



Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 04:04 am BST

PDB ID : 4B2W
Title : Pseudomonas aeruginosa RmlA in complex with allosteric inhibitor
Authors : Alphey, M.S.; Pirrie, L.; Torrie, L.; Gardiner, M.; Westwood, N.J.; Gray, D.; Naismith, J.H.
Deposited on : 2012-07-18
Resolution : 2.36 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

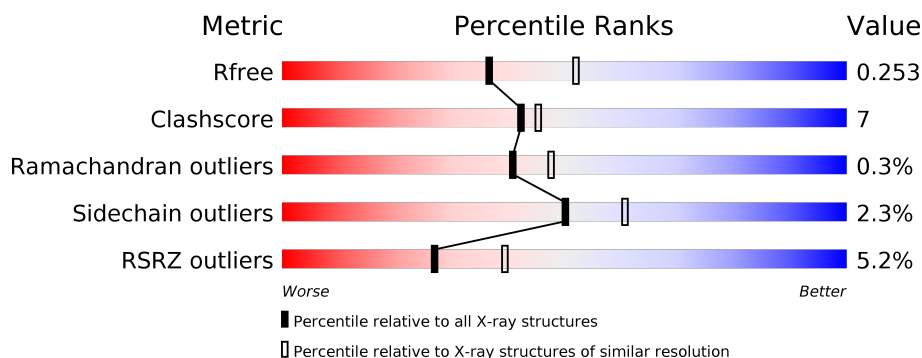
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.36 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}	130704	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (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 on 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	303	<div> <div>5%</div> <div> <div></div> <div>84%</div> <div>14%</div> <div>•</div> </div> </div>
1	B	303	<div> <div>5%</div> <div> <div></div> <div>82%</div> <div>13%</div> <div>• •</div> </div> </div>
1	C	303	<div> <div>7%</div> <div> <div></div> <div>81%</div> <div>14%</div> <div>• 5%</div> </div> </div>
1	D	303	<div> <div>4%</div> <div> <div></div> <div>81%</div> <div>14%</div> <div>• •</div> </div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9575 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 GLUCOSE-1-PHOSPHATE THYMIDYLYLTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	297	Total	C	N	O	S	0	1	0
			2339	1495	399	440	5			
1	B	290	Total	C	N	O	S	0	1	0
			2277	1458	381	433	5			
1	C	288	Total	C	N	O	S	0	1	0
			2258	1447	378	429	4			
1	D	290	Total	C	N	O	S	0	1	0
			2277	1458	381	433	5			

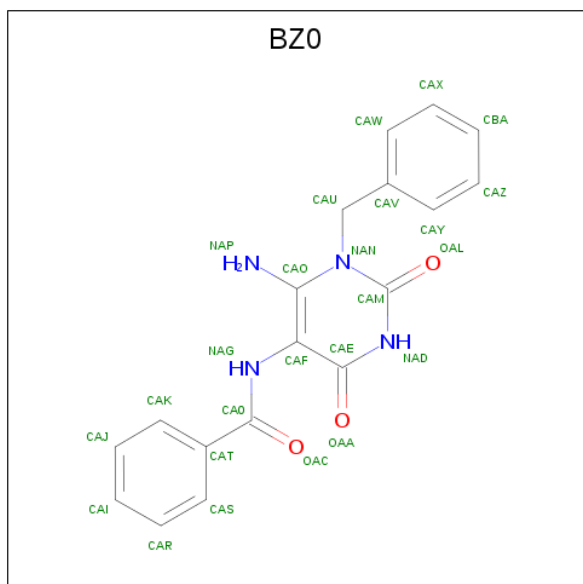
There are 39 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	HIS	-	expression tag	UNP G3XCK4
A	-8	HIS	-	expression tag	UNP G3XCK4
A	-7	HIS	-	expression tag	UNP G3XCK4
A	-6	HIS	-	expression tag	UNP G3XCK4
A	-5	HIS	-	expression tag	UNP G3XCK4
A	-4	HIS	-	expression tag	UNP G3XCK4
A	-3	GLY	-	expression tag	UNP G3XCK4
A	-2	SER	-	expression tag	UNP G3XCK4
A	-1	MET	-	expression tag	UNP G3XCK4
B	-9	HIS	-	expression tag	UNP G3XCK4
B	-8	HIS	-	expression tag	UNP G3XCK4
B	-7	HIS	-	expression tag	UNP G3XCK4
B	-6	HIS	-	expression tag	UNP G3XCK4
B	-5	HIS	-	expression tag	UNP G3XCK4
B	-4	HIS	-	expression tag	UNP G3XCK4
B	-3	GLY	-	expression tag	UNP G3XCK4
B	-2	SER	-	expression tag	UNP G3XCK4
B	-1	MET	-	expression tag	UNP G3XCK4
B	0	ALA	-	expression tag	UNP G3XCK4
C	-9	HIS	-	expression tag	UNP G3XCK4

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-8	HIS	-	expression tag	UNP G3XCK4
C	-7	HIS	-	expression tag	UNP G3XCK4
C	-6	HIS	-	expression tag	UNP G3XCK4
C	-5	HIS	-	expression tag	UNP G3XCK4
C	-4	HIS	-	expression tag	UNP G3XCK4
C	-3	GLY	-	expression tag	UNP G3XCK4
C	-2	SER	-	expression tag	UNP G3XCK4
C	-1	MET	-	expression tag	UNP G3XCK4
C	0	ALA	-	expression tag	UNP G3XCK4
D	-9	HIS	-	expression tag	UNP G3XCK4
D	-8	HIS	-	expression tag	UNP G3XCK4
D	-7	HIS	-	expression tag	UNP G3XCK4
D	-6	HIS	-	expression tag	UNP G3XCK4
D	-5	HIS	-	expression tag	UNP G3XCK4
D	-4	HIS	-	expression tag	UNP G3XCK4
D	-3	GLY	-	expression tag	UNP G3XCK4
D	-2	SER	-	expression tag	UNP G3XCK4
D	-1	MET	-	expression tag	UNP G3XCK4
D	0	ALA	-	expression tag	UNP G3XCK4

- Molecule 2 is N-(6-amino-1-benzyl-2,4-dioxo-1,2,3,4-tetrahydropyrimidin-5-yl)benzamide (three-letter code: BZ0) (formula: $C_{18}H_{16}N_4O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			25	18	4	3		

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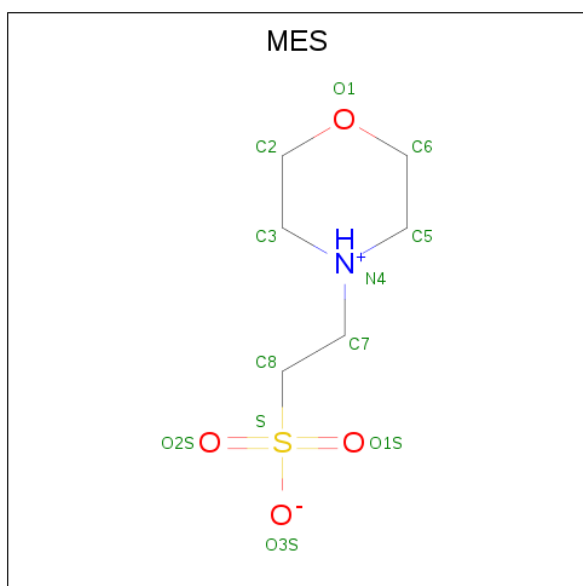
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			25	18	4	3		
2	C	1	Total	C	N	O	0	0
			25	18	4	3		
2	D	1	Total	C	N	O	0	0
			25	18	4	3		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Cl	0	0
			1	1		
3	A	1	Total	Cl	0	0
			1	1		
3	D	1	Total	Cl	0	0
			1	1		
3	C	1	Total	Cl	0	0
			1	1		

- Molecule 4 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: C₆H₁₃NO₄S).



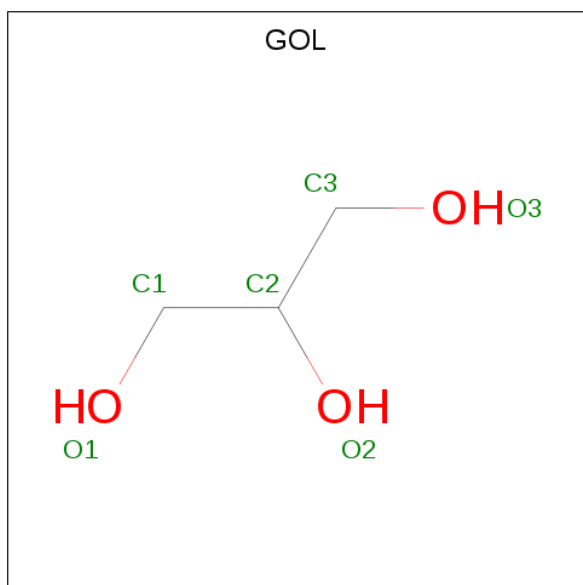
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
4	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	C	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
4	D	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	D	1	Total	C	O	0	0
			6	3	3		

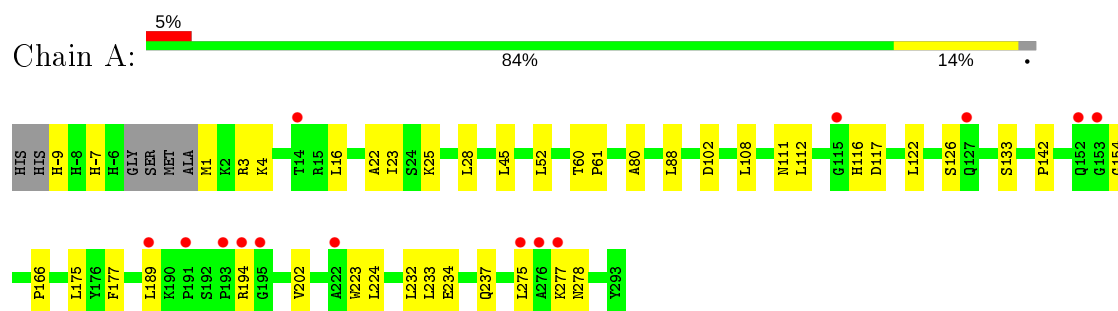
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	92	Total	O	0	0
			92	92		
6	B	66	Total	O	0	0
			66	66		
6	C	45	Total	O	0	0
			45	45		
6	D	63	Total	O	0	0
			63	63		

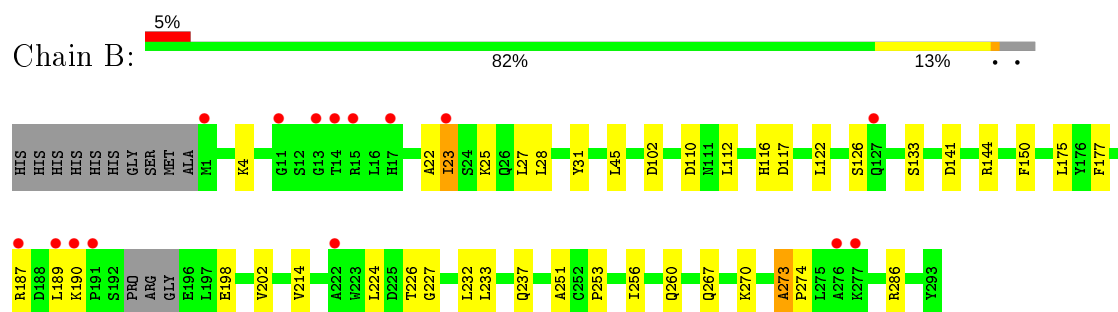
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

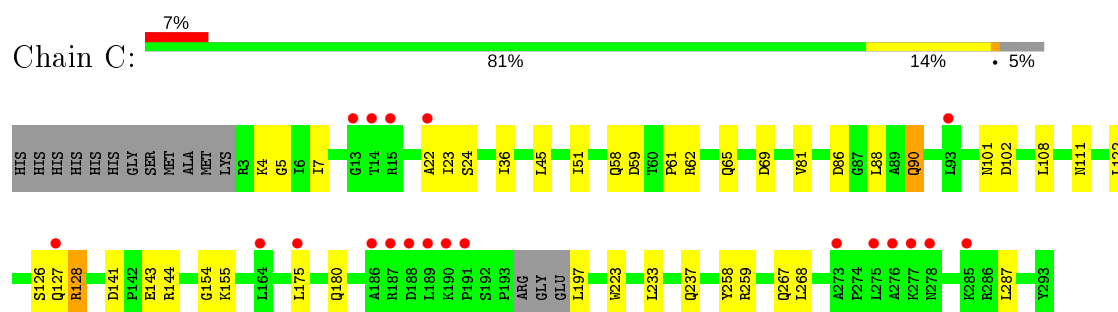
• Molecule 1: GLUCOSE-1-PHOSPHATE THYMIDYLTRANSFERASE



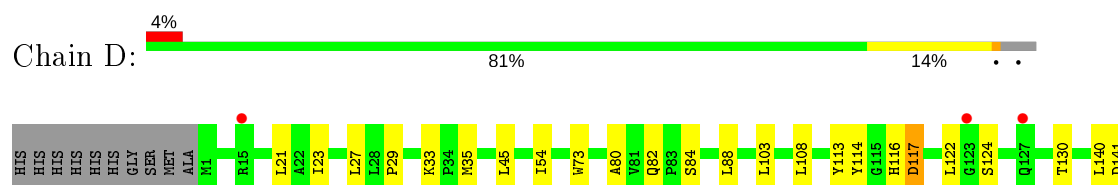
• Molecule 1: GLUCOSE-1-PHOSPHATE THYMIDYLTRANSFERASE

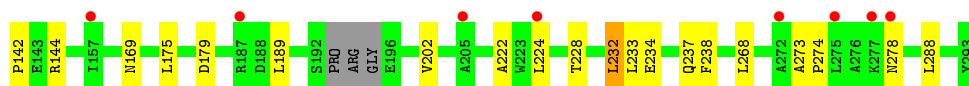


• Molecule 1: GLUCOSE-1-PHOSPHATE THYMIDYLTRANSFERASE



• Molecule 1: GLUCOSE-1-PHOSPHATE THYMIDYLTRANSFERASE





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	64.12Å 154.44Å 134.33Å 90.00° 92.46° 90.00°	Depositor
Resolution (Å)	35.02 – 2.36 35.02 – 2.35	Depositor EDS
% Data completeness (in resolution range)	99.3 (35.02-2.36) 99.3 (35.02-2.35)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.39 (at 2.36Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.197 , 0.256 0.200 , 0.253	Depositor DCC
R_{free} test set	2721 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	32.2	Xtriage
Anisotropy	0.138	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 37.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.079 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9575	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.14% 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, CL, MES, BZ0

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	0.66	0/2395	0.74	1/3247 (0.0%)
1	B	0.66	0/2328	0.75	1/3156 (0.0%)
1	C	0.60	0/2310	0.71	0/3135
1	D	0.62	1/2328 (0.0%)	0.74	0/3156
All	All	0.64	1/9361 (0.0%)	0.74	2/12694 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	73	TRP	CD2-CE2	5.11	1.47	1.41

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	286	ARG	NE-CZ-NH1	5.41	123.00	120.30
1	A	3	ARG	NE-CZ-NH2	-5.07	117.76	120.30

There are no chirality outliers.

There are no planarity outliers.

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	2339	0	2322	31	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2277	0	2268	46	0
1	C	2258	0	2244	48	0
1	D	2277	0	2268	36	0
2	A	25	0	16	0	0
2	B	25	0	16	1	0
2	C	25	0	16	1	0
2	D	25	0	16	2	0
3	A	1	0	0	1	0
3	B	1	0	0	1	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	12	0	13	1	0
4	B	12	0	13	2	0
4	C	12	0	13	2	0
4	D	12	0	13	0	0
5	D	6	0	8	0	0
6	A	92	0	0	0	0
6	B	66	0	0	1	0
6	C	45	0	0	1	0
6	D	63	0	0	0	0
All	All	9575	0	9226	128	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 7.

All (128) 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:112:LEU:HD22	1:B:224:LEU:HD12	1.22	1.14
1:A:112:LEU:HD11	1:A:224:LEU:CD1	1.80	1.12
1:A:112:LEU:CD1	1:A:224:LEU:HD12	1.79	1.10
1:B:237[A]:GLN:OE1	1:C:237[A]:GLN:OE1	1.71	1.08
1:C:128:ARG:HG3	1:C:128:ARG:HH11	0.91	1.04
1:A:23:ILE:HD13	1:D:23:ILE:HD12	1.46	0.98
1:C:128:ARG:HH11	1:C:128:ARG:CG	1.81	0.92
1:C:128:ARG:HG3	1:C:128:ARG:NH1	1.69	0.91
1:A:88:LEU:HD13	1:A:108:LEU:HD21	1.54	0.90
1:A:233:LEU:HD11	1:D:237[A]:GLN:HG3	1.57	0.85
1:B:23:ILE:HD13	1:C:23:ILE:HD13	1.62	0.82
1:B:23:ILE:CD1	1:C:23:ILE:HD13	2.10	0.81
1:C:88:LEU:HD13	1:C:108:LEU:HD21	1.62	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:154:GLY:O	4:C:1296:MES:H32	1.83	0.76
1:B:112:LEU:CD2	1:B:224:LEU:HD12	2.12	0.75
1:C:61:PRO:O	1:C:65:GLN:HG3	1.86	0.75
1:B:23:ILE:HG21	1:C:23:ILE:HG12	1.69	0.74
1:B:23:ILE:CG1	1:C:23:ILE:HD13	2.17	0.73
1:B:25:LYS:HE3	1:B:226:THR:O	1.88	0.73
1:B:237[A]:GLN:CG	1:C:237[A]:GLN:OE1	2.37	0.73
1:B:237[A]:GLN:CD	1:C:237[A]:GLN:OE1	2.29	0.71
1:B:23:ILE:HG12	1:C:23:ILE:CG2	2.20	0.70
1:B:23:ILE:HG12	1:C:23:ILE:HD13	1.75	0.69
1:B:116:HIS:ND1	3:B:1295:CL:CL	2.60	0.68
1:A:233:LEU:HD21	1:D:237[B]:GLN:HA	1.76	0.68
1:B:189:LEU:HD11	1:B:202:VAL:HG23	1.75	0.67
1:B:23:ILE:HG12	1:C:23:ILE:HG23	1.75	0.67
1:B:237[A]:GLN:OE1	1:C:237[A]:GLN:CD	2.33	0.67
1:D:45:LEU:HD22	2:D:1294:BZ0:CAY	2.26	0.66
1:B:237[A]:GLN:HG2	1:C:237[A]:GLN:OE1	1.94	0.66
1:A:233:LEU:HD21	1:D:237[A]:GLN:HA	1.78	0.66
1:C:154:GLY:O	4:C:1296:MES:C3	2.44	0.64
1:B:237[A]:GLN:OE1	1:C:237[A]:GLN:HG2	1.98	0.64
1:A:28:LEU:HD22	1:D:29:PRO:HD2	1.82	0.61
1:B:27:LEU:HD22	1:C:22:ALA:HB1	1.82	0.60
1:B:45:LEU:HD22	2:B:1294:BZ0:CAY	2.31	0.60
1:C:45:LEU:HD22	2:C:1294:BZ0:CAY	2.32	0.60
1:D:189:LEU:HD11	1:D:202:VAL:CG2	2.32	0.60
1:C:88:LEU:HD13	1:C:108:LEU:CD2	2.33	0.58
1:B:237[A]:GLN:OE1	1:C:237[A]:GLN:CG	2.52	0.58
1:C:128:ARG:NH1	1:C:128:ARG:CG	2.49	0.57
1:C:141:ASP:OD2	1:C:144:ARG:HD3	2.04	0.57
1:D:113:TYR:O	2:D:1294:BZ0:HAX	2.05	0.56
1:D:234:GLU:HA	1:D:237[A]:GLN:HE21	1.70	0.56
1:C:81:VAL:HG22	6:C:2021:HOH:O	2.05	0.56
1:D:122:LEU:CD2	1:D:175:LEU:HD21	2.36	0.55
1:A:-7:HIS:HD2	1:C:69:ASP:OD2	1.89	0.55
1:C:259:ARG:NH2	1:C:287:LEU:HD22	2.22	0.55
1:D:88:LEU:HD13	1:D:108:LEU:HD21	1.89	0.54
1:D:114:TYR:HB3	1:D:222:ALA:HB3	1.88	0.54
1:B:4:LYS:NZ	1:B:102:ASP:OD2	2.38	0.53
1:B:25:LYS:NZ	1:B:227:GLY:HA2	2.24	0.53
1:A:22:ALA:CB	1:D:27:LEU:HD22	2.40	0.52
1:B:23:ILE:HD13	1:B:28:LEU:HD23	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:PRO:HG2	1:A:166:PRO:HG2	1.91	0.51
1:A:237[B]:GLN:HA	1:D:233:LEU:HD21	1.92	0.51
1:A:111:ASN:ND2	1:A:223:TRP:HE1	2.09	0.51
1:B:25:LYS:HZ1	1:B:227:GLY:HA2	1.75	0.51
1:B:273:ALA:HB3	1:B:274:PRO:CD	2.41	0.51
1:A:189:LEU:HD11	1:A:202:VAL:CG2	2.41	0.51
1:A:112:LEU:HD11	1:A:224:LEU:HD12	0.84	0.51
1:A:116:HIS:ND1	3:A:1295:CL:CL	2.71	0.50
1:D:268:LEU:HD23	1:D:288:LEU:CD2	2.43	0.49
1:A:237[A]:GLN:HA	1:D:233:LEU:HD21	1.93	0.49
1:B:189:LEU:HD11	1:B:202:VAL:CG2	2.42	0.49
1:B:23:ILE:HG12	1:C:23:ILE:CD1	2.41	0.48
1:D:189:LEU:HD11	1:D:202:VAL:HG23	1.93	0.48
1:A:189:LEU:HD11	1:A:202:VAL:HG23	1.95	0.48
1:B:273:ALA:HB3	1:B:274:PRO:HD3	1.95	0.48
1:B:198:GLU:HG3	6:B:2046:HOH:O	2.13	0.48
1:B:233:LEU:HD21	1:C:237[B]:GLN:HA	1.95	0.48
1:A:122:LEU:CD2	1:A:175:LEU:HD21	2.44	0.48
1:B:214:VAL:O	4:B:1296:MES:H71	2.14	0.48
1:D:103:LEU:HD22	1:D:179:ASP:HA	1.96	0.48
1:A:233:LEU:HD21	1:D:237[B]:GLN:CA	2.45	0.47
1:C:258:TYR:CG	1:C:268:LEU:HD22	2.50	0.47
1:C:61:PRO:O	1:C:65:GLN:CG	2.59	0.47
1:C:4:LYS:NZ	1:C:102:ASP:OD2	2.48	0.47
1:D:54:ILE:HA	1:D:80:ALA:O	2.15	0.46
1:A:234:GLU:HA	1:A:237[A]:GLN:HE21	1.80	0.46
1:C:24:SER:HB2	1:C:59:ASP:OD1	2.15	0.46
1:D:35:MET:HE1	1:D:232:LEU:HD11	1.98	0.46
1:A:-9:HIS:NE2	1:C:267:GLN:HG2	2.31	0.46
1:C:58:GLN:OE1	1:C:58:GLN:N	2.48	0.46
1:A:154:GLY:O	4:A:1296:MES:H32	2.16	0.46
1:B:133:SER:HB2	1:B:177:PHE:HB2	1.98	0.46
1:D:21:LEU:HD23	1:D:21:LEU:O	2.16	0.46
1:B:122:LEU:CD2	1:B:175:LEU:HD21	2.46	0.45
1:A:4:LYS:NZ	1:A:102:ASP:OD2	2.42	0.45
1:B:256:ILE:O	1:B:260:GLN:HG2	2.17	0.45
1:D:116:HIS:CD2	1:D:117:ASP:HB2	2.51	0.45
1:D:141:ASP:OD2	1:D:144:ARG:NH1	2.33	0.45
1:B:233:LEU:HD21	1:C:237[A]:GLN:HA	1.98	0.44
1:A:133:SER:HB2	1:A:177:PHE:HB2	1.98	0.44
1:B:141:ASP:OD2	1:B:144:ARG:HD3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:275:LEU:HD22	1:A:278:ASN:HD21	1.83	0.44
1:A:52:LEU:HD11	1:A:80:ALA:HB2	2.00	0.44
1:B:23:ILE:HG12	1:C:23:ILE:CG1	2.48	0.44
1:D:268:LEU:HD23	1:D:288:LEU:HD23	1.99	0.43
1:A:16:LEU:CD1	1:A:25:LYS:HD3	2.48	0.43
1:D:228:THR:O	1:D:232:LEU:CD2	2.67	0.43
1:D:228:THR:O	1:D:232:LEU:HD22	2.18	0.43
1:D:82:GLN:HG2	1:D:84:SER:O	2.19	0.43
1:B:226:THR:HB	1:B:232:LEU:HD13	2.01	0.43
1:B:22:ALA:C	1:B:23:ILE:HG13	2.40	0.42
1:B:237[B]:GLN:HA	1:C:233:LEU:HD21	2.00	0.42
1:C:7:ILE:HG21	1:C:36:ILE:HD11	2.00	0.42
1:C:101:ASN:O	1:C:180:GLN:HG2	2.19	0.42
1:D:29:PRO:HA	1:D:33:LYS:O	2.19	0.42
1:A:233:LEU:HD11	1:D:237[A]:GLN:CG	2.40	0.42
1:B:267:GLN:OE1	1:B:270:LYS:HE2	2.20	0.42
1:C:5:GLY:O	1:C:51:ILE:HA	2.20	0.42
1:D:273:ALA:N	1:D:274:PRO:HD2	2.35	0.42
1:A:60:THR:N	1:A:61:PRO:CD	2.82	0.41
1:D:278:ASN:OD1	1:D:278:ASN:C	2.58	0.41
1:C:86:ASP:O	1:C:197:LEU:HB2	2.20	0.41
1:C:86:ASP:HB2	1:C:90:GLN:HE22	1.85	0.41
1:C:122:LEU:CD2	1:C:175:LEU:HD21	2.51	0.41
1:A:116:HIS:CD2	1:A:117:ASP:HB2	2.56	0.41
1:D:189:LEU:HD11	1:D:202:VAL:HG22	2.01	0.41
1:D:141:ASP:OD2	1:D:144:ARG:HD3	2.20	0.41
1:D:224:LEU:HD11	1:D:238:PHE:CD2	2.56	0.41
1:B:25:LYS:HE2	1:B:110:ASP:CB	2.51	0.41
1:C:111:ASN:ND2	1:C:223:TRP:HE1	2.19	0.41
1:D:142:PRO:HG3	1:D:169:ASN:HA	2.03	0.40
1:B:251:ALA:O	1:B:253:PRO:HD3	2.22	0.40
1:B:150:PHE:CE1	4:B:1296:MES:H51	2.57	0.40
1:B:237[A]:GLN:HA	1:C:233:LEU:HD21	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	294/303 (97%)	287 (98%)	7 (2%)	0	100	100
1	B	287/303 (95%)	281 (98%)	4 (1%)	2 (1%)	22	23
1	C	285/303 (94%)	277 (97%)	7 (2%)	1 (0%)	34	38
1	D	287/303 (95%)	281 (98%)	6 (2%)	0	100	100
All	All	1153/1212 (95%)	1126 (98%)	24 (2%)	3 (0%)	41	47

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	31	TYR
1	C	143	GLU
1	B	273	ALA

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	245/248 (99%)	239 (98%)	6 (2%)	49	59
1	B	239/248 (96%)	234 (98%)	5 (2%)	53	65
1	C	237/248 (96%)	231 (98%)	6 (2%)	47	58
1	D	239/248 (96%)	234 (98%)	5 (2%)	53	65
All	All	960/992 (97%)	938 (98%)	22 (2%)	50	61

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	45	LEU
1	A	126	SER
1	A	194	ARG
1	A	232	LEU
1	A	277	LYS
1	B	23	ILE
1	B	117	ASP
1	B	126	SER
1	B	187	ARG
1	B	190	LYS
1	C	62	ARG
1	C	90	GLN
1	C	126	SER
1	C	127	GLN
1	C	128	ARG
1	C	155	LYS
1	D	117	ASP
1	D	124	SER
1	D	130	THR
1	D	140	LEU
1	D	232	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	-7	HIS
1	A	-6	HIS
1	A	111	ASN
1	A	229	HIS
1	C	111	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 13 ligands modelled in this entry, 4 are monoatomic - leaving 9 for Mogul analysis.

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	BZ0	C	1294	-	25,27,27	1.92	4 (16%)	28,37,37	1.29	2 (7%)
4	MES	D	1297	-	12,12,12	2.10	1 (8%)	14,16,16	1.42	2 (14%)
2	BZ0	A	1294	-	25,27,27	1.74	4 (16%)	28,37,37	1.67	3 (10%)
4	MES	B	1296	-	12,12,12	2.19	1 (8%)	14,16,16	1.37	2 (14%)
4	MES	A	1296	-	12,12,12	1.88	1 (8%)	14,16,16	2.71	4 (28%)
4	MES	C	1296	-	12,12,12	1.95	1 (8%)	14,16,16	2.14	5 (35%)
2	BZ0	D	1294	-	25,27,27	1.74	4 (16%)	28,37,37	1.71	5 (17%)
2	BZ0	B	1294	-	25,27,27	1.80	4 (16%)	28,37,37	1.44	3 (10%)
5	GOL	D	1296	-	5,5,5	0.37	0	5,5,5	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
2	BZ0	C	1294	-	-	2/12/12/12	0/3/3/3
4	MES	D	1297	-	-	3/6/14/14	0/1/1/1
2	BZ0	A	1294	-	-	3/12/12/12	0/3/3/3
4	MES	B	1296	-	-	1/6/14/14	0/1/1/1
4	MES	A	1296	-	-	0/6/14/14	0/1/1/1
4	MES	C	1296	-	-	3/6/14/14	0/1/1/1
2	BZ0	D	1294	-	-	2/12/12/12	0/3/3/3
2	BZ0	B	1294	-	-	2/12/12/12	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	GOL	D	1296	-	-	4/4/4/4	-

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1296	MES	C8-S	-7.14	1.67	1.77
4	D	1297	MES	C8-S	-6.81	1.67	1.77
4	C	1296	MES	C8-S	-6.24	1.68	1.77
2	C	1294	BZ0	CAE-NAD	5.85	1.43	1.33
4	A	1296	MES	C8-S	-5.72	1.69	1.77
2	A	1294	BZ0	CAE-NAD	5.34	1.42	1.33
2	D	1294	BZ0	CAT-CA0	-4.94	1.39	1.50
2	B	1294	BZ0	CAT-CA0	-4.69	1.40	1.50
2	A	1294	BZ0	CAT-CA0	-4.40	1.41	1.50
2	C	1294	BZ0	CAU-CAV	-4.32	1.41	1.51
2	C	1294	BZ0	CAT-CA0	-4.29	1.41	1.50
2	B	1294	BZ0	CAU-CAV	-4.23	1.41	1.51
2	D	1294	BZ0	CAU-CAV	-4.23	1.41	1.51
2	B	1294	BZ0	CAE-NAD	3.96	1.39	1.33
2	D	1294	BZ0	CAE-NAD	3.87	1.39	1.33
2	B	1294	BZ0	CAE-CAF	-3.34	1.37	1.41
2	A	1294	BZ0	CAU-CAV	-3.24	1.43	1.51
2	C	1294	BZ0	CAM-NAD	3.13	1.44	1.38
2	A	1294	BZ0	CAE-CAF	-2.59	1.38	1.41
2	D	1294	BZ0	CAE-CAF	-2.56	1.38	1.41

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1294	BZ0	CAE-NAD-CAM	6.92	120.98	115.14
4	A	1296	MES	C2-C3-N4	6.69	120.25	110.10
2	A	1294	BZ0	CAE-NAD-CAM	6.55	120.67	115.14
4	A	1296	MES	O1S-S-C8	5.55	113.60	106.92
2	B	1294	BZ0	CAE-NAD-CAM	5.07	119.43	115.14
2	C	1294	BZ0	CAE-NAD-CAM	4.86	119.25	115.14
4	C	1296	MES	O2S-S-C8	4.42	112.24	106.92
4	C	1296	MES	O3S-S-C8	3.77	111.87	105.77
4	D	1297	MES	O3S-S-C8	3.43	111.31	105.77
2	B	1294	BZ0	CAF-NAG-CA0	-3.32	115.26	123.41
2	A	1294	BZ0	CAF-NAG-CA0	-3.32	115.27	123.41
4	B	1296	MES	O3S-S-C8	3.28	111.08	105.77
4	C	1296	MES	C6-C5-N4	2.98	114.62	110.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1296	MES	O1-C2-C3	2.79	117.93	111.80
2	D	1294	BZ0	CAU-NAN-CAM	2.70	121.08	117.92
2	D	1294	BZ0	CAF-NAG-CA0	-2.53	117.21	123.41
4	A	1296	MES	O2S-S-O1S	-2.46	105.43	113.95
2	A	1294	BZ0	CAF-CAE-NAD	-2.46	117.03	124.01
4	C	1296	MES	C2-C3-N4	2.31	113.60	110.10
4	B	1296	MES	O1S-S-C8	2.28	109.66	106.92
2	D	1294	BZ0	NAP-CAO-NAN	2.16	123.03	120.12
4	D	1297	MES	O3S-S-O2S	-2.15	106.02	111.27
4	C	1296	MES	O3S-S-O2S	-2.13	106.08	111.27
2	C	1294	BZ0	OAC-CA0-CAT	-2.10	117.20	120.94
2	D	1294	BZ0	CAF-CAE-NAD	-2.03	118.25	124.01
2	B	1294	BZ0	OAC-CA0-CAT	-2.02	117.34	120.94

There are no chirality outliers.

All (20) torsion outliers are listed below:

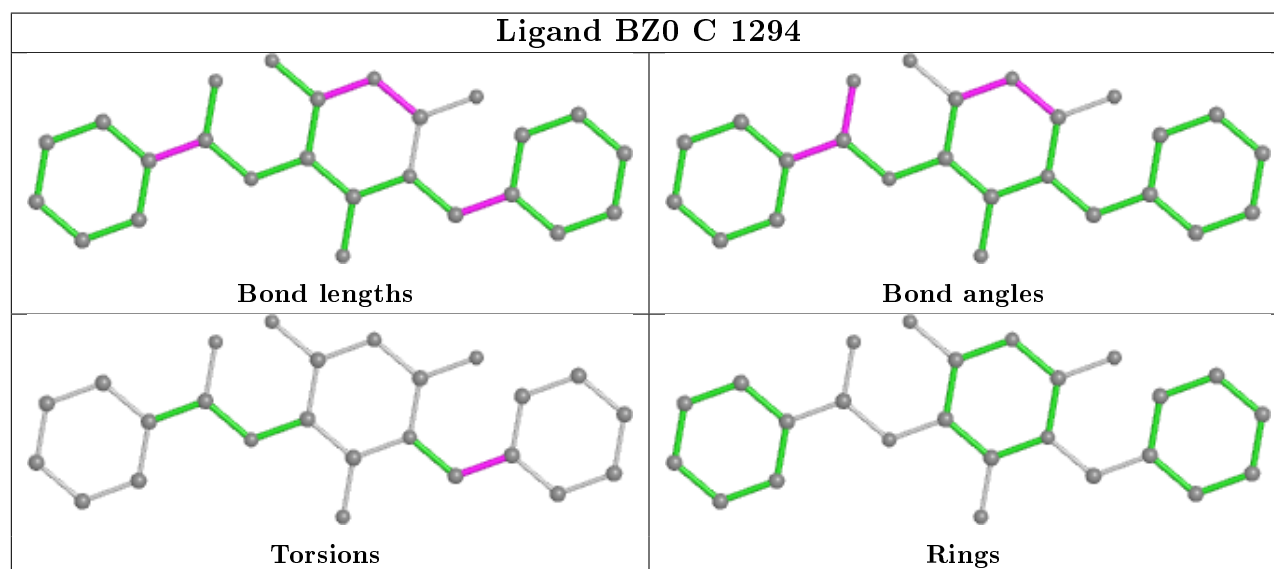
Mol	Chain	Res	Type	Atoms
4	C	1296	MES	C8-C7-N4-C3
5	D	1296	GOL	O2-C2-C3-O3
5	D	1296	GOL	O1-C1-C2-C3
5	D	1296	GOL	C1-C2-C3-O3
4	C	1296	MES	C7-C8-S-O3S
4	C	1296	MES	N4-C7-C8-S
4	D	1297	MES	C8-C7-N4-C3
4	D	1297	MES	C8-C7-N4-C5
5	D	1296	GOL	O1-C1-C2-O2
2	A	1294	BZ0	NAN-CAU-CAV-CAW
2	A	1294	BZ0	NAN-CAU-CAV-CAY
2	C	1294	BZ0	NAN-CAU-CAV-CAW
2	C	1294	BZ0	NAN-CAU-CAV-CAY
2	D	1294	BZ0	NAN-CAU-CAV-CAW
2	D	1294	BZ0	NAN-CAU-CAV-CAY
2	B	1294	BZ0	NAN-CAU-CAV-CAW
2	B	1294	BZ0	NAN-CAU-CAV-CAY
4	B	1296	MES	C8-C7-N4-C3
4	D	1297	MES	N4-C7-C8-S
2	A	1294	BZ0	CAE-CAF-NAG-CA0

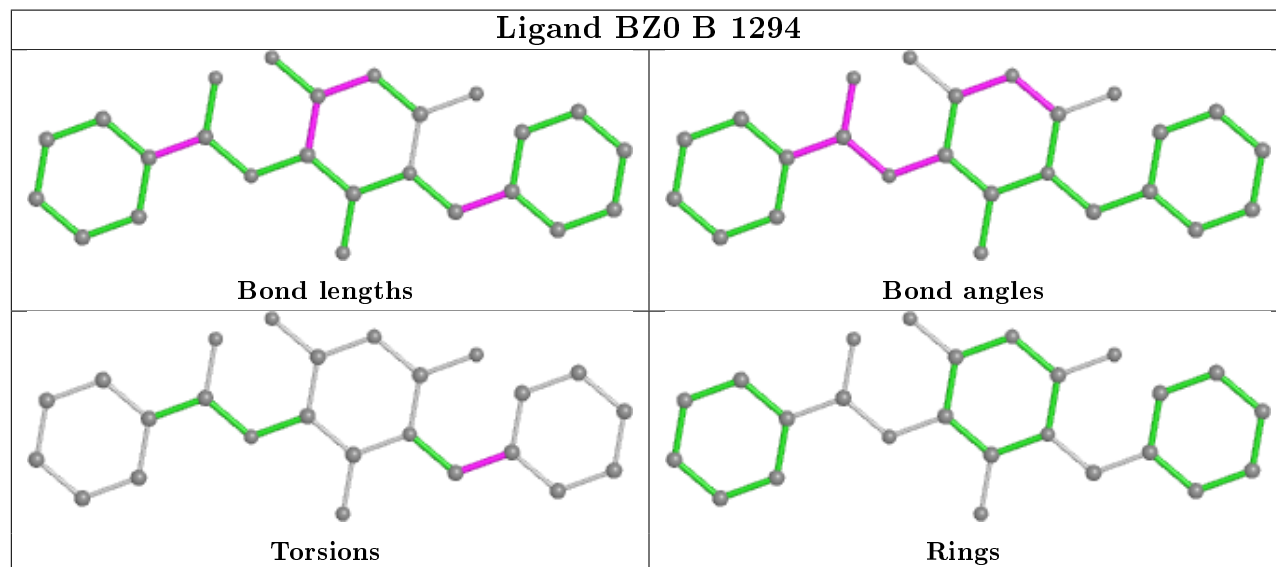
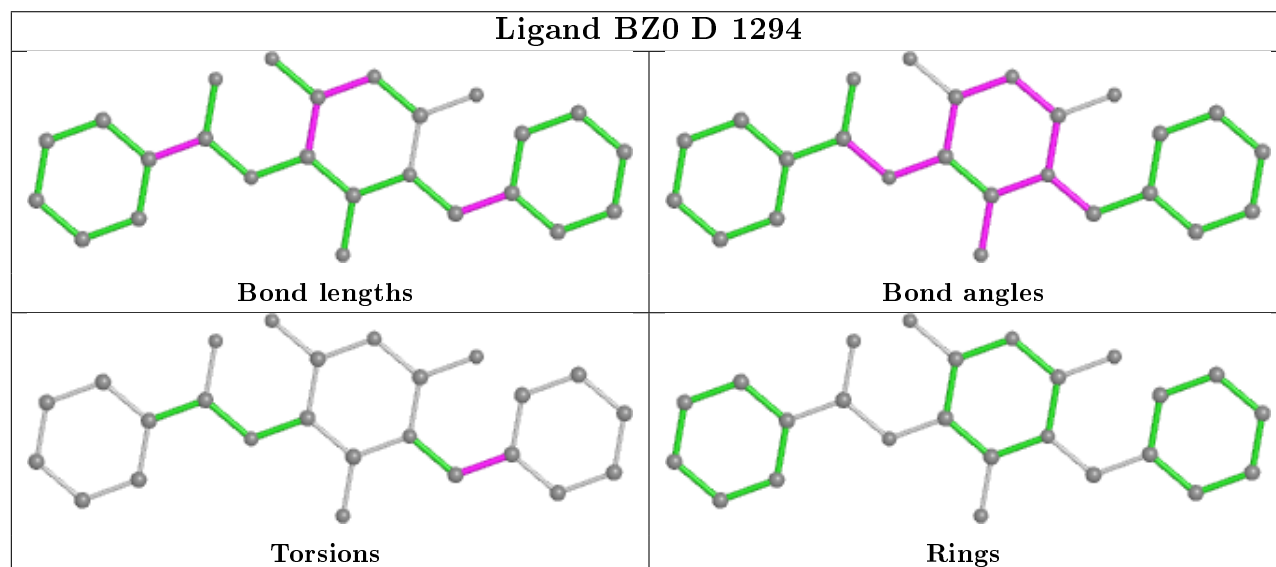
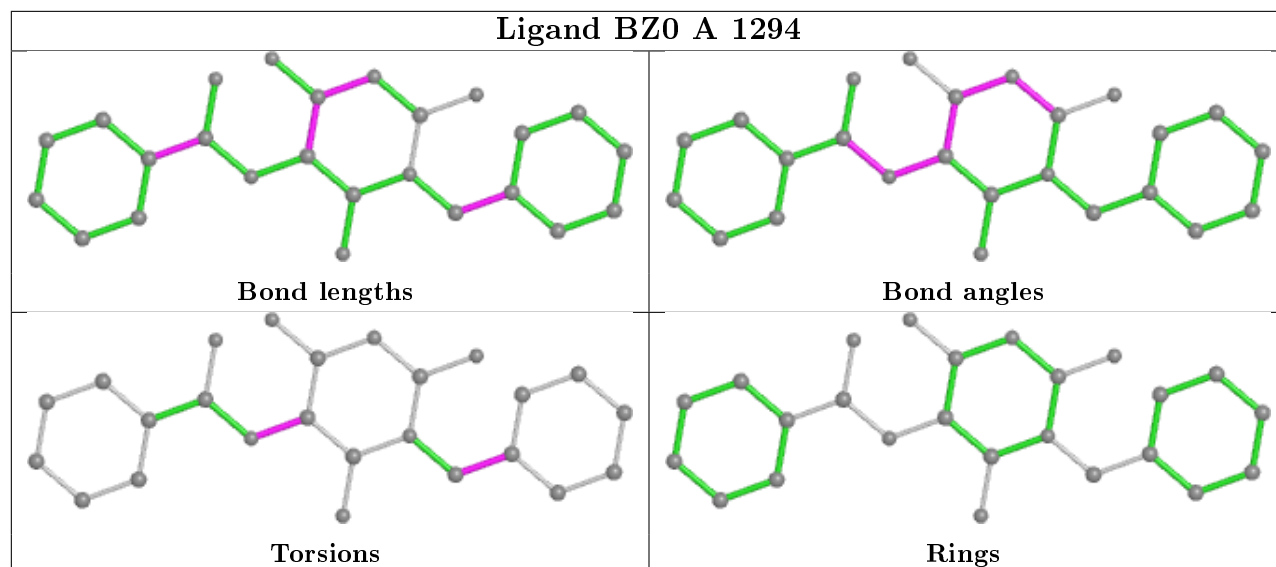
There are no ring outliers.

6 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1294	BZ0	1	0
4	B	1296	MES	2	0
4	A	1296	MES	1	0
4	C	1296	MES	2	0
2	D	1294	BZ0	2	0
2	B	1294	BZ0	1	0

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





5.7 Other polymers [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	297/303 (98%)	0.11	14 (4%)	31 44	28, 42, 72, 92	0
1	B	290/303 (95%)	0.25	15 (5%)	27 39	28, 44, 74, 113	1 (0%)
1	C	288/303 (95%)	0.33	20 (6%)	16 24	32, 49, 75, 114	1 (0%)
1	D	290/303 (95%)	0.23	11 (3%)	40 53	36, 49, 72, 88	1 (0%)
All	All	1165/1212 (96%)	0.23	60 (5%)	27 39	28, 46, 74, 114	3 (0%)

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	14	THR	5.8
1	B	189	LEU	5.3
1	C	276	ALA	4.8
1	B	13	GLY	4.4
1	C	190	LYS	4.3
1	C	277	LYS	3.9
1	B	15	ARG	3.8
1	C	187	ARG	3.5
1	B	277	LYS	3.4
1	D	277	LYS	3.2
1	B	191	PRO	3.2
1	A	193	PRO	3.1
1	B	190	LYS	3.1
1	D	127	GLN	3.0
1	A	189	LEU	3.0
1	C	164	LEU	2.9
1	B	276	ALA	2.9
1	A	276	ALA	2.9
1	C	191	PRO	2.9
1	A	127	GLN	2.8
1	C	278	ASN	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	272	ALA	2.8
1	C	189	LEU	2.7
1	C	273	ALA	2.7
1	B	127	GLN	2.7
1	C	22	ALA	2.7
1	B	187	ARG	2.7
1	A	14	THR	2.7
1	A	191	PRO	2.6
1	A	194	ARG	2.6
1	C	175	LEU	2.6
1	B	222	ALA	2.5
1	D	278	ASN	2.5
1	A	152	GLN	2.5
1	B	17	HIS	2.5
1	A	115	GLY	2.5
1	A	195	GLY	2.4
1	B	1	MET	2.4
1	A	277	LYS	2.4
1	A	275	LEU	2.4
1	C	186	ALA	2.4
1	C	275	LEU	2.3
1	C	285	LYS	2.3
1	C	15	ARG	2.3
1	B	23	ILE	2.3
1	C	93	LEU	2.2
1	D	15	ARG	2.2
1	D	123	GLY	2.2
1	C	127	GLN	2.2
1	B	11	GLY	2.2
1	C	188	ASP	2.2
1	D	205	ALA	2.1
1	D	157	ILE	2.1
1	D	275	LEU	2.1
1	C	13	GLY	2.1
1	D	187	ARG	2.1
1	D	224	LEU	2.1
1	C	14	THR	2.0
1	A	153	GLY	2.0
1	A	222	ALA	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 carbohydrates in this entry.

6.4 Ligands [i](#)

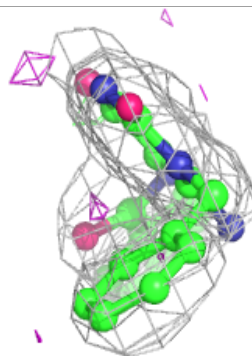
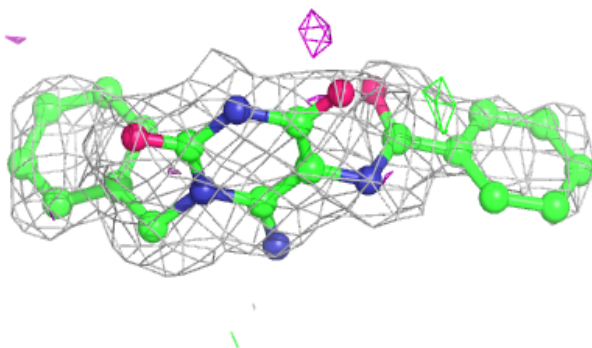
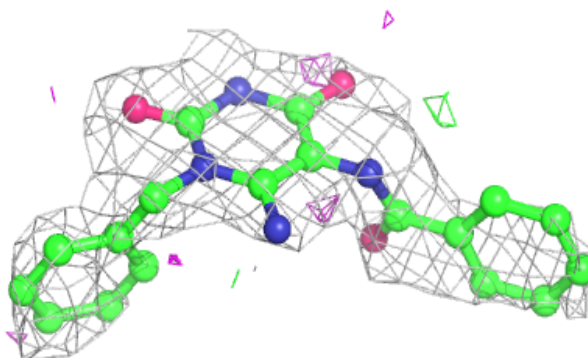
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	BZ0	C	1294	25/25	0.80	0.28	55,70,90,90	0
2	BZ0	D	1294	25/25	0.82	0.23	51,65,76,79	0
2	BZ0	B	1294	25/25	0.82	0.21	46,60,71,72	0
2	BZ0	A	1294	25/25	0.85	0.19	44,57,80,81	0
5	GOL	D	1296	6/6	0.86	0.28	55,64,66,67	0
4	MES	D	1297	12/12	0.91	0.24	61,78,81,82	0
3	CL	D	1295	1/1	0.93	0.20	54,54,54,54	0
4	MES	B	1296	12/12	0.93	0.22	45,58,63,64	0
4	MES	C	1296	12/12	0.94	0.17	57,68,73,73	0
4	MES	A	1296	12/12	0.96	0.22	51,62,68,68	0
3	CL	C	1295	1/1	0.96	0.24	57,57,57,57	0
3	CL	B	1295	1/1	0.97	0.21	50,50,50,50	0
3	CL	A	1295	1/1	0.99	0.19	43,43,43,43	0

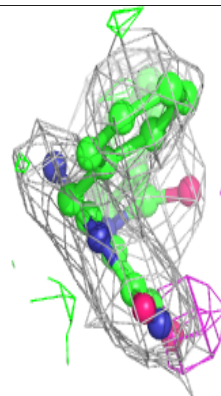
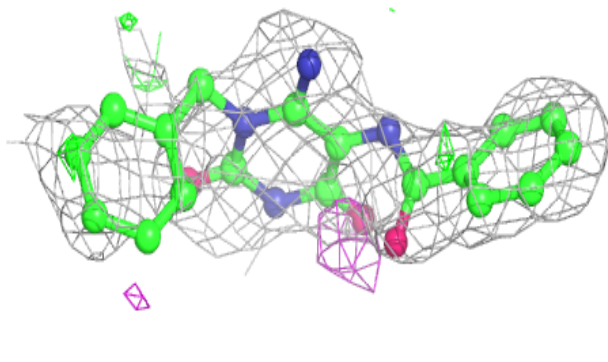
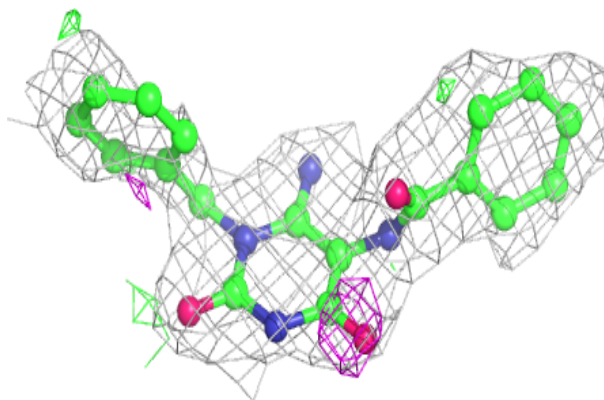
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around BZ0 C 1294:

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

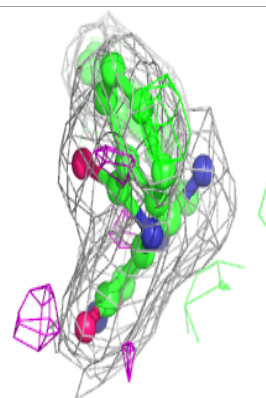
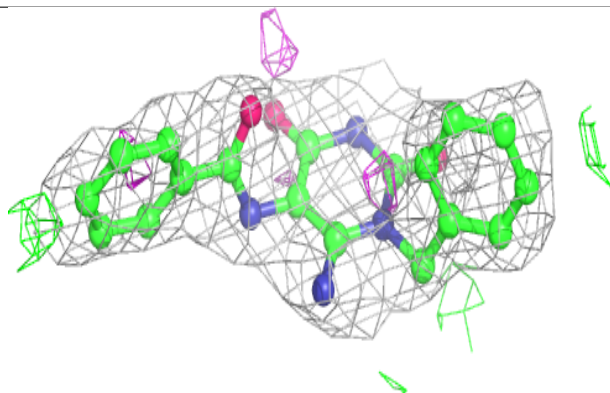
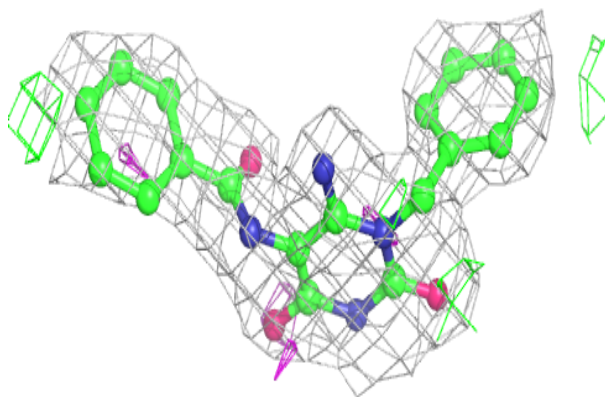
**Electron density around BZ0 D 1294:**

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

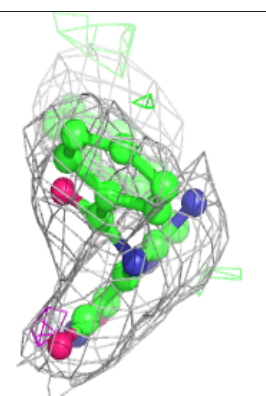
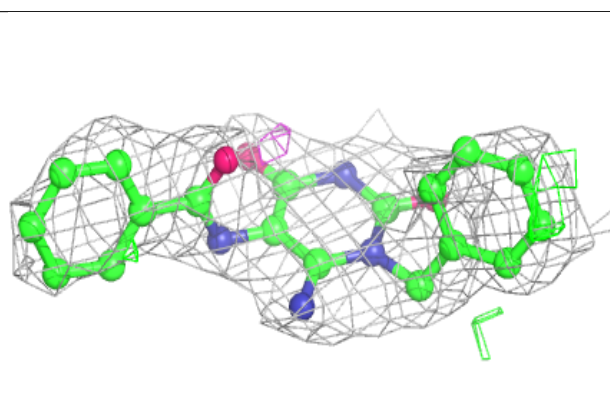
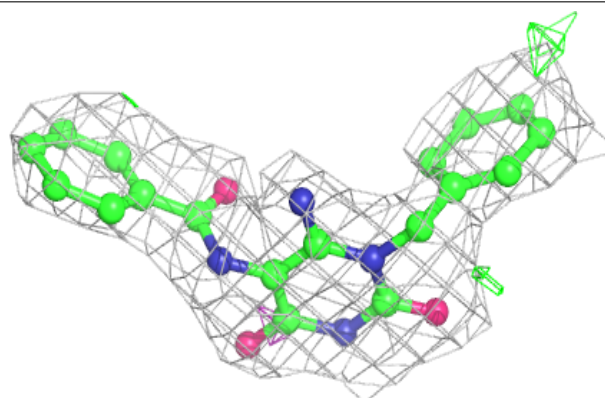


Electron density around BZ0 B 1294:

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

**Electron density around BZ0 A 1294:**

$2mF_o-DF_c$ (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.