



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 08:33 AM UTC

PDB ID : 3ZLK / pdb_00003zlk
Title : Pseudomonas aeruginosa RmlA in complex with allosteric inhibitor
Authors : Alphey, M.S.; Pirrie, L.; Torrie, L.S.; Gardiner, M.; Sarkar, A.; Brenk, R.;
Westwood, N.J.; Gray, D.; Naismith, J.H.
Deposited on : 2013-02-01
Resolution : 1.95 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

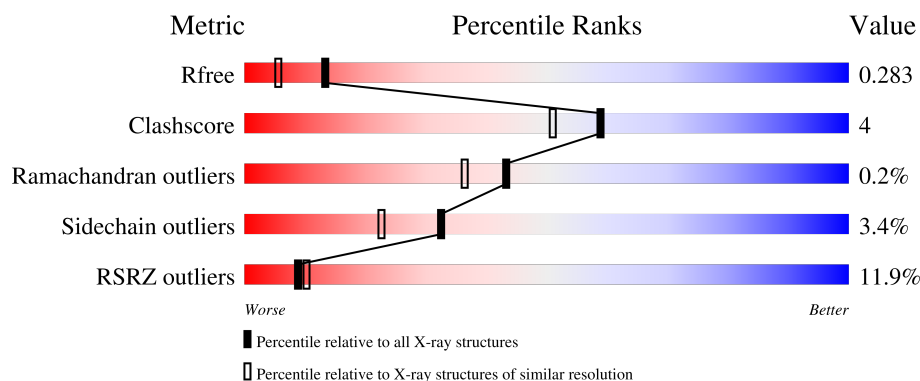
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	303	9% 89% 8% ..
1	B	303	12% 90% 6% .
1	C	303	13% 86% 10% ..
1	D	303	12% 87% 9% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10058 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	0	0
			2335	1492	400	439	4			
1	B	293	Total	C	N	O	S	0	1	0
			2299	1471	387	436	5			
1	C	292	Total	C	N	O	S	0	1	0
			2291	1466	386	435	4			
1	D	293	Total	C	N	O	S	0	1	0
			2299	1471	387	436	5			

There are 40 discrepancies between the modelled and reference sequences:

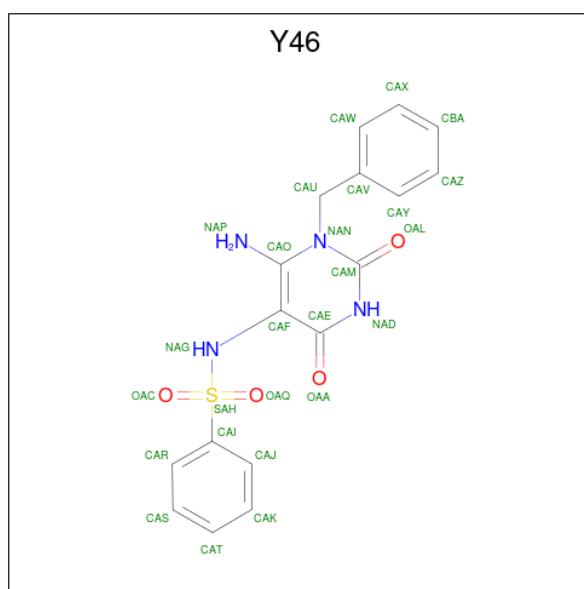
Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	HIS	-	expression tag	UNP G3XCK4
A	-8	HIS	-	expression tag	UNP G3XCK4
A	-7	HIS	-	expression tag	UNP G3XCK4
A	-6	HIS	-	expression tag	UNP G3XCK4
A	-5	HIS	-	expression tag	UNP G3XCK4
A	-4	HIS	-	expression tag	UNP G3XCK4
A	-3	GLY	-	expression tag	UNP G3XCK4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

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-9	HIS	-	expression tag	UNP G3XCK4
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-TETRAHYDOPYRIMIDIN-5-YL)BENZENESULFONAMIDE (CCD ID: Y46) (formula: C₁₇H₁₆N₄O₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			26	17	4	4	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	S	0	0
			26	17	4	4	1		
2	C	1	Total	C	N	O	S	0	0
			26	17	4	4	1		
2	D	1	Total	C	N	O	S	0	0
			26	17	4	4	1		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

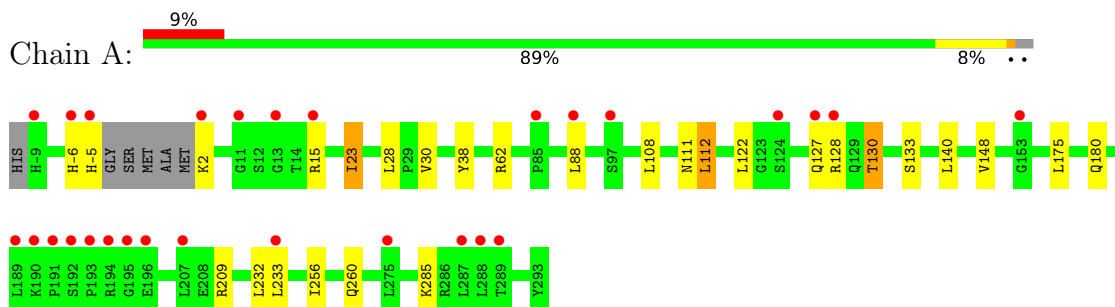
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	208	Total	O	0	0
			208	208		
4	B	189	Total	O	0	0
			189	189		
4	C	190	Total	O	0	0
			190	190		
4	D	139	Total	O	0	0
			139	139		

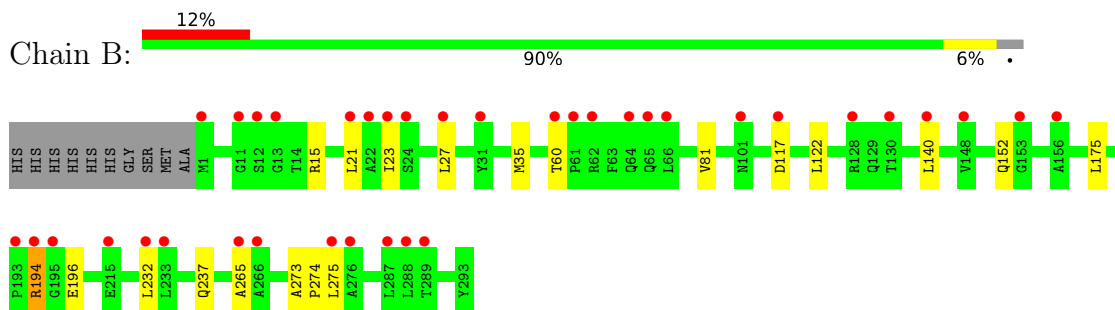
3 Residue-property plots [i](#)

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.

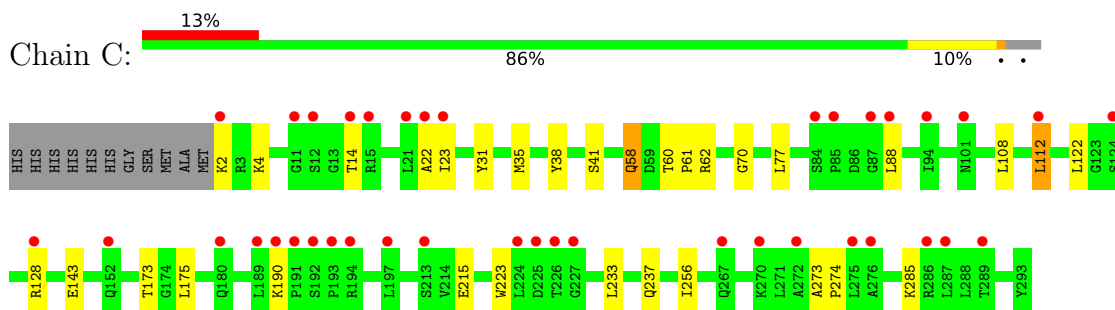
• Molecule 1: GLUCOSE-1-PHOSPHATE THYMIDYLTRANSFERASE



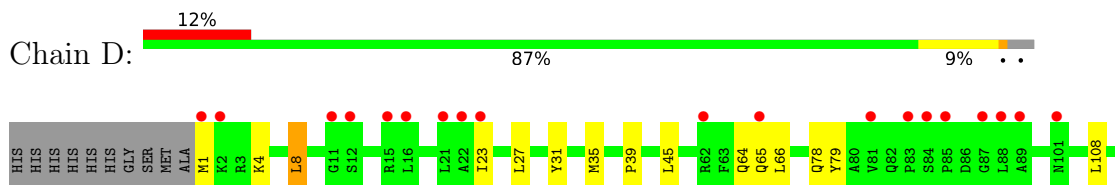
• Molecule 1: GLUCOSE-1-PHOSPHATE THYMIDYLTRANSFERASE

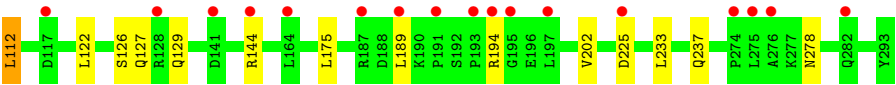


• Molecule 1: GLUCOSE-1-PHOSPHATE THYMIDYLTRANSFERASE



• 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	64.30Å 155.00Å 134.45Å 90.00° 92.03° 90.00°	Depositor
Resolution (Å)	37.23 – 1.95 37.23 – 1.95	Depositor EDS
% Data completeness (in resolution range)	84.7 (37.23-1.95) 84.7 (37.23-1.95)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.6.0017	Depositor
R, R_{free}	0.233 , 0.277 0.239 , 0.283	Depositor DCC
R_{free} test set	4049 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	21.5	Xtriage
Anisotropy	0.096	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 36.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.080 for h,-k,-l	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	10058	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.83% 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: Y46, CL

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	0.75	0/2389	0.79	0/3240
1	B	0.68	0/2352	0.78	0/3190
1	C	0.69	0/2344	0.79	0/3180
1	D	0.67	0/2352	0.78	0/3190
All	All	0.70	0/9437	0.78	0/12800

There are no bond length outliers.

There are no bond angle outliers.

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	2335	0	2308	30	0
1	B	2299	0	2292	20	0
1	C	2291	0	2280	26	0
1	D	2299	0	2292	24	0
2	A	26	0	16	0	0
2	B	26	0	16	0	0
2	C	26	0	16	2	0
2	D	26	0	16	2	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	208	0	0	9	0
4	B	189	0	0	4	0
4	C	190	0	0	9	0
4	D	139	0	0	1	0
All	All	10058	0	9236	82	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 4.

All (82) 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:C:22:ALA:HB3	4:C:2014:HOH:O	1.38	1.21
1:A:130:THR:HG23	4:A:2102:HOH:O	1.59	1.03
1:A:23:ILE:HG23	1:B:23:ILE:HD13	1.46	0.95
1:A:233:LEU:HD21	1:B:237[A]:GLN:HG2	1.53	0.91
1:A:23:ILE:CG2	1:B:23:ILE:HD13	2.02	0.90
1:A:233:LEU:HD21	1:B:237[A]:GLN:CG	2.11	0.79
1:C:23:ILE:HD13	1:D:23:ILE:HD12	1.69	0.75
1:A:23:ILE:HD11	1:A:28:LEU:HD21	1.70	0.72
1:A:233:LEU:HD11	1:B:237[A]:GLN:HG3	1.73	0.70
1:A:23:ILE:CD1	1:A:28:LEU:HD21	2.23	0.68
1:C:128:ARG:NH2	4:C:2089:HOH:O	2.29	0.64
4:A:2084:HOH:O	1:D:65:GLN:HB2	1.97	0.64
1:B:15:ARG:NH1	4:B:2013:HOH:O	2.32	0.63
1:B:273:ALA:HB3	1:B:274:PRO:HD3	1.80	0.62
1:A:23:ILE:HG23	1:B:23:ILE:CD1	2.25	0.61
1:B:23:ILE:HD11	1:B:27:LEU:HD13	1.82	0.60
2:C:400:Y46:CAS	4:C:2028:HOH:O	2.50	0.60
1:C:273:ALA:HB3	1:C:274:PRO:HD3	1.83	0.59
1:A:23:ILE:CG2	1:B:23:ILE:CD1	2.80	0.58
1:A:2:LYS:HB3	4:A:2007:HOH:O	2.02	0.58
4:A:2084:HOH:O	1:D:65:GLN:CB	2.52	0.57
1:A:23:ILE:HD12	1:A:23:ILE:O	2.04	0.57
1:D:45:LEU:HD22	2:D:400:Y46:CAJ	2.35	0.57
1:C:233:LEU:HD21	1:D:237[B]:GLN:HA	1.86	0.56
1:D:45:LEU:HD22	2:D:400:Y46:HAJ	1.88	0.56
1:C:233:LEU:HD21	1:D:237[A]:GLN:HA	1.86	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:ILE:CD1	1:A:28:LEU:CD2	2.84	0.55
1:C:35:MET:HG2	4:C:2024:HOH:O	2.07	0.54
1:A:88:LEU:HD13	1:A:108:LEU:HD21	1.90	0.54
1:D:1:MET:HE3	1:D:129:GLN:NE2	2.23	0.53
1:A:23:ILE:HD13	1:A:28:LEU:CD2	2.41	0.50
1:D:27:LEU:HD22	1:D:66:LEU:HD22	1.93	0.50
1:B:60:THR:HG23	4:B:2045:HOH:O	2.10	0.50
1:A:23:ILE:HG21	1:B:23:ILE:HD13	1.92	0.50
1:D:112:LEU:HD12	1:D:112:LEU:C	2.36	0.50
1:A:62:ARG:NH1	4:A:2047:HOH:O	2.44	0.50
1:A:38:TYR:HB2	1:A:112:LEU:HD22	1.93	0.50
1:C:237[B]:GLN:HA	1:D:233:LEU:HD21	1.94	0.50
1:D:108:LEU:HD12	4:D:2053:HOH:O	2.11	0.49
1:C:233:LEU:HD11	1:D:237[A]:GLN:HG3	1.96	0.48
1:D:122:LEU:CD2	1:D:175:LEU:HD21	2.43	0.48
1:A:-5:HIS:CE1	1:D:78:GLN:NE2	2.82	0.48
1:C:273:ALA:HB3	1:C:274:PRO:CD	2.44	0.48
1:C:237[A]:GLN:HA	1:D:233:LEU:HD21	1.96	0.48
1:A:23:ILE:HG21	1:B:23:ILE:HG21	1.96	0.48
1:C:237[A]:GLN:HG3	1:D:233:LEU:HD11	1.97	0.47
1:B:122:LEU:CD2	1:B:175:LEU:HD21	2.45	0.46
1:A:30:VAL:HG11	1:A:112:LEU:HD11	1.96	0.46
1:A:233:LEU:HD21	1:B:237[B]:GLN:HB2	1.98	0.46
1:C:122:LEU:HD21	1:C:175:LEU:HD21	1.96	0.46
1:A:128:ARG:HD2	1:A:133:SER:OG	2.16	0.46
1:B:81:VAL:HG22	4:B:2062:HOH:O	2.14	0.46
2:C:400:Y46:CAR	4:C:2028:HOH:O	2.64	0.46
1:C:190:LYS:N	4:C:2130:HOH:O	2.48	0.45
1:B:35:MET:HE1	1:B:232:LEU:HD21	1.99	0.45
1:C:23:ILE:HD13	1:D:23:ILE:CD1	2.43	0.45
4:A:2084:HOH:O	1:D:65:GLN:CG	2.65	0.45
1:A:130:THR:CG2	4:A:2102:HOH:O	2.39	0.44
1:C:273:ALA:HB3	4:C:2179:HOH:O	2.17	0.44
1:C:215:GLU:HG2	4:C:2029:HOH:O	2.17	0.44
1:C:58:GLN:CD	1:C:58:GLN:H	2.25	0.44
1:D:64:GLN:HG2	1:D:79:TYR:CD1	2.53	0.44
1:A:2:LYS:HE3	4:A:2007:HOH:O	2.17	0.44
1:B:194:ARG:NH1	1:B:196:GLU:OE1	2.51	0.43
1:C:88:LEU:HD13	1:C:108:LEU:HD21	2.01	0.43
1:A:122:LEU:CD2	1:A:175:LEU:HD21	2.49	0.43
1:D:35:MET:O	1:D:39:PRO:HD2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:38:TYR:CB	1:A:112:LEU:HD22	2.49	0.42
1:C:70:GLY:HA3	1:C:77:LEU:HG	2.02	0.42
1:D:8:LEU:C	1:D:8:LEU:HD12	2.44	0.42
1:C:38:TYR:HB2	1:C:112:LEU:HD22	2.01	0.42
1:C:60:THR:N	1:C:61:PRO:HD2	2.35	0.42
1:C:108:LEU:HD12	4:C:2074:HOH:O	2.19	0.42
1:C:41:SER:HB3	1:C:256:ILE:CD1	2.50	0.41
1:D:189:LEU:HD11	1:D:202:VAL:CG2	2.50	0.41
1:A:256:ILE:O	1:A:260:GLN:HG2	2.21	0.41
1:A:122:LEU:HD21	1:A:175:LEU:HD21	2.03	0.41
1:C:14:THR:O	1:D:278:ASN:ND2	2.54	0.41
1:C:173:THR:HG22	1:C:223:TRP:CE3	2.56	0.41
1:B:140:LEU:HD12	1:B:140:LEU:O	2.21	0.41
1:B:265:ALA:HB3	4:B:2173:HOH:O	2.21	0.40
1:A:128:ARG:HG3	4:A:2101:HOH:O	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	293/303 (97%)	286 (98%)	7 (2%)	0	100	100
1	B	292/303 (96%)	286 (98%)	6 (2%)	0	100	100
1	C	291/303 (96%)	286 (98%)	4 (1%)	1 (0%)	36	28
1	D	292/303 (96%)	287 (98%)	4 (1%)	1 (0%)	36	28
All	All	1168/1212 (96%)	1145 (98%)	21 (2%)	2 (0%)	43	36

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	31	TYR
1	D	31	TYR

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	244/248 (98%)	231 (95%)	13 (5%)	20	9
1	B	241/248 (97%)	236 (98%)	5 (2%)	47	41
1	C	240/248 (97%)	233 (97%)	7 (3%)	37	29
1	D	241/248 (97%)	233 (97%)	8 (3%)	33	24
All	All	966/992 (97%)	933 (97%)	33 (3%)	32	23

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-6	HIS
1	A	15	ARG
1	A	23	ILE
1	A	111	ASN
1	A	112	LEU
1	A	127	GLN
1	A	130	THR
1	A	140	LEU
1	A	148	VAL
1	A	180	GLN
1	A	209	ARG
1	A	232	LEU
1	A	285	LYS
1	B	21	LEU
1	B	117	ASP
1	B	152	GLN
1	B	194	ARG
1	B	275	LEU
1	C	2	LYS
1	C	4	LYS

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Mol	Chain	Res	Type
1	C	58	GLN
1	C	62	ARG
1	C	112	LEU
1	C	143	GLU
1	C	285	LYS
1	D	4	LYS
1	D	8	LEU
1	D	112	LEU
1	D	126	SER
1	D	127	GLN
1	D	144	ARG
1	D	194	ARG
1	D	225	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	-5	HIS
1	B	152	GLN
1	B	169	ASN
1	B	181	GLN
1	C	82	GLN
1	C	90	GLN
1	C	127	GLN
1	D	78	GLN

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	Y46	B	400	-	28,28,28	2.86	6 (21%)	35,40,40	2.07	9 (25%)
2	Y46	A	400	-	28,28,28	3.09	7 (25%)	35,40,40	1.96	6 (17%)
2	Y46	C	400	-	28,28,28	3.03	7 (25%)	35,40,40	1.85	6 (17%)
2	Y46	D	400	-	28,28,28	2.81	6 (21%)	35,40,40	2.10	8 (22%)

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	Y46	B	400	-	-	0/13/15/15	0/3/3/3
2	Y46	A	400	-	-	0/13/15/15	0/3/3/3
2	Y46	C	400	-	-	0/13/15/15	0/3/3/3
2	Y46	D	400	-	-	0/13/15/15	0/3/3/3

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	400	Y46	OAQ-SAH	9.51	1.54	1.43
2	D	400	Y46	OAQ-SAH	8.99	1.53	1.43
2	A	400	Y46	OAQ-SAH	8.62	1.53	1.43
2	A	400	Y46	OAC-SAH	8.41	1.53	1.43
2	B	400	Y46	OAC-SAH	8.07	1.52	1.43
2	B	400	Y46	OAQ-SAH	7.98	1.52	1.43
2	C	400	Y46	OAC-SAH	7.77	1.52	1.43
2	D	400	Y46	OAC-SAH	7.27	1.52	1.43
2	A	400	Y46	CAF-NAG	-7.18	1.34	1.42
2	B	400	Y46	CAF-NAG	-6.84	1.35	1.42
2	C	400	Y46	CAF-NAG	-6.54	1.35	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	400	Y46	CAF-NAG	-6.22	1.35	1.42
2	C	400	Y46	CAU-CAV	-6.16	1.40	1.51
2	D	400	Y46	CAU-CAV	-5.66	1.41	1.51
2	A	400	Y46	CAU-CAV	-5.52	1.41	1.51
2	B	400	Y46	CAU-CAV	-5.01	1.42	1.51
2	A	400	Y46	CAI-SAH	3.76	1.82	1.76
2	B	400	Y46	CAM-NAN	-2.87	1.35	1.39
2	A	400	Y46	CAU-NAN	2.80	1.52	1.47
2	A	400	Y46	CAM-NAN	-2.52	1.35	1.39
2	C	400	Y46	CAM-NAN	-2.40	1.35	1.39
2	C	400	Y46	SAH-NAG	2.35	1.69	1.64
2	D	400	Y46	CAM-NAN	-2.34	1.35	1.39
2	D	400	Y46	CAU-NAN	2.17	1.51	1.47
2	B	400	Y46	CAU-NAN	2.14	1.51	1.47
2	C	400	Y46	CAU-NAN	2.09	1.50	1.47

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	400	Y46	OAQ-SAH-OAC	-7.03	110.98	119.52
2	D	400	Y46	OAQ-SAH-OAC	-6.80	111.27	119.52
2	B	400	Y46	OAQ-SAH-OAC	-6.41	111.74	119.52
2	D	400	Y46	CAE-NAD-CAM	-5.93	119.56	127.34
2	C	400	Y46	CAE-NAD-CAM	-5.64	119.95	127.34
2	C	400	Y46	OAQ-SAH-OAC	-5.40	112.96	119.52
2	B	400	Y46	CAE-NAD-CAM	-5.17	120.56	127.34
2	A	400	Y46	CAE-NAD-CAM	-4.98	120.81	127.34
2	B	400	Y46	CAI-SAH-NAG	4.81	113.73	106.06
2	A	400	Y46	NAD-CAM-NAN	3.55	121.82	115.45
2	D	400	Y46	NAD-CAM-NAN	3.51	121.76	115.45
2	D	400	Y46	CAF-CAO-NAP	-3.46	116.58	123.34
2	D	400	Y46	NAP-CAO-NAN	3.44	123.02	116.29
2	C	400	Y46	NAD-CAM-NAN	3.40	121.56	115.45
2	B	400	Y46	NAD-CAM-NAN	3.07	120.96	115.45
2	C	400	Y46	NAP-CAO-NAN	2.83	121.83	116.29
2	B	400	Y46	CAF-CAO-NAP	-2.75	117.97	123.34
2	C	400	Y46	CAF-CAO-NAP	-2.52	118.43	123.34
2	B	400	Y46	NAP-CAO-NAN	2.45	121.08	116.29
2	B	400	Y46	CAU-NAN-CAM	2.44	119.84	117.05
2	D	400	Y46	CAV-CAU-NAN	2.43	117.16	113.09
2	A	400	Y46	CAU-NAN-CAM	2.42	119.82	117.05
2	D	400	Y46	OAC-SAH-CAI	2.39	110.99	107.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	400	Y46	CAF-CAO-NAP	-2.29	118.86	123.34
2	C	400	Y46	CAF-CAE-NAD	2.27	118.37	112.13
2	D	400	Y46	CAF-CAE-NAD	2.21	118.21	112.13
2	A	400	Y46	NAP-CAO-NAN	2.20	120.58	116.29
2	B	400	Y46	CAF-CAE-NAD	2.03	117.70	112.13
2	B	400	Y46	CAV-CAU-NAN	2.00	116.44	113.09

There are no chirality outliers.

There are no torsion outliers.

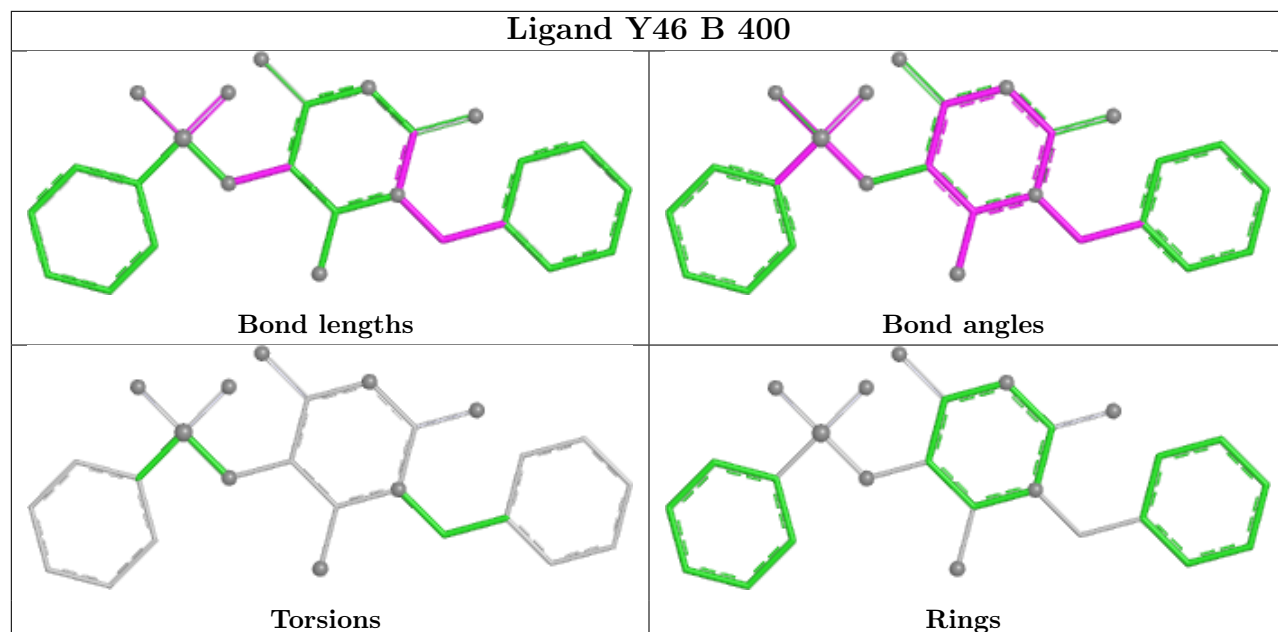
There are no ring outliers.

2 monomers are involved in 4 short contacts:

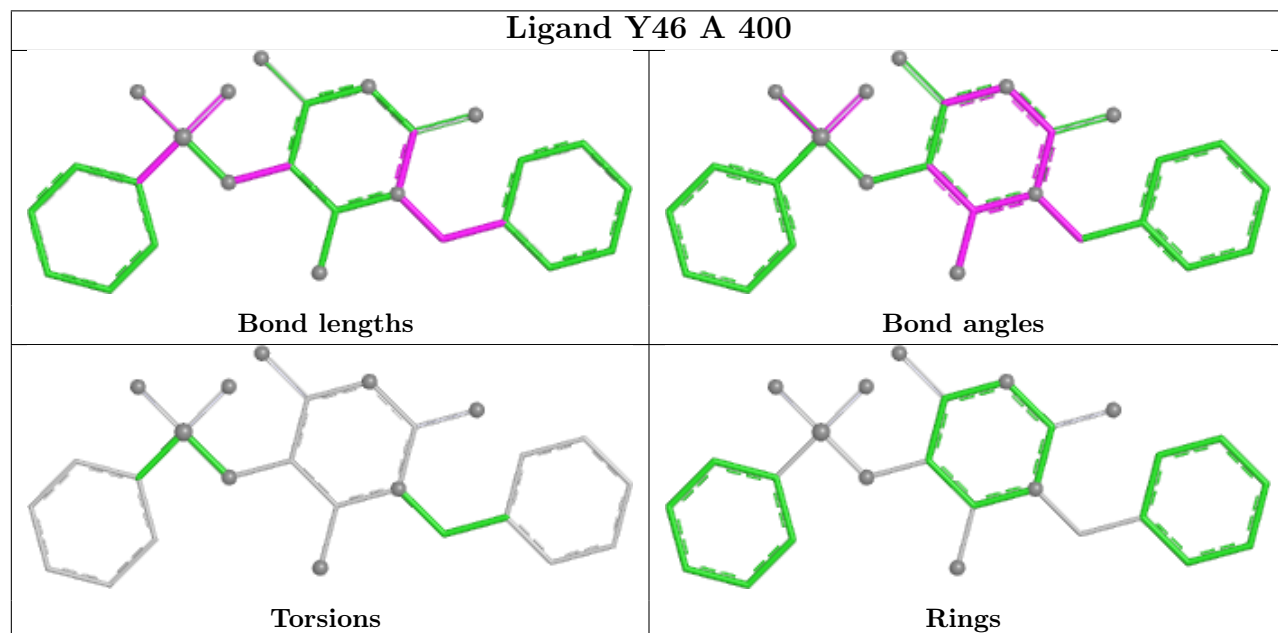
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	400	Y46	2	0
2	D	400	Y46	2	0

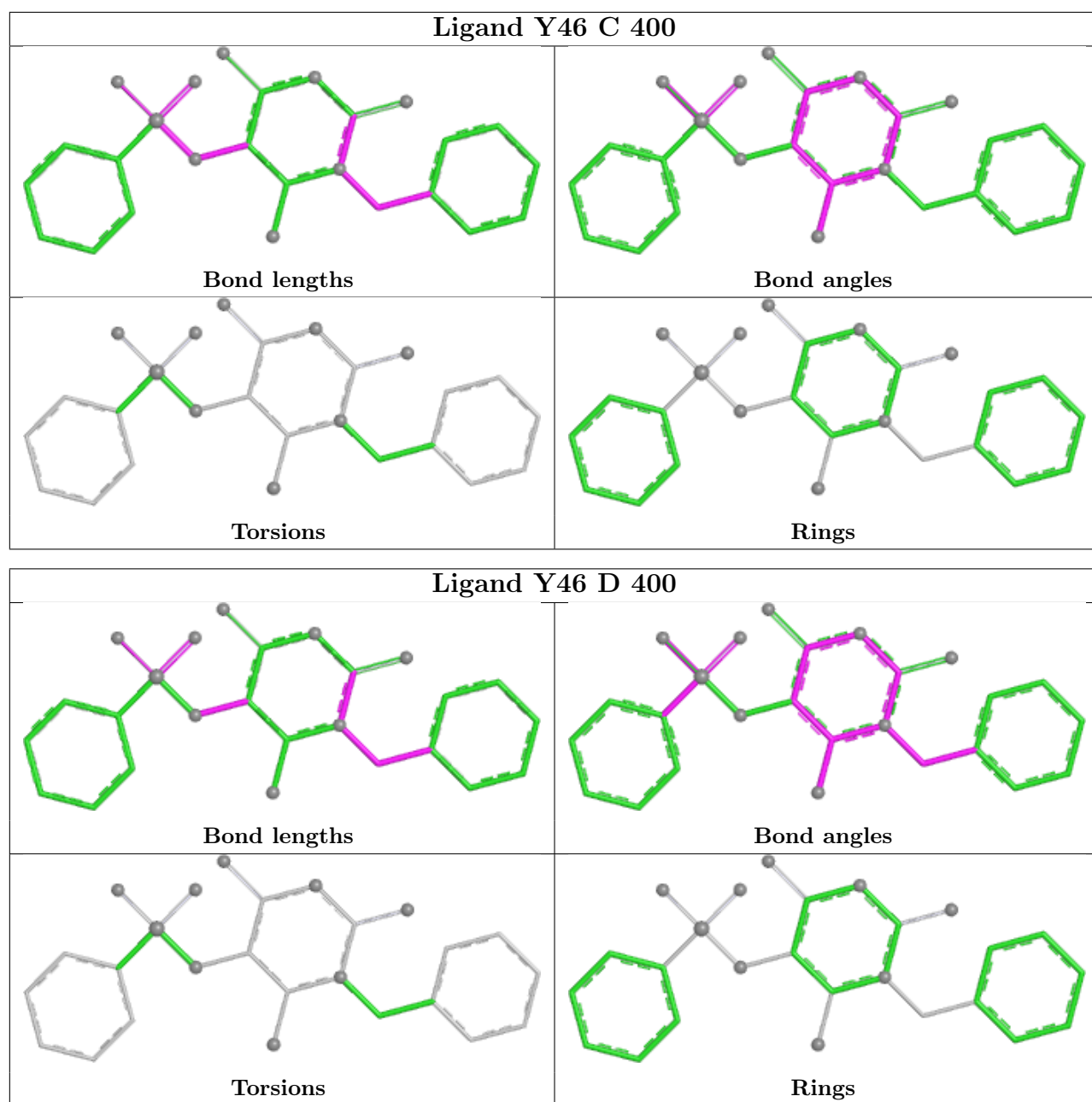
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 Y46 B 400



Ligand Y46 A 400





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%)	0.87	28 (9%) 14 16	16, 31, 53, 79	2 (0%)
1	B	293/303 (96%)	1.05	37 (12%) 8 9	18, 34, 49, 62	2 (0%)
1	C	292/303 (96%)	1.00	39 (13%) 7 8	13, 33, 57, 90	2 (0%)
1	D	293/303 (96%)	1.09	36 (12%) 8 9	14, 35, 53, 83	2 (0%)
All	All	1175/1212 (96%)	1.00	140 (11%) 9 10	13, 33, 53, 90	8 (0%)

All (140) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	22	ALA	6.1
1	A	-6	HIS	5.3
1	C	189	LEU	5.3
1	C	88	LEU	5.2
1	C	11	GLY	5.1
1	D	23	ILE	4.6
1	C	23	ILE	4.4
1	A	11	GLY	4.2
1	A	-5	HIS	4.1
1	C	22	ALA	4.0
1	D	65	GLN	3.9
1	D	11	GLY	3.9
1	C	191	PRO	3.9
1	D	128	ARG	3.7
1	B	24	SER	3.6
1	A	88	LEU	3.6
1	B	65	GLN	3.5
1	D	84	SER	3.5
1	B	22	ALA	3.5
1	C	193	PRO	3.4
1	D	195	GLY	3.4

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Mol	Chain	Res	Type	RSRZ
1	C	190	LYS	3.4
1	C	287	LEU	3.4
1	C	87	GLY	3.3
1	B	1	MET	3.3
1	B	64	GLN	3.2
1	A	275	LEU	3.2
1	C	275	LEU	3.2
1	D	81	VAL	3.2
1	A	192	SER	3.2
1	B	12	SER	3.2
1	B	23	ILE	3.1
1	A	287	LEU	3.1
1	C	194	ARG	3.1
1	D	194	ARG	3.1
1	A	193	PRO	3.0
1	D	225	ASP	2.9
1	A	288	LEU	2.9
1	D	141	ASP	2.9
1	A	195	GLY	2.9
1	C	226	THR	2.9
1	A	15	ARG	2.9
1	A	128	ARG	2.9
1	A	194	ARG	2.9
1	D	21	LEU	2.9
1	A	2	LYS	2.8
1	D	197	LEU	2.8
1	A	289	THR	2.8
1	B	128	ARG	2.8
1	C	128	ARG	2.8
1	B	13	GLY	2.8
1	A	85	PRO	2.8
1	B	194	ARG	2.7
1	B	232	LEU	2.7
1	B	11	GLY	2.7
1	D	187	ARG	2.7
1	C	85	PRO	2.7
1	D	85	PRO	2.7
1	B	153	GLY	2.6
1	D	276	ALA	2.6
1	C	192	SER	2.6
1	A	191	PRO	2.6
1	B	193	PRO	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	189	LEU	2.6
1	B	27	LEU	2.6
1	D	16	LEU	2.6
1	C	14	THR	2.6
1	B	61	PRO	2.6
1	D	164	LEU	2.6
1	C	15	ARG	2.6
1	D	1	MET	2.6
1	C	227	GLY	2.6
1	B	287	LEU	2.6
1	A	127	GLN	2.6
1	C	225	ASP	2.5
1	C	94	ILE	2.5
1	B	130	THR	2.5
1	C	289	THR	2.5
1	D	191	PRO	2.5
1	C	2	LYS	2.5
1	B	140	LEU	2.5
1	C	21	LEU	2.5
1	A	13	GLY	2.4
1	C	197	LEU	2.4
1	C	124	SER	2.4
1	B	276	ALA	2.4
1	B	289	THR	2.4
1	C	12	SER	2.4
1	D	275	LEU	2.4
1	D	101	ASN	2.4
1	C	112	LEU	2.4
1	C	224	LEU	2.4
1	D	88	LEU	2.4
1	B	215	GLU	2.4
1	D	83	PRO	2.4
1	B	148	VAL	2.4
1	B	60	THR	2.3
1	B	101	ASN	2.3
1	B	233	LEU	2.3
1	B	266	ALA	2.3
1	C	272	ALA	2.3
1	C	84	SER	2.3
1	B	275	LEU	2.3
1	D	2	LYS	2.3
1	B	265	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	195	GLY	2.3
1	D	15	ARG	2.3
1	B	21	LEU	2.3
1	D	144	ARG	2.2
1	A	124	SER	2.2
1	C	180	GLN	2.2
1	C	267	GLN	2.2
1	A	-9	HIS	2.2
1	B	66	LEU	2.2
1	A	153	GLY	2.2
1	D	193	PRO	2.2
1	C	276	ALA	2.2
1	D	62	ARG	2.2
1	D	189	LEU	2.2
1	B	117	ASP	2.2
1	C	213	SER	2.2
1	D	89	ALA	2.2
1	C	152	GLN	2.2
1	D	274	PRO	2.2
1	A	196	GLU	2.1
1	C	270	LYS	2.1
1	A	233	LEU	2.1
1	A	190	LYS	2.1
1	B	288	LEU	2.1
1	D	117	ASP	2.1
1	B	31	TYR	2.0
1	D	87	GLY	2.0
1	C	101	ASN	2.0
1	D	282	GLN	2.0
1	A	207	LEU	2.0
1	B	62	ARG	2.0
1	B	156	ALA	2.0
1	A	97	SER	2.0
1	D	12	SER	2.0
1	C	286	ARG	2.0

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

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

6.3 Carbohydrates

There are no oligosaccharides in this entry.

6.4 Ligands

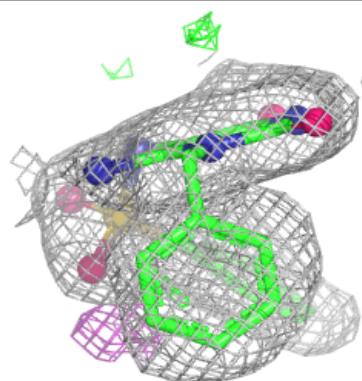
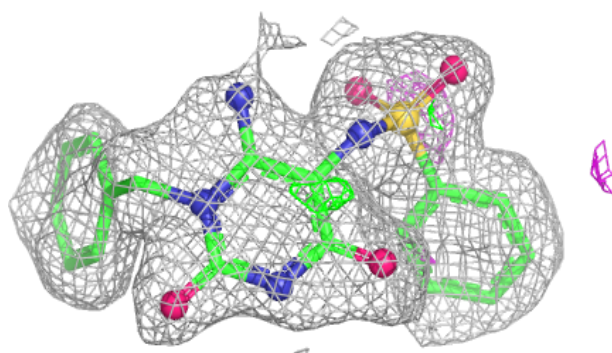
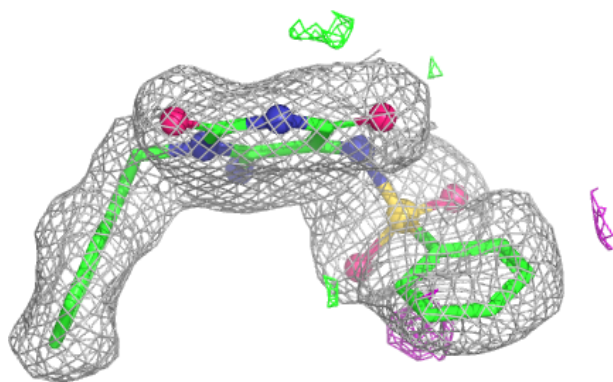
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	Y46	B	400	26/26	0.94	0.09	30,31,34,34	0
2	Y46	A	400	26/26	0.95	0.08	25,27,34,36	0
2	Y46	D	400	26/26	0.95	0.09	31,34,37,39	0
2	Y46	C	400	26/26	0.96	0.07	24,25,30,31	0
3	CL	C	1294	1/1	0.98	0.04	27,27,27,27	0
3	CL	B	1294	1/1	0.99	0.03	29,29,29,29	0
3	CL	A	1294	1/1	0.99	0.04	27,27,27,27	0
3	CL	D	1294	1/1	0.99	0.04	34,34,34,34	0

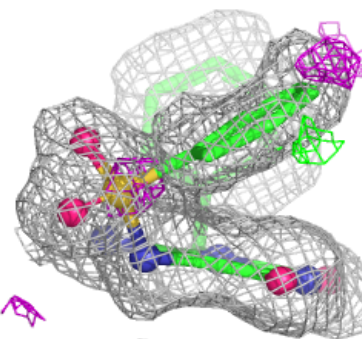
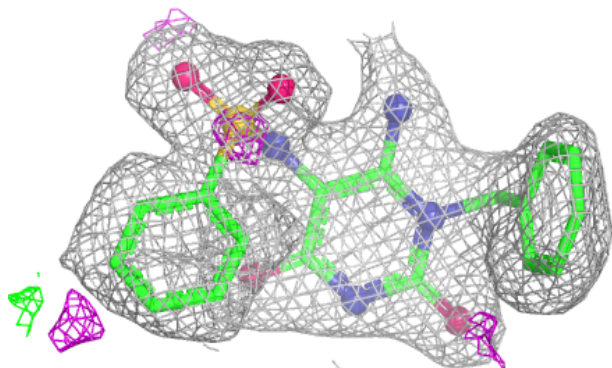
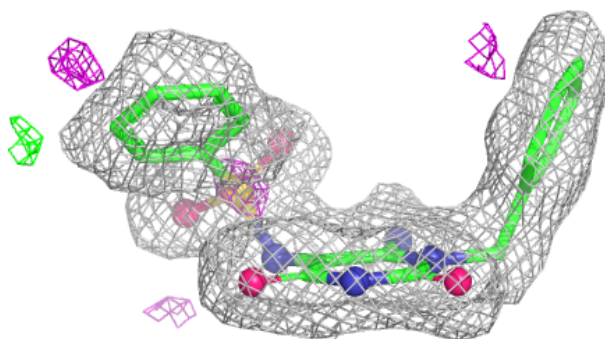
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around Y46 B 400:

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

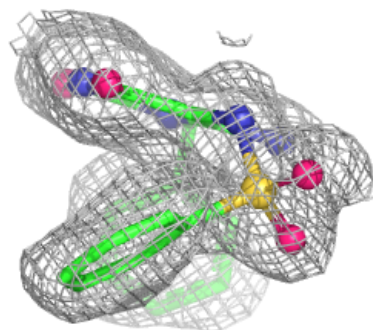
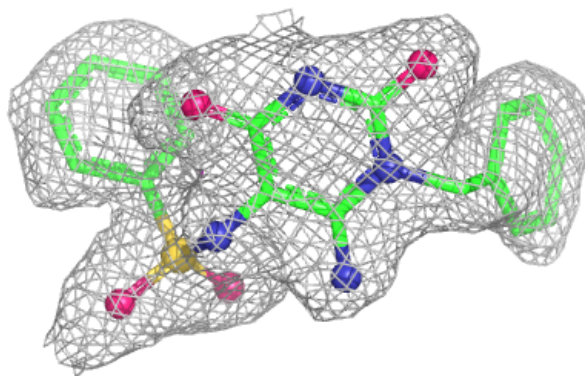
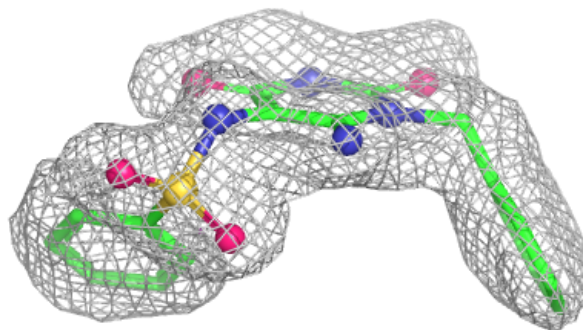
**Electron density around Y46 A 400:**

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

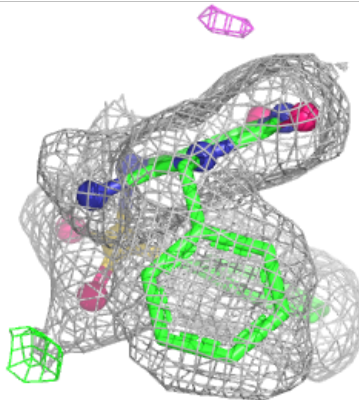
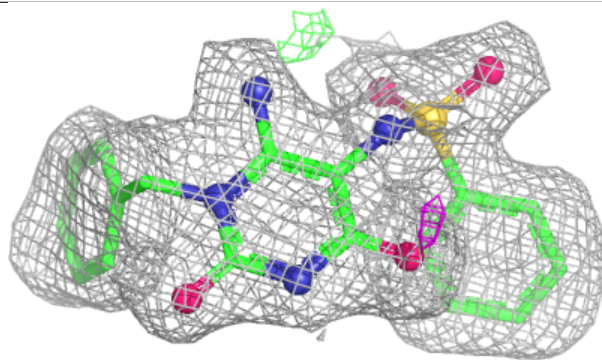
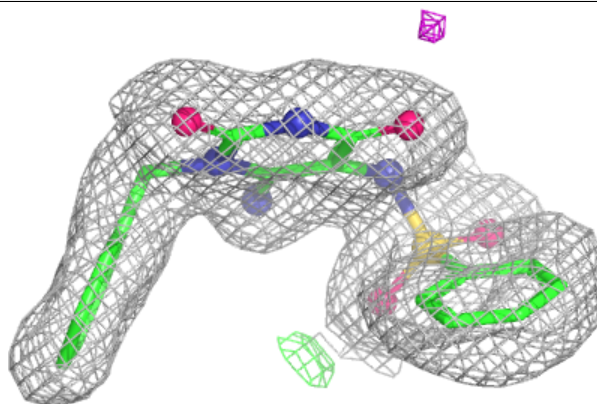


Electron density around Y46 D 400:

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

**Electron density around Y46 C 400:**

$2mF_o-DF_c$ (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.