



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:17 PM UTC

PDB ID : 4R3A / pdb_00004r3a
Title : Erythrobacter litoralis EL346 blue-light activated histidine kinase
Authors : Tomchick, D.R.; Rivera-Cancel, G.; Gardner, K.H.
Deposited on : 2014-08-14
Resolution : 2.92 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

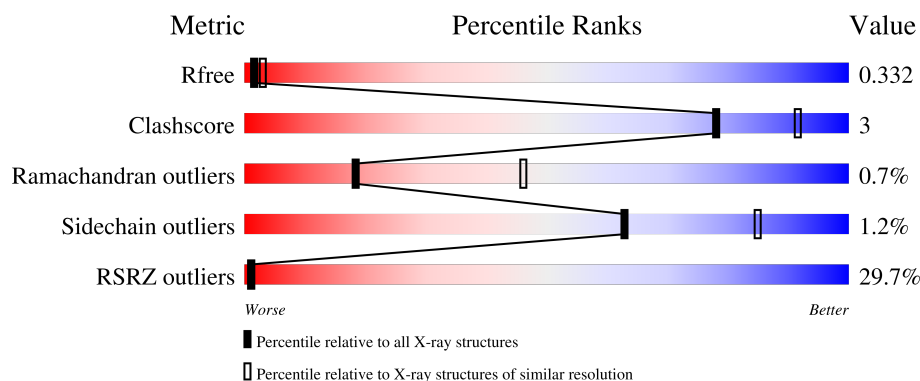
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.92 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	2995 (2.94-2.90)
Clashscore	190562	3213 (2.94-2.90)
Ramachandran outliers	187476	3128 (2.94-2.90)
Sidechain outliers	187428	3130 (2.94-2.90)
RSRZ outliers	180081	2995 (2.94-2.90)

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	352	 15% 81% 9% 10%
1	B	352	 35% 74% 5% 21%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9372 atoms, of which 4639 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 Blue-light-activated histidine kinase 2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	318	Total	C	H	N	O	S	0	0	0
			4903	1525	2437	449	481	11			
1	B	278	Total	C	H	N	O	S	0	0	0
			4285	1336	2136	386	416	11			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	GLY	-	expression tag	UNP Q2NB77
A	-4	GLU	-	expression tag	UNP Q2NB77
A	-3	PHE	-	expression tag	UNP Q2NB77
A	-2	LYS	-	expression tag	UNP Q2NB77
A	-1	GLY	-	expression tag	UNP Q2NB77
A	0	LEU	-	expression tag	UNP Q2NB77
B	-5	GLY	-	expression tag	UNP Q2NB77
B	-4	GLU	-	expression tag	UNP Q2NB77
B	-3	PHE	-	expression tag	UNP Q2NB77
B	-2	LYS	-	expression tag	UNP Q2NB77
B	-1	GLY	-	expression tag	UNP Q2NB77
B	0	LEU	-	expression tag	UNP Q2NB77

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

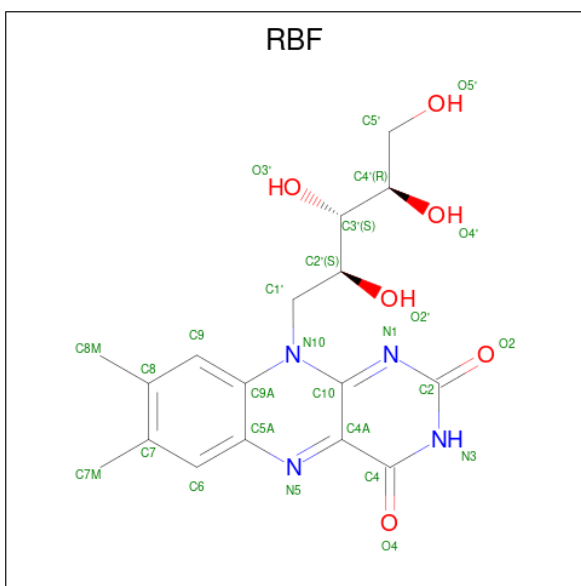
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		
2	B	1	Total	Mg	0	0
			1	1		

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total 44	C 10	H 13	N 6	O 12	P 3	0	0
3	B	1	Total 44	C 10	H 13	N 6	O 12	P 3	0	0

- Molecule 4 is RIBOFLAVIN (CCD ID: RBF) (formula: $C_{17}H_{20}N_4O_6$).

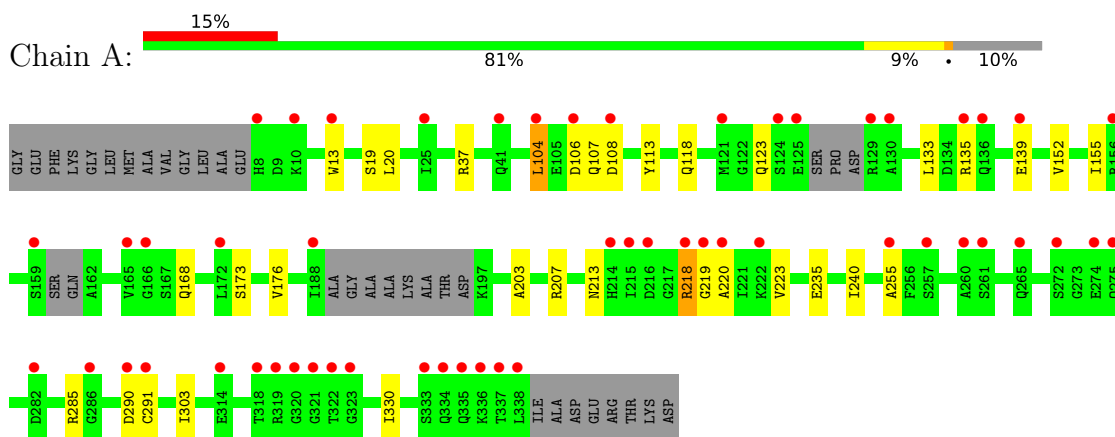


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 47	C 17	H 20	N 4	O 6	0	0
4	B	1	Total 47	C 17	H 20	N 4	O 6	0	0

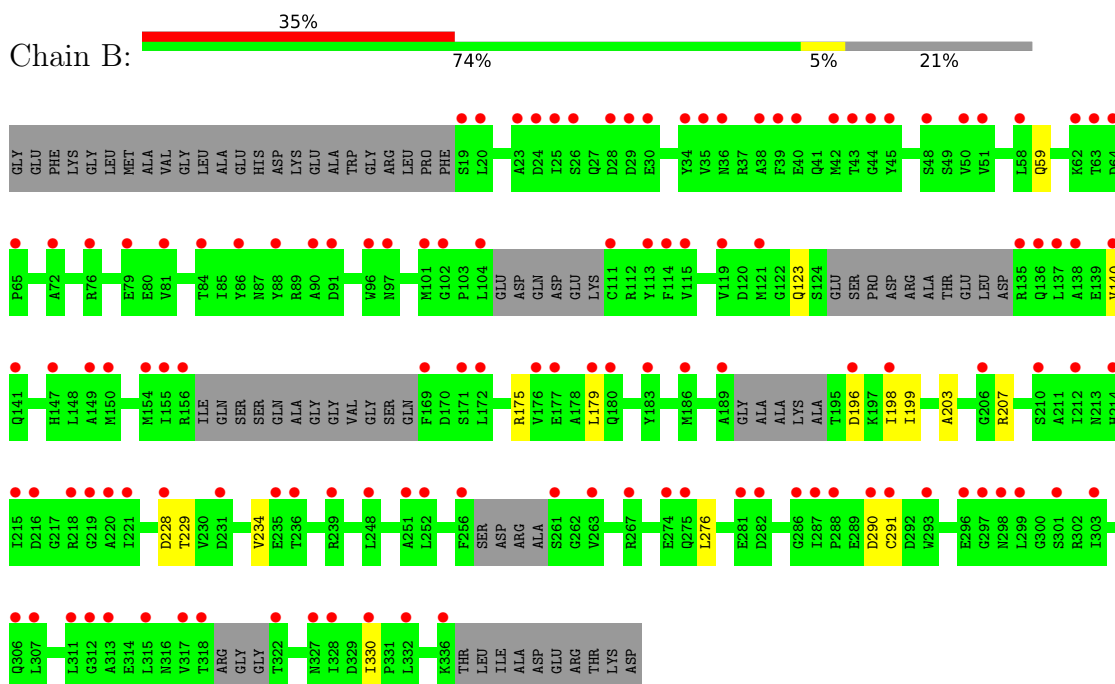
3 Residue-property plots [i](#)

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: Blue-light-activated histidine kinase 2



• Molecule 1: Blue-light-activated histidine kinase 2



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	98.88Å 98.88Å 422.25Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.60 – 2.92 46.60 – 2.92	Depositor EDS
% Data completeness (in resolution range)	94.7 (46.60-2.92) 94.8 (46.60-2.92)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.42 (at 2.91Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.290 , 0.330 0.294 , 0.332	Depositor DCC
R_{free} test set	1693 reflections (9.47%)	wwPDB-VP
Wilson B-factor (Å ²)	53.1	Xtriage
Anisotropy	0.073	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 65.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	9372	wwPDB-VP
Average B, all atoms (Å ²)	98.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.62% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, ANP, RBF

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.40	0/2497	0.91	4/3370 (0.1%)
1	B	0.34	0/2172	0.83	3/2931 (0.1%)
All	All	0.37	0/4669	0.87	7/6301 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	107	GLN	N-CA-C	-9.29	99.53	112.45
1	B	199	ILE	N-CA-C	8.17	116.92	107.77
1	B	140	VAL	N-CA-C	-6.66	107.38	113.71
1	A	104	LEU	N-CA-C	-6.32	101.13	110.48
1	A	133	LEU	N-CA-C	-6.22	104.04	111.69
1	A	168	GLN	N-CA-C	-5.69	105.83	112.89
1	B	228	ASP	N-CA-C	-5.66	103.06	110.53

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	2466	2437	2437	20	0
1	B	2149	2136	2134	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	31	13	13	1	0
3	B	31	13	13	1	0
4	A	27	20	20	0	0
4	B	27	20	20	1	0
All	All	4733	4639	4637	26	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 3.

All (26) 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:123:GLN:O	1:B:123:GLN:NE2	2.40	0.55
1:A:135:ARG:NH1	1:A:139:GLU:OE2	2.44	0.51
1:A:13:TRP:HB3	1:A:37:ARG:HG3	1.94	0.50
1:A:218:ARG:N	1:A:218:ARG:HD3	2.27	0.49
1:A:104:LEU:HG	1:A:152:VAL:HG11	1.94	0.48
1:A:37:ARG:CD	1:A:37:ARG:H	2.27	0.48
1:A:255:ALA:O	1:A:285:ARG:N	2.45	0.48
1:A:207:ARG:HG2	1:B:203:ALA:HB1	1.97	0.46
1:B:276:LEU:N	1:B:330:ILE:O	2.44	0.46
1:A:290:ASP:OD1	1:A:291:CYS:N	2.49	0.46
1:A:152:VAL:O	1:A:155:ILE:HG13	2.16	0.45
1:B:290:ASP:OD1	1:B:291:CYS:N	2.49	0.45
1:A:19:SER:C	1:A:20:LEU:HD12	2.42	0.45
1:B:59:GLN:OE1	4:B:402:RBF:O4'	2.35	0.45
3:A:401:ANP:O2G	3:A:401:ANP:O3A	2.33	0.45
1:A:203:ALA:HB1	1:B:207:ARG:HG2	1.98	0.45
1:B:175:ARG:NH2	3:B:401:ANP:O1G	2.47	0.44
1:A:223:VAL:HB	1:B:229:THR:CG2	2.47	0.44
1:A:104:LEU:N	1:A:113:TYR:O	2.45	0.44
1:A:240:ILE:HG12	1:A:330:ILE:HD13	1.99	0.43
1:A:37:ARG:H	1:A:37:ARG:HD3	1.84	0.43
1:A:37:ARG:HH12	1:A:173:SER:HB3	1.83	0.42
1:A:104:LEU:HD12	1:A:104:LEU:HA	1.95	0.42
1:A:19:SER:HB3	1:A:118:GLN:HB2	2.02	0.42
1:B:198:ILE:HD11	1:B:234:VAL:CG2	2.51	0.41
1:A:176:VAL:O	1:A:176:VAL:HG12	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	310/352 (88%)	295 (95%)	11 (4%)	4 (1%)	9	30
1	B	264/352 (75%)	257 (97%)	7 (3%)	0	100	100
All	All	574/704 (82%)	552 (96%)	18 (3%)	4 (1%)	18	46

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	106	ASP
1	A	219	GLY
1	A	108	ASP
1	A	220	ALA

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	266/289 (92%)	262 (98%)	4 (2%)	57	82
1	B	234/289 (81%)	232 (99%)	2 (1%)	70	88
All	All	500/578 (86%)	494 (99%)	6 (1%)	63	85

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	213	ASN
1	A	218	ARG
1	A	235	GLU
1	A	303	ILE
1	B	179	LEU
1	B	196	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	253	GLN
1	B	77	ASN
1	B	327	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 2 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	$\# Z > 2$
3	ANP	A	401	2	33,33,33	0.93	4 (12%)	45,52,52	0.65	1 (2%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	RBF	A	402	-	29,29,29	1.29	2 (6%)	42,43,43	1.31	8 (19%)
3	ANP	B	401	2	33,33,33	0.98	4 (12%)	45,52,52	0.65	1 (2%)
4	RBF	B	402	-	29,29,29	1.32	3 (10%)	42,43,43	1.43	9 (21%)

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	ANP	A	401	2	-	6/18/38/38	0/3/3/3
4	RBF	A	402	-	-	3/14/14/14	0/3/3/3
3	ANP	B	401	2	-	4/18/38/38	0/3/3/3
4	RBF	B	402	-	-	0/14/14/14	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	402	RBF	C5A-N5	-3.15	1.33	1.39
4	A	402	RBF	C5A-N5	-3.13	1.33	1.39
3	A	401	ANP	PB-O1B	2.96	1.50	1.46
3	B	401	ANP	PG-N3B	2.63	1.70	1.63
3	B	401	ANP	PG-O1G	2.62	1.50	1.46
3	B	401	ANP	PB-O1B	2.58	1.50	1.46
3	A	401	ANP	PG-N3B	2.24	1.69	1.63
3	A	401	ANP	PG-O1G	2.20	1.49	1.46
3	B	401	ANP	PB-O3A	-2.13	1.56	1.59
3	A	401	ANP	PB-O3A	-2.10	1.56	1.59
4	B	402	RBF	C7M-C7	2.06	1.54	1.51
4	A	402	RBF	O2'-C2'	-2.04	1.39	1.43
4	B	402	RBF	O2'-C2'	-2.01	1.39	1.43

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	402	RBF	C4-N3-C2	-2.99	120.33	125.64
4	B	402	RBF	C4A-C4-N3	2.84	120.49	113.25
4	A	402	RBF	C4-N3-C2	-2.80	120.67	125.64
4	A	402	RBF	C4A-C4-N3	2.70	120.11	113.25
4	B	402	RBF	O4-C4-C4A	-2.60	119.68	126.53
4	B	402	RBF	C4A-C10-N1	-2.54	118.36	124.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	ANP	O1G-PG-N3B	-2.53	108.04	111.77
4	A	402	RBF	O4-C4-C4A	-2.51	119.90	126.53
4	A	402	RBF	C4A-C10-N1	-2.51	118.44	124.59
4	A	402	RBF	C5A-C9A-N10	2.40	120.14	117.97
4	B	402	RBF	C4-C4A-N5	2.32	121.41	118.21
4	B	402	RBF	C10-C4A-N5	-2.26	120.19	124.81
4	B	402	RBF	C10-N1-C2	2.24	121.71	116.85
4	B	402	RBF	C4A-C10-N10	2.22	119.66	116.48
4	A	402	RBF	C4A-C10-N10	2.21	119.65	116.48
4	A	402	RBF	C9A-N10-C10	-2.17	117.44	120.75
4	A	402	RBF	C10-N1-C2	2.16	121.53	116.85
4	B	402	RBF	C5A-C9A-N10	2.07	119.84	117.97
3	B	401	ANP	O1G-PG-N3B	-2.05	108.75	111.77

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	401	ANP	PB-N3B-PG-O1G
3	A	401	ANP	PB-O3A-PA-O5'
3	A	401	ANP	C5'-O5'-PA-O2A
3	B	401	ANP	PB-N3B-PG-O1G
3	B	401	ANP	PG-N3B-PB-O1B
3	B	401	ANP	PA-O3A-PB-O2B
4	A	402	RBF	N10-C1'-C2'-O2'
4	A	402	RBF	O4'-C4'-C5'-O5'
3	A	401	ANP	C5'-O5'-PA-O1A
3	A	401	ANP	C5'-O5'-PA-O3A
3	A	401	ANP	PG-N3B-PB-O1B
4	A	402	RBF	C3'-C4'-C5'-O5'
3	B	401	ANP	PG-N3B-PB-O3A

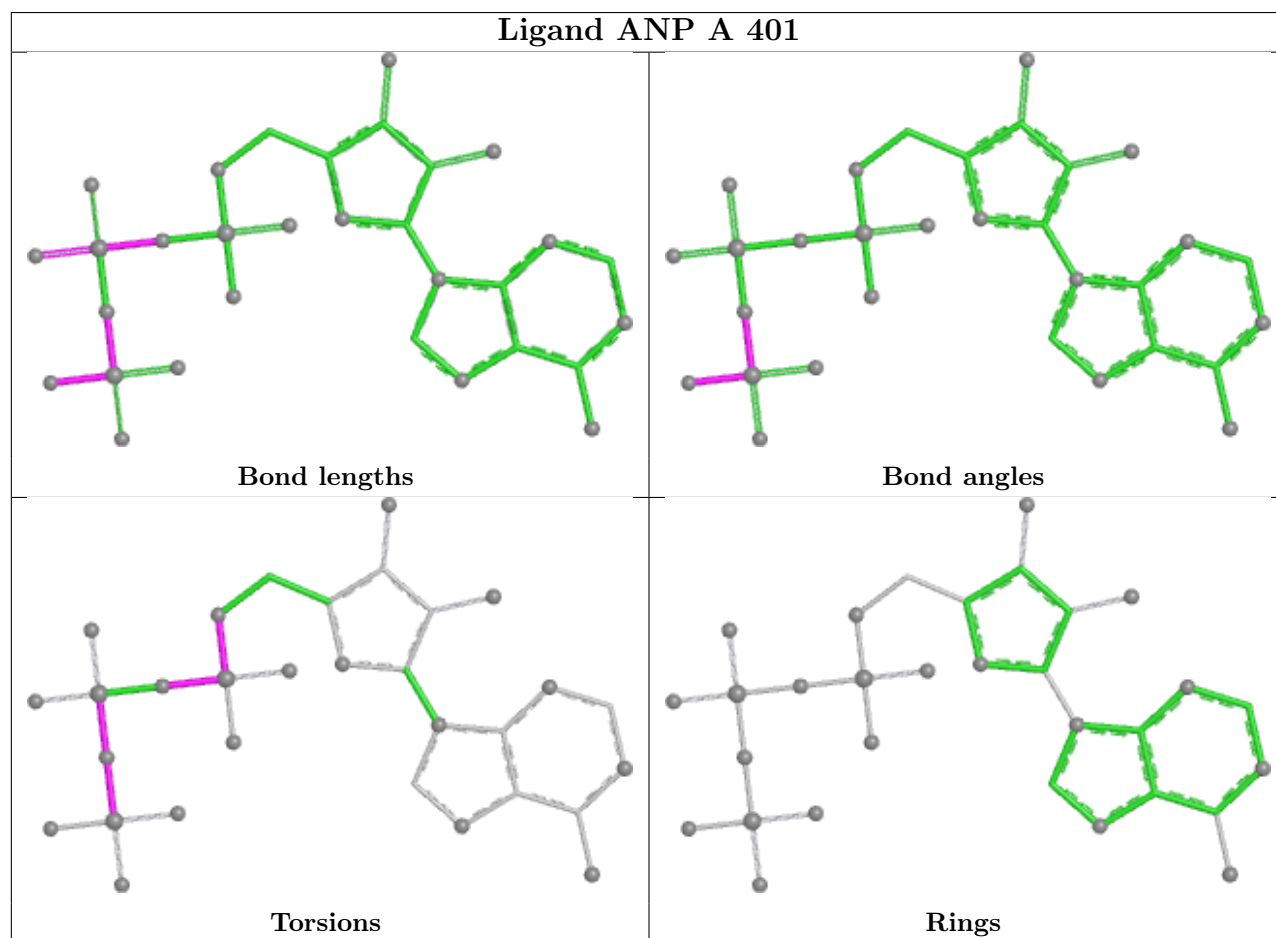
There are no ring outliers.

3 monomers are involved in 3 short contacts:

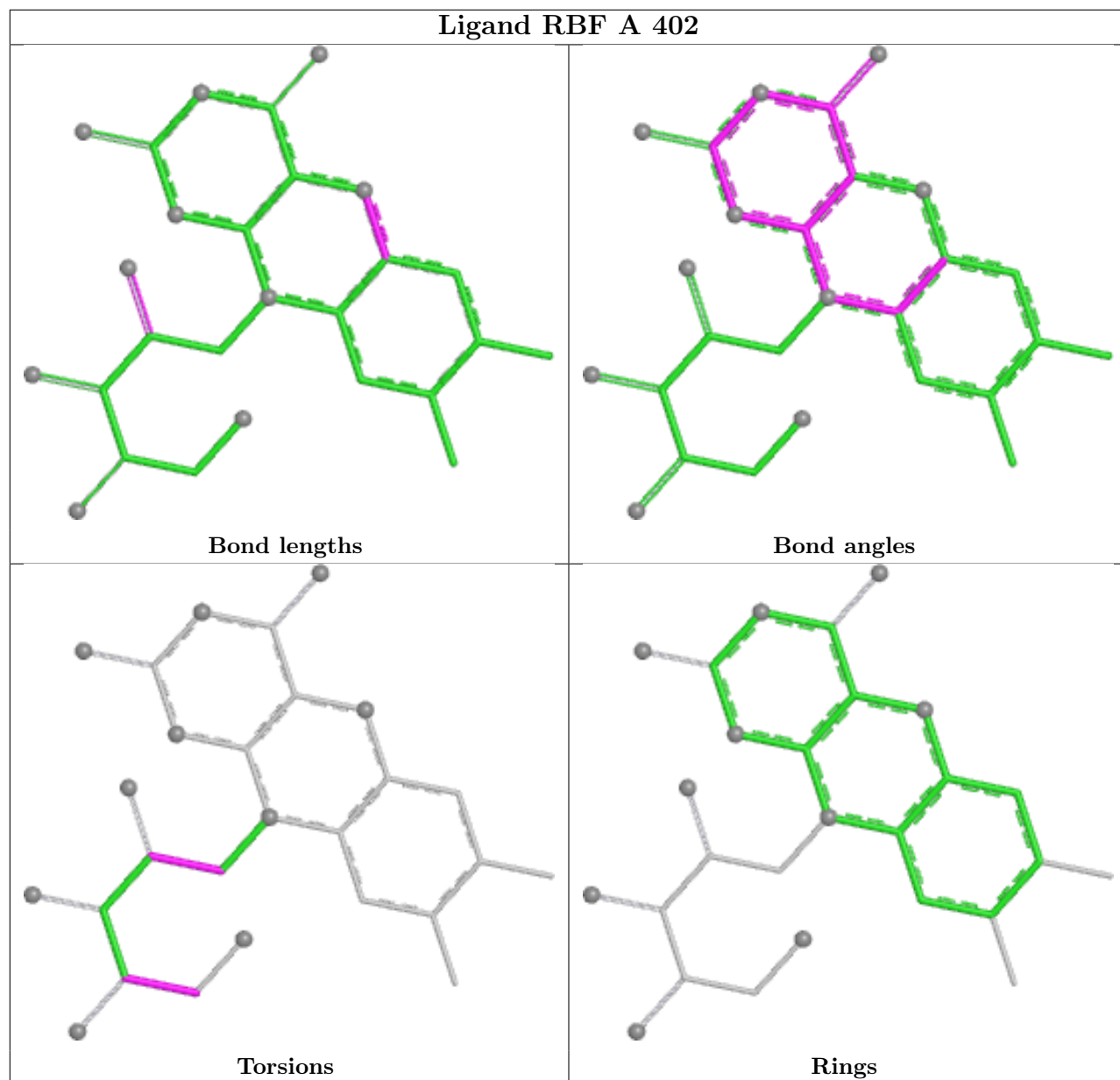
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	401	ANP	1	0
3	B	401	ANP	1	0
4	B	402	RBF	1	0

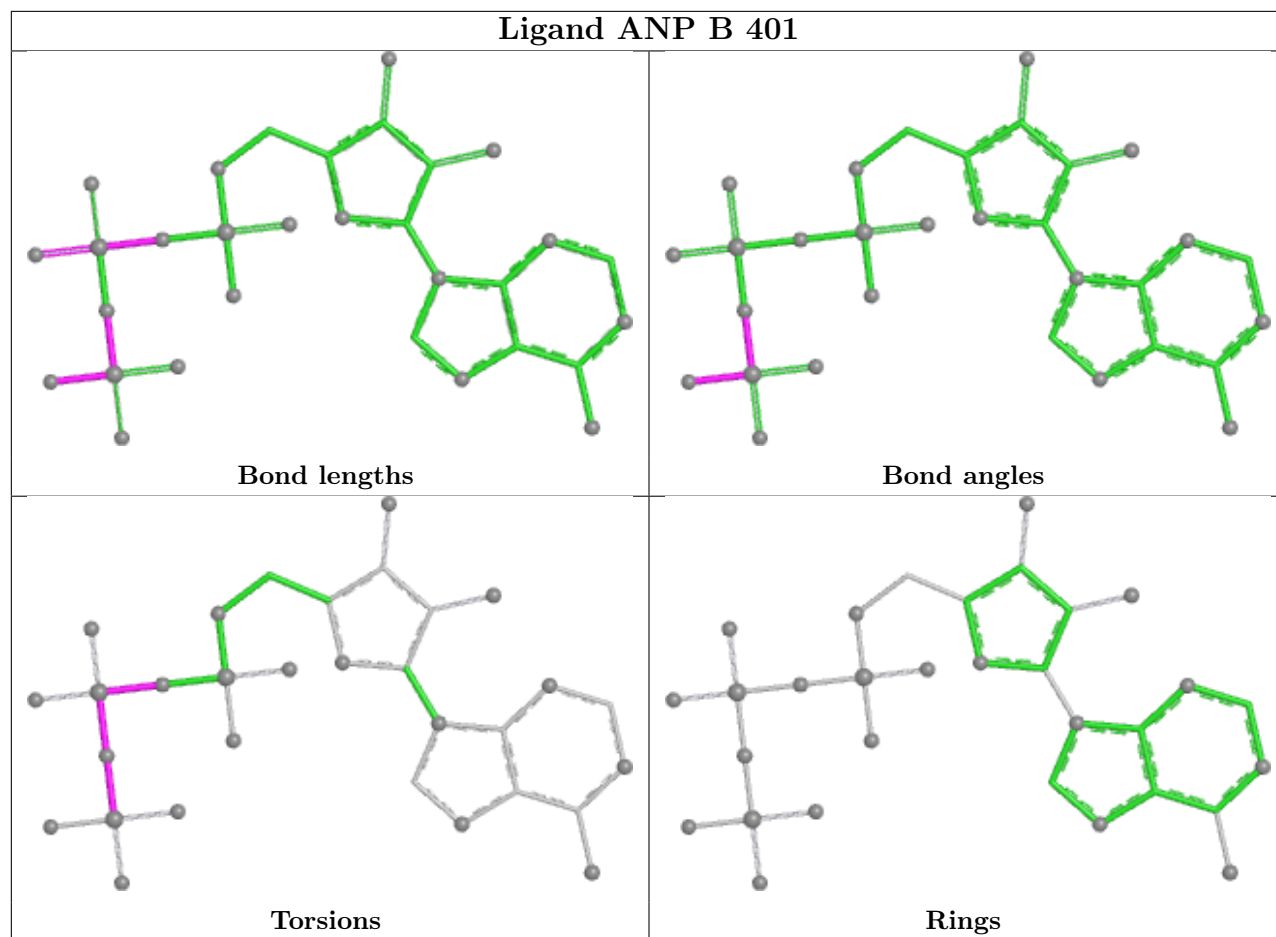
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

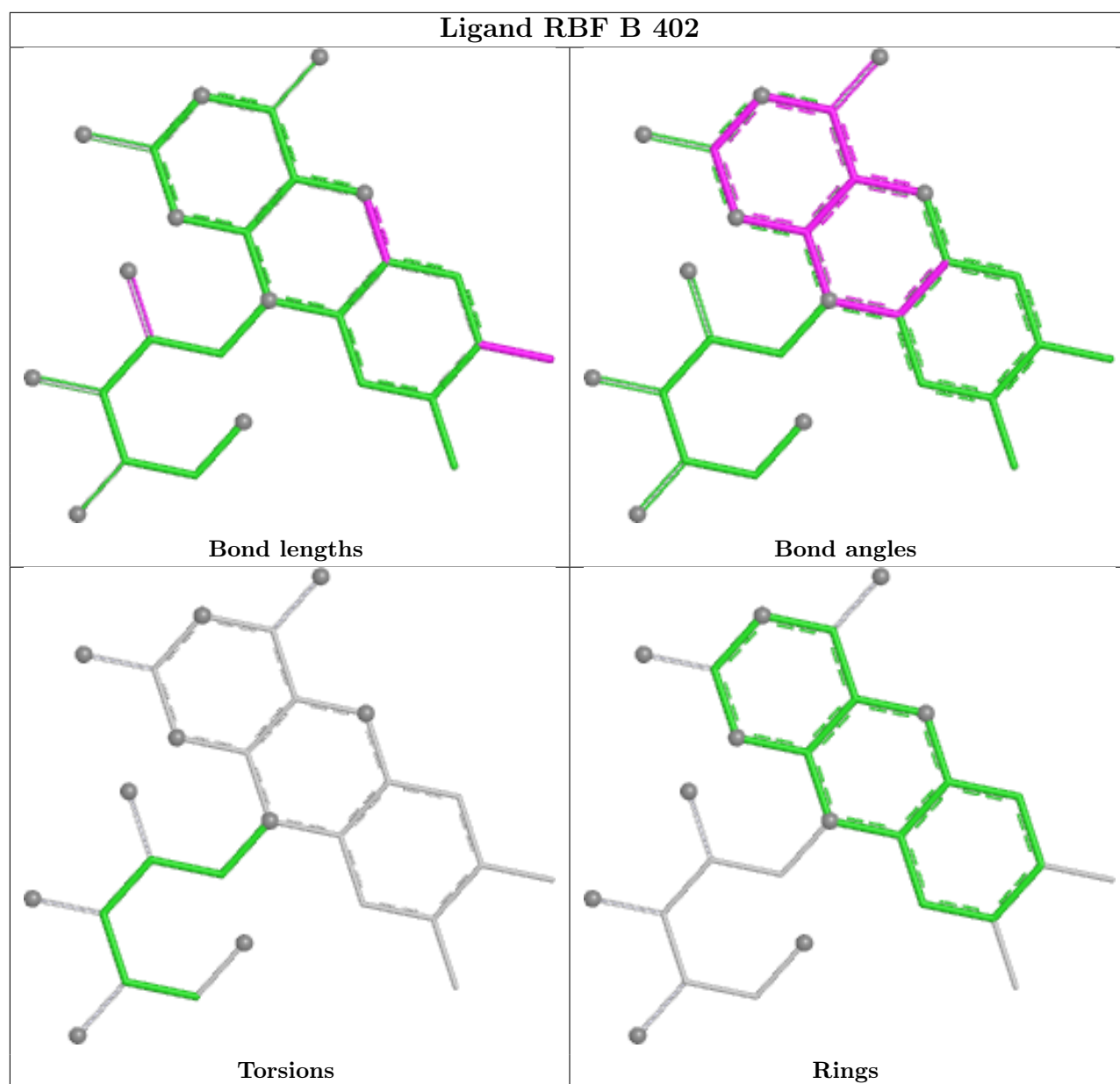
bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 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.



Ligand RBF A 402







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	318/352 (90%)	1.04	54 (16%) 4 3	22, 59, 135, 194	0
1	B	278/352 (78%)	2.07	123 (44%) 0 0	76, 130, 187, 212	0
All	All	596/704 (84%)	1.52	177 (29%) 1 1	22, 95, 175, 212	0

All (177) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	186	MET	6.9
1	A	261	SER	6.2
1	B	38	ALA	6.0
1	A	319	ARG	5.9
1	A	165	VAL	5.8
1	B	156	ARG	5.3
1	A	159	SER	4.8
1	B	220	ALA	4.6
1	A	220	ALA	4.6
1	B	147	HIS	4.6
1	B	104	LEU	4.6
1	B	63	THR	4.5
1	B	313	ALA	4.4
1	B	136	GLN	4.4
1	B	44	GLY	4.3
1	B	137	LEU	4.3
1	A	318	THR	4.2
1	B	307	LEU	4.1
1	A	275	GLN	4.1
1	B	19	SER	4.1
1	A	320	GLY	4.0
1	B	261	SER	4.0
1	B	169	PHE	4.0
1	B	141	GLN	4.0

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Mol	Chain	Res	Type	RSRZ
1	B	50	VAL	3.9
1	B	287	ILE	3.9
1	A	257	SER	3.9
1	B	176	VAL	3.9
1	B	155	ILE	3.9
1	A	219	GLY	3.8
1	A	334	GLN	3.7
1	B	330	ILE	3.7
1	A	321	GLY	3.7
1	A	282	ASP	3.7
1	B	65	PRO	3.6
1	B	39	PHE	3.6
1	B	62	LYS	3.6
1	B	274	GLU	3.6
1	A	260	ALA	3.6
1	B	51	VAL	3.6
1	B	311	LEU	3.6
1	B	111	CYS	3.5
1	A	124	SER	3.5
1	B	172	LEU	3.4
1	B	76	ARG	3.4
1	B	140	VAL	3.4
1	B	34	TYR	3.4
1	B	35	VAL	3.4
1	B	218	ARG	3.4
1	B	43	THR	3.4
1	B	88	TYR	3.4
1	B	20	LEU	3.3
1	B	36	ASN	3.3
1	B	336	LYS	3.3
1	A	13	TRP	3.3
1	A	338	LEU	3.3
1	B	306	GLN	3.2
1	B	248	LEU	3.2
1	A	322	THR	3.1
1	B	113	TYR	3.1
1	B	30	GLU	3.1
1	B	322	THR	3.1
1	B	45	TYR	3.1
1	B	114	PHE	3.1
1	B	327	ASN	3.1
1	B	115	VAL	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	135	ARG	3.0
1	B	293	TRP	3.0
1	B	317	VAL	3.0
1	B	48	SER	3.0
1	B	215	ILE	3.0
1	B	275	GLN	3.0
1	B	303	ILE	3.0
1	B	179	LEU	3.0
1	A	166	GLY	3.0
1	B	252	LEU	3.0
1	B	312	GLY	3.0
1	A	106	ASP	3.0
1	A	291	CYS	3.0
1	A	130	ALA	2.9
1	B	239	ARG	2.9
1	B	297	GLY	2.9
1	B	267	ARG	2.9
1	A	314	GLU	2.8
1	B	23	ALA	2.8
1	B	64	ASP	2.8
1	A	337	THR	2.8
1	B	212	ILE	2.8
1	B	315	LEU	2.8
1	A	222	LYS	2.8
1	B	235	GLU	2.8
1	B	177	GLU	2.8
1	B	25	ILE	2.7
1	B	90	ALA	2.7
1	B	138	ALA	2.7
1	B	290	ASP	2.7
1	A	10	LYS	2.7
1	A	41	GLN	2.7
1	B	198	ILE	2.7
1	A	135	ARG	2.7
1	A	218	ARG	2.7
1	B	286	GLY	2.7
1	A	290	ASP	2.7
1	B	86	TYR	2.6
1	A	136	GLN	2.6
1	B	263	VAL	2.6
1	B	96	TRP	2.6
1	B	196	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	251	ALA	2.6
1	B	288	PRO	2.6
1	A	25	ILE	2.6
1	B	216	ASP	2.6
1	B	206	GLY	2.6
1	B	72	ALA	2.6
1	B	81	VAL	2.5
1	B	154	MET	2.5
1	A	265	GLN	2.5
1	B	84	THR	2.5
1	A	188	ILE	2.5
1	B	282	ASP	2.5
1	B	210	SER	2.5
1	B	58	LEU	2.5
1	B	318	THR	2.5
1	B	91	ASP	2.5
1	B	228	ASP	2.5
1	A	121	MET	2.4
1	B	28	ASP	2.4
1	B	119	VAL	2.4
1	A	274	GLU	2.4
1	A	8	HIS	2.4
1	B	299	LEU	2.4
1	A	214	HIS	2.4
1	A	286	GLY	2.4
1	B	29	ASP	2.4
1	A	215	ILE	2.4
1	A	323	GLY	2.4
1	B	24	ASP	2.4
1	B	171	SER	2.4
1	B	189	ALA	2.3
1	A	335	GLN	2.3
1	A	272	SER	2.3
1	B	256	PHE	2.3
1	B	40	GLU	2.3
1	A	172	LEU	2.3
1	B	180	GLN	2.3
1	B	332	LEU	2.2
1	A	216	ASP	2.2
1	A	129	ARG	2.2
1	B	236	THR	2.2
1	B	121	MET	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	221	ILE	2.2
1	B	150	MET	2.2
1	A	139	GLU	2.2
1	B	298	ASN	2.2
1	B	281	GLU	2.2
1	B	183	TYR	2.2
1	A	156	ARG	2.1
1	B	214	HIS	2.1
1	B	102	GLY	2.1
1	B	291	CYS	2.1
1	B	149	ALA	2.1
1	B	296	GLU	2.1
1	B	101	MET	2.1
1	A	108	ASP	2.1
1	A	333	SER	2.1
1	B	26	SER	2.1
1	B	79	GLU	2.1
1	B	301	SER	2.1
1	B	97	ASN	2.1
1	A	255	ALA	2.1
1	A	336	LYS	2.1
1	B	231	ASP	2.1
1	A	125	GLU	2.1
1	B	42	MET	2.0
1	A	104	LEU	2.0
1	B	219	GLY	2.0
1	B	328	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,

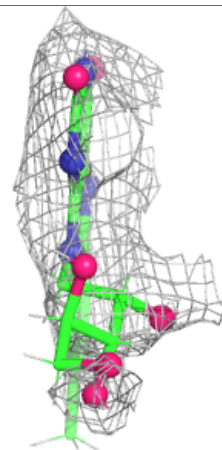
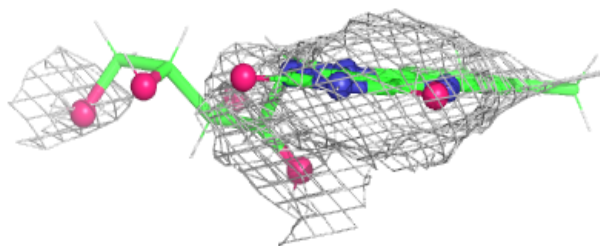
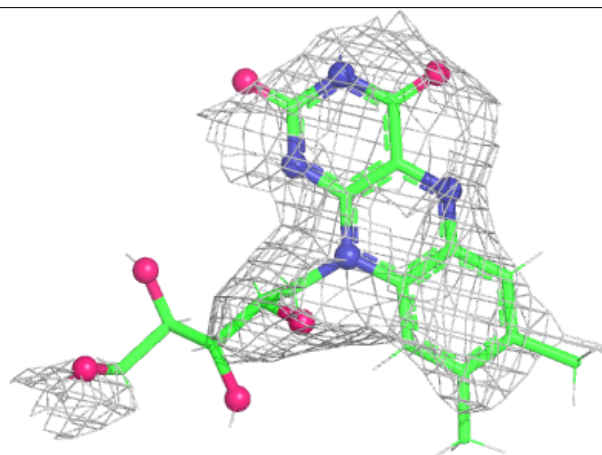
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	RBF	B	402	27/27	0.82	0.19	121,148,228,236	0
3	ANP	B	401	31/31	0.85	0.14	42,106,139,160	0
3	ANP	A	401	31/31	0.85	0.22	23,52,185,217	0
4	RBF	A	402	27/27	0.92	0.11	12,37,74,86	0
2	MG	A	400	1/1	0.95	0.10	24,24,24,24	0
2	MG	B	400	1/1	0.96	0.06	79,79,79,79	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.

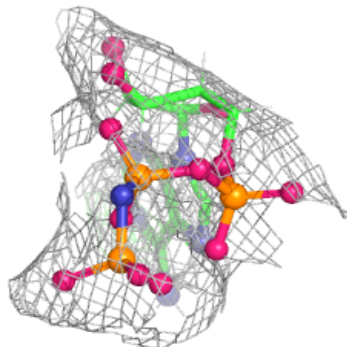
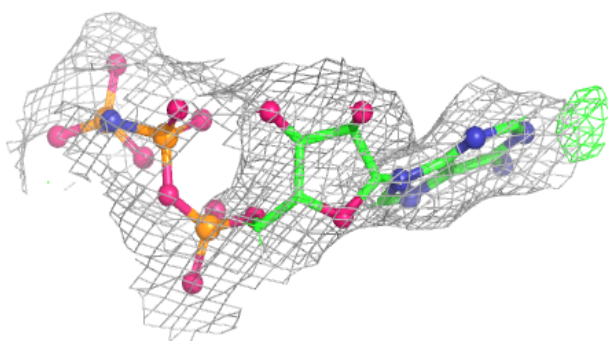
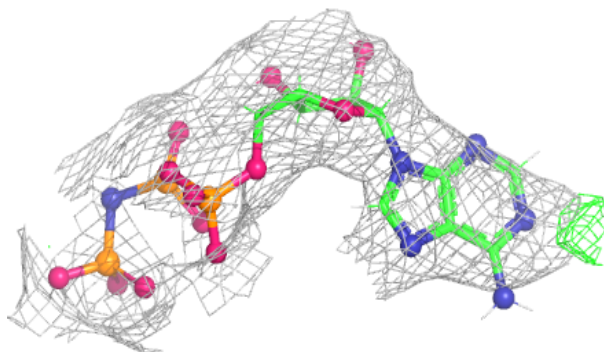
Electron density around RBF B 402:

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

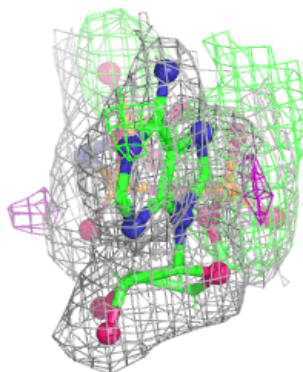
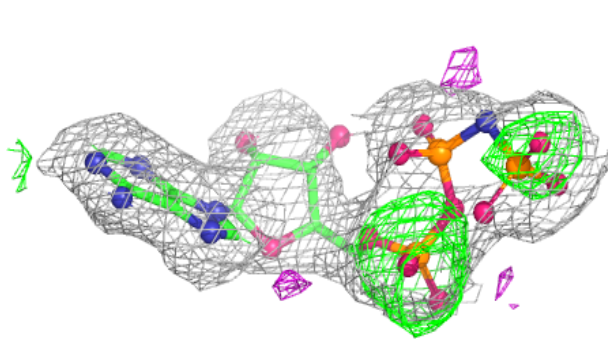
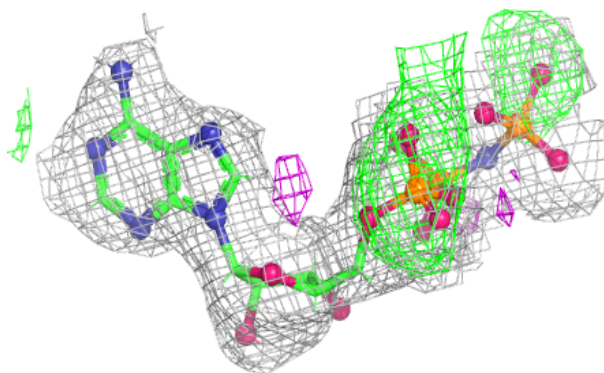


Electron density around ANP B 401:

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

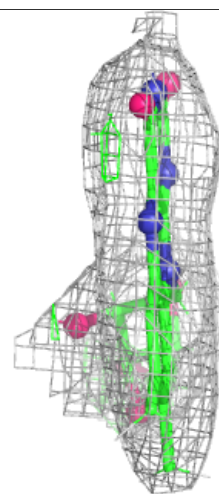
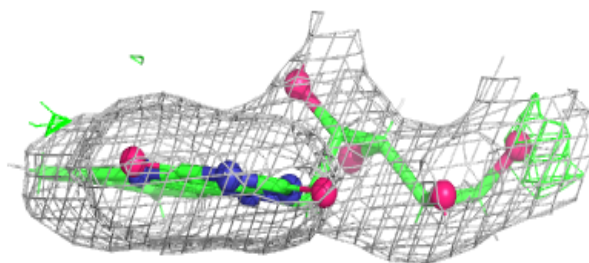
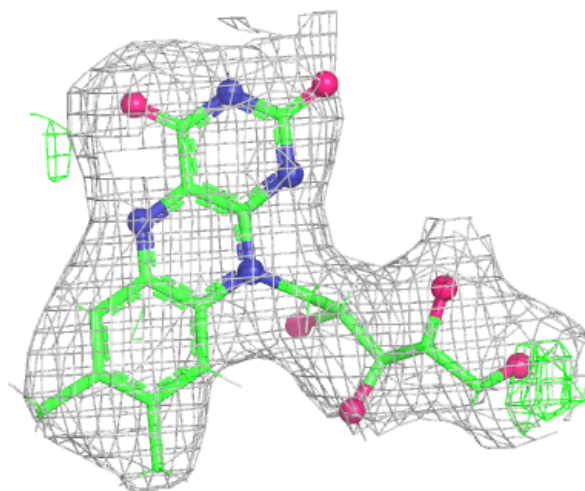
**Electron density around ANP A 401:**

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



Electron density around RBF A 402:

$2mF_o-DF_c$ (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.