



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2023 – 11:54 PM EDT

PDB ID : 1OX5
Title : TOWARDS UNDERSTANDING THE MECHANISM OF THE COMPLEX CYCLIZATION REACTION CATALYZED BY IMIDAZOLE GLYCEROPHOSPHATE SYNTHASE
Authors : Chaudhuri, B.N.; Smith, J.L.
Deposited on : 2003-04-01
Resolution : 2.50 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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.35
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) : 2.35

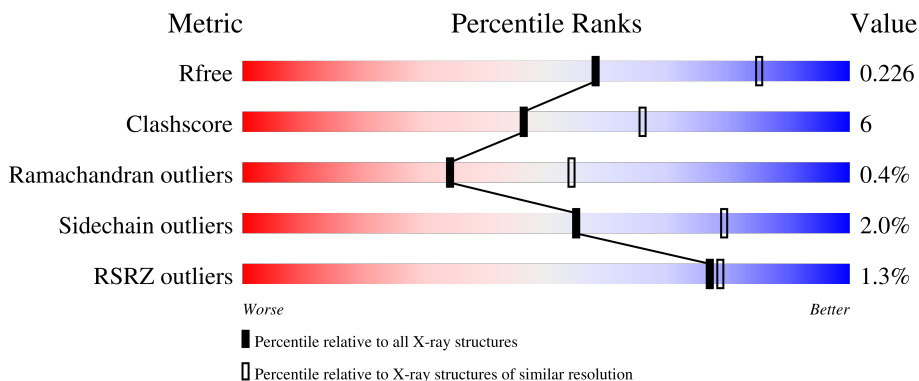
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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	555	 81% 13% ••
1	B	555	 82% 12% • 5%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
1	5CS	A	83	X	-	-	-
1	5CS	B	83	X	-	-	-
3	1PR	A	991	X	-	-	-
3	1PR	B	891	X	-	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 8528 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 Imidazole glycerol phosphate synthase hisHF.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	532	4146	2639	694	798	15	0	0	0
1	B	527	4111	2618	687	792	14	0	0	0

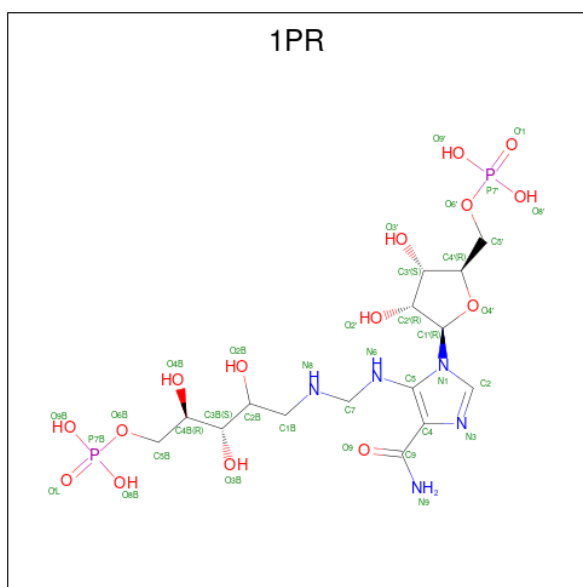
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	cloning artifact	UNP P33734
A	-2	SER	-	cloning artifact	UNP P33734
A	-1	HIS	-	cloning artifact	UNP P33734
A	83	5CS	CYS	modified residue	UNP P33734
B	-2	GLY	-	cloning artifact	UNP P33734
B	-1	SER	-	cloning artifact	UNP P33734
B	0	HIS	-	cloning artifact	UNP P33734
B	83	5CS	CYS	modified residue	UNP P33734

- Molecule 2 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ni	0	0
			1	1		

- Molecule 3 is PHOSPHORIC ACID MONO-[5-({[5-CARBAMOYL-3-(5-PHOSPHONOXY-5-DEOXY-RIBOFURANOSYL)-3H-IMIDAZOL-4-YLAMINO]-METHYL}-AMINO)-2,3,4-TRIHYDROXY-PENTYL] ESTER (three-letter code: 1PR) (formula: C₁₅H₂₉N₅O₁₅P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
3	A	1	Total 37	15	5	15	2	0	0
3	B	1	Total 37	15	5	15	2	0	0

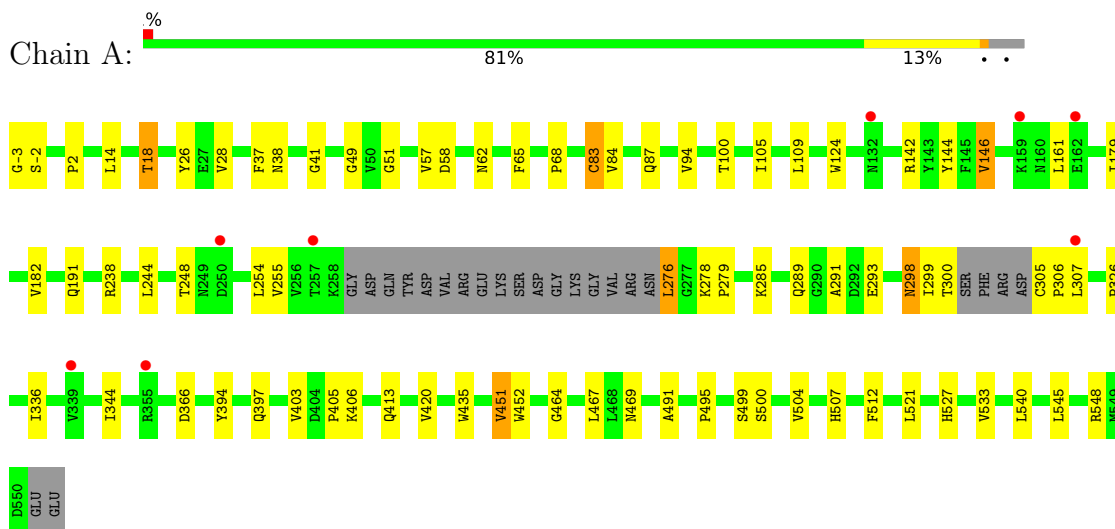
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	118	Total 118	118	0	0
4	B	78	Total 78	78	0	0

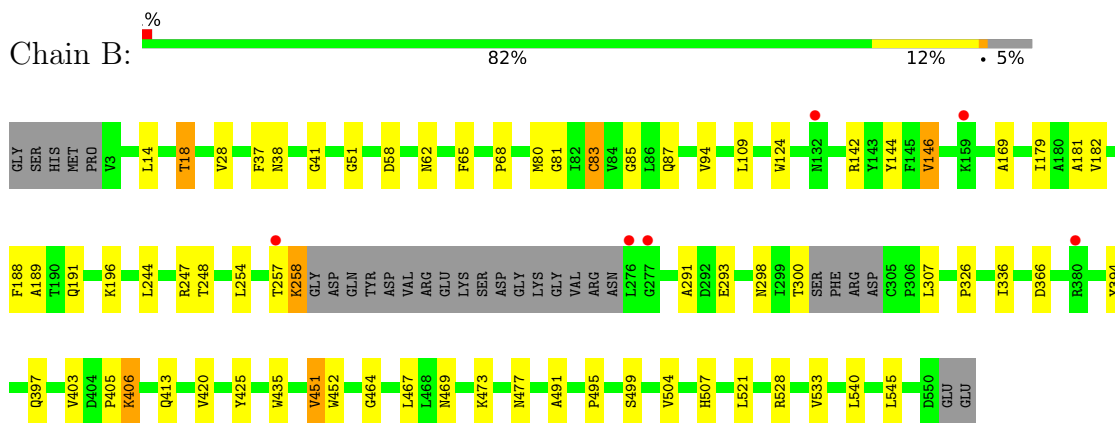
3 Residue-property plots [i](#)

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: Imidazole glycerol phosphate synthase hisHF



- Molecule 1: Imidazole glycerol phosphate synthase hisHF



4 Data and refinement statistics i

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	98.80Å 111.50Å 117.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.50 80.78 – 2.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.50) 99.9 (80.78-2.50)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.48 (at 2.51Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.223 , 0.246 0.218 , 0.226	Depositor DCC
R_{free} test set	2282 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	34.1	Xtrriage
Anisotropy	0.433	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 35.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.016 for -h,l,k	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8528	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.11% 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: 1PR, 5CS, NI

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	0.42	0/4216	0.66	0/5702
1	B	0.42	0/4179	0.67	0/5652
All	All	0.42	0/8395	0.66	0/11354

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	0
1	B	1	0
All	All	2	0

There are no bond length outliers.

There are no bond angle outliers.

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	83	5CS	CF
1	B	83	5CS	CF

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	4146	0	4091	55	0
1	B	4111	0	4059	48	0
2	A	1	0	0	0	0
3	A	37	0	22	1	0
3	B	37	0	22	1	0
4	A	118	0	0	1	0
4	B	78	0	0	2	0
All	All	8528	0	8194	104	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 6.

All (104) 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:293:GLU:HG3	1:A:326:PRO:HG2	1.58	0.85
1:B:293:GLU:HG3	1:B:326:PRO:HG2	1.60	0.81
1:B:477:ASN:HD21	1:B:528:ARG:NH2	1.83	0.76
1:A:298:ASN:HD22	1:A:299:ILE:N	1.82	0.76
1:B:257:THR:HG22	1:B:258:LYS:H	1.52	0.72
1:B:247:ARG:HE	1:B:257:THR:HB	1.56	0.70
1:B:477:ASN:HD21	1:B:528:ARG:HH21	1.38	0.70
1:B:51:GLY:HA2	1:B:394:TYR:O	1.94	0.67
1:A:298:ASN:HD22	1:A:299:ILE:H	1.42	0.66
1:B:257:THR:HG22	1:B:258:LYS:N	2.11	0.65
1:A:298:ASN:ND2	1:A:300:THR:H	1.95	0.64
1:A:298:ASN:C	1:A:299:ILE:HG13	2.18	0.64
1:B:403:VAL:HG12	1:B:405:PRO:HD3	1.80	0.63
1:A:403:VAL:HG12	1:A:405:PRO:HD3	1.81	0.63
1:B:473:LYS:NZ	4:B:938:HOH:O	2.31	0.62
1:B:146:VAL:HG22	1:B:397:GLN:HB2	1.82	0.62
1:B:467:LEU:HD11	1:B:521:LEU:HD13	1.82	0.61
1:A:83:5CS:HE2	1:A:191:GLN:HE22	1.66	0.59
1:A:57:VAL:HG11	1:A:100:THR:O	2.02	0.59
1:A:298:ASN:ND2	1:A:299:ILE:N	2.51	0.58
1:B:83:5CS:HE2	1:B:191:GLN:HE22	1.69	0.57
1:A:254:LEU:HD22	1:A:276:LEU:HB2	1.88	0.56
1:A:527:HIS:ND1	1:A:527:HIS:O	2.38	0.56
1:A:248:THR:HG23	1:A:300:THR:HB	1.86	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:255:VAL:HG12	1:A:278:LYS:HD2	1.87	0.55
1:B:80:MET:HA	1:B:188:PHE:O	2.08	0.54
1:B:540:LEU:HD22	1:B:545:LEU:HD12	1.89	0.54
1:A:18:THR:HG23	1:A:28:VAL:HG11	1.91	0.53
1:A:51:GLY:HA2	1:A:394:TYR:O	2.08	0.53
1:A:179:ILE:HG21	1:A:182:VAL:HG13	1.90	0.53
1:A:298:ASN:ND2	1:A:300:THR:HG23	2.23	0.53
1:A:540:LEU:HD22	1:A:545:LEU:HD12	1.90	0.53
1:A:467:LEU:HD11	1:A:521:LEU:HD13	1.90	0.53
1:A:307:LEU:HD22	1:A:336:ILE:HD11	1.91	0.53
1:B:406:LYS:HA	4:B:961:HOH:O	2.08	0.53
1:B:247:ARG:NE	1:B:257:THR:HB	2.24	0.52
1:B:179:ILE:HG21	1:B:182:VAL:HG13	1.91	0.52
1:B:307:LEU:HD22	1:B:336:ILE:HD11	1.91	0.52
1:A:58:ASP:O	1:A:62:ASN:HB2	2.10	0.51
1:B:38:ASN:ND2	1:B:41:GLY:H	2.09	0.51
1:A:65:PHE:C	1:A:68:PRO:HD2	2.31	0.51
1:B:65:PHE:C	1:B:68:PRO:HD2	2.31	0.50
1:A:38:ASN:ND2	1:A:41:GLY:H	2.10	0.50
1:B:452:TRP:CE3	1:B:491:ALA:HB2	2.47	0.50
1:B:18:THR:HG23	1:B:28:VAL:HG11	1.94	0.50
1:A:285:LYS:HE2	1:A:289:GLN:NE2	2.27	0.49
3:A:991:1PR:H72	3:A:991:1PR:H1'	1.94	0.49
1:A:18:THR:CG2	1:A:28:VAL:HG11	2.42	0.49
3:B:891:1PR:H1'	3:B:891:1PR:H72	1.95	0.49
1:A:144:TYR:CZ	1:A:326:PRO:HG3	2.47	0.49
1:B:18:THR:CG2	1:B:28:VAL:HG11	2.42	0.49
1:A:452:TRP:CE3	1:A:491:ALA:HB2	2.47	0.49
1:A:124:TRP:CE3	1:A:548:ARG:HD2	2.47	0.49
1:A:299:ILE:HG22	1:A:299:ILE:O	2.13	0.48
1:B:124:TRP:HB3	1:B:142:ARG:HB3	1.95	0.48
1:B:146:VAL:HG13	1:B:146:VAL:O	2.14	0.48
1:B:87:GLN:NE2	1:B:109:LEU:HD22	2.29	0.48
1:B:58:ASP:O	1:B:62:ASN:HB2	2.13	0.48
1:B:146:VAL:CG2	1:B:397:GLN:HB2	2.44	0.48
1:B:403:VAL:HG11	1:B:451:VAL:HG22	1.95	0.47
1:A:14:LEU:O	1:A:18:THR:OG1	2.32	0.47
1:B:248:THR:HG23	1:B:300:THR:HB	1.97	0.47
1:A:2:PRO:HB2	1:A:26:TYR:CD2	2.49	0.47
1:A:146:VAL:HG22	1:A:397:GLN:HB2	1.95	0.47
1:A:146:VAL:O	1:A:146:VAL:HG13	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:403:VAL:HG11	1:A:451:VAL:HG22	1.96	0.46
1:B:469:ASN:ND2	1:B:499:SER:OG	2.48	0.46
1:B:254:LEU:HD12	1:B:254:LEU:N	2.31	0.46
1:A:124:TRP:HB3	1:A:142:ARG:HB3	1.96	0.46
1:B:81:GLY:O	1:B:189:ALA:HA	2.16	0.45
1:A:298:ASN:C	1:A:299:ILE:CG1	2.83	0.45
1:A:298:ASN:HD21	1:A:300:THR:HG23	1.81	0.45
1:A:464:GLY:O	1:A:495:PRO:HD2	2.17	0.45
1:B:420:VAL:HG12	1:B:435:TRP:HB3	1.99	0.45
1:B:521:LEU:C	1:B:521:LEU:HD23	2.38	0.45
1:A:521:LEU:C	1:A:521:LEU:HD23	2.37	0.44
1:A:87:GLN:NE2	1:A:109:LEU:HD22	2.31	0.44
1:B:413:GLN:OE1	1:B:420:VAL:HG22	2.17	0.44
1:B:464:GLY:O	1:B:495:PRO:HD2	2.18	0.44
1:B:14:LEU:O	1:B:18:THR:OG1	2.36	0.43
1:A:49:GLY:HA3	1:A:84:VAL:HB	2.01	0.43
1:B:477:ASN:ND2	1:B:528:ARG:NH2	2.59	0.43
1:B:247:ARG:HD2	1:B:257:THR:OG1	2.19	0.42
1:A:-3:GLY:HA3	1:B:425:TYR:CE2	2.55	0.42
1:A:504:VAL:HG22	1:A:507:HIS:CE1	2.54	0.42
1:B:504:VAL:HG22	1:B:507:HIS:CE1	2.55	0.42
1:A:305:CYS:HA	1:A:306:PRO:HD3	1.92	0.42
1:A:278:LYS:N	1:A:279:PRO:HD2	2.34	0.42
1:B:257:THR:CG2	1:B:258:LYS:H	2.27	0.42
1:A:105:ILE:HG23	1:A:161:LEU:HD21	2.02	0.41
1:B:196:LYS:HD2	1:B:326:PRO:HA	2.02	0.41
1:A:254:LEU:HB2	1:A:276:LEU:HA	2.01	0.41
1:B:83:5CS:C	1:B:85:GLY:N	2.83	0.41
1:A:420:VAL:HG12	1:A:435:TRP:HB3	2.01	0.41
1:A:469:ASN:ND2	1:A:499:SER:OG	2.54	0.41
1:A:238:ARG:HB3	1:A:512:PHE:CE2	2.56	0.40
1:A:326:PRO:HB2	4:A:1065:HOH:O	2.20	0.40
1:A:291:ALA:HA	1:A:533:VAL:HG11	2.03	0.40
1:B:169:ALA:HB3	1:B:181:ALA:HB3	2.04	0.40
1:A:307:LEU:HD23	1:A:344:ILE:HD13	2.04	0.40
1:A:179:ILE:HG21	1:A:182:VAL:CG1	2.51	0.40
1:A:413:GLN:OE1	1:A:420:VAL:HG22	2.21	0.40
1:B:144:TYR:CZ	1:B:326:PRO:HG3	2.57	0.40
1:B:291:ALA:HA	1:B:533:VAL:HG11	2.03	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	525/555 (95%)	504 (96%)	18 (3%)	3 (1%)	25	43
1	B	520/555 (94%)	501 (96%)	18 (4%)	1 (0%)	47	68
All	All	1045/1110 (94%)	1005 (96%)	36 (3%)	4 (0%)	34	54

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	-2	SER
1	A	500	SER
1	A	146	VAL
1	B	146	VAL

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	446/466 (96%)	437 (98%)	9 (2%)	55	79
1	B	442/466 (95%)	433 (98%)	9 (2%)	55	79
All	All	888/932 (95%)	870 (98%)	18 (2%)	55	79

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	THR
1	A	37	PHE

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Mol	Chain	Res	Type
1	A	94	VAL
1	A	244	LEU
1	A	276	LEU
1	A	298	ASN
1	A	366	ASP
1	A	406	LYS
1	A	451	VAL
1	B	18	THR
1	B	37	PHE
1	B	94	VAL
1	B	244	LEU
1	B	258	LYS
1	B	298	ASN
1	B	366	ASP
1	B	406	LYS
1	B	451	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	ASN
1	A	87	GLN
1	A	191	GLN
1	A	298	ASN
1	A	469	ASN
1	B	38	ASN
1	B	55	HIS
1	B	87	GLN
1	B	191	GLN
1	B	298	ASN
1	B	469	ASN
1	B	527	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
1	5CS	B	83	1	10,16,17	1.84	2 (20%)	10,21,23	4.52	5 (50%)
1	5CS	A	83	1	10,16,17	1.74	2 (20%)	10,21,23	4.51	5 (50%)

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
1	5CS	B	83	1	1/1/5/7	3/11/24/26	0/1/1/1
1	5CS	A	83	1	1/1/5/7	3/11/24/26	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	83	5CS	OF-CF	-4.07	1.38	1.44
1	A	83	5CS	OF-CF	-4.05	1.38	1.44
1	B	83	5CS	CB-SG	-3.31	1.78	1.82
1	A	83	5CS	CB-SG	-2.67	1.79	1.82

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	83	5CS	OF-CF-CE	8.87	113.15	103.56
1	A	83	5CS	OF-CF-CE	8.82	113.11	103.56
1	B	83	5CS	CD-CE-CF	-6.85	97.41	107.78
1	A	83	5CS	CD-CE-CF	-6.84	97.42	107.78
1	B	83	5CS	OF-CF-CI	6.45	121.63	109.23
1	A	83	5CS	OF-CF-CI	6.31	121.36	109.23
1	A	83	5CS	CE-CF-CI	5.57	124.80	115.08
1	B	83	5CS	CE-CF-CI	5.38	124.47	115.08
1	A	83	5CS	CA-CB-SG	2.70	123.24	113.58
1	B	83	5CS	CA-CB-SG	2.68	123.17	113.58

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	83	5CS	CF
1	B	83	5CS	CF

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	83	5CS	OF-CF-CI-CJ
1	A	83	5CS	NI-CI-CJ-OJ1
1	B	83	5CS	NI-CI-CJ-OJ1
1	A	83	5CS	OF-CF-CI-CJ
1	A	83	5CS	CA-CB-SG-CD
1	B	83	5CS	CA-CB-SG-CD

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	83	5CS	2	0
1	A	83	5CS	1	0

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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	1PR	B	891	-	32,38,38	2.85	9 (28%)	42,56,56	2.23	13 (30%)
3	1PR	A	991	-	32,38,38	2.84	9 (28%)	42,56,56	2.30	11 (26%)

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	1PR	B	891	-	1/1/10/12	13/24/51/51	0/2/2/2
3	1PR	A	991	-	1/1/10/12	11/24/51/51	0/2/2/2

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	891	1PR	O2B-C2B	-8.39	1.25	1.43
3	A	991	1PR	O2B-C2B	-8.31	1.25	1.43
3	A	991	1PR	O3'-C3'	-7.47	1.25	1.43
3	B	891	1PR	O3'-C3'	-7.15	1.26	1.43
3	B	891	1PR	C7-N6	-6.85	1.33	1.45
3	A	991	1PR	C7-N6	-6.41	1.33	1.45
3	B	891	1PR	C7-N8	-4.43	1.33	1.45
3	A	991	1PR	C7-N8	-4.21	1.34	1.45
3	B	891	1PR	C9-N9	3.98	1.40	1.33
3	A	991	1PR	C9-N9	3.95	1.40	1.33
3	B	891	1PR	O6B-C5B	-3.20	1.32	1.44
3	A	991	1PR	C2-N3	-3.11	1.29	1.35
3	A	991	1PR	O6B-C5B	-3.04	1.33	1.44
3	B	891	1PR	P7B-O6B	-3.01	1.50	1.60
3	B	891	1PR	C2-N3	-2.82	1.30	1.35
3	A	991	1PR	P7B-O6B	-2.74	1.51	1.60
3	A	991	1PR	C4-C9	-2.50	1.44	1.51
3	B	891	1PR	C4-C9	-2.30	1.44	1.51

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	991	1PR	C2B-C1B-N8	7.36	122.06	111.97
3	B	891	1PR	C2B-C1B-N8	6.69	121.15	111.97
3	A	991	1PR	C7-N8-C1B	5.14	127.32	112.86
3	A	991	1PR	O2B-C2B-C1B	4.89	119.56	109.69
3	B	891	1PR	C7-N8-C1B	4.71	126.12	112.86
3	B	891	1PR	O2B-C2B-C1B	4.62	119.01	109.69
3	B	891	1PR	O2B-C2B-C3B	4.24	119.42	109.10
3	A	991	1PR	O2B-C2B-C3B	4.17	119.23	109.10
3	B	891	1PR	O9-C9-C4	3.85	129.13	119.89
3	A	991	1PR	O9-C9-C4	3.84	129.12	119.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	891	1PR	C1B-C2B-C3B	3.36	121.58	110.93
3	B	891	1PR	O9B-P7B-O6B	-3.30	97.95	106.73
3	A	991	1PR	O9B-P7B-O6B	-3.27	98.04	106.73
3	A	991	1PR	P7B-O6B-C5B	3.25	127.26	118.30
3	A	991	1PR	C1B-C2B-C3B	3.24	121.21	110.93
3	B	891	1PR	P7B-O6B-C5B	2.91	126.30	118.30
3	A	991	1PR	C4-C9-N9	-2.86	111.24	115.75
3	B	891	1PR	C4-C9-N9	-2.69	111.51	115.75
3	A	991	1PR	C5B-C4B-C3B	-2.36	107.64	112.20
3	B	891	1PR	C5B-C4B-C3B	-2.36	107.64	112.20
3	B	891	1PR	O9-C9-N9	-2.27	119.35	122.58
3	B	891	1PR	C3'-C2'-C1'	2.08	104.11	100.98
3	A	991	1PR	O9-C9-N9	-2.06	119.65	122.58
3	B	891	1PR	O9'-P7'-O6'	2.01	112.08	106.73

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	991	1PR	C2B
3	B	891	1PR	C2B

All (24) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	991	1PR	C5'-O6'-P7'-O9'
3	A	991	1PR	C5'-O6'-P7'-O8'
3	A	991	1PR	C5'-O6'-P7'-O'1
3	A	991	1PR	N8-C7-N6-C5
3	A	991	1PR	N8-C1B-C2B-O2B
3	A	991	1PR	C2B-C1B-N8-C7
3	A	991	1PR	C1B-C2B-C3B-C4B
3	A	991	1PR	C1B-C2B-C3B-O3B
3	A	991	1PR	C3B-C4B-C5B-O6B
3	A	991	1PR	O4B-C4B-C5B-O6B
3	B	891	1PR	C5'-O6'-P7'-O9'
3	B	891	1PR	C5'-O6'-P7'-O8'
3	B	891	1PR	C5'-O6'-P7'-O'1
3	B	891	1PR	N8-C7-N6-C5
3	B	891	1PR	N8-C1B-C2B-O2B
3	B	891	1PR	C2B-C1B-N8-C7
3	B	891	1PR	C1B-C2B-C3B-C4B
3	B	891	1PR	C1B-C2B-C3B-O3B

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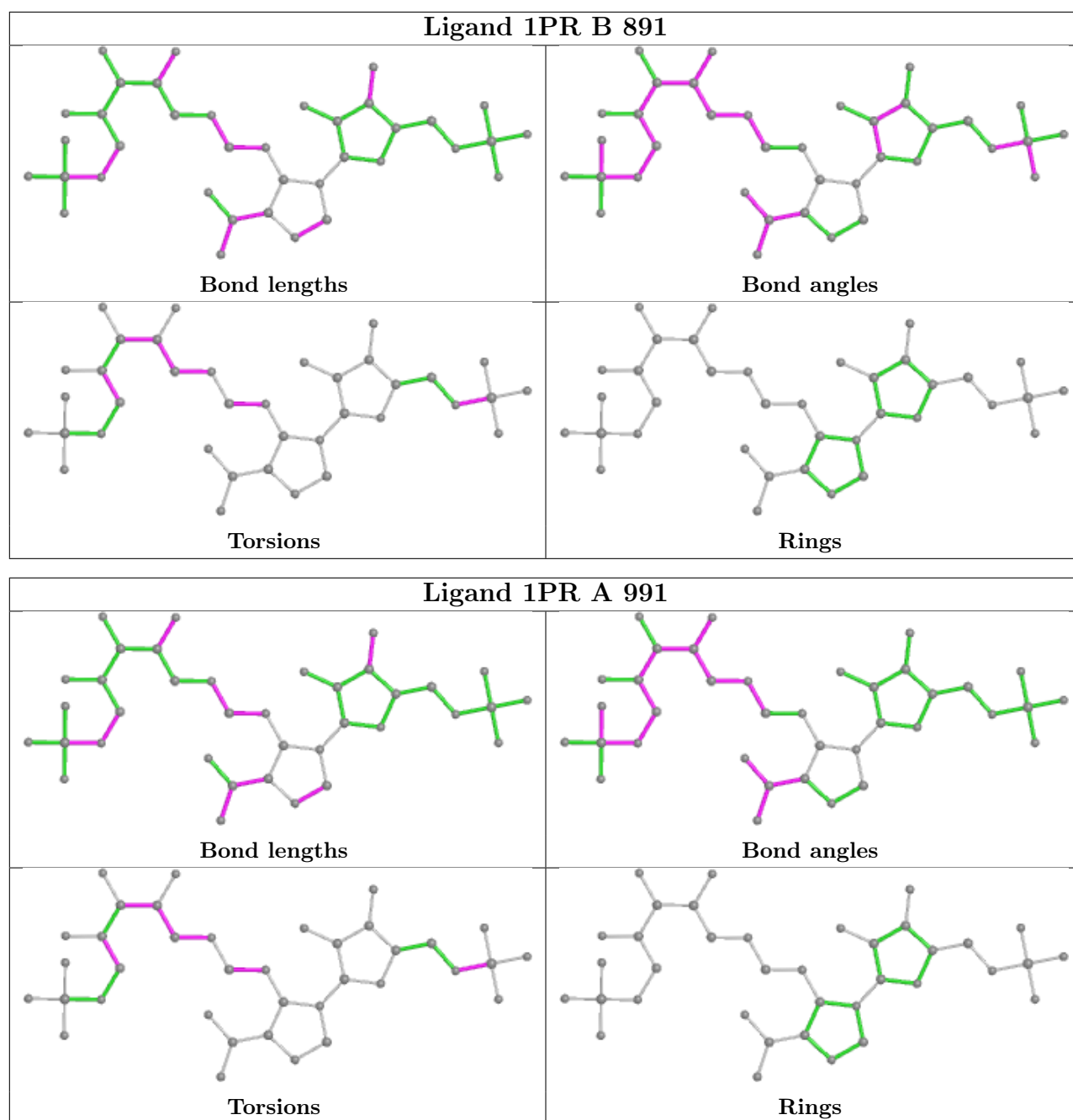
Mol	Chain	Res	Type	Atoms
3	B	891	1PR	C3B-C4B-C5B-O6B
3	B	891	1PR	O4B-C4B-C5B-O6B
3	A	991	1PR	N8-C1B-C2B-C3B
3	B	891	1PR	N8-C1B-C2B-C3B
3	B	891	1PR	O2B-C2B-C3B-O3B
3	B	891	1PR	O2B-C2B-C3B-C4B

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	891	1PR	1	0
3	A	991	1PR	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	531/555 (95%)	0.02	8 (1%) 73 75	15, 30, 54, 73	0
1	B	526/555 (94%)	-0.01	6 (1%) 80 82	16, 32, 54, 74	0
All	All	1057/1110 (95%)	0.00	14 (1%) 77 79	15, 31, 54, 74	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	276	LEU	4.5
1	A	162	GLU	2.7
1	B	380	ARG	2.7
1	B	132	ASN	2.7
1	A	307	LEU	2.7
1	A	257	THR	2.6
1	A	132	ASN	2.6
1	A	159	LYS	2.4
1	A	355	ARG	2.2
1	B	277	GLY	2.2
1	A	250	ASP	2.2
1	B	257	THR	2.2
1	B	159	LYS	2.1
1	A	339	VAL	2.1

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(\AA^2)	Q<0.9
1	5CS	B	83	16/17	0.85	0.25	30,52,56,57	0
1	5CS	A	83	16/17	0.91	0.27	29,53,56,57	0

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

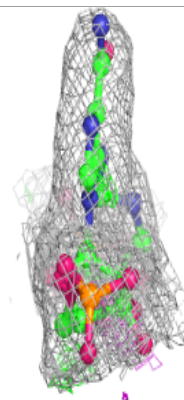
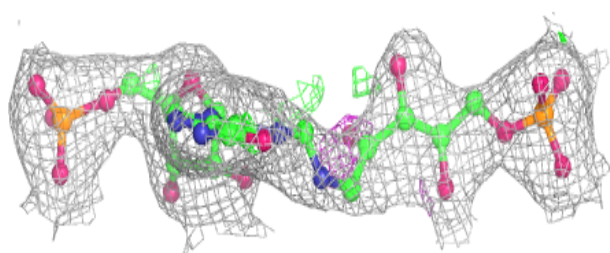
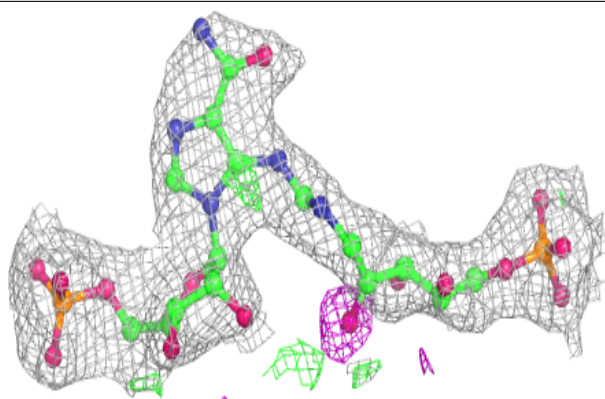
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	1PR	A	991	37/37	0.94	0.16	30,43,48,53	0
3	1PR	B	891	37/37	0.96	0.18	27,42,50,51	0
2	NI	A	902	1/1	0.99	0.13	26,26,26,26	0

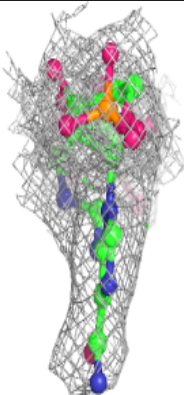
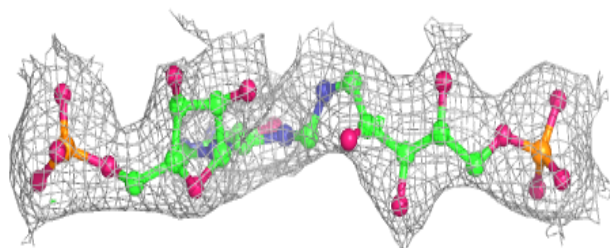
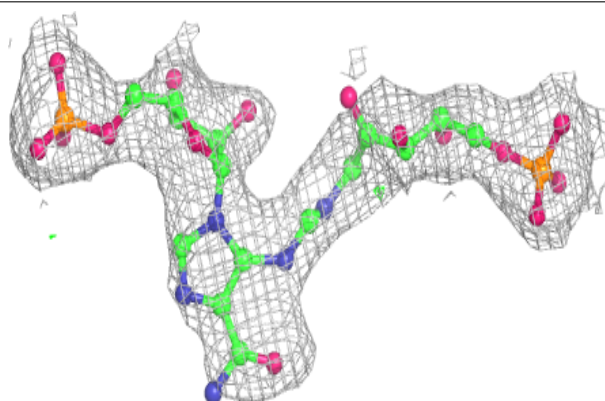
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 1PR A 991:

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

**Electron density around 1PR B 891:**

$2mF_o-DF_c$ (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.