



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:52 PM UTC

PDB ID : 8VXL / pdb_00008vxl
Title : Crystal Structure of the External Aldimine Complex of Pyridoxal-5'-Tetrazole and (S)-4-amino-5-phenoxy-pentanoate with the Bacillus subtilis GabR C-terminal Effector-binding and Oligomerization Domain
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Deposited on : 2024-02-05
Resolution : 2.24 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

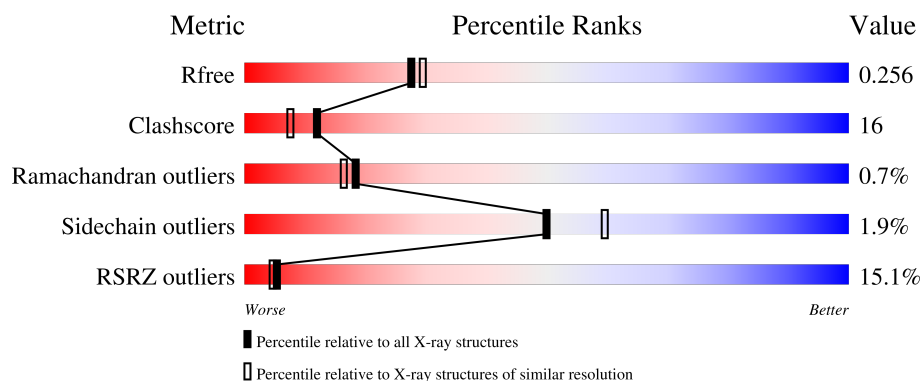
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.24 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	3416 (2.26-2.22)
Clashscore	190562	3556 (2.26-2.22)
Ramachandran outliers	187476	3500 (2.26-2.22)
Sidechain outliers	187428	3501 (2.26-2.22)
RSRZ outliers	180081	3415 (2.26-2.22)

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	383	13% 70% 22% 5%
1	B	383	16% 71% 23% 5%

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	GOL	A	501	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6089 atoms, of which 60 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 HTH-type transcriptional regulatory protein GabR.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	363	Total	C	H	N	O	S	0	4	0
			2960	1851	46	502	545	16			
1	B	364	Total	C	H	N	O	S	0	2	0
			2898	1833	14	494	539	18			

There are 14 discrepancies between the modelled and reference sequences:

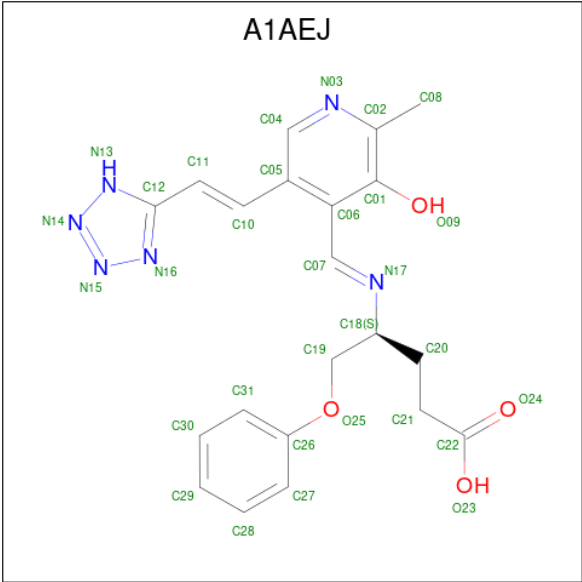
Chain	Residue	Modelled	Actual	Comment	Reference
A	103	MET	-	initiating methionine	UNP P94426
A	480	HIS	-	expression tag	UNP P94426
A	481	HIS	-	expression tag	UNP P94426
A	482	HIS	-	expression tag	UNP P94426
A	483	HIS	-	expression tag	UNP P94426
A	484	HIS	-	expression tag	UNP P94426
A	485	HIS	-	expression tag	UNP P94426
B	103	MET	-	initiating methionine	UNP P94426
B	480	HIS	-	expression tag	UNP P94426
B	481	HIS	-	expression tag	UNP P94426
B	482	HIS	-	expression tag	UNP P94426
B	483	HIS	-	expression tag	UNP P94426
B	484	HIS	-	expression tag	UNP P94426
B	485	HIS	-	expression tag	UNP P94426

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



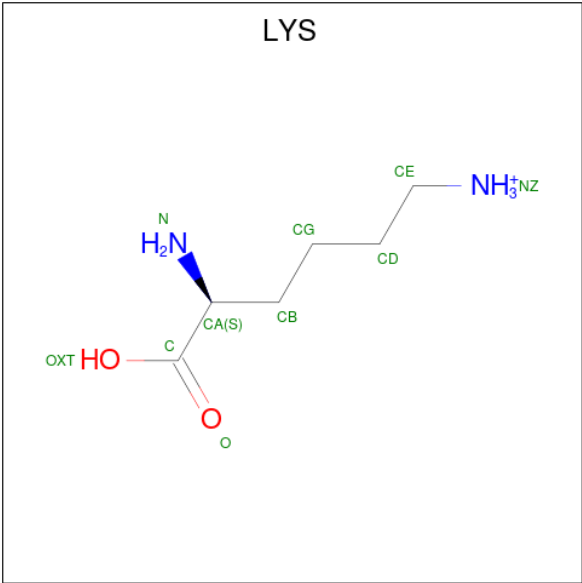
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is (4S)-4-[(E)-({3-hydroxy-2-methyl-5-[(E)-2-(1H-tetrazol-5-yl)ethenyl]pyridin-4-yl}methylidene)amino]-5-phenoxy-pentanoic acid (CCD ID: A1AEJ) (formula: C₂₁H₂₂N₆O₄) (labeled as "Ligand of Interest" by depositor).



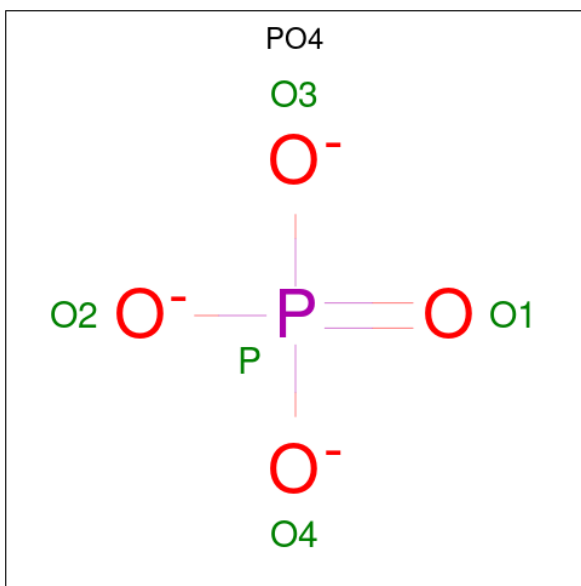
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			31	21	6	4		
3	B	1	Total	C	N	O	0	0
			31	21	6	4		

- Molecule 4 is LYSINE (CCD ID: LYS) (formula: C₆H₁₅N₂O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	N	O	0	0
			5	3	1	1		
4	B	1	Total	C	N	O	0	0
			9	6	2	1		

- Molecule 5 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	O	P	0	0
			5	4	1		

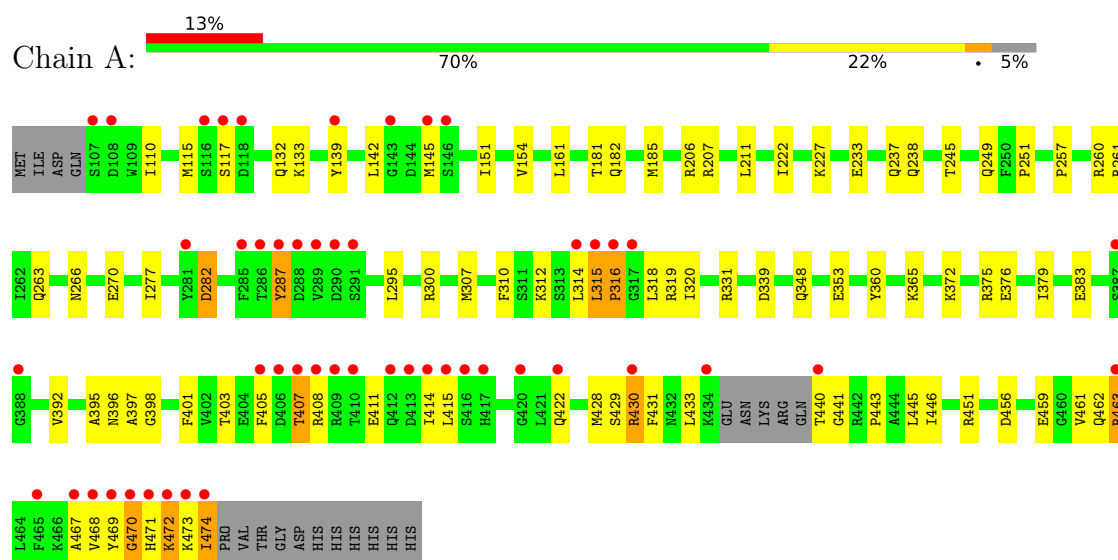
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	57	Total	O	0	0
			57	57		
6	B	63	Total	O	0	0
			63	63		

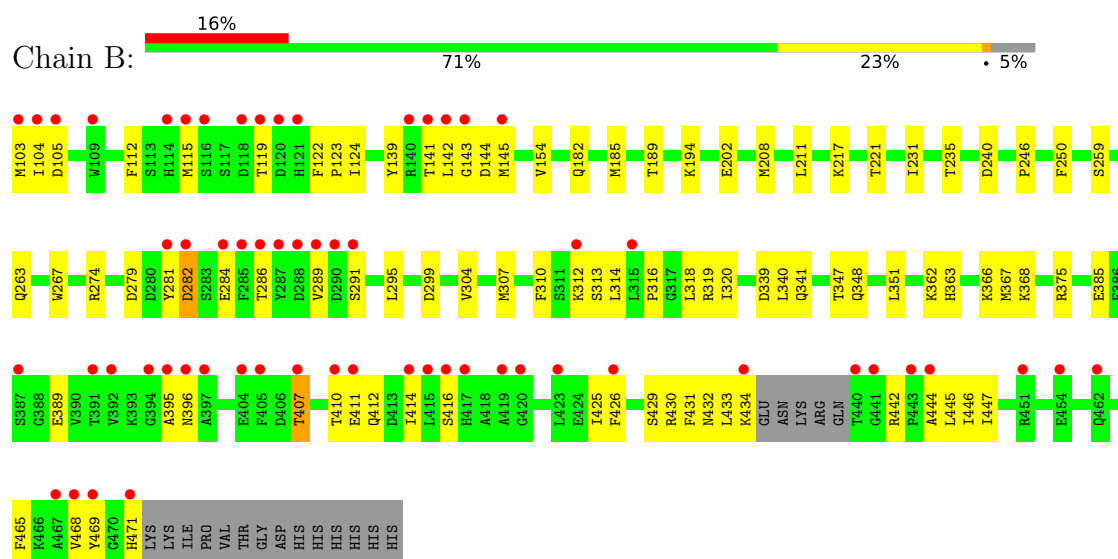
3 Residue-property plots

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: HTH-type transcriptional regulatory protein GabR



- Molecule 1: HTH-type transcriptional regulatory protein GabR



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	118.32Å 118.32Å 75.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.44 – 2.24 39.44 – 2.24	Depositor EDS
% Data completeness (in resolution range)	99.7 (39.44-2.24) 99.7 (39.44-2.24)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.29 (at 2.24Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.213 , 0.256 0.213 , 0.256	Depositor DCC
R_{free} test set	2537 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	37.9	Xtriage
Anisotropy	0.244	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 46.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.040 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6089	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.39% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, PO4, A1AEJ

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	0.40	0/2972	0.58	0/4013
1	B	0.40	0/2947	0.56	0/3980
All	All	0.40	0/5919	0.57	0/7993

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.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	430	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	2914	46	2895	104	1
1	B	2884	14	2849	90	1
2	A	12	0	16	6	0
2	B	18	0	24	4	0
3	A	31	0	0	3	0
3	B	31	0	0	2	0
4	B	14	0	14	5	0
5	B	5	0	0	0	0
6	A	57	0	0	2	0
6	B	63	0	0	3	0
All	All	6029	60	5798	184	1

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 (184) 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:341:GLN:HA	2:B:505:GOL:H31	1.24	1.16
1:B:367:MET:HA	1:B:367:MET:HE2	1.35	1.05
1:A:431:PHE:CE2	1:A:446:ILE:HD11	1.94	1.01
1:A:318:LEU:HD21	1:A:320:ILE:HD11	1.50	0.92
1:A:182:GLN:NE2	6:A:601:HOH:O	2.06	0.89
1:B:320:ILE:HD11	1:B:347:THR:HG22	1.58	0.85
1:A:315[A]:LEU:HD23	1:A:318:LEU:HB2	1.59	0.84
1:A:472:LYS:HD2	1:A:474:ILE:HG12	1.61	0.82
1:A:266:ASN:O	1:A:270:GLU:HG3	1.80	0.81
1:B:410:THR:HG22	1:B:412:GLN:H	1.45	0.81
1:A:316[A]:PRO:HB3	1:B:143:GLY:HA2	1.64	0.79
1:B:375:ARG:HE	1:B:395:ALA:HB1	1.47	0.79
1:A:227:LYS:HD3	2:A:501:GOL:H12	1.67	0.77
1:A:318:LEU:HD12	1:A:319:ARG:H	1.49	0.77
1:B:367:MET:HA	1:B:367:MET:CE	2.15	0.76
1:A:315[A]:LEU:HD12	1:A:316[A]:PRO:HD2	1.69	0.75
1:A:462:GLN:O	1:A:463:ARG:HB2	1.87	0.72
1:A:318:LEU:HD12	1:A:319:ARG:N	2.04	0.72
1:A:431:PHE:CE2	1:A:446:ILE:CD1	2.71	0.72
1:A:431:PHE:HE2	1:A:446:ILE:HD11	1.50	0.71
1:A:472:LYS:CD	1:A:474:ILE:HG12	2.21	0.71
1:A:287:TYR:HB3	1:A:372:LYS:HD3	1.73	0.71
1:A:227:LYS:CD	2:A:501:GOL:H12	2.21	0.71
1:A:115:MET:HE2	1:B:145:MET:SD	2.31	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:451:ARG:O	1:A:451:ARG:HG3	1.91	0.69
1:A:263:GLN:OE1	2:A:501:GOL:H11	1.92	0.69
1:A:405:PHE:O	1:A:443:PRO:HD2	1.91	0.69
1:B:425:ILE:HD11	1:B:447:ILE:CD1	2.23	0.68
1:A:451:ARG:NH2	1:B:144:ASP:OD1	2.28	0.66
1:B:284:GLU:H	1:B:284:GLU:CD	2.04	0.66
1:B:231:ILE:O	1:B:235:THR:HG23	1.95	0.66
1:B:246:PRO:HD2	6:B:634:HOH:O	1.95	0.66
1:B:320:ILE:HD11	1:B:347:THR:CG2	2.27	0.64
1:A:472:LYS:HD3	1:A:472:LYS:C	2.24	0.63
1:A:469:TYR:HD2	1:A:470:GLY:H	1.44	0.62
4:B:501:LYS:O	4:B:502:LYS:HB2	1.98	0.62
1:A:316[A]:PRO:HB3	1:B:143:GLY:CA	2.29	0.62
1:B:154:VAL:HG21	1:B:348:GLN:HB3	1.83	0.61
1:B:471:HIS:C	4:B:501:LYS:HA	2.26	0.61
1:A:206:ARG:NH2	1:A:433:LEU:O	2.34	0.60
1:A:468:VAL:HG23	1:A:469:TYR:CD1	2.36	0.60
1:A:316[A]:PRO:HB2	1:B:142:LEU:HB3	1.82	0.59
1:A:468:VAL:HG23	1:A:469:TYR:CE1	2.37	0.59
1:A:133:LYS:HE3	1:A:353:GLU:CD	2.28	0.59
1:B:411:GLU:OE2	1:B:429:SER:OG	2.17	0.59
1:A:185:MET:HE1	1:A:277:ILE:HG21	1.85	0.58
1:B:295:LEU:C	1:B:295:LEU:HD23	2.27	0.58
1:A:181:THR:HG23	1:A:307:MET:HE1	1.86	0.58
1:B:240:ASP:O	1:B:274:ARG:HA	2.04	0.57
1:B:286:THR:H	1:B:368:LYS:NZ	2.01	0.57
1:B:341:GLN:CA	2:B:505:GOL:H31	2.16	0.57
1:A:372:LYS:HE2	1:A:376:GLU:OE2	2.05	0.56
1:A:375:ARG:NE	1:A:395:ALA:O	2.39	0.55
1:A:473:LYS:O	1:A:474:ILE:HG23	2.06	0.55
1:A:440:THR:HG23	1:A:441:GLY:N	2.22	0.55
1:B:122:PHE:CD1	1:B:123:PRO:HD2	2.42	0.55
1:B:185:MET:HE1	1:B:307:MET:HE3	1.89	0.55
1:A:472:LYS:HD2	1:A:474:ILE:CG1	2.34	0.54
1:A:315[B]:LEU:HD22	1:B:142:LEU:HB3	1.90	0.54
1:A:462:GLN:O	1:A:463:ARG:CB	2.56	0.54
1:A:407:THR:HG21	1:A:468:VAL:O	2.08	0.54
1:A:372:LYS:HD2	1:A:372:LYS:O	2.08	0.54
1:A:407:THR:CG2	1:A:408:ARG:N	2.71	0.53
1:B:202:GLU:OE2	1:B:432:ASN:HA	2.08	0.53
1:B:341:GLN:HA	2:B:505:GOL:C3	2.17	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:251:PRO:HG2	1:A:428:MET:HE1	1.90	0.53
1:A:318:LEU:HD21	1:A:320:ILE:CD1	2.33	0.53
1:A:110:ILE:HD13	1:A:456:ASP:HB3	1.91	0.53
1:B:286:THR:H	1:B:368:LYS:CE	2.21	0.52
1:A:314[A]:LEU:HD21	1:A:360:TYR:CD1	2.44	0.52
1:A:314[A]:LEU:HD21	1:A:360:TYR:HD1	1.74	0.52
1:B:112:PHE:CB	1:B:425:ILE:HD12	2.41	0.51
1:A:139:TYR:CE1	1:B:124:ILE:HD13	2.45	0.51
1:B:185:MET:HE3	1:B:208:MET:CE	2.40	0.51
1:B:286:THR:O	1:B:368:LYS:HE2	2.11	0.51
1:B:425:ILE:HD11	1:B:447:ILE:HD12	1.91	0.51
1:B:115:MET:HE2	1:B:430:ARG:NH1	2.26	0.51
1:B:185:MET:HE3	1:B:208:MET:HE2	1.93	0.51
1:B:385:GLU:HG3	1:B:465:PHE:CD2	2.47	0.50
1:B:139:TYR:C	1:B:141:THR:H	2.20	0.50
1:B:318:LEU:HD12	1:B:319:ARG:N	2.27	0.49
1:B:444:ALA:O	1:B:445:LEU:HD23	2.12	0.49
1:B:104:ILE:O	1:B:105:ASP:C	2.55	0.49
1:B:433:LEU:O	1:B:434:LYS:CB	2.60	0.49
1:A:145:MET:HB3	1:A:151:ILE:HG12	1.95	0.49
1:A:440:THR:CG2	1:A:441:GLY:N	2.75	0.48
1:B:314:LEU:HD11	1:B:351:LEU:HD13	1.93	0.48
1:B:295:LEU:HD23	1:B:295:LEU:O	2.14	0.48
1:A:372:LYS:HD2	1:A:372:LYS:C	2.38	0.48
1:A:379:ILE:O	1:A:383:GLU:HG3	2.13	0.48
1:A:238:GLN:HA	1:A:238:GLN:OE1	2.13	0.48
1:A:430:ARG:NH1	3:A:502:A1AEJ:O23	2.47	0.48
1:A:407:THR:HG23	1:A:408:ARG:H	1.78	0.48
1:B:250:PHE:HB2	1:B:281:TYR:CZ	2.49	0.48
1:B:221:THR:HG22	1:B:433:LEU:HD12	1.94	0.48
1:B:318:LEU:HD12	1:B:319:ARG:H	1.78	0.48
1:A:472:LYS:HD3	1:A:472:LYS:O	2.14	0.48
1:A:316[A]:PRO:HB2	1:B:142:LEU:CB	2.43	0.47
1:B:279:ASP:OD2	3:B:503:A1AEJ:N03	2.48	0.47
1:B:431:PHE:CE2	1:B:446:ILE:HG13	2.50	0.47
1:B:414:ILE:HD13	1:B:468:VAL:HG22	1.96	0.47
1:A:139:TYR:HE1	1:B:124:ILE:HD13	1.80	0.47
1:B:284:GLU:CD	1:B:284:GLU:N	2.73	0.47
1:B:410:THR:O	1:B:414:ILE:HG13	2.14	0.47
1:A:261:ARG:HB3	1:A:295:LEU:HD12	1.96	0.47
1:A:227:LYS:HG2	1:A:257:PRO:HG3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:103:MET:SD	1:B:104:ILE:N	2.87	0.47
1:B:471:HIS:C	4:B:501:LYS:CA	2.88	0.46
1:B:259:SER:O	1:B:263:GLN:HG3	2.16	0.46
1:A:154:VAL:HG21	1:A:348:GLN:HB3	1.97	0.46
1:B:267:TRP:CH2	1:B:274:ARG:HD3	2.51	0.46
1:B:444:ALA:C	1:B:445:LEU:HD23	2.41	0.46
1:B:103:MET:SD	1:B:105:ASP:N	2.89	0.46
1:A:414:ILE:HG23	1:A:467:ALA:O	2.15	0.46
1:A:287:TYR:HB3	1:A:372:LYS:CD	2.43	0.46
1:A:375:ARG:HE	1:A:395:ALA:HB1	1.81	0.45
1:A:431:PHE:HE2	1:A:446:ILE:CD1	2.18	0.45
1:B:362:LYS:HG2	1:B:366:LYS:HZ3	1.80	0.45
1:B:442:ARG:HB2	6:B:620:HOH:O	2.15	0.45
1:A:233:GLU:HG3	1:A:237:GLN:NE2	2.31	0.45
1:A:339:ASP:HA	1:B:182:GLN:HE22	1.82	0.45
1:A:459:GLU:O	1:A:462:GLN:O	2.35	0.45
1:A:310:PHE:HB3	1:A:314[A]:LEU:HD12	1.98	0.45
1:A:139:TYR:CD1	1:A:142:LEU:HD12	2.52	0.45
4:B:501:LYS:O	4:B:502:LYS:CB	2.65	0.45
1:B:286:THR:H	1:B:368:LYS:HE3	1.82	0.45
1:A:315[A]:LEU:HD11	1:B:142:LEU:HD23	1.98	0.44
1:A:375:ARG:CZ	1:A:395:ALA:O	2.65	0.44
1:B:362:LYS:HG2	1:B:366:LYS:NZ	2.31	0.44
1:B:112:PHE:HB2	1:B:425:ILE:HD12	2.00	0.44
1:B:281:TYR:CD2	1:B:282[B]:ASP:HB2	2.52	0.44
1:B:407:THR:HG21	1:B:414:ILE:HD11	1.99	0.44
1:B:471:HIS:C	4:B:501:LYS:N	2.76	0.44
1:A:411:GLU:OE1	1:A:429:SER:HB3	2.17	0.44
1:A:260:ARG:HH21	2:A:501:GOL:C3	2.30	0.44
1:B:312:LYS:HB2	6:B:638:HOH:O	2.17	0.44
1:A:207:ARG:NH1	6:A:608:HOH:O	2.51	0.43
1:A:472:LYS:O	1:A:472:LYS:CG	2.66	0.43
1:A:261:ARG:HG2	1:A:295:LEU:CD1	2.48	0.43
1:A:312:LYS:HD3	3:A:502:A1AEJ:C28	2.48	0.43
1:B:123:PRO:HD3	1:B:363:HIS:ND1	2.33	0.43
1:B:339:ASP:HB2	2:B:504:GOL:C1	2.48	0.43
1:B:119:THR:HG22	1:B:316:PRO:CG	2.48	0.43
1:A:211:LEU:C	1:A:211:LEU:HD23	2.44	0.43
1:A:392:VAL:HG22	1:A:403:THR:HG22	2.00	0.43
1:A:117:SER:OG	1:A:316[B]:PRO:HD2	2.19	0.43
1:A:207:ARG:NH2	1:A:430:ARG:HH22	2.16	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:222:ILE:HG23	1:A:233:GLU:HB3	2.01	0.43
1:A:245:THR:OG1	1:A:249:GLN:HA	2.19	0.43
1:A:461:VAL:HG23	1:A:462:GLN:N	2.33	0.43
1:B:267:TRP:CZ2	1:B:274:ARG:HD3	2.54	0.43
1:B:194:LYS:O	1:B:217:LYS:NZ	2.50	0.42
1:A:139:TYR:HD1	1:A:142:LEU:HD12	1.83	0.42
1:A:365:LYS:HD3	1:A:365:LYS:HA	1.70	0.42
1:B:211:LEU:C	1:B:211:LEU:HD23	2.44	0.42
1:A:282:ASP:OD2	1:A:398:GLY:HA3	2.19	0.42
1:B:103:MET:SD	1:B:105:ASP:HA	2.60	0.42
1:B:289:VAL:CG2	1:B:396:ASN:HB3	2.49	0.42
1:A:181:THR:HG23	1:A:307:MET:CE	2.49	0.42
1:A:395:ALA:HB2	1:A:401:PHE:CB	2.50	0.42
1:B:310:PHE:HB3	1:B:314:LEU:HG	2.01	0.42
1:B:389:GLU:CB	1:B:469:TYR:OH	2.67	0.42
1:B:425:ILE:HG13	1:B:426:PHE:H	1.85	0.42
1:A:287:TYR:CD2	1:A:375:ARG:HD2	2.55	0.42
1:A:407:THR:CG2	1:A:408:ARG:H	2.32	0.41
1:A:415:LEU:HD21	1:A:445:LEU:HD22	2.02	0.41
1:A:461:VAL:CG2	1:A:462:GLN:N	2.83	0.41
1:B:281:TYR:HD2	3:B:503:A1AEJ:C28	2.32	0.41
1:A:331:ARG:NH1	1:A:331:ARG:HG2	2.36	0.41
1:A:396:ASN:O	1:A:397:ALA:HB2	2.21	0.41
1:A:115:MET:HE3	3:A:502:A1AEJ:C20	2.51	0.41
1:A:446:ILE:HD13	1:A:446:ILE:HG21	1.81	0.41
1:B:185:MET:O	1:B:189:THR:HG23	2.20	0.41
1:A:117:SER:OG	1:A:316[B]:PRO:CD	2.69	0.41
1:A:260:ARG:NH2	2:A:501:GOL:H31	2.36	0.41
1:A:422:GLN:O	1:A:463:ARG:NH2	2.49	0.41
1:B:139:TYR:C	1:B:141:THR:N	2.79	0.41
1:A:260:ARG:HH21	2:A:501:GOL:H31	1.86	0.41
1:A:206:ARG:HH21	1:A:433:LEU:C	2.28	0.40
1:B:299:ASP:CG	1:B:304:VAL:HG23	2.46	0.40
1:B:425:ILE:HG13	1:B:426:PHE:N	2.37	0.40
1:A:182:GLN:NE2	1:B:340:LEU:H	2.19	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:GLN:OE1	1:B:103:MET:N[2_454]	2.18	0.02

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	363/383 (95%)	342 (94%)	15 (4%)	6 (2%)	7	3
1	B	362/383 (94%)	342 (94%)	20 (6%)	0	100	100
All	All	725/766 (95%)	684 (94%)	35 (5%)	6 (1%)	18	13

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	463	ARG
1	A	472	LYS
1	A	471	HIS
1	A	315[A]	LEU
1	A	315[B]	LEU
1	A	470	GLY

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	313/338 (93%)	305 (97%)	8 (3%)	40	48
1	B	308/338 (91%)	302 (98%)	6 (2%)	50	59
All	All	621/676 (92%)	607 (98%)	14 (2%)	50	53

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

Mol	Chain	Res	Type
1	A	161	LEU
1	A	282	ASP
1	A	287	TYR
1	A	300	ARG
1	A	316[A]	PRO
1	A	316[B]	PRO
1	A	407	THR
1	A	474	ILE
1	B	282[A]	ASP
1	B	282[B]	ASP
1	B	291	SER
1	B	313	SER
1	B	407	THR
1	B	416	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	149	GLN
1	A	182	GLN
1	A	237	GLN
1	A	302	GLN
1	A	417	HIS
1	A	422	GLN
1	A	462	GLN
1	B	132	GLN
1	B	149	GLN
1	B	182	GLN
1	B	186	GLN
1	B	303	ASN
1	B	335	GLN
1	B	369	GLN
1	B	396	ASN
1	B	432	ASN
1	B	462	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](#)

10 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	A1AEJ	B	503	-	32,33,33	1.75	5 (15%)	36,43,43	2.06	7 (19%)
5	PO4	B	507	-	4,4,4	0.69	0	6,6,6	0.35	0
2	GOL	A	503	-	5,5,5	0.52	0	5,5,5	1.33	0
3	A1AEJ	A	502	-	32,33,33	1.76	4 (12%)	36,43,43	1.99	5 (13%)
4	LYS	B	501	-	3,4,9	0.67	0	2,4,10	0.91	0
2	GOL	A	501	-	5,5,5	1.06	0	5,5,5	0.91	0
2	GOL	B	505	-	5,5,5	1.34	0	5,5,5	0.90	0
2	GOL	B	506	-	5,5,5	0.97	0	5,5,5	1.19	1 (20%)
2	GOL	B	504	-	5,5,5	0.95	0	5,5,5	1.16	0
4	LYS	B	502	-	7,8,9	0.51	0	3,8,10	0.30	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
3	A1AEJ	B	503	-	-	9/20/20/20	0/3/3/3
2	GOL	A	503	-	-	0/4/4/4	-
3	A1AEJ	A	502	-	-	11/20/20/20	0/3/3/3
4	LYS	B	501	-	-	1/1/2/9	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	A	501	-	-	2/4/4/4	-
2	GOL	B	505	-	-	4/4/4/4	-
2	GOL	B	506	-	-	2/4/4/4	-
2	GOL	B	504	-	-	2/4/4/4	-
4	LYS	B	502	-	-	2/6/7/9	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502	A1AEJ	C11-C12	5.01	1.52	1.45
3	B	503	A1AEJ	C11-C12	4.97	1.52	1.45
3	A	502	A1AEJ	C06-C07	4.52	1.56	1.46
3	A	502	A1AEJ	C05-C10	4.30	1.54	1.47
3	B	503	A1AEJ	C06-C07	4.21	1.55	1.46
3	B	503	A1AEJ	C05-C10	3.97	1.53	1.47
3	B	503	A1AEJ	C08-C02	2.60	1.54	1.50
3	A	502	A1AEJ	C08-C02	2.56	1.54	1.50
3	B	503	A1AEJ	C06-C05	-2.38	1.39	1.42

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	A1AEJ	C10-C11-C12	8.24	143.81	124.44
3	B	503	A1AEJ	C10-C11-C12	7.06	141.03	124.44
3	A	502	A1AEJ	C05-C10-C11	-4.96	117.18	126.91
3	B	503	A1AEJ	C06-C05-C10	-4.52	112.42	121.09
3	B	503	A1AEJ	C19-O25-C26	-4.20	108.60	117.85
3	B	503	A1AEJ	C05-C10-C11	-3.05	120.93	126.91
3	A	502	A1AEJ	C19-O25-C26	-2.78	111.72	117.85
3	B	503	A1AEJ	C06-C01-C02	-2.75	118.59	120.14
3	A	502	A1AEJ	O09-C01-C02	2.57	122.91	117.58
3	A	502	A1AEJ	C06-C05-C10	-2.57	116.17	121.09
3	B	503	A1AEJ	C06-C07-N17	2.53	127.70	122.73
3	B	503	A1AEJ	O09-C01-C02	2.49	122.74	117.58
2	B	506	GOL	C3-C2-C1	-2.10	104.09	111.80

There are no chirality outliers.

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	504	GOL	O1-C1-C2-C3
2	B	505	GOL	C1-C2-C3-O3
2	B	505	GOL	O2-C2-C3-O3
2	B	506	GOL	C1-C2-C3-O3
3	A	502	A1AEJ	N17-C18-C20-C21
3	A	502	A1AEJ	C19-C18-N17-C07
3	A	502	A1AEJ	C20-C18-N17-C07
3	B	503	A1AEJ	C10-C11-C12-N13
3	B	503	A1AEJ	C10-C11-C12-N16
3	B	503	A1AEJ	C18-C19-O25-C26
4	B	501	LYS	O-C-CA-CB
3	A	502	A1AEJ	C19-C18-C20-C21
3	A	502	A1AEJ	C27-C26-O25-C19
3	A	502	A1AEJ	C31-C26-O25-C19
3	B	503	A1AEJ	C27-C26-O25-C19
3	B	503	A1AEJ	C31-C26-O25-C19
2	B	505	GOL	O1-C1-C2-C3
2	B	505	GOL	O1-C1-C2-O2
2	B	504	GOL	O1-C1-C2-O2
3	A	502	A1AEJ	C01-C06-C07-N17
2	B	506	GOL	O2-C2-C3-O3
3	B	503	A1AEJ	N17-C18-C20-C21
3	A	502	A1AEJ	C18-C20-C21-C22
2	A	501	GOL	O2-C2-C3-O3
3	B	503	A1AEJ	C19-C18-C20-C21
4	B	502	LYS	CE-CD-CG-CB
3	A	502	A1AEJ	C20-C21-C22-O23
3	B	503	A1AEJ	C20-C21-C22-O24
3	B	503	A1AEJ	C20-C21-C22-O23
3	A	502	A1AEJ	C05-C06-C07-N17
3	A	502	A1AEJ	C20-C21-C22-O24
2	A	501	GOL	C1-C2-C3-O3
4	B	502	LYS	CG-CD-CE-NZ

There are no ring outliers.

7 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	503	A1AEJ	2	0
3	A	502	A1AEJ	3	0
4	B	501	LYS	5	0
2	A	501	GOL	6	0
2	B	505	GOL	3	0

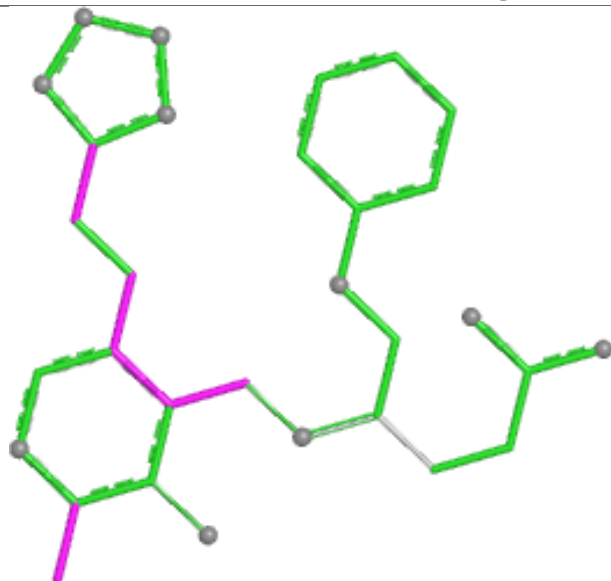
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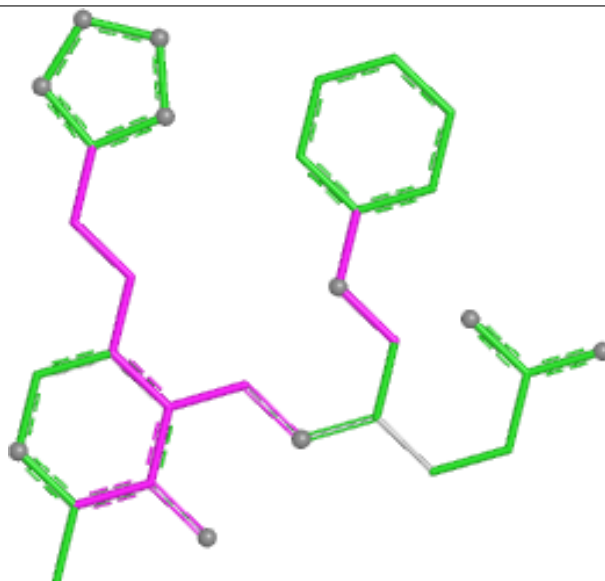
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	504	GOL	1	0
4	B	502	LYS	2	0

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.

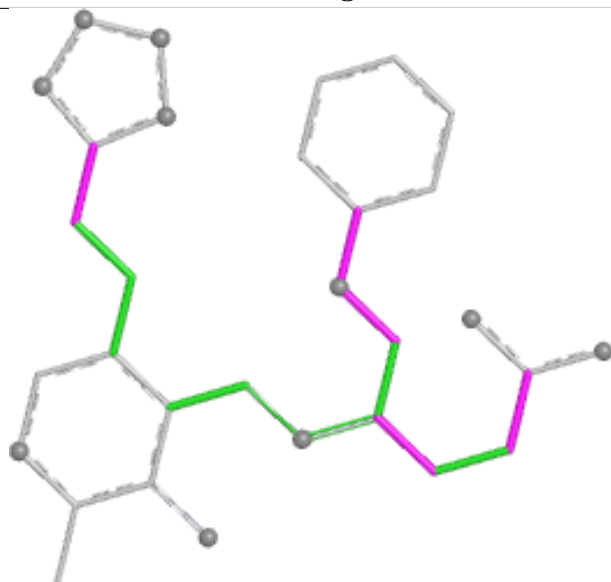
Ligand A1AEJ B 503



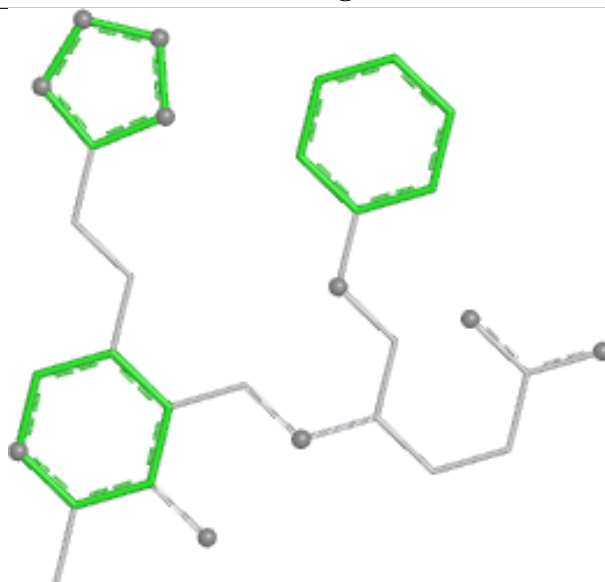
Bond lengths



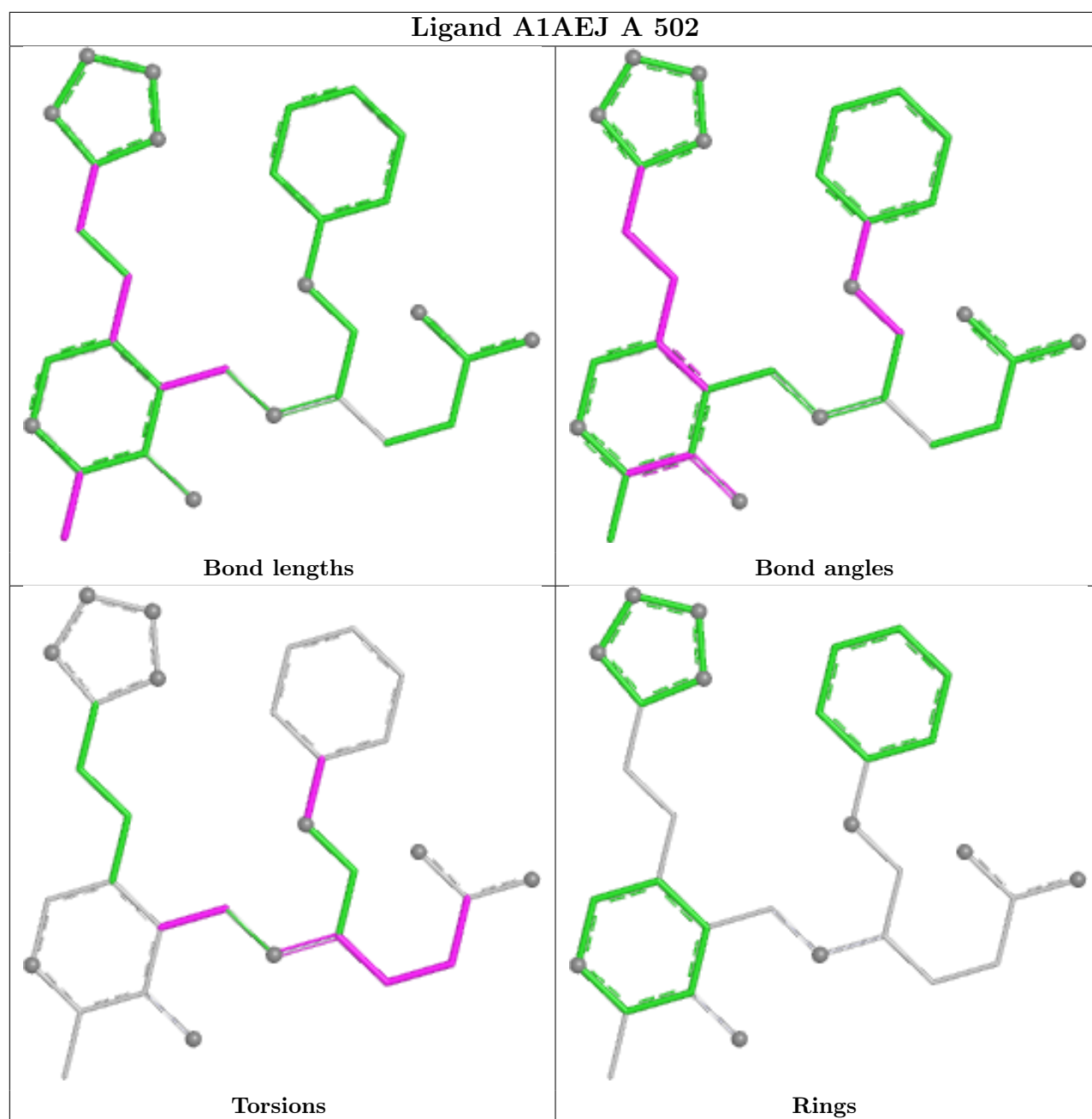
Bond angles



Torsions



Rings



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	363/383 (94%)	0.62	50 (13%) 6 6	26, 43, 80, 123	4 (1%)
1	B	364/383 (95%)	0.78	60 (16%) 4 3	25, 45, 85, 139	2 (0%)
All	All	727/766 (94%)	0.70	110 (15%) 5 5	25, 44, 82, 139	6 (0%)

All (110) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	315[A]	LEU	9.3
1	B	103	MET	7.1
1	B	289	VAL	7.0
1	A	469	TYR	6.9
1	B	290	ASP	6.6
1	B	287	TYR	5.6
1	A	474	ILE	5.6
1	A	468	VAL	5.5
1	A	145	MET	5.3
1	B	440	THR	5.1
1	B	118	ASP	5.1
1	B	281	TYR	5.0
1	A	434	LYS	4.8
1	B	396	ASN	4.7
1	A	467	ALA	4.6
1	B	395	ALA	4.4
1	A	289	VAL	4.4
1	A	316[A]	PRO	4.4
1	A	473	LYS	4.4
1	B	434	LYS	4.3
1	A	314[A]	LEU	4.2
1	B	145	MET	4.2
1	B	121	HIS	4.0
1	B	116	SER	4.0

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Mol	Chain	Res	Type	RSRZ
1	A	139	TYR	4.0
1	B	288	ASP	4.0
1	B	405	PHE	3.9
1	B	104	ILE	3.9
1	A	407	THR	3.9
1	B	120	ASP	3.8
1	A	417	HIS	3.7
1	A	416	SER	3.6
1	B	471	HIS	3.6
1	A	317[A]	GLY	3.6
1	A	440	THR	3.5
1	B	142	LEU	3.5
1	B	469	TYR	3.4
1	A	288	ASP	3.4
1	B	115	MET	3.4
1	B	286	THR	3.3
1	B	467	ALA	3.3
1	B	140	ARG	3.3
1	A	387	SER	3.3
1	A	420	GLY	3.2
1	B	410	THR	3.1
1	B	387	SER	3.1
1	B	141	THR	3.1
1	A	409	ARG	3.1
1	B	415	LEU	3.1
1	B	468	VAL	3.0
1	B	411	GLU	2.9
1	B	443	PRO	2.9
1	B	315	LEU	2.9
1	B	444	ALA	2.9
1	A	290	ASP	2.9
1	B	282[A]	ASP	2.8
1	A	388	GLY	2.8
1	A	470	GLY	2.8
1	B	397	ALA	2.8
1	B	391	THR	2.7
1	B	143	GLY	2.7
1	B	407	THR	2.7
1	A	291	SER	2.7
1	A	108	ASP	2.6
1	A	471	HIS	2.6
1	A	287	TYR	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	143	GLY	2.6
1	A	286	THR	2.6
1	B	441	GLY	2.6
1	B	114	HIS	2.6
1	B	417	HIS	2.6
1	B	284	GLU	2.5
1	A	118	ASP	2.5
1	A	413	ASP	2.5
1	A	408	ARG	2.5
1	A	414	ILE	2.5
1	B	119	THR	2.4
1	B	285	PHE	2.4
1	B	394	GLY	2.4
1	B	416	SER	2.4
1	A	412	GLN	2.4
1	A	415	LEU	2.4
1	A	410	THR	2.4
1	B	291	SER	2.4
1	A	465	PHE	2.4
1	A	472	LYS	2.3
1	B	420	GLY	2.3
1	B	419	ALA	2.3
1	A	117	SER	2.3
1	B	426	PHE	2.3
1	B	451	ARG	2.3
1	B	404	GLU	2.3
1	B	414	ILE	2.3
1	A	406	ASP	2.2
1	A	430	ARG	2.2
1	A	422	GLN	2.2
1	B	312	LYS	2.2
1	A	463	ARG	2.2
1	B	423	LEU	2.2
1	B	462	GLN	2.2
1	A	405	PHE	2.1
1	A	107	SER	2.1
1	B	109	TRP	2.1
1	A	285	PHE	2.1
1	A	116	SER	2.1
1	A	281	TYR	2.1
1	B	392	VAL	2.1
1	B	454	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	105	ASP	2.0
1	A	146	SER	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

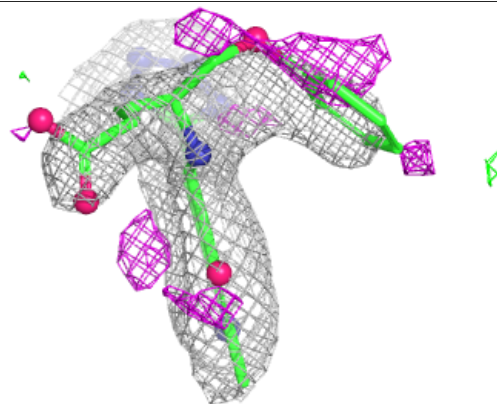
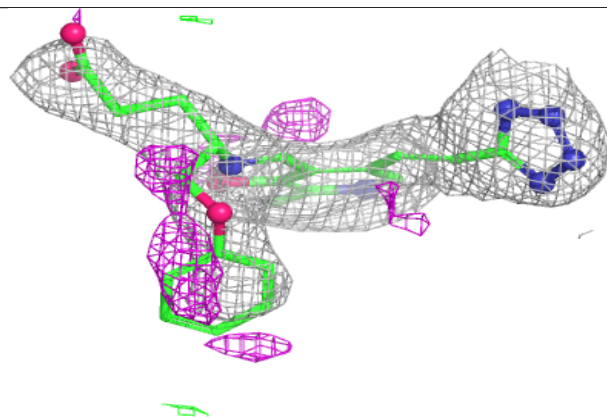
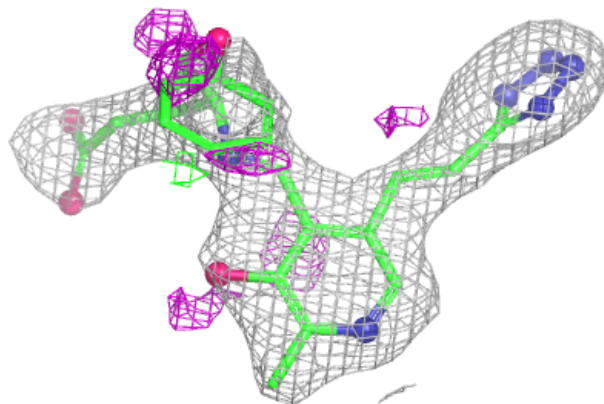
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	GOL	A	501	6/6	0.75	0.18	62,70,72,74	0
2	GOL	B	506	6/6	0.75	0.22	69,73,75,76	0
4	LYS	B	502	9/10	0.76	0.20	64,75,76,76	0
3	A1AEJ	B	503	31/31	0.85	0.18	53,62,82,85	0
4	LYS	B	501	5/10	0.85	0.18	77,81,83,84	0
3	A1AEJ	A	502	31/31	0.85	0.18	50,62,77,82	0
2	GOL	B	505	6/6	0.88	0.16	37,38,41,50	0
5	PO4	B	507	5/5	0.91	0.14	50,54,61,66	0
2	GOL	A	503	6/6	0.94	0.08	28,32,33,33	0
2	GOL	B	504	6/6	0.95	0.09	30,36,36,38	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.

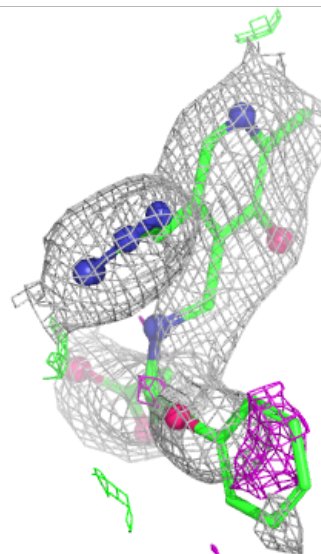
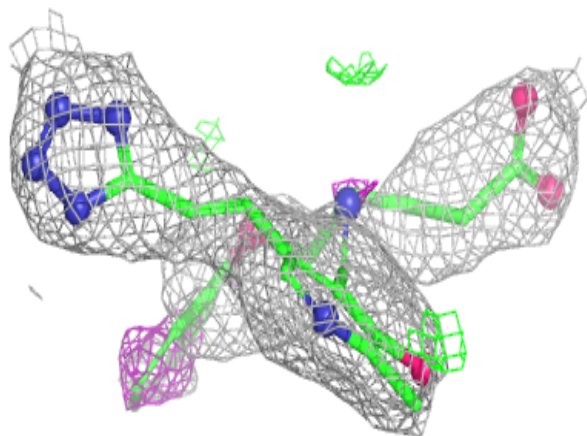
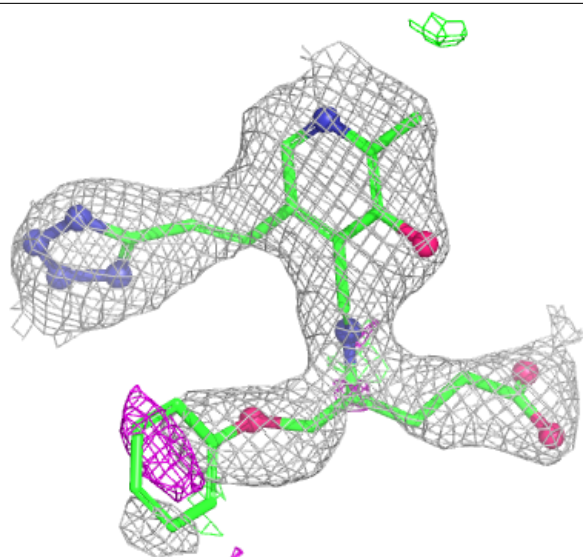
Electron density around A1AEJ B 503:

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



Electron density around A1AEJ A 502:

$2mF_o-DF_c$ (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.