Full wwPDB X-ray Structure Validation Report (i)

Mar 5, 2026 – 01:52 PM UTC

PDB ID	:	7C92 / pdb_00007c92
Title	:	Structure of Serratia marcescens chitinase B complexed with compound 6k
Authors	:	Jiang, X.; Yuan, P.T.
Deposited on	:	2020-06-04
Resolution	:	2.32 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 (i) 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 (i)) were used in the production of this report:

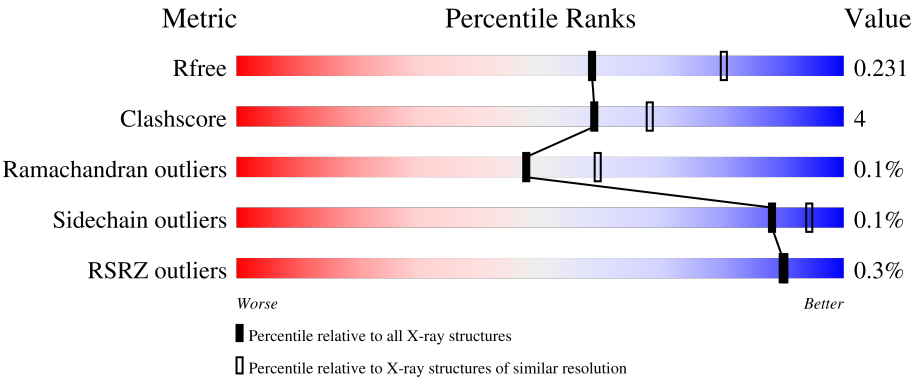
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal 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 (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.32 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	7754 (2.34-2.30)
Clashscore	190562	8383 (2.34-2.30)
Ramachandran outliers	187476	8303 (2.34-2.30)
Sidechain outliers	187428	8303 (2.34-2.30)
RSRZ outliers	180081	7760 (2.34-2.30)

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 <=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	509	% 88% 10% ·
1	B	509	87% 10% ·

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 8325 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 Chitinase.

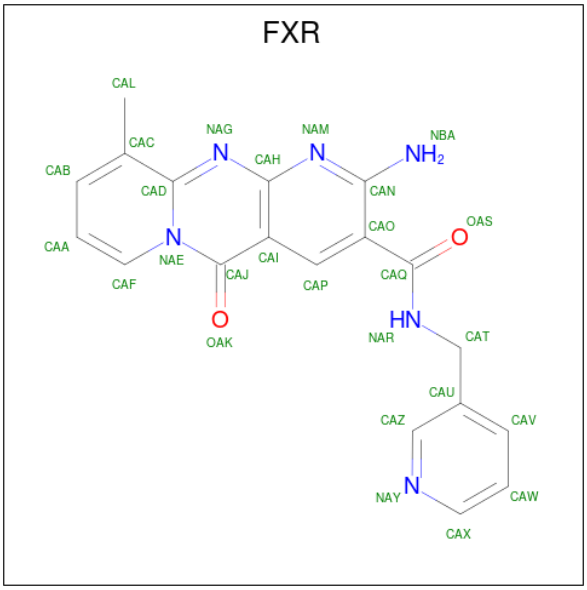
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	495	Total	C	N	O	S	0	0	0
			3889	2488	654	733	14			
1	B	495	Total	C	N	O	S	0	0	0
			3889	2488	654	733	14			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	ASP	-	expression tag	UNP Q54276
A	-2	PRO	-	expression tag	UNP Q54276
A	-1	SER	-	expression tag	UNP Q54276
A	0	SER	-	expression tag	UNP Q54276
A	1	ARG	-	expression tag	UNP Q54276
A	500	HIS	-	expression tag	UNP Q54276
A	501	HIS	-	expression tag	UNP Q54276
A	502	HIS	-	expression tag	UNP Q54276
A	503	HIS	-	expression tag	UNP Q54276
A	504	HIS	-	expression tag	UNP Q54276
A	505	HIS	-	expression tag	UNP Q54276
B	-3	ASP	-	expression tag	UNP Q54276
B	-2	PRO	-	expression tag	UNP Q54276
B	-1	SER	-	expression tag	UNP Q54276
B	0	SER	-	expression tag	UNP Q54276
B	1	ARG	-	expression tag	UNP Q54276
B	500	HIS	-	expression tag	UNP Q54276
B	501	HIS	-	expression tag	UNP Q54276
B	502	HIS	-	expression tag	UNP Q54276
B	503	HIS	-	expression tag	UNP Q54276
B	504	HIS	-	expression tag	UNP Q54276
B	505	HIS	-	expression tag	UNP Q54276

- Molecule 2 is 6-azanyl-11-methyl-2-oxidanylidene-N-(pyridin-3-ylmethyl)-1,7,9-triazatricyclo[8.4.0.0^3,8]tetradeca-3(8),4,6,9,11,13-hexaene-5-carboxamide (CCD ID: FXR) (formula:

$C_{19}H_{16}N_6O_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C N O 54 38 12 4	0	1
2	B	1	Total C N O 27 19 6 2	0	0
2	B	1	Total C N O 27 19 6 2	0	0

- Molecule 3 is water.

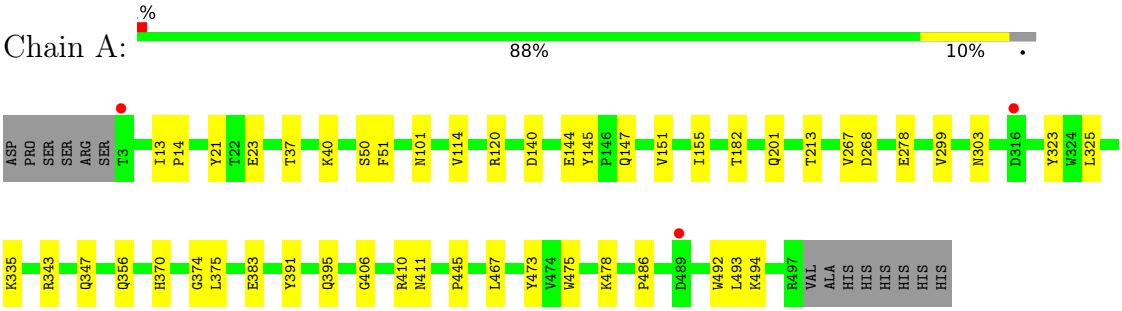
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	225	Total O 225 225	0	0
3	B	214	Total O 214 214	0	0

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.

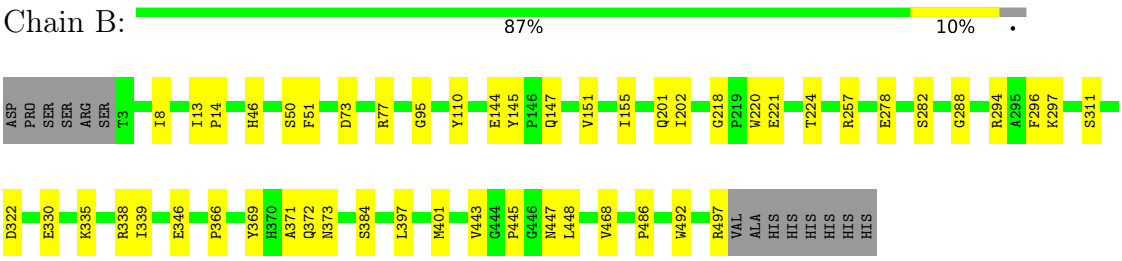
• Molecule 1: Chitinase

Chain A:



• Molecule 1: Chitinase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	57.36Å 103.64Å 186.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.88 – 2.32 48.88 – 2.32	Depositor EDS
% Data completeness (in resolution range)	93.6 (48.88-2.32) 93.8 (48.88-2.32)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.10 (at 2.32Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.173 , 0.230 0.173 , 0.231	Depositor DCC
R_{free} test set	2000 reflections (4.05%)	wwPDB-VP
Wilson B-factor (Å^2)	33.7	Xtriage
Anisotropy	0.264	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 37.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	8325	wwPDB-VP
Average B, all atoms (Å^2)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.94% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: FXR

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	0.32	0/4000	0.51	0/5453
1	B	0.33	0/4000	0.50	0/5453
All	All	0.33	0/8000	0.50	0/10906

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

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	3889	0	3716	34	0
1	B	3889	0	3716	33	0
2	A	54	0	0	1	0
2	B	54	0	0	0	0
3	A	225	0	0	3	0
3	B	214	0	0	3	0
All	All	8325	0	7432	67	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (67) 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:347:GLN:HE22	1:A:410:ARG:HH21	1.28	0.80
1:B:278:GLU:HG3	1:B:445:PRO:HB2	1.69	0.74
1:A:478:LYS:HE2	1:A:493:LEU:HB2	1.72	0.70
1:B:151:VAL:O	1:B:155:ILE:HG12	1.96	0.66
1:A:155:ILE:HD13	1:A:201:GLN:OE1	1.96	0.65
1:B:155:ILE:HD11	1:B:202:ILE:HG13	1.80	0.62
1:A:278:GLU:HG3	1:A:445:PRO:HB2	1.81	0.62
1:A:335:LYS:NZ	3:A:702:HOH:O	2.27	0.60
1:B:335:LYS:NZ	3:B:703:HOH:O	2.30	0.60
1:B:448:LEU:O	1:B:497:ARG:NH1	2.35	0.59
1:B:155:ILE:CD1	1:B:202:ILE:HG13	2.33	0.59
1:B:257:ARG:NH1	3:B:702:HOH:O	2.29	0.55
1:A:144:GLU:HA	1:A:145:TYR:CG	2.41	0.54
1:A:356:GLN:OE1	3:A:701:HOH:O	2.18	0.54
1:B:294:ARG:NH1	1:B:339:ILE:HD11	2.23	0.53
1:A:343:ARG:HH11	1:A:410:ARG:HG3	1.73	0.53
1:B:330:GLU:OE2	3:B:701:HOH:O	2.18	0.53
1:B:144:GLU:HA	1:B:145:TYR:CG	2.44	0.53
1:A:213:THR:HB	1:A:267:VAL:HG22	1.91	0.52
1:A:323:TYR:OH	1:A:335:LYS:HE3	2.10	0.52
1:A:114:VAL:HA	1:A:120:ARG:HG2	1.92	0.51
1:A:383:GLU:HG2	3:A:905:HOH:O	2.08	0.51
1:B:486:PRO:HA	1:B:492:TRP:CD1	2.46	0.51
1:A:151:VAL:HG12	1:A:155:ILE:HD11	1.93	0.50
1:A:343:ARG:NH2	1:A:406:GLY:O	2.44	0.49
1:A:21:TYR:OH	1:A:23:GLU:OE2	2.29	0.48
1:A:475:TRP:CD2	1:A:486:PRO:HB3	2.49	0.48
1:B:50:SER:HB2	1:B:51:PHE:CG	2.49	0.48
1:A:299:VAL:HG22	1:A:374:GLY:C	2.39	0.47
1:B:338:ARG:C	1:B:339:ILE:HD12	2.40	0.47
1:A:37:THR:OG1	1:A:40:LYS:HG3	2.14	0.47
1:A:268:ASP:OD1	1:A:391:TYR:OH	2.25	0.47
1:B:297:LYS:HE2	1:B:311:SER:O	2.15	0.47
1:A:101:ASN:HA	1:A:144:GLU:O	2.14	0.46
1:B:95:GLY:HA2	1:B:110:TYR:OH	2.15	0.46
1:B:155:ILE:HG13	1:B:201:GLN:OE1	2.15	0.46
2:A:601[A]:FXR:OAS	2:A:601[A]:FXR:NBA	2.47	0.46
1:A:467:LEU:HA	1:A:475:TRP:O	2.16	0.46
1:A:473:TYR:CD1	1:A:494:LYS:HD3	2.51	0.45
1:B:147:GLN:O	1:B:151:VAL:HG23	2.17	0.45
1:A:50:SER:HB2	1:A:51:PHE:CG	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:8:ILE:HG12	1:B:46:HIS:HB2	1.99	0.44
1:B:443:VAL:HA	1:B:447:ASN:HD21	1.83	0.43
1:B:288:GLY:HA2	1:B:401:MET:O	2.19	0.43
1:B:369:TYR:CZ	1:B:371:ALA:HA	2.54	0.43
1:A:50:SER:HA	1:A:51:PHE:HA	1.74	0.43
1:B:50:SER:HA	1:B:51:PHE:HA	1.79	0.42
1:A:325:LEU:HD12	1:A:325:LEU:HA	1.86	0.42
1:A:140:ASP:HA	1:A:182:THR:O	2.20	0.42
1:B:218:GLY:HA3	1:B:220:TRP:CH2	2.55	0.42
1:B:144:GLU:HG2	1:B:145:TYR:CZ	2.54	0.42
1:B:372:GLN:HG2	1:B:373:ASN:OD1	2.18	0.42
1:B:282:SER:O	1:B:397:LEU:HA	2.20	0.42
1:A:151:VAL:HG12	1:A:155:ILE:CD1	2.50	0.41
1:B:73:ASP:O	1:B:77:ARG:HG3	2.20	0.41
1:A:370:HIS:HB3	1:A:375:LEU:HB2	2.02	0.41
1:A:147:GLN:O	1:A:151:VAL:HG23	2.20	0.41
1:B:220:TRP:CD1	1:B:221:GLU:HG2	2.56	0.41
1:B:366:PRO:HD3	1:B:384:SER:HB3	2.03	0.41
1:A:411:ASN:HD21	1:B:346:GLU:CD	2.29	0.41
1:A:486:PRO:HG3	1:A:492:TRP:CD2	2.56	0.41
1:B:224:THR:HG22	1:B:296:PHE:CD2	2.55	0.41
1:B:50:SER:HB2	1:B:51:PHE:CD1	2.56	0.41
1:B:13:ILE:HA	1:B:14:PRO:HD3	1.96	0.40
1:A:13:ILE:HA	1:A:14:PRO:HD3	1.94	0.40
1:A:303:ASN:HD22	1:A:303:ASN:HA	1.68	0.40
1:A:391:TYR:CE1	1:A:395:GLN:HG3	2.56	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	493/509 (97%)	480 (97%)	13 (3%)	0	100 100
1	B	493/509 (97%)	481 (98%)	11 (2%)	1 (0%)	43 53
All	All	986/1018 (97%)	961 (98%)	24 (2%)	1 (0%)	48 59

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	322	ASP

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	401/415 (97%)	401 (100%)	0	100 100
1	B	401/415 (97%)	400 (100%)	1 (0%)	87 94
All	All	802/830 (97%)	801 (100%)	1 (0%)	88 94

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	468	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	ASN
1	A	57	ASN
1	A	84	HIS
1	A	147	GLN
1	A	303	ASN
1	A	347	GLN
1	A	423	ASN
1	B	272	GLN
1	B	394	GLN

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Mol	Chain	Res	Type
1	B	464	GLN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

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

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

4 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	FXR	A	601[B]	-	29,30,30	2.91	9 (31%)	33,43,43	2.63	12 (36%)
2	FXR	A	601[A]	-	29,30,30	2.91	8 (27%)	33,43,43	2.93	12 (36%)
2	FXR	B	601	-	29,30,30	2.89	9 (31%)	33,43,43	1.83	7 (21%)
2	FXR	B	602	-	29,30,30	2.80	9 (31%)	33,43,43	2.02	9 (27%)

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	FXR	A	601[B]	-	-	6/9/9/9	0/4/4/4
2	FXR	A	601[A]	-	-	0/9/9/9	0/4/4/4
2	FXR	B	601	-	-	0/9/9/9	0/4/4/4
2	FXR	B	602	-	-	1/9/9/9	0/4/4/4

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	602	FXR	OAK-CAJ	9.36	1.41	1.22
2	A	601[A]	FXR	OAK-CAJ	9.28	1.41	1.22
2	A	601[B]	FXR	OAK-CAJ	9.24	1.41	1.22
2	B	601	FXR	OAK-CAJ	9.23	1.41	1.22
2	A	601[B]	FXR	CAQ-NAR	6.43	1.47	1.33
2	B	601	FXR	CAQ-NAR	6.28	1.47	1.33
2	A	601[B]	FXR	CAD-NAG	6.26	1.38	1.30
2	A	601[A]	FXR	CAQ-NAR	6.20	1.47	1.33
2	B	602	FXR	CAQ-NAR	6.05	1.47	1.33
2	B	601	FXR	CAD-NAG	6.00	1.38	1.30
2	A	601[A]	FXR	CAD-NAG	5.85	1.38	1.30
2	B	602	FXR	CAD-NAG	5.37	1.37	1.30
2	B	602	FXR	CAN-NBA	4.95	1.46	1.34
2	A	601[B]	FXR	CAN-NBA	4.78	1.46	1.34
2	A	601[A]	FXR	CAN-NBA	4.75	1.46	1.34
2	B	601	FXR	CAN-NBA	4.58	1.45	1.34
2	A	601[A]	FXR	CAD-NAE	-3.94	1.31	1.40
2	A	601[B]	FXR	CAD-NAE	-3.92	1.31	1.40
2	B	602	FXR	CAD-NAE	-3.81	1.31	1.40
2	B	601	FXR	CAD-NAE	-3.77	1.31	1.40
2	A	601[A]	FXR	CAP-CAI	3.53	1.45	1.39
2	B	601	FXR	CAP-CAI	2.96	1.44	1.39
2	A	601[A]	FXR	CAO-CAQ	2.69	1.55	1.50
2	A	601[B]	FXR	CAO-CAQ	2.60	1.55	1.50
2	B	601	FXR	CAH-NAM	2.59	1.38	1.34
2	A	601[B]	FXR	CAH-NAM	2.50	1.38	1.34
2	A	601[B]	FXR	CAP-CAI	2.45	1.43	1.39
2	B	602	FXR	CAF-CAA	2.40	1.40	1.35
2	B	601	FXR	CAO-CAQ	2.39	1.55	1.50
2	A	601[A]	FXR	CAF-CAA	2.38	1.40	1.35
2	B	601	FXR	CAF-CAA	2.33	1.40	1.35
2	A	601[B]	FXR	CAF-CAA	2.26	1.40	1.35
2	B	602	FXR	CAH-NAM	2.11	1.37	1.34
2	B	602	FXR	CAO-CAQ	2.04	1.54	1.50
2	B	602	FXR	CAP-CAI	2.01	1.43	1.39

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	601[A]	FXR	CAN-CAO-CAQ	-10.88	116.70	120.66
2	A	601[B]	FXR	CAN-CAO-CAQ	9.50	124.11	120.66
2	A	601[A]	FXR	CAI-CAJ-NAE	6.34	118.78	114.20
2	B	602	FXR	CAN-CAO-CAQ	5.79	122.76	120.66
2	A	601[B]	FXR	CAI-CAJ-NAE	5.67	118.29	114.20
2	B	601	FXR	CAI-CAJ-NAE	5.56	118.22	114.20
2	B	602	FXR	CAI-CAJ-NAE	4.73	117.61	114.20
2	A	601[A]	FXR	CAO-CAN-NBA	-4.46	118.98	122.52
2	A	601[B]	FXR	CAO-CAQ-NAR	4.22	125.52	117.32
2	B	601	FXR	CAO-CAN-NBA	-4.06	119.30	122.52
2	A	601[B]	FXR	CAU-CAT-NAR	-3.78	105.11	113.07
2	A	601[A]	FXR	CAA-CAB-CAC	-3.72	117.79	122.03
2	B	602	FXR	OAK-CAJ-NAE	3.69	122.37	119.93
2	A	601[A]	FXR	CAX-NAY-CAZ	3.28	122.60	116.85
2	A	601[B]	FXR	CAA-CAB-CAC	-3.20	118.39	122.03
2	B	602	FXR	CAA-CAB-CAC	-3.19	118.40	122.03
2	A	601[A]	FXR	CAP-CAI-CAJ	3.10	124.52	118.37
2	B	601	FXR	CAA-CAB-CAC	-2.97	118.65	122.03
2	A	601[A]	FXR	CAO-CAQ-NAR	2.95	123.06	117.32
2	B	602	FXR	CAC-CAD-NAE	2.93	121.25	115.92
2	A	601[A]	FXR	NAE-CAD-NAG	-2.87	120.78	123.65
2	B	601	FXR	CAX-NAY-CAZ	2.87	121.88	116.85
2	A	601[B]	FXR	CAO-CAN-NBA	-2.84	120.27	122.52
2	A	601[A]	FXR	CAP-CAO-CAQ	2.80	125.16	117.21
2	A	601[A]	FXR	CAC-CAD-NAE	2.62	120.69	115.92
2	A	601[A]	FXR	CAU-CAT-NAR	2.59	118.53	113.07
2	B	602	FXR	CAO-CAN-NBA	-2.58	120.47	122.52
2	B	601	FXR	CAN-CAO-CAQ	-2.47	119.76	120.66
2	A	601[B]	FXR	OAS-CAQ-NAR	-2.44	117.83	122.59
2	B	601	FXR	CAC-CAD-NAE	2.43	120.35	115.92
2	A	601[B]	FXR	OAS-CAQ-CAO	-2.41	116.62	121.03
2	A	601[B]	FXR	CAX-NAY-CAZ	2.37	121.01	116.85
2	A	601[B]	FXR	NAE-CAD-NAG	-2.34	121.31	123.65
2	B	602	FXR	CAX-NAY-CAZ	2.32	120.92	116.85
2	A	601[A]	FXR	OAS-CAQ-NAR	-2.28	118.14	122.59
2	B	602	FXR	CAT-NAR-CAQ	2.28	127.22	121.89
2	A	601[B]	FXR	CAP-CAI-CAJ	2.19	122.72	118.37
2	B	601	FXR	CAP-CAI-CAJ	2.15	122.64	118.37
2	A	601[B]	FXR	CAC-CAD-NAE	2.15	119.84	115.92
2	B	602	FXR	CAV-CAU-CAZ	2.12	120.21	117.10

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	602	FXR	CAU-CAT-NAR-CAQ
2	A	601[B]	FXR	OAS-CAQ-NAR-CAT
2	A	601[B]	FXR	CAO-CAQ-NAR-CAT
2	A	601[B]	FXR	CAN-CAO-CAQ-OAS
2	A	601[B]	FXR	CAP-CAO-CAQ-OAS
2	A	601[B]	FXR	CAP-CAO-CAQ-NAR
2	A	601[B]	FXR	CAN-CAO-CAQ-NAR

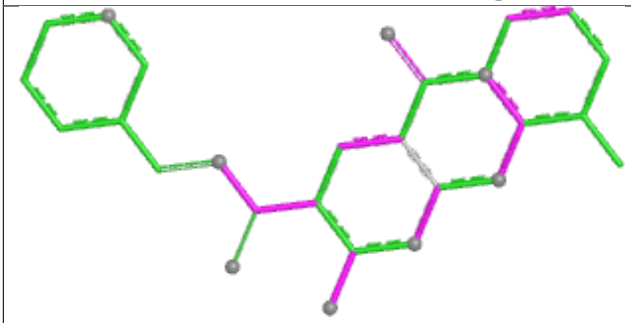
There are no ring outliers.

1 monomer is involved in 1 short contact:

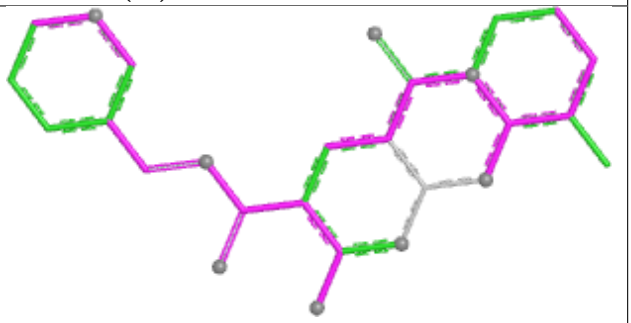
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601[A]	FXR	1	0

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 then 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.

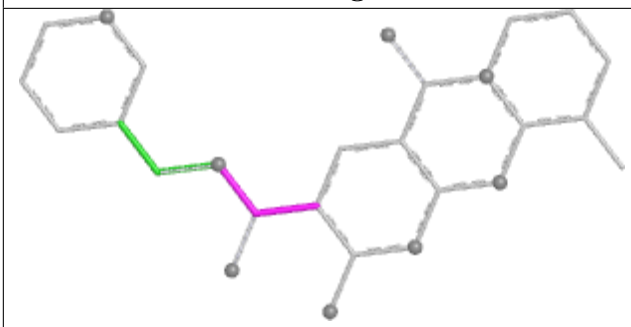
Ligand FXR A 601 (B)



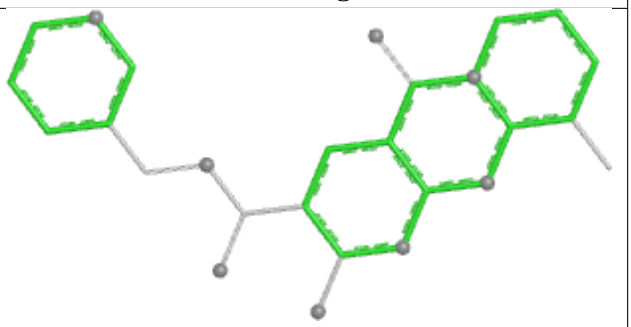
Bond lengths



Bond angles

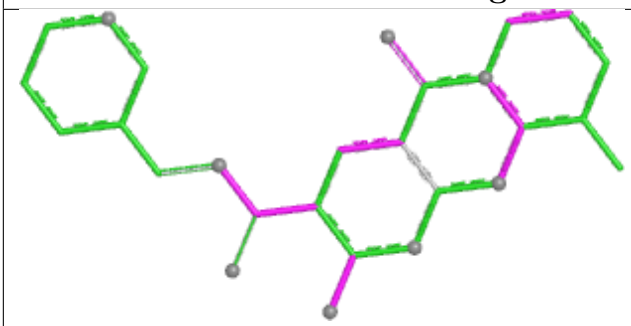


Torsions

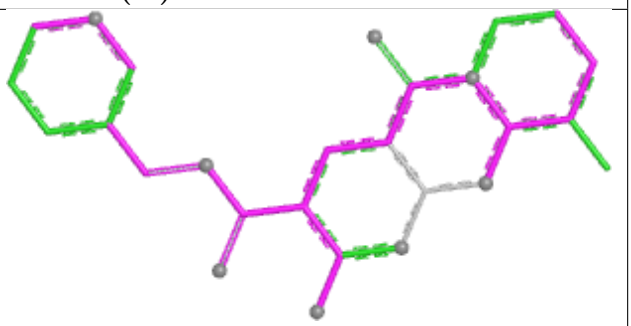


Rings

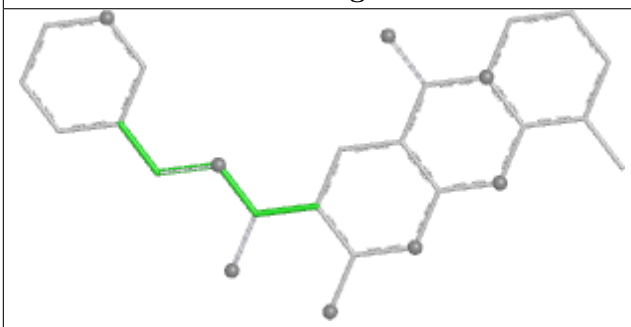
Ligand FXR A 601 (A)



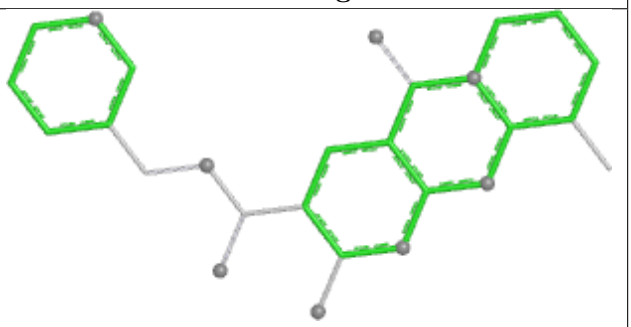
Bond lengths



Bond angles

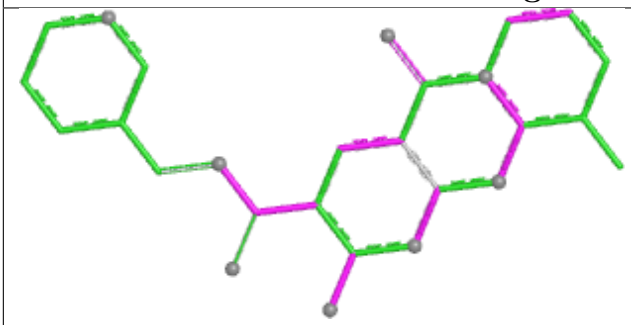


Torsions

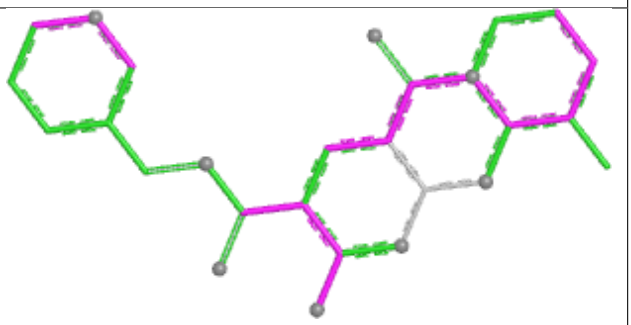


Rings

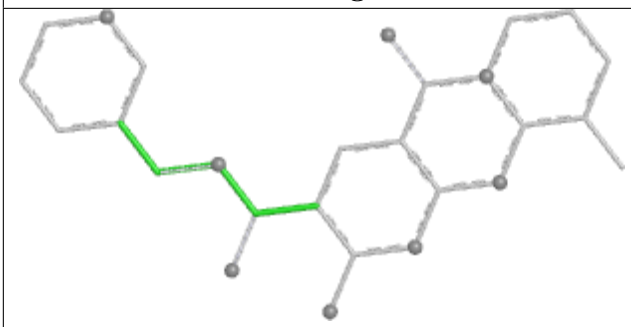
Ligand FXR B 601



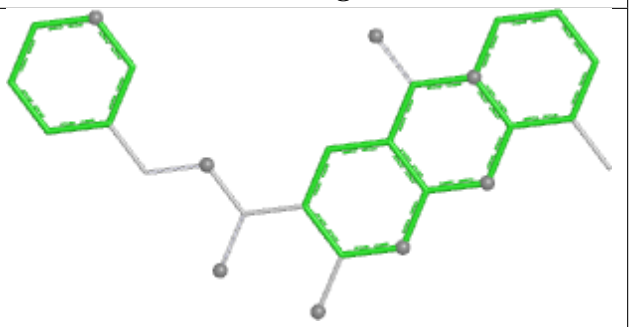
Bond lengths



Bond angles

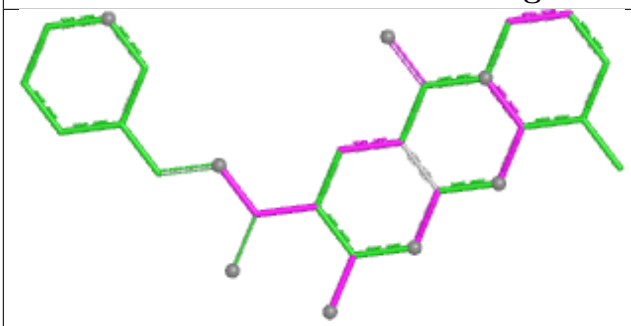


Torsions

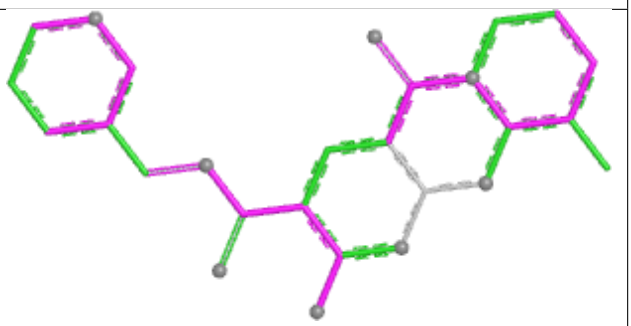


Rings

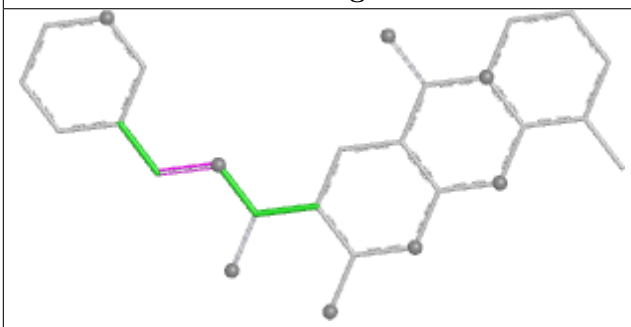
Ligand FXR B 602



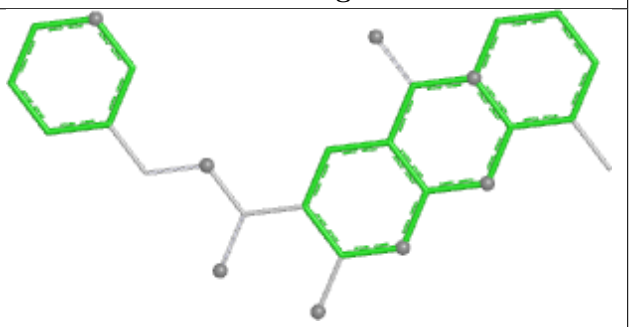
Bond lengths



Bond angles



Torsions



Rings

5.7 Other polymers

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, 95^{th} 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	495/509 (97%)	-0.40	3 (0%) 85 86	22, 32, 46, 58	0
1	B	495/509 (97%)	-0.35	0 100 100	23, 33, 48, 66	0
All	All	990/1018 (97%)	-0.38	3 (0%) 90 90	22, 33, 47, 66	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	3	THR	3.3
1	A	316	ASP	2.4
1	A	489	ASP	2.2

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

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

6.3 Carbohydrates

There are no oligosaccharides in this entry.

6.4 Ligands

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} 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	FXR	A	601[A]	27/27	0.59	0.24	30,50,57,59	27
2	FXR	A	601[B]	27/27	0.59	0.24	30,50,57,59	27

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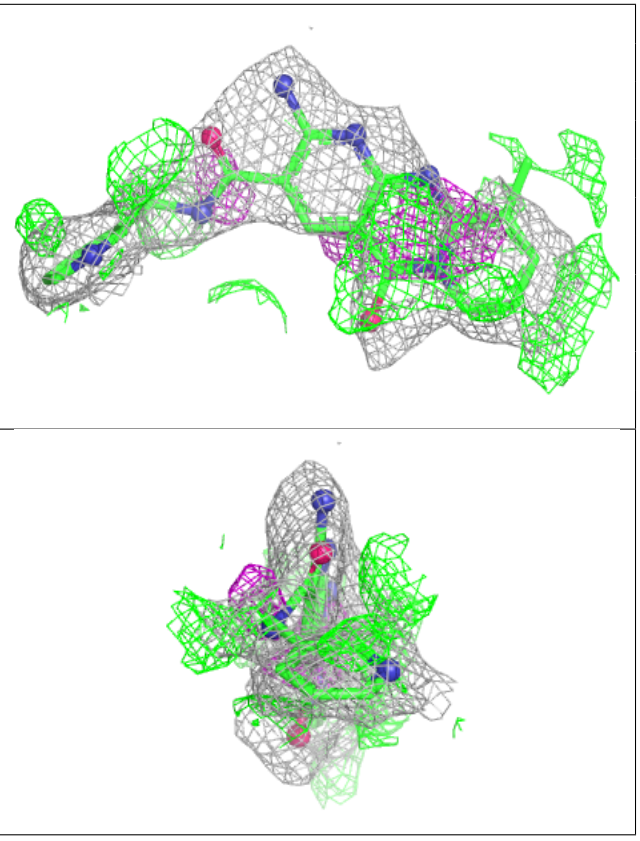
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	FXR	B	602	27/27	0.75	0.16	46,53,62,63	0
2	FXR	B	601	27/27	0.88	0.10	34,41,45,48	0

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.

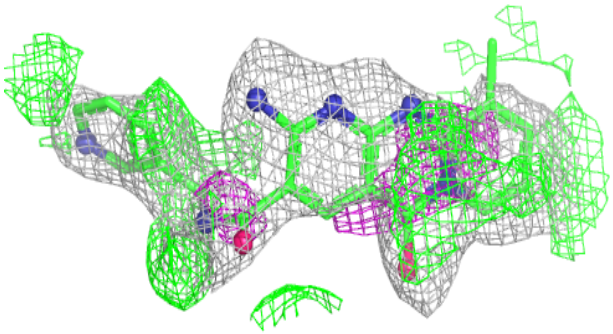
Electron density around FXR A 601 (A):

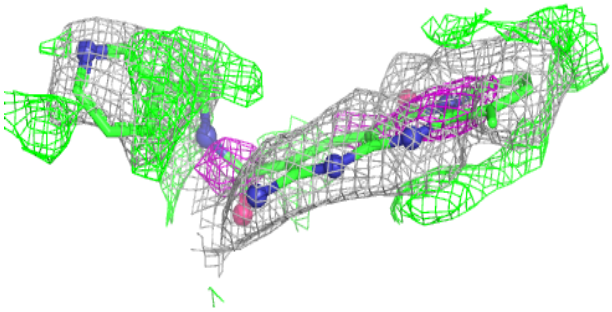
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

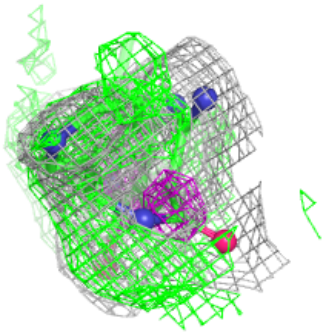


Electron density around FXR A 601 (B):

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

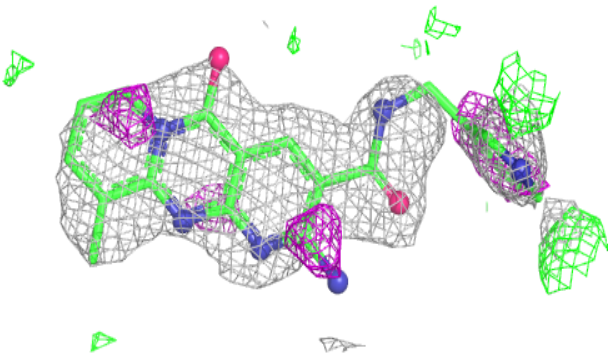


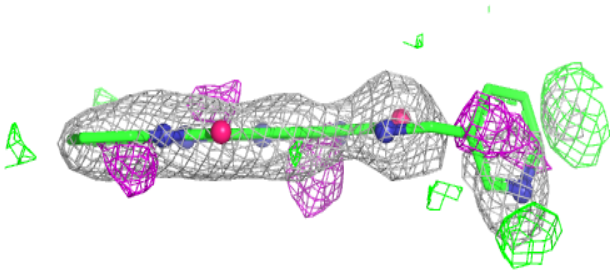


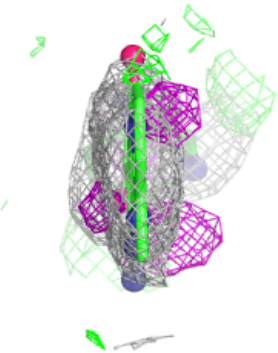


Electron density around FXR B 602:

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

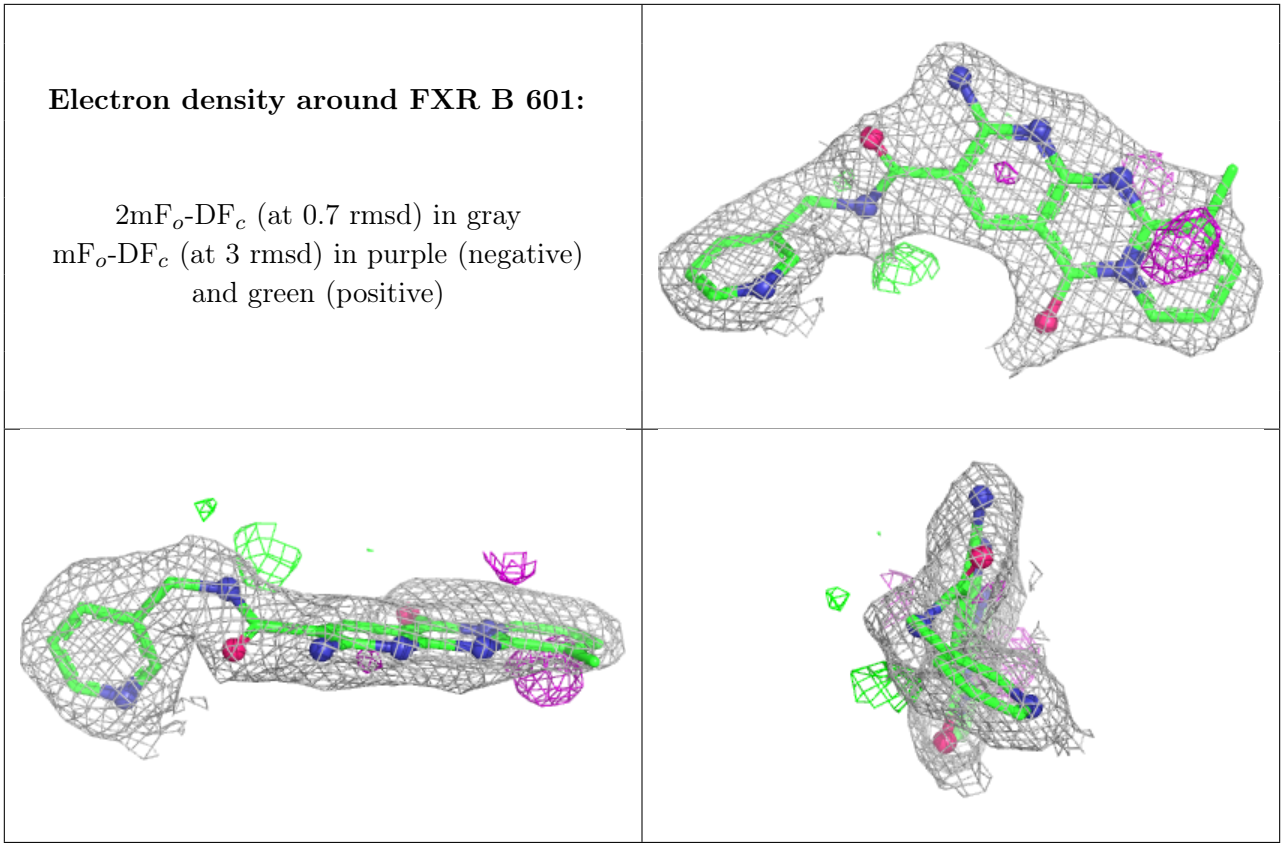






Electron density around FXR B 601:

$2mF_o$-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

There are no such residues in this entry.