



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 06:04 pm BST

PDB ID : 3A35
Title : Crystal structure of LumP complexed with riboflavin
Authors : Sato, Y.
Deposited on : 2009-06-09
Resolution : 1.42 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

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

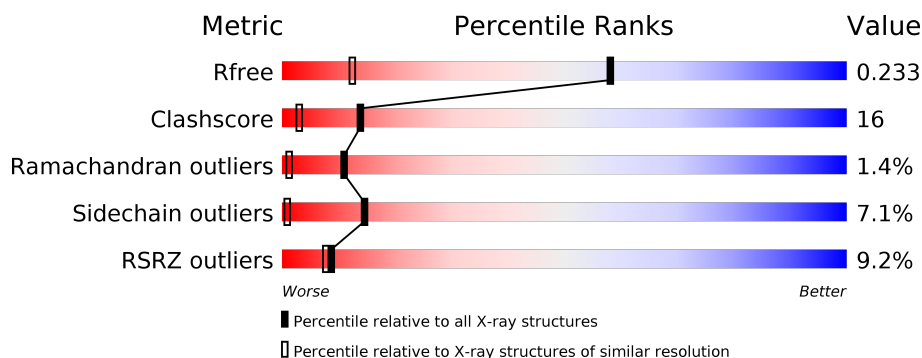
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.42 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	2579 (1.44-1.40)
Clashscore	141614	2696 (1.44-1.40)
Ramachandran outliers	138981	2632 (1.44-1.40)
Sidechain outliers	138945	2631 (1.44-1.40)
RSRZ outliers	127900	2528 (1.44-1.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	190	<div> <div>8%</div> <div>66%</div> <div>24%</div> <div>• • •</div> </div>
1	B	190	<div> <div>9%</div> <div>68%</div> <div>22%</div> <div>5%</div> <div>• • •</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3327 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 Lumazine protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	184	Total	C	N	O	S	0	3	0
			1442	906	242	291	3			
1	B	184	Total	C	N	O	S	0	1	0
			1425	898	238	286	3			

There are 26 discrepancies between the modelled and reference sequences:

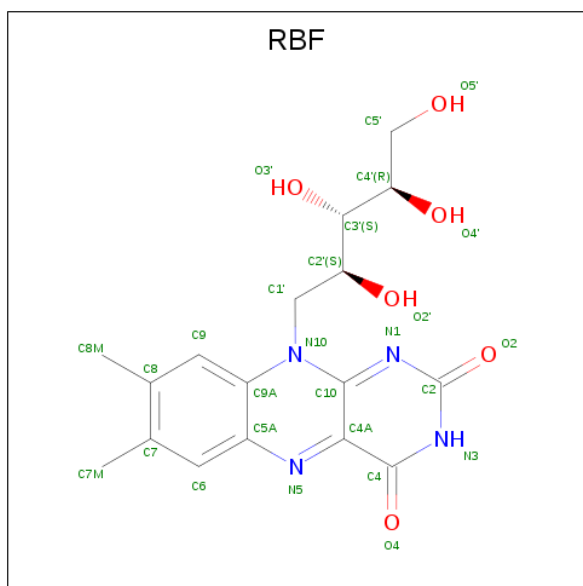
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	SEE REMARK 999	UNP C4TPG1
A	2	PHE	-	SEE REMARK 999	UNP C4TPG1
A	3	LYS	-	SEE REMARK 999	UNP C4TPG1
A	4	GLY	-	SEE REMARK 999	UNP C4TPG1
A	5	ILE	-	SEE REMARK 999	UNP C4TPG1
A	6	VAL	-	SEE REMARK 999	UNP C4TPG1
A	7	GLN	-	SEE REMARK 999	UNP C4TPG1
A	185	VAL	-	SEE REMARK 999	UNP C4TPG1
A	186	SER	-	SEE REMARK 999	UNP C4TPG1
A	187	ASN	-	SEE REMARK 999	UNP C4TPG1
A	188	GLU	-	SEE REMARK 999	UNP C4TPG1
A	189	TRP	-	SEE REMARK 999	UNP C4TPG1
A	190	ASP	-	SEE REMARK 999	UNP C4TPG1
B	1	MET	-	SEE REMARK 999	UNP C4TPG1
B	2	PHE	-	SEE REMARK 999	UNP C4TPG1
B	3	LYS	-	SEE REMARK 999	UNP C4TPG1
B	4	GLY	-	SEE REMARK 999	UNP C4TPG1
B	5	ILE	-	SEE REMARK 999	UNP C4TPG1
B	6	VAL	-	SEE REMARK 999	UNP C4TPG1
B	7	GLN	-	SEE REMARK 999	UNP C4TPG1
B	185	VAL	-	SEE REMARK 999	UNP C4TPG1
B	186	SER	-	SEE REMARK 999	UNP C4TPG1
B	187	ASN	-	SEE REMARK 999	UNP C4TPG1
B	188	GLU	-	SEE REMARK 999	UNP C4TPG1
B	189	TRP	-	SEE REMARK 999	UNP C4TPG1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	190	ASP	-	SEE REMARK 999	UNP C4TPG1

- Molecule 2 is RIBOFLAVIN (three-letter code: RBF) (formula: $C_{17}H_{20}N_4O_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			27	17	4	6		
2	B	1	Total	C	N	O	0	0
			27	17	4	6		

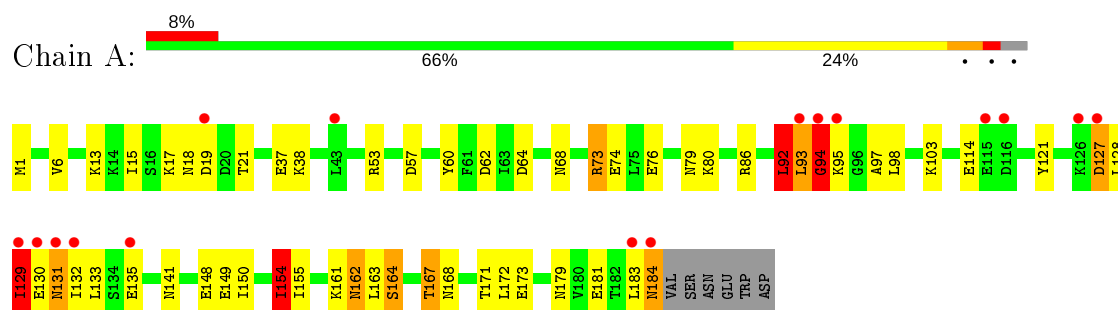
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	201	Total	O	0	1
			202	202		
3	B	204	Total	O	0	0
			204	204		

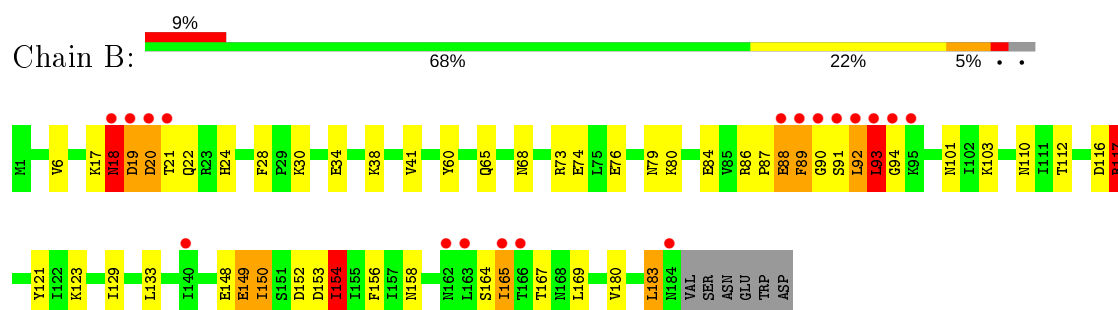
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Lumazine protein



• Molecule 1: Lumazine protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	46.29Å 46.58Å 161.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.15 – 1.42 20.14 – 1.42	Depositor EDS
% Data completeness (in resolution range)	97.1 (20.15-1.42) 97.1 (20.14-1.42)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.05 (at 1.42Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.192 , 0.234 0.190 , 0.233	Depositor DCC
R_{free} test set	6533 reflections (10.10%)	wwPDB-VP
Wilson B-factor (Å ²)	16.3	Xtriage
Anisotropy	0.046	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.46 , 65.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.029 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3327	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.43% 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: 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	1.45	6/1457 (0.4%)	1.46	18/1968 (0.9%)
1	B	1.50	5/1440 (0.3%)	1.43	13/1945 (0.7%)
All	All	1.48	11/2897 (0.4%)	1.45	31/3913 (0.8%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	1	1
1	B	2	1
All	All	3	2

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	164	SER	CB-OG	10.06	1.55	1.42
1	B	180	VAL	CB-CG2	-8.51	1.34	1.52
1	A	37	GLU	CB-CG	-7.66	1.37	1.52
1	A	181	GLU	CG-CD	-7.34	1.41	1.51
1	A	38	LYS	CG-CD	5.89	1.72	1.52
1	B	65	GLN	CB-CG	-5.71	1.37	1.52
1	A	149	GLU	CG-CD	5.61	1.60	1.51
1	A	167	THR	C-O	5.48	1.33	1.23
1	B	84	GLU	CD-OE1	-5.34	1.19	1.25
1	B	123	LYS	CE-NZ	5.16	1.61	1.49
1	A	148	GLU	CB-CG	-5.12	1.42	1.52

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	53	ARG	NE-CZ-NH2	-10.07	115.26	120.30
1	A	64	ASP	CB-CG-OD1	9.23	126.61	118.30
1	B	117	ARG	NE-CZ-NH2	-8.96	115.82	120.30
1	A	93	LEU	N-CA-C	7.25	130.58	111.00
1	A	154	ILE	CG1-CB-CG2	7.09	126.99	111.40
1	B	38	LYS	CD-CE-NZ	-6.88	95.87	111.70
1	A	38	LYS	CD-CE-NZ	-6.81	96.04	111.70
1	A	92	LEU	N-CA-C	-6.78	92.71	111.00
1	B	117	ARG	NE-CZ-NH1	6.68	123.64	120.30
1	B	116	ASP	CB-CG-OD2	-6.65	112.31	118.30
1	A	94	GLY	N-CA-C	-6.60	96.61	113.10
1	B	88	GLU	N-CA-C	6.49	128.51	111.00
1	B	183	LEU	CB-CG-CD2	6.42	121.91	111.00
1	A	60	TYR	CE1-CZ-CE2	6.40	130.04	119.80
1	B	150	ILE	CG1-CB-CG2	6.15	124.92	111.40
1	A	129	ILE	N-CA-C	-6.13	94.45	111.00
1	A	172	LEU	CB-CG-CD1	-6.10	100.63	111.00
1	B	28	PHE	CB-CG-CD2	-6.04	116.57	120.80
1	A	57	ASP	CB-CG-OD2	-5.82	113.06	118.30
1	A	13	LYS	CD-CE-NZ	-5.81	98.34	111.70
1	A	60	TYR	CD1-CE1-CZ	-5.68	114.69	119.80
1	B	60	TYR	CB-CG-CD2	5.67	124.40	121.00
1	A	62	ASP	CB-CG-OD2	5.53	123.28	118.30
1	B	154	ILE	CG1-CB-CG2	5.38	123.23	111.40
1	B	93	LEU	CA-CB-CG	5.25	127.38	115.30
1	B	133	LEU	CB-CG-CD2	-5.25	102.07	111.00
1	A	73[A]	ARG	NE-CZ-NH1	-5.14	117.73	120.30
1	A	73[B]	ARG	NE-CZ-NH1	-5.14	117.73	120.30
1	A	86	ARG	CG-CD-NE	-5.11	101.07	111.80
1	B	86	ARG	N-CA-CB	-5.11	101.41	110.60
1	A	163	LEU	CB-CG-CD1	-5.02	102.46	111.00

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	93	LEU	CA
1	B	112	THR	CB
1	B	154	ILE	CB

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	92	LEU	Peptide

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Mol	Chain	Res	Type	Group
1	B	88	GLU	Peptide

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	1442	0	1469	49	0
1	B	1425	0	1459	43	0
2	A	27	0	20	0	0
2	B	27	0	19	2	0
3	A	202	0	0	9	1
3	B	204	0	0	8	1
All	All	3327	0	2967	93	1

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 16.

All (93) 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:89:PHE:HB2	1:B:90:GLY:CA	1.64	1.23
1:A:94:GLY:CA	1:A:95:LYS:HB2	1.79	1.13
1:B:169:LEU:HD21	3:B:239:HOH:O	1.48	1.10
1:B:87:PRO:O	1:B:89:PHE:HA	1.51	1.09
1:A:131:ASN:HB3	1:A:132:ILE:HA	1.33	1.09
1:A:94:GLY:HA2	1:A:95:LYS:CB	1.81	1.08
1:B:89:PHE:CB	1:B:90:GLY:HA3	1.92	0.99
1:B:89:PHE:HB2	1:B:90:GLY:HA3	0.97	0.96
1:B:121:TYR:HB3	1:B:154:ILE:HD11	1.47	0.95
1:B:129:ILE:HD11	1:B:150:ILE:HD11	1.49	0.93
1:A:128:LEU:HA	1:A:129:ILE:HB	1.52	0.90
1:A:94:GLY:HA2	1:A:95:LYS:HB2	0.91	0.88
1:B:117:ARG:HD3	1:B:158:ASN:OD1	1.76	0.85
1:A:131:ASN:CB	1:A:132:ILE:HA	2.06	0.85
1:A:162:ASN:OD1	3:A:657:HOH:O	1.95	0.82
1:B:21:THR:HG22	3:B:473:HOH:O	1.84	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:ASN:HB3	1:A:132:ILE:CA	2.14	0.77
1:B:21:THR:CG2	3:B:473:HOH:O	2.33	0.77
1:B:129:ILE:HD11	1:B:150:ILE:CD1	2.16	0.75
1:B:110:ASN:HD21	1:B:112:THR:HG23	1.51	0.74
1:A:18:ASN:HB2	1:A:21:THR:HG23	1.70	0.73
1:A:15:ILE:CG2	1:A:17:LYS:HZ3	2.01	0.73
1:B:89:PHE:CB	1:B:90:GLY:CA	2.55	0.73
1:B:169:LEU:HD23	3:B:311:HOH:O	1.87	0.73
1:B:92:LEU:HD23	1:B:94:GLY:N	2.03	0.72
1:A:161:LYS:O	1:A:164:SER:HB2	1.90	0.72
1:A:128:LEU:HA	1:A:129:ILE:CB	2.22	0.70
1:A:131:ASN:HD21	1:A:184:ASN:C	1.93	0.70
1:A:15:ILE:HG21	1:A:17:LYS:HZ3	1.57	0.69
1:B:19:ASP:N	1:B:19:ASP:OD1	2.25	0.69
1:A:128:LEU:CA	1:A:129:ILE:HB	2.23	0.69
1:B:80:LYS:HE3	3:B:652:HOH:O	1.93	0.68
1:B:30:LYS:NZ	1:B:34:GLU:OE2	2.23	0.67
1:A:121:TYR:HB3	1:A:154:ILE:HD11	1.76	0.67
1:B:87:PRO:C	1:B:89:PHE:HA	2.15	0.67
1:A:121:TYR:HB3	1:A:154:ILE:CD1	2.27	0.64
1:B:101:ASN:HD22	1:B:103:LYS:NZ	1.96	0.63
1:B:80:LYS:CE	3:B:652:HOH:O	2.49	0.61
1:A:127:ASP:O	1:A:129:ILE:HD13	2.02	0.60
1:B:117:ARG:CD	1:B:158:ASN:OD1	2.49	0.59
1:B:148:GLU:C	1:B:149:GLU:HG2	2.23	0.58
3:A:337:HOH:O	1:B:93:LEU:HG	2.06	0.55
1:A:94:GLY:CA	1:A:95:LYS:CB	2.54	0.55
1:B:110:ASN:ND2	1:B:112:THR:HG23	2.20	0.55
1:B:129:ILE:CD1	1:B:150:ILE:HD11	2.30	0.55
1:A:131:ASN:CB	1:A:132:ILE:CA	2.80	0.55
1:B:76:GLU:H	1:B:79:ASN:ND2	2.04	0.55
1:A:133:LEU:HB3	1:A:135:GLU:OE2	2.07	0.54
1:A:80:LYS:NZ	1:A:168:ASN:HA	2.23	0.54
1:B:110:ASN:HD21	1:B:112:THR:CG2	2.21	0.54
1:A:133:LEU:HD12	1:A:135:GLU:OE1	2.07	0.54
1:A:1:MET:HB2	1:A:98:LEU:O	2.08	0.53
1:A:80:LYS:NZ	1:A:171:THR:HG23	2.23	0.52
1:A:93:LEU:HD12	1:A:93:LEU:N	2.25	0.52
1:A:141:ASN:ND2	1:A:179:ASN:H	2.09	0.51
1:A:76:GLU:H	1:A:79:ASN:ND2	2.09	0.50
1:B:92:LEU:HB3	1:B:93:LEU:HA	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:VAL:HB	1:A:167:THR:HG22	1.93	0.50
1:B:17:LYS:O	1:B:18:ASN:HB2	2.11	0.50
1:A:129:ILE:O	1:A:130:GLU:HG2	2.12	0.49
1:A:80:LYS:NZ	1:A:171:THR:CG2	2.76	0.49
1:A:93:LEU:HD21	3:A:606:HOH:O	2.13	0.49
1:B:22:GLN:OE1	1:B:24:HIS:HE1	1.95	0.49
1:B:41:VAL:HG11	2:B:191:RBF:HC72	1.94	0.49
1:B:149:GLU:HG3	1:B:156:PHE:HD2	1.76	0.48
1:A:103:LYS:NZ	3:A:655[B]:HOH:O	2.42	0.48
1:A:15:ILE:HG22	1:A:17:LYS:HZ3	1.74	0.48
1:B:18:ASN:C	1:B:19:ASP:OD1	2.51	0.48
1:A:93:LEU:HB3	1:A:97:ALA:HB2	1.96	0.47
1:A:133:LEU:CD1	1:A:135:GLU:OE1	2.64	0.46
1:A:150:ILE:HG12	1:A:155:ILE:HG22	1.99	0.45
1:A:173:GLU:HB3	3:A:593:HOH:O	2.16	0.45
1:B:19:ASP:HB2	1:B:20:ASP:H	1.59	0.45
1:A:133:LEU:HB2	1:A:135:GLU:OE1	2.16	0.45
1:A:133:LEU:CB	1:A:135:GLU:OE1	2.65	0.44
1:A:98:LEU:HD23	1:A:183:LEU:HD21	2.00	0.44
1:A:154:ILE:HG12	3:A:259:HOH:O	2.17	0.44
1:A:15:ILE:HG21	1:A:17:LYS:NZ	2.28	0.44
1:A:68:ASN:HD22	1:A:73[A]:ARG:HH12	1.66	0.43
1:B:20:ASP:OD1	1:B:20:ASP:N	2.50	0.43
1:A:154:ILE:HG13	3:A:236:HOH:O	2.17	0.43
1:A:68:ASN:ND2	1:A:73[B]:ARG:HH12	2.17	0.42
1:A:93:LEU:HB2	3:A:335:HOH:O	2.19	0.42
2:B:191:RBF:HC11	2:B:191:RBF:HC9	1.90	0.42
1:A:93:LEU:CB	1:A:97:ALA:HB2	2.49	0.42
1:B:169:LEU:CD2	3:B:311:HOH:O	2.57	0.41
3:A:248:HOH:O	1:B:165:ILE:HD11	2.20	0.41
1:B:73:ARG:HG3	1:B:74:GLU:HG3	2.01	0.41
1:B:153:ASP:N	1:B:153:ASP:OD1	2.53	0.41
1:B:6:VAL:HB	1:B:167:THR:HG22	2.03	0.41
1:A:93:LEU:HD12	1:A:93:LEU:H	1.86	0.40
1:B:101:ASN:HD22	1:B:103:LYS:HZ3	1.66	0.40
1:B:167:THR:HA	3:B:199:HOH:O	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:608:HOH:O	3:B:293:HOH:O[4_455]	2.15	0.05

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	185/190 (97%)	176 (95%)	5 (3%)	4 (2%)	6	0
1	B	183/190 (96%)	172 (94%)	10 (6%)	1 (0%)	29	8
All	All	368/380 (97%)	348 (95%)	15 (4%)	5 (1%)	11	1

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	129	ILE
1	B	18	ASN
1	A	131	ASN
1	A	19	ASP
1	A	94	GLY

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	166/169 (98%)	157 (95%)	9 (5%)	22	2
1	B	164/169 (97%)	150 (92%)	14 (8%)	10	0
All	All	330/338 (98%)	307 (93%)	23 (7%)	14	1

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	74	GLU
1	A	92	LEU
1	A	114	GLU
1	A	127	ASP
1	A	129	ILE
1	A	154	ILE
1	A	162	ASN
1	A	164	SER
1	A	184	ASN
1	B	18	ASN
1	B	19	ASP
1	B	20	ASP
1	B	68	ASN
1	B	89	PHE
1	B	91	SER
1	B	92	LEU
1	B	93	LEU
1	B	117	ARG
1	B	149	GLU
1	B	152	ASP
1	B	154	ILE
1	B	165	ILE
1	B	183	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	68	ASN
1	A	79	ASN
1	A	110	ASN
1	A	131	ASN
1	A	141	ASN
1	A	162	ASN
1	A	184	ASN
1	B	24	HIS
1	B	79	ASN
1	B	101	ASN
1	B	110	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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	RBF	B	191	-	27,29,29	2.86	10 (37%)	33,43,43	2.72	15 (45%)
2	RBF	A	191	-	27,29,29	3.06	10 (37%)	33,43,43	2.59	13 (39%)

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	RBF	B	191	-	-	1/14/14/14	0/3/3/3
2	RBF	A	191	-	-	0/14/14/14	0/3/3/3

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	191	RBF	C4A-N5	8.33	1.45	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	191	RBF	C2-N3	7.44	1.52	1.38
2	A	191	RBF	C1'-N10	-7.12	1.40	1.48
2	B	191	RBF	C4A-C10	7.03	1.45	1.38
2	B	191	RBF	C2-N1	-6.18	1.25	1.38
2	A	191	RBF	C4-N3	5.19	1.42	1.33
2	A	191	RBF	C2'-C3'	-4.92	1.44	1.53
2	A	191	RBF	O3'-C3'	4.35	1.53	1.43
2	B	191	RBF	C1'-N10	4.18	1.52	1.48
2	A	191	RBF	C9A-N10	4.08	1.44	1.38
2	B	191	RBF	C4A-N5	3.25	1.38	1.33
2	B	191	RBF	C7M-C7	-3.22	1.44	1.51
2	B	191	RBF	C9A-C5A	3.01	1.48	1.42
2	A	191	RBF	C5A-N5	2.74	1.39	1.35
2	A	191	RBF	C9-C8	2.60	1.44	1.37
2	A	191	RBF	C7M-C7	2.54	1.56	1.51
2	B	191	RBF	C5'-C4'	2.27	1.58	1.52
2	B	191	RBF	O3'-C3'	2.19	1.48	1.43
2	A	191	RBF	C10-N1	2.11	1.36	1.33
2	B	191	RBF	C8M-C8	2.07	1.55	1.51

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	191	RBF	C4A-C4-N3	-6.59	114.42	123.43
2	A	191	RBF	C4-N3-C2	6.59	120.71	115.14
2	B	191	RBF	C1'-N10-C9A	-5.96	113.61	118.29
2	A	191	RBF	C1'-N10-C9A	5.58	122.69	118.29
2	B	191	RBF	O5'-C5'-C4'	-5.44	99.22	111.07
2	B	191	RBF	C4-N3-C2	5.32	119.64	115.14
2	B	191	RBF	C5'-C4'-C3'	-4.89	101.80	112.41
2	B	191	RBF	C4-C4A-C10	-4.61	116.90	119.95
2	A	191	RBF	C7-C6-C5A	-4.52	114.83	121.22
2	B	191	RBF	C9-C9A-C5A	-3.81	113.37	119.88
2	A	191	RBF	C5A-C9A-N10	3.64	120.35	117.72
2	B	191	RBF	C4A-C10-N10	-3.41	116.79	120.30
2	B	191	RBF	C1'-N10-C10	3.19	121.27	118.41
2	A	191	RBF	C9A-N10-C10	-3.19	117.73	121.91
2	B	191	RBF	C5A-C9A-N10	-3.14	115.44	117.72
2	B	191	RBF	C7-C6-C5A	3.10	125.60	121.22
2	A	191	RBF	C6-C5A-C9A	3.02	123.01	119.05
2	A	191	RBF	C6-C5A-N5	-2.92	115.83	119.05
2	B	191	RBF	C4A-N5-C5A	2.78	119.55	116.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	191	RBF	C7M-C7-C6	-2.56	114.21	120.34
2	A	191	RBF	O3'-C3'-C4'	-2.55	102.65	108.81
2	A	191	RBF	C6-C7-C8	2.54	124.19	119.91
2	B	191	RBF	C4A-C4-N3	-2.44	120.10	123.43
2	A	191	RBF	C4-C4A-C10	2.43	121.56	119.95
2	B	191	RBF	C4-C4A-N5	2.32	121.24	118.60
2	B	191	RBF	C9A-N10-C10	2.30	124.92	121.91
2	B	191	RBF	C8-C9-C9A	2.21	126.05	119.19
2	A	191	RBF	C4-C4A-N5	-2.01	116.30	118.60

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	191	RBF	O4'-C4'-C5'-O5'

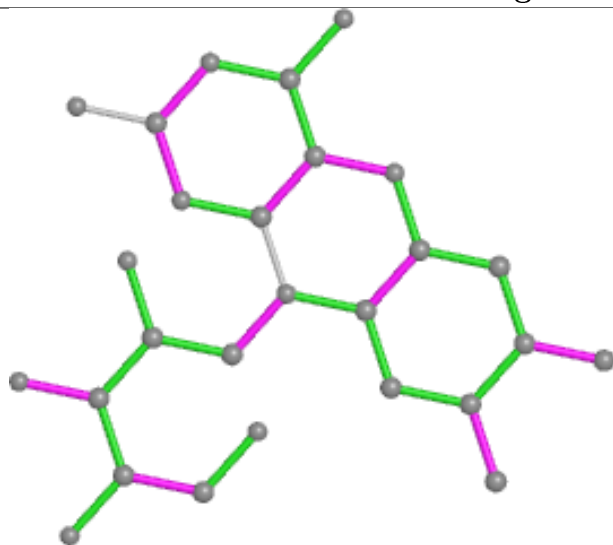
There are no ring outliers.

1 monomer is involved in 2 short contacts:

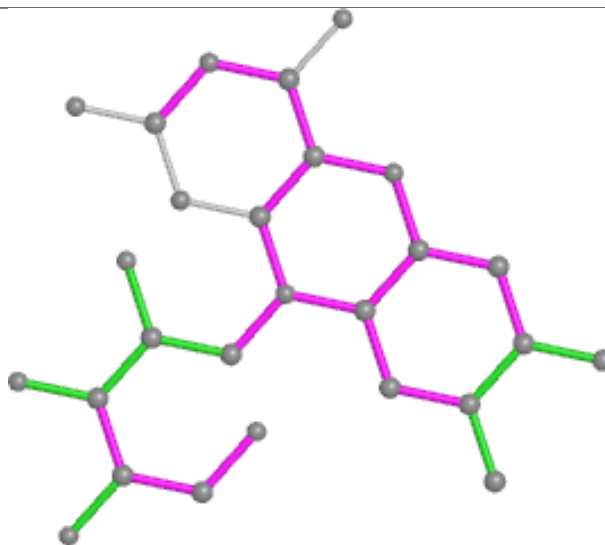
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	191	RBF	2	0

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.

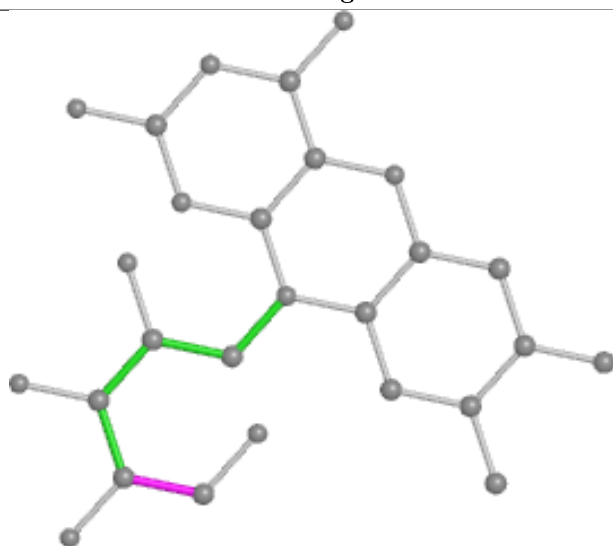
Ligand RBF B 191



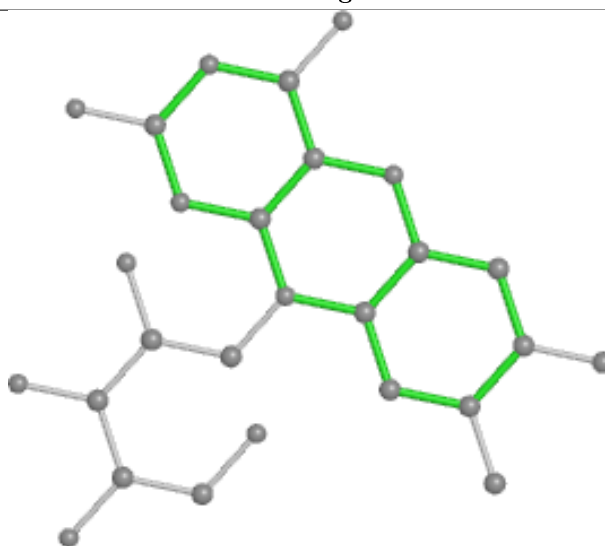
Bond lengths



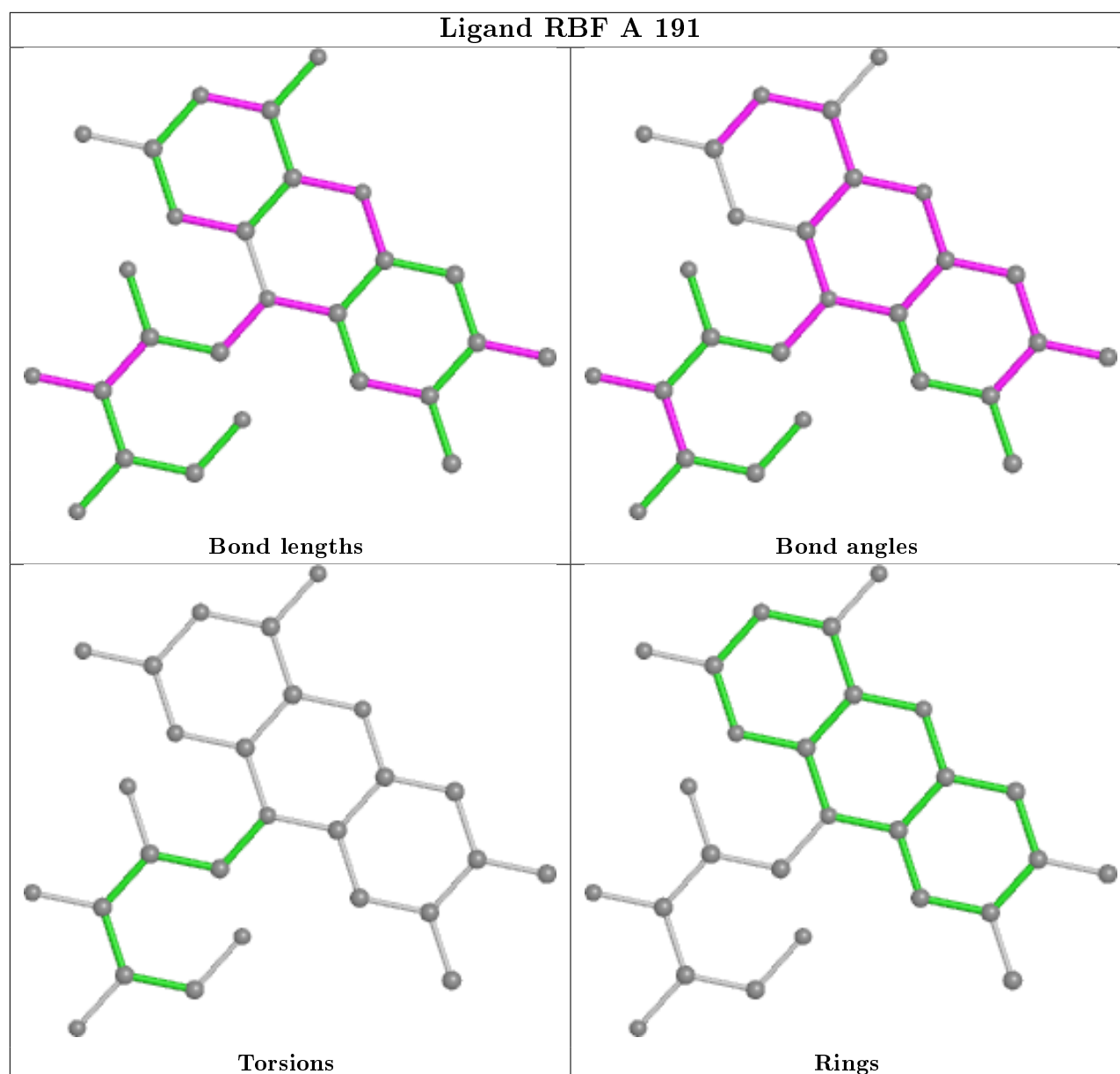
Bond angles



Torsions



Rings



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	184/190 (96%)	0.37	16 (8%)	10 8	10, 18, 45, 60	0
1	B	184/190 (96%)	0.41	18 (9%)	7 6	12, 19, 47, 68	0
All	All	368/380 (96%)	0.39	34 (9%)	9 8	10, 19, 46, 68	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	92	LEU	8.6
1	B	89	PHE	8.2
1	A	93	LEU	8.1
1	A	129	ILE	7.8
1	B	18	ASN	7.7
1	B	91	SER	7.6
1	B	165	ILE	6.5
1	B	94	GLY	5.5
1	B	93	LEU	5.3
1	A	184	ASN	4.9
1	A	127	ASP	4.6
1	A	130	GLU	4.6
1	B	90	GLY	4.5
1	A	131	ASN	4.5
1	A	126	LYS	4.2
1	A	135	GLU	3.7
1	B	163	LEU	3.6
1	B	19	ASP	3.4
1	B	184	ASN	3.4
1	A	19	ASP	3.3
1	A	94	GLY	3.1
1	B	20	ASP	2.9
1	B	162	ASN	2.7
1	B	21	THR	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	95	LYS	2.5
1	B	95	LYS	2.5
1	A	115	GLU	2.5
1	B	166	THR	2.5
1	A	132	ILE	2.4
1	A	183	LEU	2.3
1	B	88	GLU	2.2
1	A	116	ASP	2.1
1	A	43	LEU	2.1
1	B	140	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 carbohydrates in this entry.

6.4 Ligands [i](#)

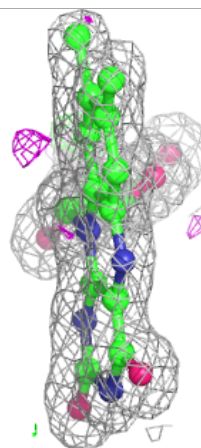
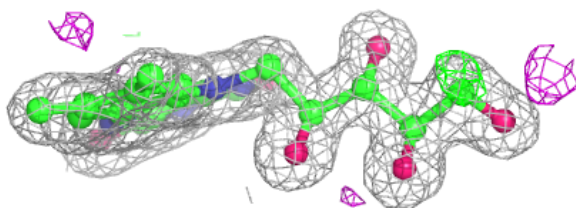
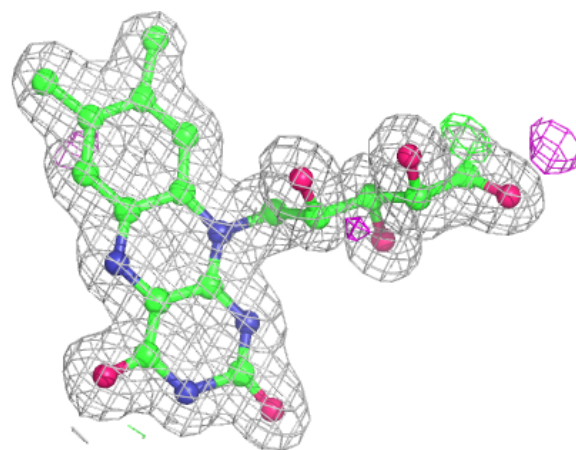
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

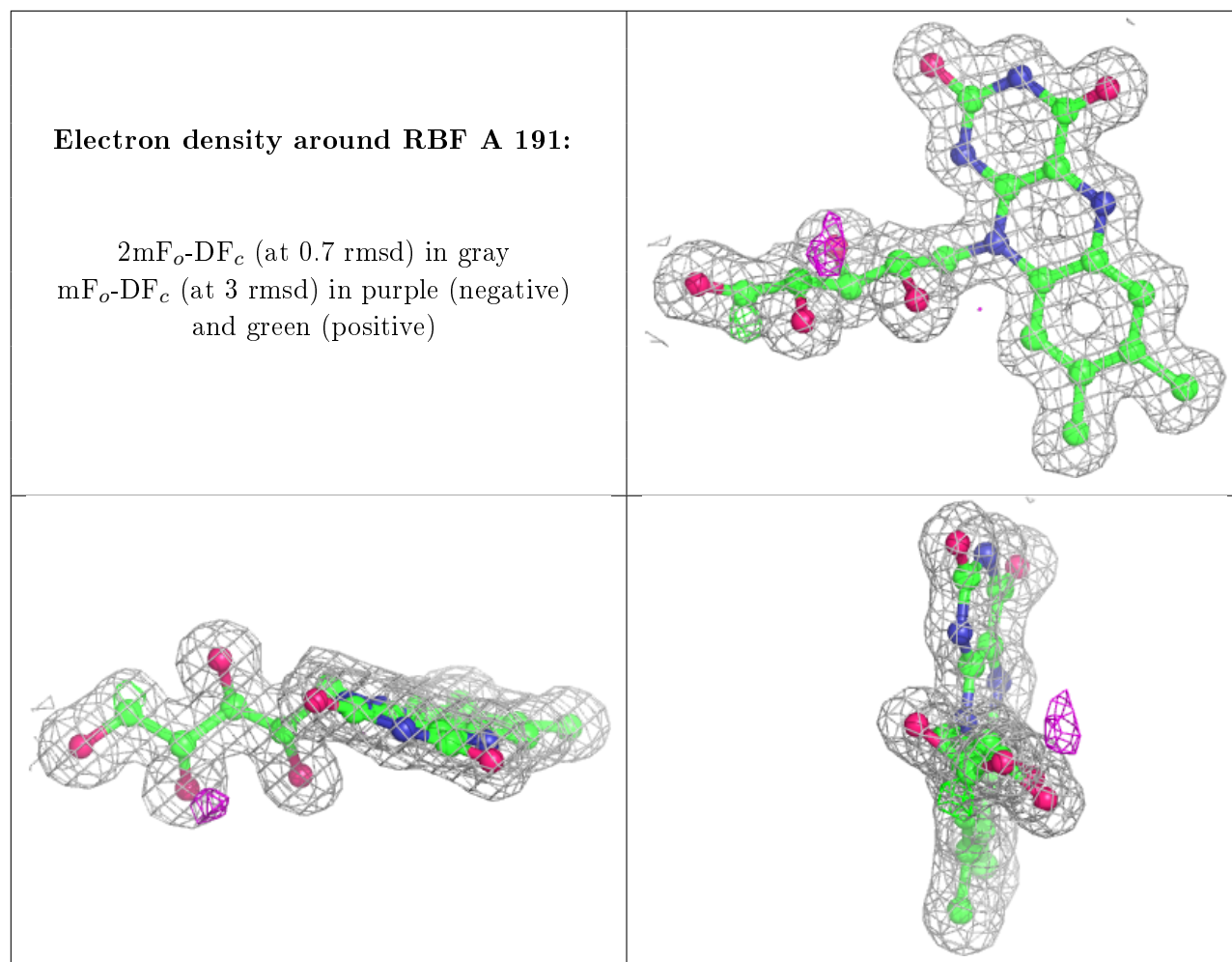
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	RBF	B	191	27/27	0.96	0.07	12,15,19,23	0
2	RBF	A	191	27/27	0.97	0.07	8,10,13,14	0

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 191:

$2mF_o-DF_c$ (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.