



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 20, 2022 – 03:29 pm GMT

PDB ID : 8A9G
Title : Binary complex of 14-3-3 zeta Glucocorticoid Receptor (GR) pT524 peptide stabilised by (R)-para chloropyrrolidone1
Authors : Munier, C.C.; Edman, K.; Perry, M.W.D.; Ottmann, C.
Deposited on : 2022-06-28
Resolution : 1.96 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 : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.31.3
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) : 2.31.3

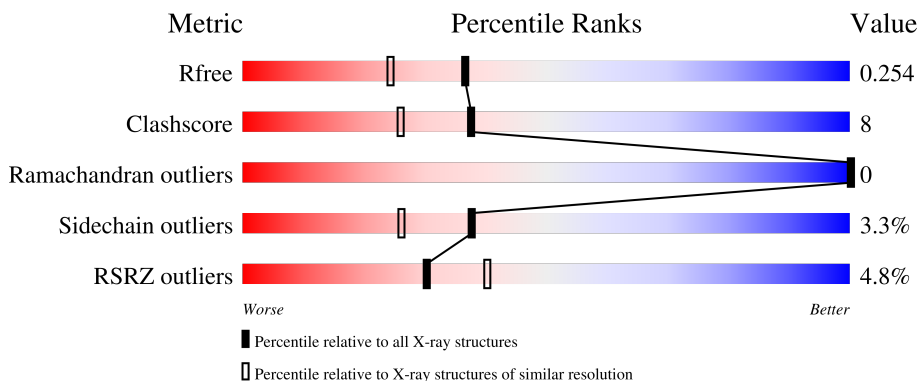
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.96 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	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (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	235	
1	B	235	
2	C	13	
2	D	13	

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 4269 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 14-3-3 protein zeta/delta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	230	1851	1158	310	373	10	0	1	0
1	B	230	1843	1154	309	370	10	0	0	0

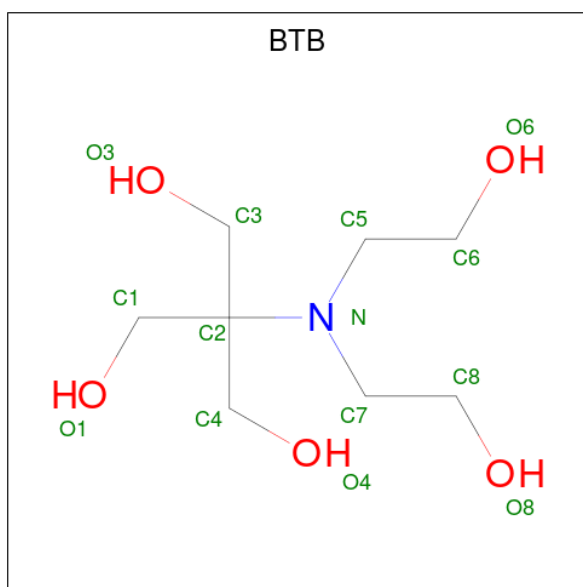
There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	expression tag	UNP P63104
A	-3	ALA	-	expression tag	UNP P63104
A	-2	MET	-	expression tag	UNP P63104
A	-1	GLY	-	expression tag	UNP P63104
A	0	SER	-	expression tag	UNP P63104
B	-4	GLY	-	expression tag	UNP P63104
B	-3	ALA	-	expression tag	UNP P63104
B	-2	MET	-	expression tag	UNP P63104
B	-1	GLY	-	expression tag	UNP P63104
B	0	SER	-	expression tag	UNP P63104

- Molecule 2 is a protein called Glucocorticoid receptor.

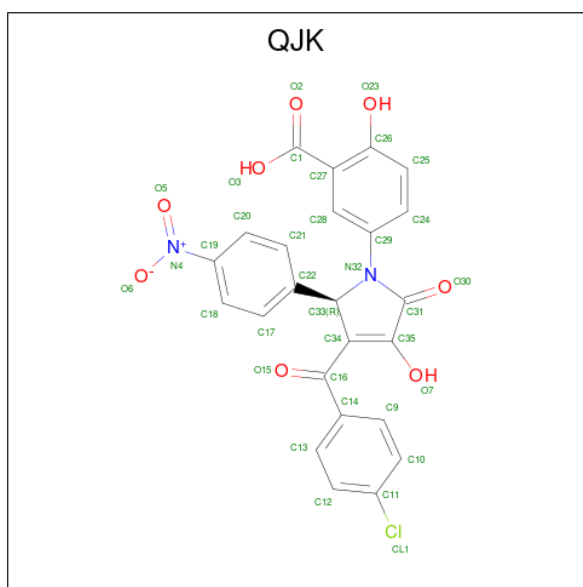
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	C	12	92	58	13	20	1	0	0	0
2	D	13	101	64	15	21	1	0	0	0

- Molecule 3 is 2-[BIS-(2-HYDROXY-ETHYL)-AMINO]-2-HYDROXYMETHYL-PROPAN E-1,3-DIOL (three-letter code: BTB) (formula: C₈H₁₉NO₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is 5-[(2 {R})-3-(4-chlorophenyl)carbonyl-2-(4-nitrophenyl)-4-oxidanyl-5-oxidanylidene-2 {H}-pyrrol-1-yl]-2-oxidanyl-benzoic acid (three-letter code: QJK) (formula: C₂₄H₁₅ClN₂O₈) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	Cl	N	O	0	0
			35	24	1	2	8		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
4	B	1	Total	C	Cl	N	O	0	0
			35	24	1	2	8		

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Mg	0	0
			1	1		

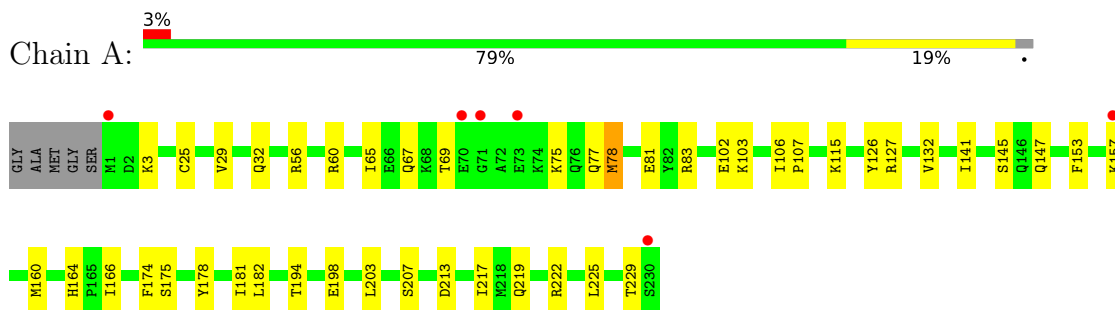
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	123	Total	O	0	0
			123	123		
6	B	135	Total	O	0	0
			135	135		
6	C	14	Total	O	0	0
			14	14		
6	D	11	Total	O	0	0
			11	11		

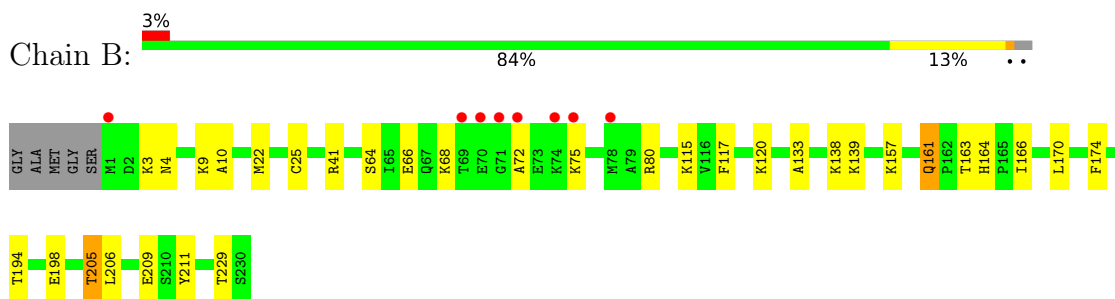
3 Residue-property plots [i](#)

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.

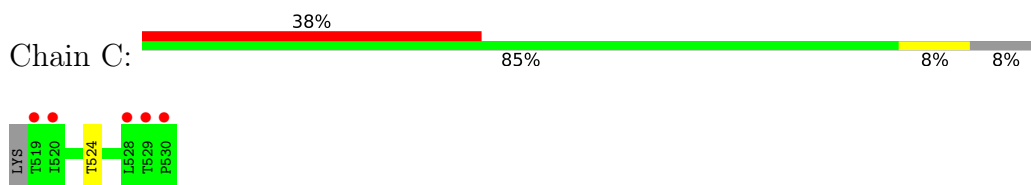
- Molecule 1: 14-3-3 protein zeta/delta



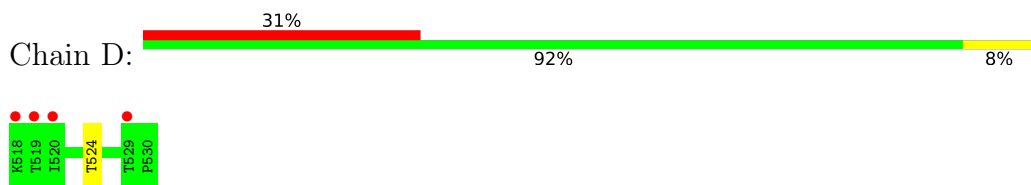
- Molecule 1: 14-3-3 protein zeta/delta



- Molecule 2: Glucocorticoid receptor



- Molecule 2: Glucocorticoid receptor



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	57.35Å 78.07Å 122.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	65.85 – 1.96 65.85 – 1.96	Depositor EDS
% Data completeness (in resolution range)	78.7 (65.85-1.96) 78.8 (65.85-1.96)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.74 (at 1.95Å)	Xtrriage
Refinement program	BUSTER 2.11.8 (24-FEB-2021)	Depositor
R, R_{free}	0.250 , 0.261 0.236 , 0.254	Depositor DCC
R_{free} test set	1587 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	39.3	Xtrriage
Anisotropy	0.028	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 57.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4269	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 27.95 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.0096e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: TPO, QJK, BTB, MG

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	1.38	0/1876	0.80	0/2521
1	B	1.39	0/1868	0.80	0/2510
2	C	0.37	0/82	0.51	0/112
2	D	0.42	0/91	0.56	0/123
All	All	1.36	0/3917	0.79	0/5266

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	1851	0	1836	36	0
1	B	1843	0	1833	24	0
2	C	92	0	94	0	0
2	D	101	0	107	0	0
3	A	14	0	19	1	0
3	B	14	0	19	0	0
4	A	35	0	0	0	0
4	B	35	0	0	0	0
5	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	123	0	0	0	0
6	B	135	0	0	0	0
6	C	14	0	0	0	0
6	D	11	0	0	0	0
All	All	4269	0	3908	61	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 8.

All (61) 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:64:SER:O	1:B:68:LYS:HD3	1.66	0.94
1:B:161:GLN:HG2	1:B:205:THR:CG2	1.99	0.92
1:B:161:GLN:HG2	1:B:205:THR:HG23	1.51	0.90
1:B:163:THR:OG1	1:B:205:THR:HG22	1.73	0.89
1:A:56:ARG:O	1:A:60:ARG:HG3	1.71	0.88
1:A:153:PHE:CZ	1:A:157:LYS:HE3	2.14	0.83
1:A:225:LEU:O	1:A:229:THR:HG23	1.79	0.81
1:A:69:THR:HG21	1:A:75:LYS:HE2	1.68	0.76
1:A:65:ILE:O	1:A:69:THR:HG23	1.87	0.74
1:A:77:GLN:O	1:A:81:GLU:HG3	1.87	0.73
1:A:106:ILE:O	1:A:115:LYS:NZ	2.22	0.72
1:A:164:HIS:HD2	1:A:166:ILE:H	1.38	0.72
1:B:163:THR:OG1	1:B:205:THR:CG2	2.38	0.71
1:A:153:PHE:CE2	1:A:157:LYS:HE3	2.28	0.69
1:A:78:MET:HA	1:A:78:MET:CE	2.24	0.68
1:A:102:GLU:HG3	1:A:103:LYS:HD3	1.76	0.67
1:B:163:THR:HG1	1:B:205:THR:HG22	1.58	0.66
1:B:194:THR:O	1:B:198:GLU:HG3	1.96	0.64
1:A:106:ILE:N	1:A:107:PRO:HD2	2.15	0.62
1:A:194:THR:O	1:A:198:GLU:HG3	2.01	0.60
1:A:178:TYR:CD1	1:A:182:LEU:HD12	2.36	0.60
1:A:153:PHE:CZ	1:A:157:LYS:CE	2.83	0.60
1:B:161:GLN:HG2	1:B:205:THR:HG21	1.82	0.55
1:B:66:GLU:OE2	1:B:80:ARG:HG3	2.07	0.54
1:A:160:MET:CE	1:A:164:HIS:CG	2.91	0.53
1:B:164:HIS:HD2	1:B:166:ILE:H	1.57	0.52
1:A:213:ASP:O	1:A:217:ILE:HG13	2.09	0.52
1:A:25:CYS:O	1:A:29:VAL:HG23	2.10	0.51
1:B:9:LYS:HD3	1:B:25:CYS:SG	2.51	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:MET:HA	1:A:78:MET:HE3	1.93	0.51
1:B:209:GLU:O	1:B:209:GLU:HG2	2.11	0.51
1:A:3:LYS:HB2	1:A:32:GLN:OE1	2.11	0.50
1:A:3:LYS:HG3	1:A:29:VAL:HG13	1.95	0.48
1:B:72:ALA:HB1	1:B:75:LYS:HG3	1.96	0.48
1:A:178:TYR:HD1	1:A:182:LEU:HD12	1.78	0.48
1:B:133:ALA:HB3	1:B:138:LYS:HG3	1.95	0.48
1:A:83:ARG:HH11	1:A:83:ARG:HG2	1.79	0.47
3:A:301:BTB:H41	3:A:301:BTB:H71	1.52	0.47
1:A:83:ARG:HG2	1:A:83:ARG:NH1	2.30	0.47
1:B:161:GLN:CG	1:B:205:THR:HG21	2.45	0.46
1:A:102:GLU:CG	1:A:103:LYS:HD3	2.43	0.46
1:B:120:LYS:HD2	1:B:170:LEU:HA	1.97	0.46
1:A:160:MET:CE	1:A:164:HIS:CD2	2.98	0.46
1:A:160:MET:HE3	1:A:164:HIS:CG	2.51	0.46
1:B:139:LYS:N	1:B:139:LYS:HD3	2.31	0.45
1:B:41:ARG:HD2	1:B:117:PHE:CD1	2.52	0.45
1:A:106:ILE:N	1:A:107:PRO:CD	2.80	0.45
1:B:115:LYS:HB2	1:B:115:LYS:HE2	1.70	0.45
1:B:163:THR:HB	1:B:206:LEU:HD23	1.99	0.44
1:B:157:LYS:NZ	1:B:198:GLU:OE2	2.51	0.44
1:A:78:MET:HA	1:A:78:MET:HE2	1.97	0.44
1:B:206:LEU:HD13	1:B:211:TYR:HB2	1.99	0.43
1:A:203:LEU:HD23	1:A:203:LEU:HA	1.76	0.42
1:B:10:ALA:O	1:B:22:MET:HG3	2.20	0.42
1:A:132:VAL:O	1:A:132:VAL:HG23	2.20	0.42
1:A:103:LYS:HA	1:A:103:LYS:HD2	1.87	0.41
1:A:164:HIS:CD2	1:A:166:ILE:H	2.26	0.41
1:A:126:TYR:HB3	1:A:145:SER:HB2	2.03	0.41
1:A:127:ARG:HG3	1:A:181:ILE:HG13	2.03	0.41
1:B:161:GLN:CG	1:B:205:THR:CG2	2.87	0.40
1:A:219:GLN:O	1:A:219:GLN:HG3	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	229/235 (97%)	225 (98%)	4 (2%)	0	100	100
1	B	228/235 (97%)	222 (97%)	6 (3%)	0	100	100
2	C	9/13 (69%)	9 (100%)	0	0	100	100
2	D	10/13 (77%)	9 (90%)	1 (10%)	0	100	100
All	All	476/496 (96%)	465 (98%)	11 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	201/202 (100%)	193 (96%)	8 (4%)	31	19
1	B	200/202 (99%)	194 (97%)	6 (3%)	41	30
2	C	10/11 (91%)	10 (100%)	0	100	100
2	D	11/11 (100%)	11 (100%)	0	100	100
All	All	422/426 (99%)	408 (97%)	14 (3%)	38	26

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

Mol	Chain	Res	Type
1	A	67	GLN
1	A	78	MET
1	A	141	ILE
1	A	147	GLN
1	A	174	PHE
1	A	175	SER
1	A	207	SER
1	A	222	ARG

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Mol	Chain	Res	Type
1	B	3	LYS
1	B	4	ASN
1	B	161	GLN
1	B	174	PHE
1	B	205	THR
1	B	229	THR

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

Mol	Chain	Res	Type
1	A	164	HIS
1	B	15	GLN
1	B	67	GLN
1	B	147	GLN
1	B	164	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	TPO	C	524	2	8,10,11	1.22	1 (12%)	10,14,16	1.26	2 (20%)
2	TPO	D	524	2	8,10,11	0.96	1 (12%)	10,14,16	1.51	4 (40%)

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	TPO	C	524	2	-	2/9/11/13	-
2	TPO	D	524	2	-	3/9/11/13	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	524	TPO	P-OG1	-3.05	1.53	1.59
2	D	524	TPO	P-OG1	-2.38	1.54	1.59

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	524	TPO	O3P-P-OG1	2.47	117.05	105.99
2	D	524	TPO	P-OG1-CB	-2.11	116.84	123.21
2	D	524	TPO	O-C-CA	-2.10	119.28	124.78
2	C	524	TPO	OG1-P-O1P	2.05	117.29	109.39
2	C	524	TPO	O-C-CA	-2.04	119.42	124.78
2	D	524	TPO	OG1-P-O1P	-2.00	101.66	109.39

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	524	TPO	CB-OG1-P-O1P
2	D	524	TPO	N-CA-CB-CG2
2	C	524	TPO	CB-OG1-P-O2P
2	C	524	TPO	O-C-CA-CB
2	D	524	TPO	O-C-CA-CB

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 1 is 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
3	BTB	B	301	-	13,13,13	1.28	1 (7%)	7,16,16	1.27	0
3	BTB	A	301	-	13,13,13	1.27	1 (7%)	7,16,16	1.23	1 (14%)
4	QJK	B	303	5	37,38,38	0.76	2 (5%)	53,56,56	2.18	9 (16%)
4	QJK	A	302	-	37,38,38	0.74	0	53,56,56	2.12	7 (13%)

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	BTB	B	301	-	-	4/21/21/21	-
3	BTB	A	301	-	-	9/21/21/21	-
4	QJK	B	303	5	-	3/22/44/44	0/4/4/4
4	QJK	A	302	-	-	8/22/44/44	0/4/4/4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	301	BTB	C2-N	3.61	1.55	1.48
3	B	301	BTB	C2-N	3.49	1.55	1.48
4	B	303	QJK	C35-C34	2.27	1.36	1.34
4	B	303	QJK	C31-N32	-2.03	1.35	1.37

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	303	QJK	C33-C34-C35	-8.37	105.18	109.90
4	A	302	QJK	C33-C34-C35	-7.72	105.55	109.90
4	A	302	QJK	C34-C33-N32	7.55	107.28	102.02
4	B	303	QJK	C34-C33-N32	7.41	107.19	102.02
4	A	302	QJK	C35-C31-N32	5.81	110.17	105.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	303	QJK	C35-C31-N32	5.64	110.04	105.86
4	A	302	QJK	C33-N32-C31	-5.36	107.16	112.11
4	B	303	QJK	C33-N32-C31	-5.09	107.40	112.11
4	B	303	QJK	C29-N32-C33	4.63	126.87	121.76
4	A	302	QJK	C29-N32-C33	3.63	125.77	121.76
4	A	302	QJK	O7-C35-C34	-3.04	125.37	131.71
4	B	303	QJK	O7-C35-C34	-2.89	125.68	131.71
4	B	303	QJK	C14-C16-C34	2.29	125.04	120.84
3	A	301	BTB	O4-C4-C2	2.23	117.54	111.44
4	B	303	QJK	O3-C1-C27	2.23	121.72	115.31
4	B	303	QJK	O3-C1-O2	-2.14	118.60	123.35
4	A	302	QJK	C28-C27-C26	2.13	120.75	118.70

There are no chirality outliers.

All (24) torsion outliers are listed below:

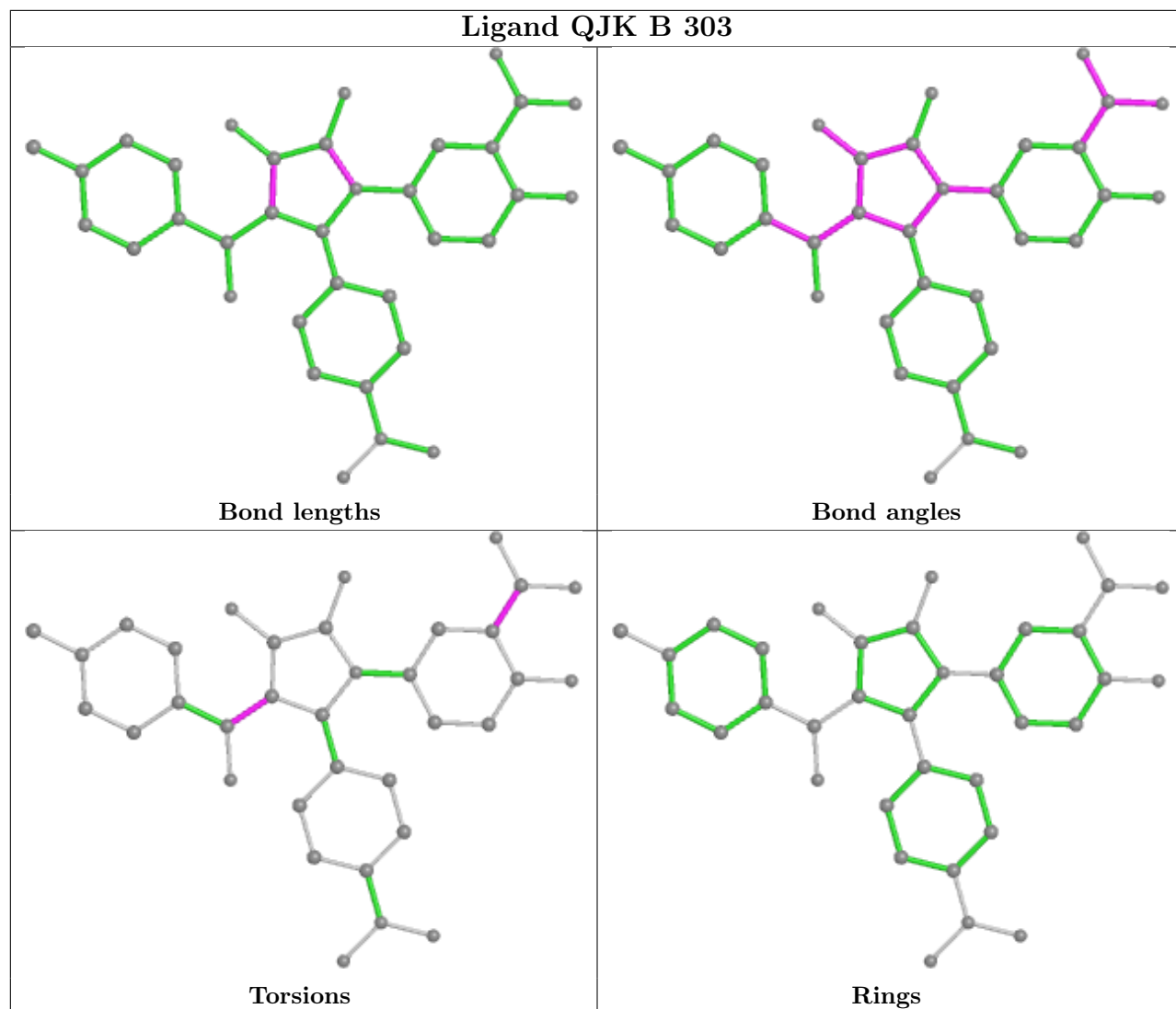
Mol	Chain	Res	Type	Atoms
3	A	301	BTB	C1-C2-C4-O4
3	A	301	BTB	C3-C2-C4-O4
3	A	301	BTB	N-C2-C4-O4
3	A	301	BTB	C1-C2-N-C5
3	A	301	BTB	C3-C2-N-C5
3	A	301	BTB	C4-C2-N-C5
3	A	301	BTB	C4-C2-N-C7
3	A	301	BTB	N-C5-C6-O6
3	B	301	BTB	O1-C1-C2-C4
3	B	301	BTB	N-C7-C8-O8
4	A	302	QJK	O3-C1-C27-C26
3	B	301	BTB	N-C5-C6-O6
4	A	302	QJK	O2-C1-C27-C26
4	A	302	QJK	C20-C19-N4-O5
4	A	302	QJK	O15-C16-C34-C35
4	A	302	QJK	C14-C16-C34-C35
4	B	303	QJK	C14-C16-C34-C35
4	A	302	QJK	C18-C19-N4-O5
4	A	302	QJK	C14-C16-C34-C33
3	B	301	BTB	O1-C1-C2-C3
3	A	301	BTB	C3-C2-N-C7
4	A	302	QJK	O15-C16-C34-C33
4	B	303	QJK	O15-C16-C34-C35
4	B	303	QJK	O3-C1-C27-C26

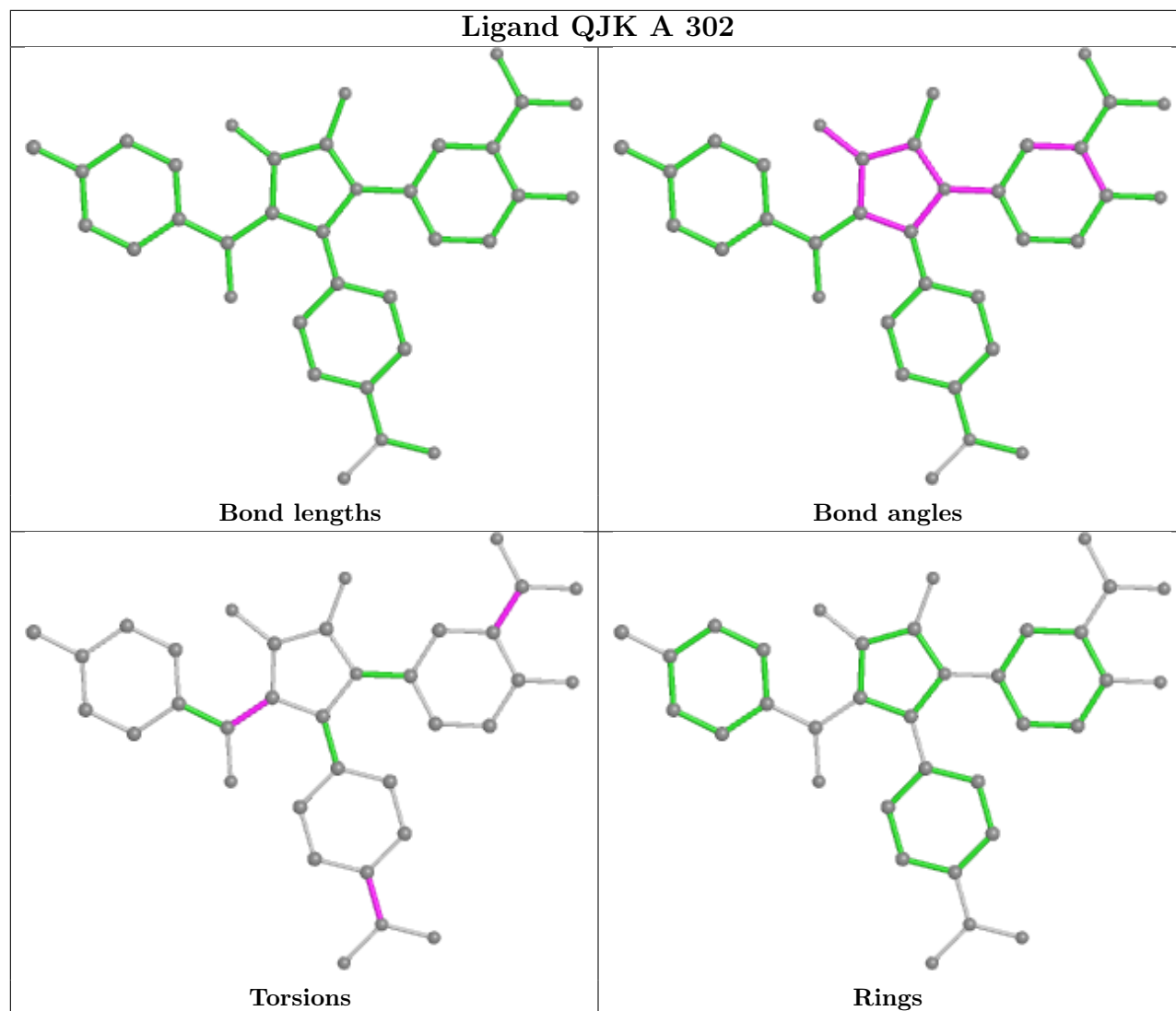
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	301	BTB	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	230/235 (97%)	0.27	6 (2%) 56 65	18, 42, 70, 82	0
1	B	230/235 (97%)	0.19	8 (3%) 44 53	17, 39, 76, 95	0
2	C	11/13 (84%)	2.26	5 (45%) 0 0	44, 61, 72, 76	0
2	D	12/13 (92%)	2.42	4 (33%) 0 0	41, 57, 79, 91	0
All	All	483/496 (97%)	0.33	23 (4%) 30 40	17, 41, 74, 95	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	519	THR	9.5
1	A	71	GLY	7.7
1	B	70	GLU	6.6
1	B	72	ALA	6.1
2	D	520	ILE	5.5
1	A	1	MET	5.1
2	C	529	THR	5.0
2	C	528	LEU	4.9
2	C	520	ILE	4.1
1	B	71	GLY	4.1
1	A	73	GLU	3.3
2	C	519	THR	3.1
2	D	518	LYS	3.0
2	C	530	PRO	2.9
1	B	75	LYS	2.9
1	B	1	MET	2.8
2	D	529	THR	2.7
1	A	70	GLU	2.4
1	B	74	LYS	2.3
1	B	78	MET	2.3
1	A	230	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	157	LYS	2.1
1	B	69	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
2	TPO	C	524	11/12	0.96	0.17	29,34,42,43	0
2	TPO	D	524	11/12	0.97	0.14	32,39,42,44	0

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

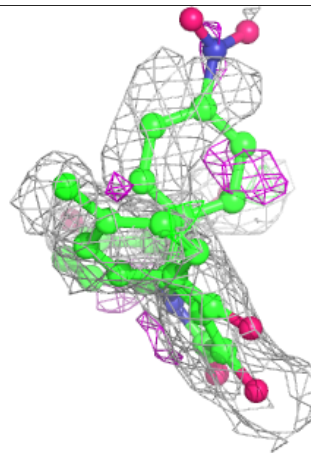
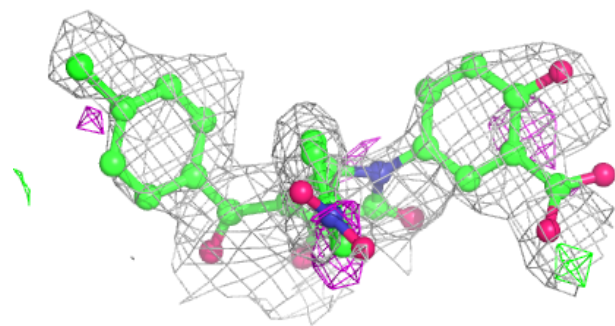
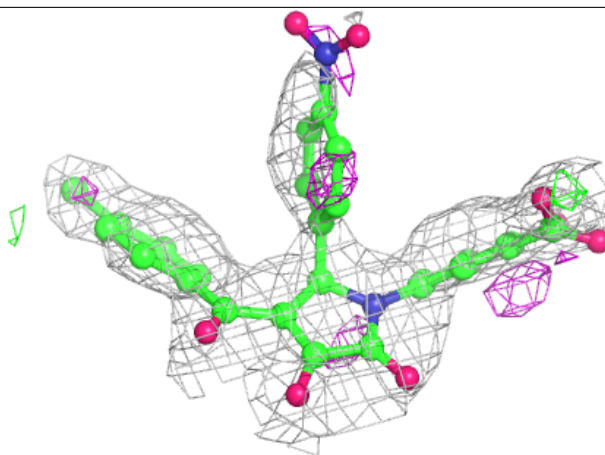
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

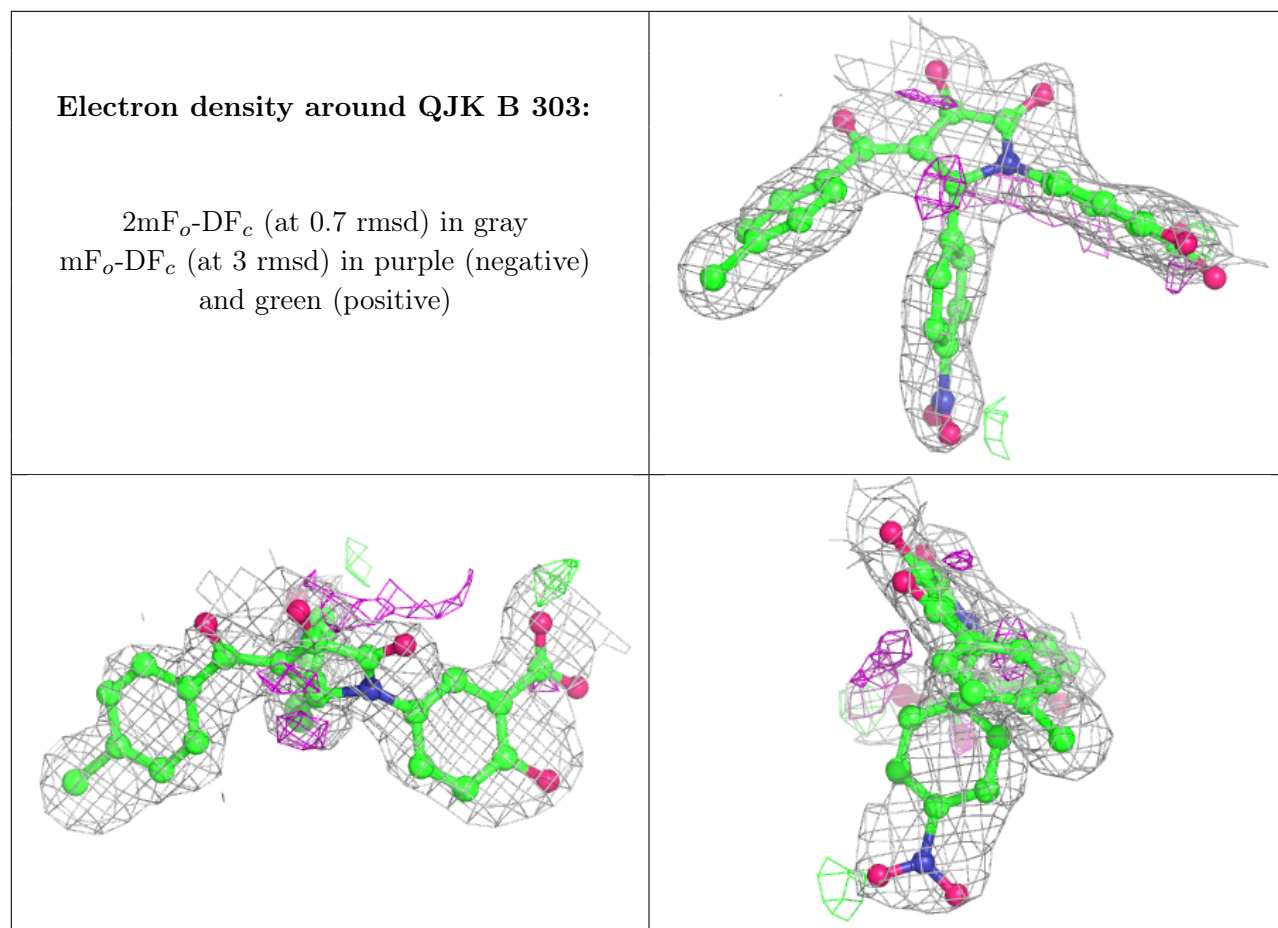
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	QJK	A	302	35/35	0.54	0.34	91,92,93,93	0
3	BTB	B	301	14/14	0.63	0.38	87,87,88,88	0
5	MG	B	302	1/1	0.63	0.09	88,88,88,88	0
3	BTB	A	301	14/14	0.64	0.39	73,73,74,74	0
4	QJK	B	303	35/35	0.74	0.23	78,78,79,79	0

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 QJK A 302:

$2mF_o-DF_c$ (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.