



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:21 AM UTC

PDB ID : 4L3Q / pdb_00004l3q
Title : Crystal structure of glucokinase-activator complex
Authors : Gajiwala, K.S.; Filipski, K.J.
Deposited on : 2013-06-06
Resolution : 2.70 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

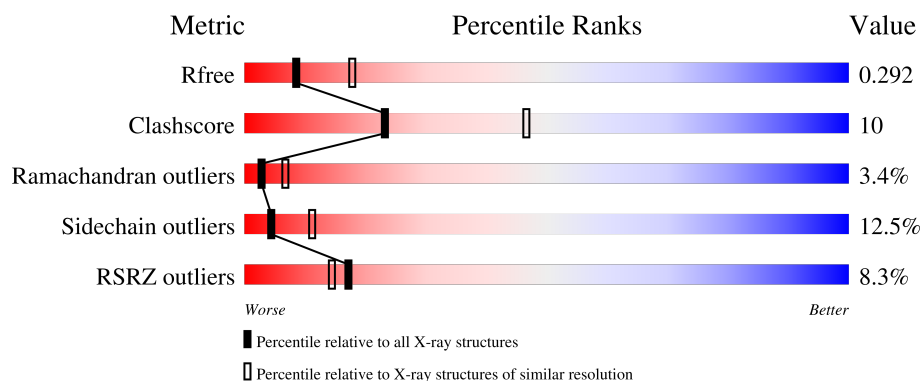
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	457	8% 67% 25% 5%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3537 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.

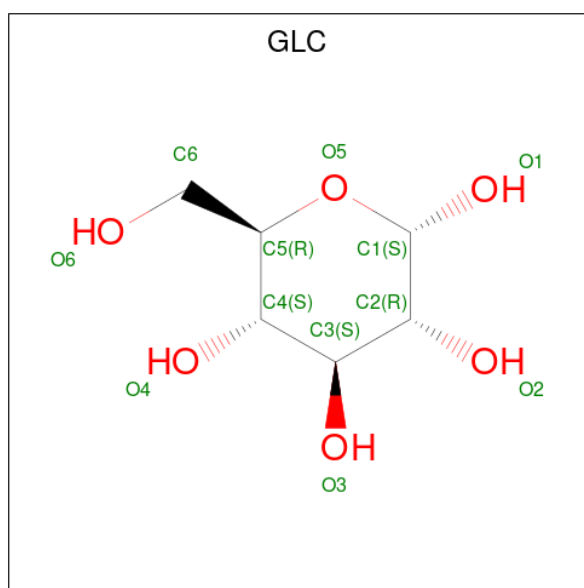
- Molecule 1 is a protein called Glucokinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	447	3499	2175	608	685	31	0	0	0

There are 7 discrepancies between the modelled and reference sequences:

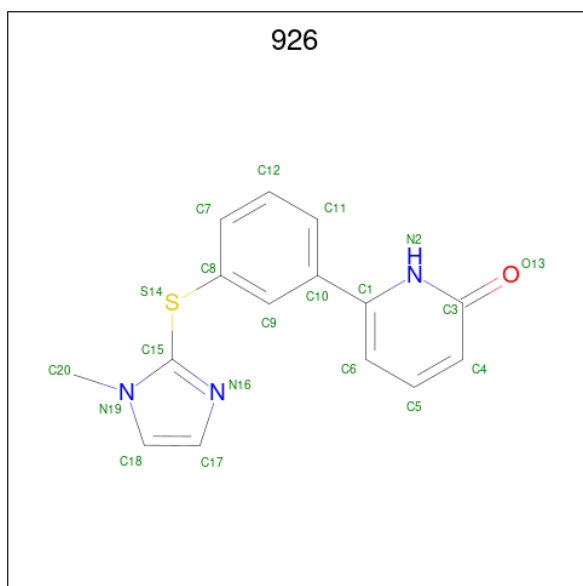
Chain	Residue	Modelled	Actual	Comment	Reference
A	9	GLY	-	expression tag	UNP P35557
A	10	PRO	-	expression tag	UNP P35557
A	11	MET	-	expression tag	UNP P35557
A	12	ALA	-	expression tag	UNP P35557
A	13	LEU	-	expression tag	UNP P35557
A	14	THR	-	expression tag	UNP P35557
A	15	LEU	-	expression tag	UNP P35557

- Molecule 2 is alpha-D-glucopyranose (CCD ID: GLC) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			12	6	6		

- Molecule 3 is 6-{3-[(1-methyl-1H-imidazol-2-yl)sulfanyl]phenyl}pyridin-2(1H)-one (CCD ID: 926) (formula: C₁₅H₁₃N₃OS).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			20	15	3	1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	6	Total	O	0	0
			6	6		

4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	79.51Å 79.51Å 321.13Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.84 – 2.70 39.84 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.6 (39.84-2.70) 99.5 (39.84-2.70)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	10.05 (at 2.69Å)	Xtriage
Refinement program	CNX	Depositor
R, R_{free}	0.254 , 0.304 0.243 , 0.292	Depositor DCC
R_{free} test set	1769 reflections (10.10%)	wwPDB-VP
Wilson B-factor (Å ²)	51.5	Xtriage
Anisotropy	0.006	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 30.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3537	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.96% 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: 926, GLC

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	0.78	2/3552 (0.1%)	1.06	3/4775 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	89	VAL	CA-CB	6.20	1.63	1.54
1	A	210	MET	SD-CE	5.38	1.93	1.79

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	335	VAL	N-CA-C	-7.30	103.42	110.42
1	A	97	GLY	N-CA-C	-5.18	100.89	113.18
1	A	296	LYS	N-CA-C	-5.01	107.12	113.18

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	3499	0	3438	72	0
2	A	12	0	12	0	0
3	A	20	0	13	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	6	0	0	0	0
All	All	3537	0	3463	72	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 10.

All (72) 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:263:SER:HB2	1:A:265:GLU:OE2	1.79	0.81
1:A:457:CYS:C	1:A:459:LYS:H	1.99	0.68
1:A:75:LEU:HD23	1:A:146:LEU:HD13	1.80	0.63
1:A:111:PRO:HD2	1:A:114:ALA:HB3	1.78	0.63
1:A:126:ILE:O	1:A:130:ILE:HG13	2.00	0.61
1:A:73:ASP:HB3	1:A:88:LEU:HD11	1.83	0.59
1:A:92:GLY:HA3	1:A:99:TRP:CE2	2.40	0.57
1:A:103:THR:O	1:A:104:LYS:HG3	2.04	0.56
1:A:238:MET:HB2	1:A:249:GLY:O	2.06	0.56
1:A:111:PRO:O	1:A:112:GLU:HB3	2.05	0.56
1:A:36:ARG:HG2	1:A:36:ARG:HH11	1.69	0.56
1:A:166:ASN:HA	1:A:177:GLU:HG2	1.88	0.55
1:A:409:ASP:OD1	1:A:414:LYS:HE3	2.09	0.53
1:A:259:ALA:HB2	1:A:287:GLN:OE1	2.10	0.52
1:A:64:SER:O	1:A:65:THR:C	2.52	0.52
1:A:85:ARG:HD3	1:A:107:MET:HE3	1.91	0.52
1:A:447:ARG:O	1:A:451:LEU:HD12	2.09	0.51
1:A:209:THR:HG23	1:A:441:SER:HB3	1.94	0.50
1:A:452:VAL:HG13	3:A:501:926:H2	1.93	0.50
1:A:92:GLY:HA3	1:A:99:TRP:CZ2	2.46	0.50
1:A:100:SER:HB2	1:A:447:ARG:HG2	1.93	0.50
1:A:111:PRO:HD2	1:A:114:ALA:CB	2.42	0.50
1:A:277:VAL:HG23	1:A:304:LEU:HD22	1.94	0.49
1:A:369:ARG:NH1	1:A:369:ARG:HA	2.27	0.49
1:A:119:ALA:HB2	1:A:176:ALA:HB2	1.93	0.49
1:A:165:LEU:O	1:A:166:ASN:HB3	2.13	0.49
1:A:41:MET:HE2	1:A:392:ARG:HD3	1.95	0.48
1:A:74:PHE:CD2	1:A:145:PRO:HB2	2.48	0.48
1:A:457:CYS:C	1:A:459:LYS:N	2.67	0.47
1:A:22:GLU:O	1:A:271:LEU:HD21	2.15	0.47
1:A:154:VAL:HG12	1:A:156:HIS:HD2	1.80	0.47
1:A:368:ARG:O	1:A:372:GLU:HG3	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:ARG:H	3:A:501:926:C3	2.28	0.47
1:A:77:LEU:HG	1:A:146:LEU:HD11	1.98	0.46
1:A:192:ARG:C	1:A:194:ASP:H	2.23	0.46
1:A:85:ARG:HG2	1:A:107:MET:HG2	1.98	0.46
1:A:152:PHE:HB3	1:A:153:PRO:HD2	1.96	0.45
1:A:260:PHE:O	1:A:289:TYR:HB2	2.17	0.45
1:A:14:THR:HG23	1:A:14:THR:O	2.17	0.45
1:A:163:ILE:HA	1:A:179:ASN:O	2.17	0.45
1:A:139:MET:HG2	1:A:142:LYS:NZ	2.32	0.44
1:A:194:ASP:HB2	1:A:195:PHE:HD2	1.82	0.44
1:A:258:GLY:O	1:A:287:GLN:HA	2.17	0.44
1:A:57:MET:CE	1:A:234:TYR:OH	2.65	0.44
1:A:84:PHE:CZ	1:A:126:ILE:HG23	2.53	0.44
1:A:224:MET:HE1	1:A:423:PHE:CZ	2.53	0.44
1:A:36:ARG:HG2	1:A:36:ARG:NH1	2.32	0.43
1:A:110:ILE:H	1:A:110:ILE:HG13	1.60	0.43
1:A:83:ASN:OD1	1:A:107:MET:HE2	2.17	0.43
1:A:262:ASP:OD2	1:A:286:GLN:HA	2.18	0.43
1:A:265:GLU:H	1:A:265:GLU:CD	2.27	0.43
1:A:165:LEU:HA	1:A:165:LEU:HD23	1.73	0.43
1:A:15:LEU:H	1:A:15:LEU:HD12	1.84	0.43
1:A:398:SER:C	1:A:399:GLU:HG2	2.42	0.43
1:A:447:ARG:CB	1:A:447:ARG:HH11	2.31	0.43
1:A:222:VAL:HB	1:A:406:VAL:HG22	2.00	0.43
1:A:101:VAL:HG22	1:A:102:LYS:O	2.18	0.42
1:A:21:ALA:HB1	1:A:24:GLN:NE2	2.34	0.42
1:A:410:GLY:O	1:A:414:LYS:HG3	2.20	0.42
1:A:94:GLY:N	1:A:97:GLY:O	2.52	0.42
1:A:369:ARG:HH12	1:A:372:GLU:HB2	1.85	0.42
1:A:56:LYS:O	1:A:57:MET:C	2.62	0.41
1:A:21:ALA:CB	1:A:24:GLN:HE21	2.32	0.41
1:A:21:ALA:HB1	1:A:24:GLN:HE21	1.86	0.41
1:A:43:ARG:HH11	1:A:43:ARG:HG2	1.86	0.41
1:A:160:ASP:HB3	1:A:200:VAL:O	2.21	0.41
1:A:237:GLU:OE2	1:A:237:GLU:HA	2.21	0.41
1:A:323:GLN:C	1:A:325:ARG:N	2.74	0.41
1:A:309:LEU:HD12	1:A:309:LEU:HA	1.85	0.40
1:A:320:ALA:HB3	1:A:325:ARG:HH11	1.86	0.40
1:A:186:ARG:O	1:A:190:LYS:HD2	2.21	0.40
1:A:55:VAL:O	1:A:57:MET:HG2	2.20	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	445/457 (97%)	402 (90%)	28 (6%)	15 (3%)	3 7

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	112	GLU
1	A	101	VAL
1	A	111	PRO
1	A	116	THR
1	A	136	LYS
1	A	166	ASN
1	A	454	ALA
1	A	458	LYS
1	A	22	GLU
1	A	67	GLU
1	A	135	ASP
1	A	115	MET
1	A	65	THR
1	A	68	GLY
1	A	159	ILE

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	383/390 (98%)	335 (88%)	48 (12%)	4 11

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	THR
1	A	15	LEU
1	A	16	VAL
1	A	26	GLN
1	A	32	LYS
1	A	36	ARG
1	A	39	LYS
1	A	64	SER
1	A	65	THR
1	A	82	THR
1	A	83	ASN
1	A	89	VAL
1	A	96	GLU
1	A	100	SER
1	A	102	LYS
1	A	110	ILE
1	A	120	GLU
1	A	127	SER
1	A	140	LYS
1	A	142	LYS
1	A	144	LEU
1	A	187	ASP
1	A	198	ASP
1	A	200	VAL
1	A	216	GLU
1	A	243	LEU
1	A	248	GLU
1	A	271	LEU
1	A	277	VAL
1	A	304	LEU
1	A	309	LEU
1	A	310	VAL
1	A	315	LEU
1	A	317	HIS
1	A	327	ARG
1	A	340	SER
1	A	347	GLN
1	A	361	THR
1	A	368	ARG
1	A	381	MET
1	A	386	LEU
1	A	397	ARG

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Mol	Chain	Res	Type
1	A	399	GLU
1	A	400	ASP
1	A	421	GLU
1	A	447	ARG
1	A	455	VAL
1	A	458	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	38	GLN
1	A	98	GLN
1	A	156	HIS
1	A	313	ASN
1	A	337	GLN

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	GLC	A	500	-	12,12,12	0.79	0	17,17,17	3.10	11 (64%)
3	926	A	501	-	22,22,22	2.31	7 (31%)	29,30,30	2.36	14 (48%)

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	GLC	A	500	-	-	0/2/22/22	0/1/1/1
3	926	A	501	-	-	4/8/8/8	0/3/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	501	926	C15-N19	4.91	1.40	1.36
3	A	501	926	C6-C1	4.55	1.40	1.36
3	A	501	926	C15-N16	3.59	1.38	1.31
3	A	501	926	C12-C7	3.48	1.44	1.38
3	A	501	926	C12-C11	3.32	1.44	1.38
3	A	501	926	C7-C8	2.50	1.43	1.39
3	A	501	926	C10-C1	2.27	1.51	1.47

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	GLC	O5-C1-C2	5.55	120.05	110.30
2	A	500	GLC	C4-C3-C2	5.11	119.79	110.83
3	A	501	926	C8-S14-C15	4.60	109.48	101.87
3	A	501	926	S14-C15-N16	4.46	127.18	124.96
2	A	500	GLC	C1-O5-C5	4.28	121.93	113.65
2	A	500	GLC	O5-C5-C4	4.19	117.25	109.70
2	A	500	GLC	O4-C4-C5	4.02	119.23	109.32
2	A	500	GLC	C3-C4-C5	3.73	116.99	110.23
3	A	501	926	C9-C10-C1	-3.68	114.74	120.43
3	A	501	926	C11-C10-C1	3.66	126.36	120.81
3	A	501	926	C1-N2-C3	3.42	128.35	124.23
3	A	501	926	C5-C4-C3	-3.41	116.57	120.54
3	A	501	926	C10-C9-C8	3.39	123.20	119.97
3	A	501	926	C7-C8-S14	3.09	128.62	120.14
2	A	500	GLC	O1-C1-O5	-3.09	101.25	110.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	926	C9-C8-S14	-3.05	112.69	119.39
2	A	500	GLC	O2-C2-C1	2.86	115.85	109.25
3	A	501	926	C10-C1-C6	2.74	127.22	123.59
2	A	500	GLC	C1-C2-C3	2.45	115.36	110.36
3	A	501	926	C5-C6-C1	-2.26	117.31	118.88
3	A	501	926	C6-C5-C4	2.26	124.71	121.14
3	A	501	926	C20-N19-C15	-2.25	123.31	127.02
3	A	501	926	C20-N19-C18	2.20	128.78	125.11
2	A	500	GLC	O1-C1-C2	-2.12	102.83	108.98
2	A	500	GLC	O3-C3-C4	2.04	115.18	110.38

There are no chirality outliers.

All (4) torsion outliers are listed below:

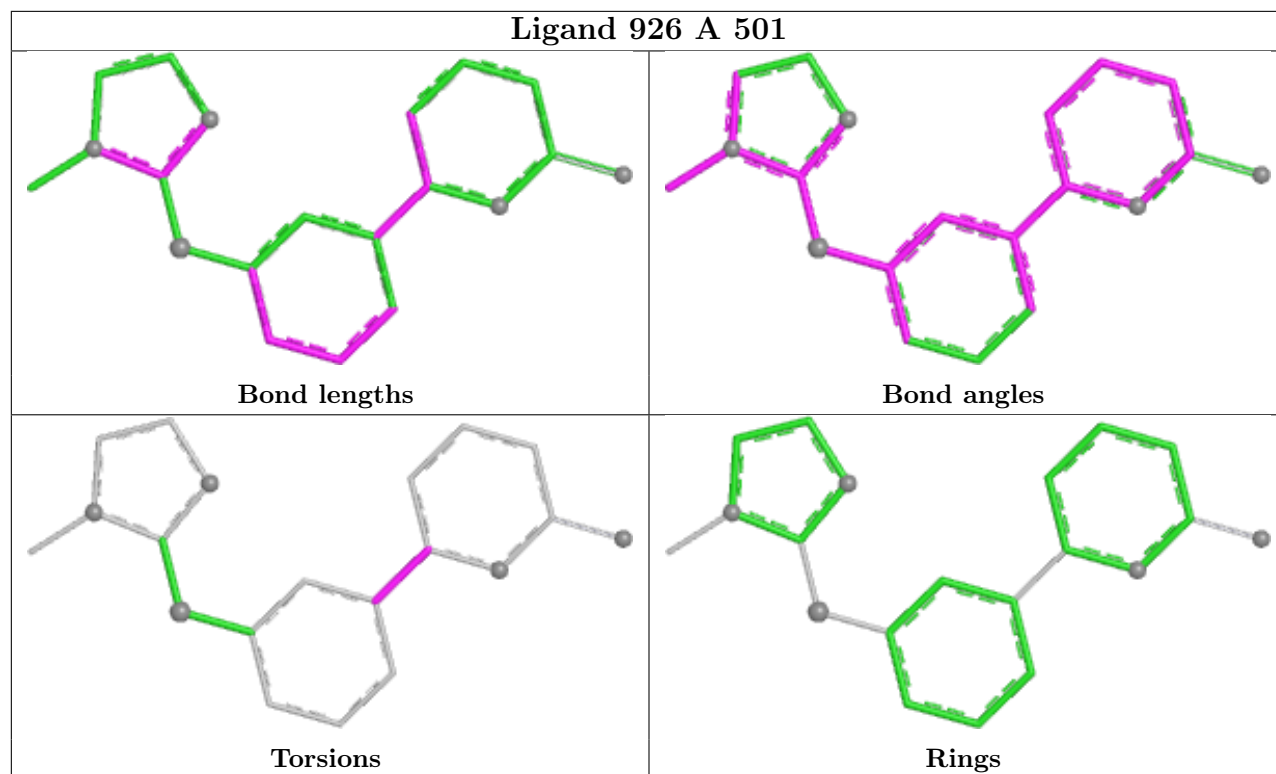
Mol	Chain	Res	Type	Atoms
3	A	501	926	C6-C1-C10-C11
3	A	501	926	N2-C1-C10-C11
3	A	501	926	N2-C1-C10-C9
3	A	501	926	C6-C1-C10-C9

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	501	926	2	0

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



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	447/457 (97%)	0.49	37 (8%) 17 14	25, 49, 90, 101	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	460	ALA	5.0
1	A	139	MET	3.7
1	A	134	LEU	3.6
1	A	190	LYS	3.5
1	A	65	THR	3.4
1	A	130	ILE	3.2
1	A	110	ILE	3.1
1	A	114	ALA	3.1
1	A	197	MET	3.0
1	A	116	THR	3.0
1	A	317	HIS	2.9
1	A	195	PHE	2.9
1	A	147	GLY	2.9
1	A	82	THR	2.7
1	A	175	GLY	2.6
1	A	109	SER	2.5
1	A	15	LEU	2.5
1	A	144	LEU	2.5
1	A	320	ALA	2.5
1	A	133	PHE	2.5
1	A	99	TRP	2.3
1	A	91	VAL	2.2
1	A	198	ASP	2.2
1	A	89	VAL	2.2
1	A	173	ALA	2.2
1	A	142	LYS	2.2
1	A	14	THR	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	454	ALA	2.1
1	A	272	GLU	2.1
1	A	81	GLY	2.1
1	A	137	HIS	2.1
1	A	118	THR	2.1
1	A	66	PRO	2.1
1	A	141	HIS	2.1
1	A	191	ARG	2.0
1	A	92	GLY	2.0
1	A	189	ILE	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](#)

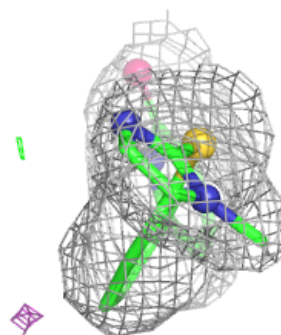
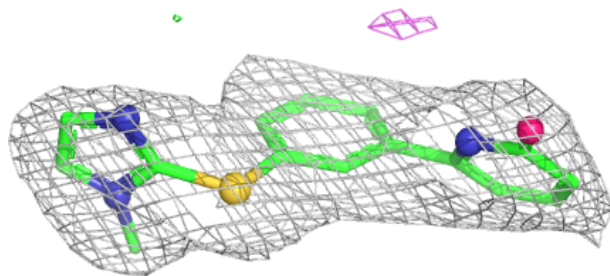
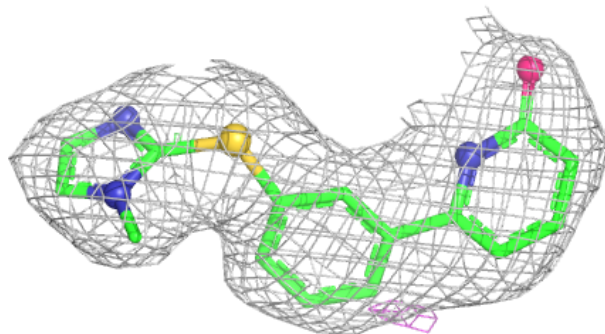
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	GLC	A	500	12/12	0.91	0.11	32,36,39,40	0
3	926	A	501	20/20	0.94	0.09	37,40,44,44	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 926 A 501:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.