LOR	C28-C23-C24	2.12	120.94	118.30

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	961	LOR	C1-C2-C3-C4
2	A	961	LOR	N2-C2-C3-C4
2	A	961	LOR	N8-C7-C8-O82
2	B	961	LOR	N2-C2-C3-C4
2	B	961	LOR	N8-C7-C8-O82
2	A	961	LOR	C3-C2-N2-C21
2	B	961	LOR	N2-C21-C22-N22
2	B	961	LOR	O21-C21-C22-N22
2	A	961	LOR	O1-C1-C2-C3
2	B	961	LOR	O1-C1-C2-C3
2	A	961	LOR	N8-C7-C8-O81
2	B	961	LOR	N8-C7-C8-O81

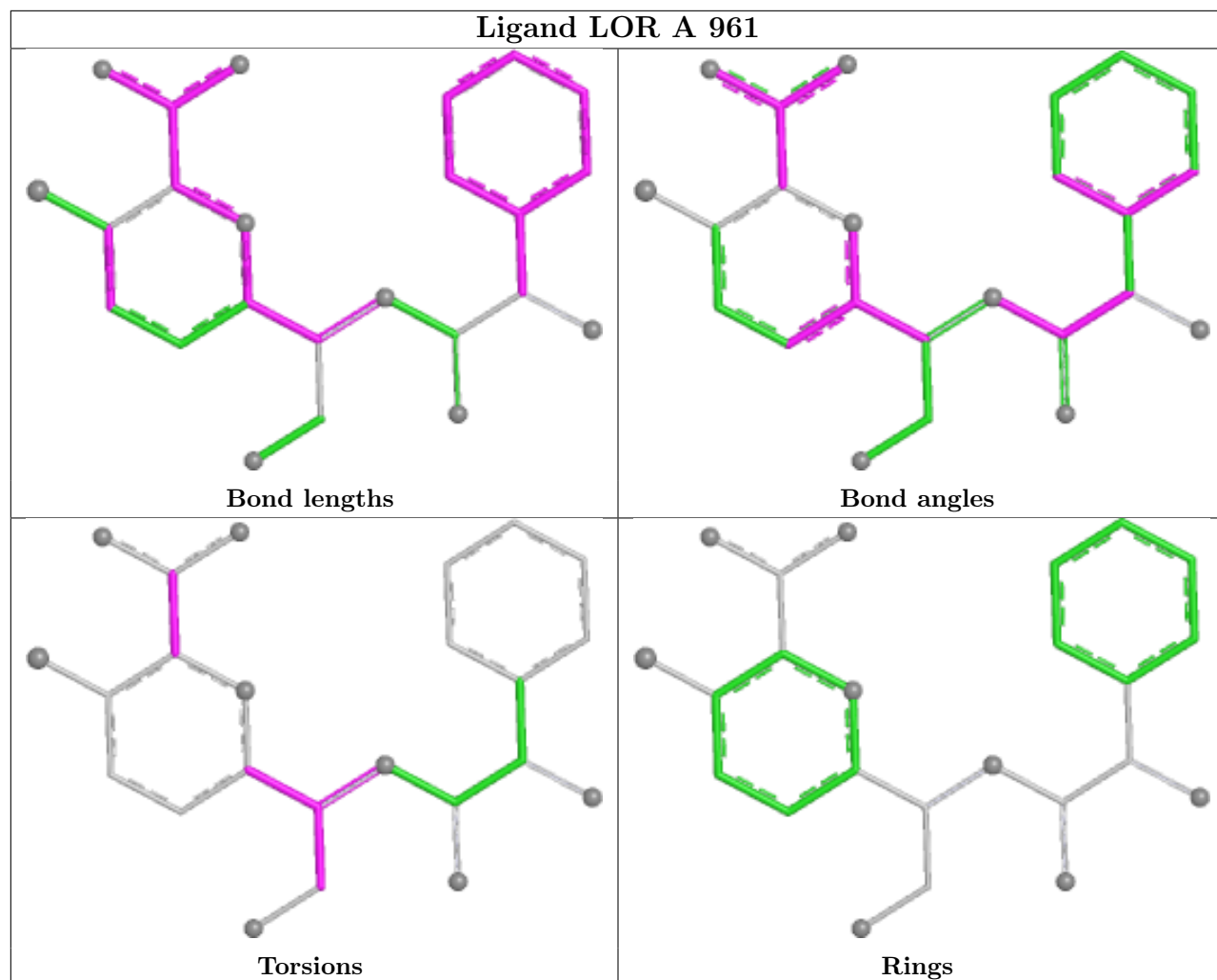
There are no ring outliers.

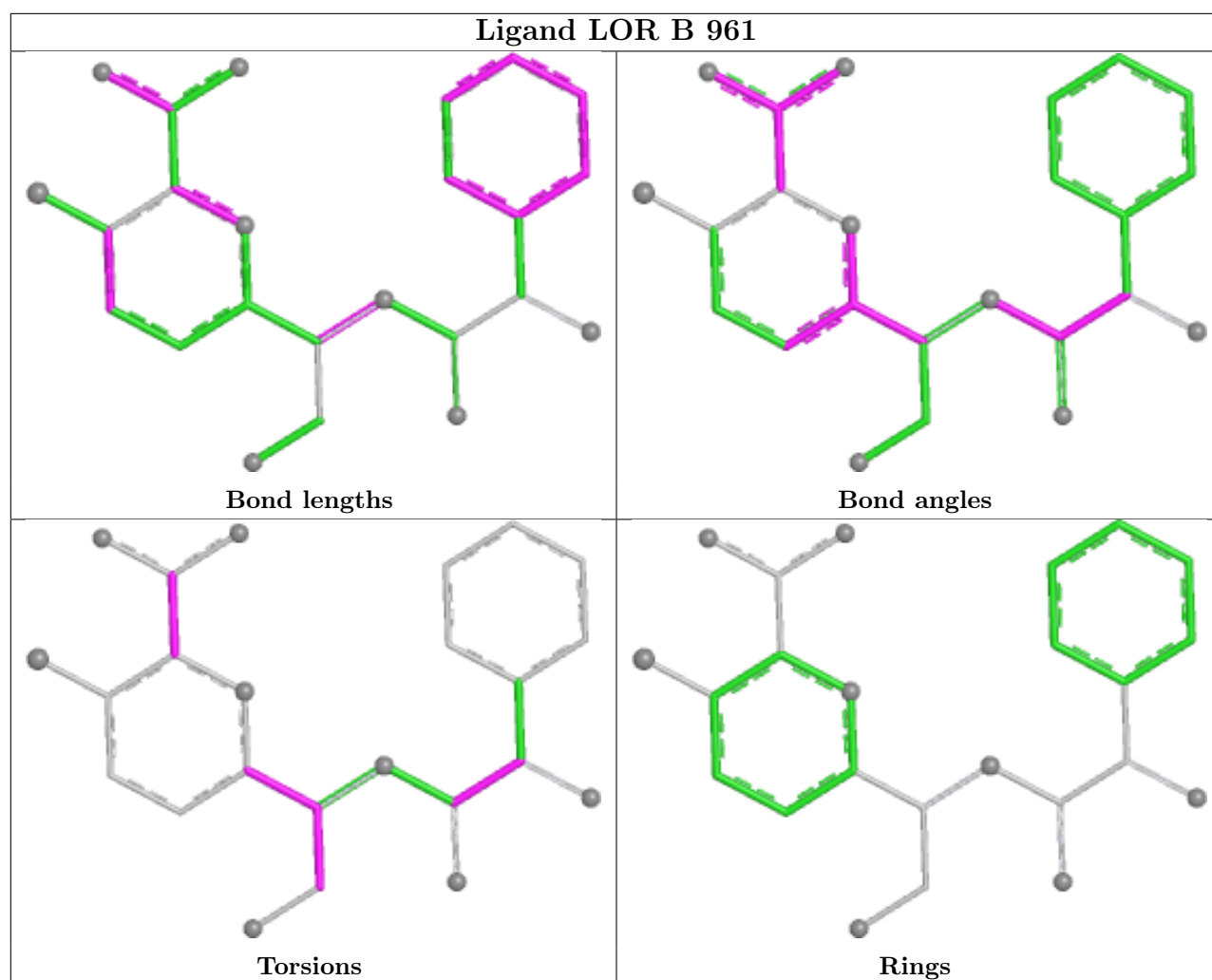
2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	961	LOR	7	0
2	B	961	LOR	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.

Ligand LOR A 961





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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.