

wwPDB X-ray Structure Validation Summary Report (i)

Oct 9, 2023 – 11:15 PM EDT

PDB ID : 7N9O

Title: Estrogen Receptor Alpha Ligand Binding Domain in Complex with Aliphatic

SERD S-C10(15)

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Deposited on : 2021-06-18

Resolution : 2.00 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (1)) were used in the production of this report:

 $Mol Probity \quad : \quad 4.02b\text{--}467$

Mogul : 1.8.5 (274361), CSD as541be (2020)

 $Xtriage\ (Phenix) \quad : \quad 1.13$

EDS : 2.35.1buster-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

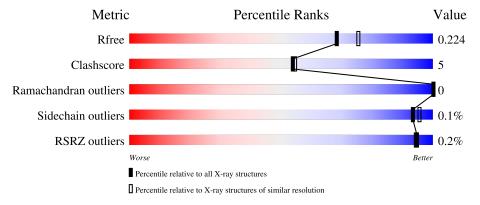


1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.00 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	Similar resolution $(\#\text{Entries, resolution range}(\text{\AA}))$
R_{free}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	249	88%	8%	.
1	В	249	83%	12%	6%
1	С	249	87%	9%	.
1	D	249	84%	11%	5%



2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 8242 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 Estrogen receptor.

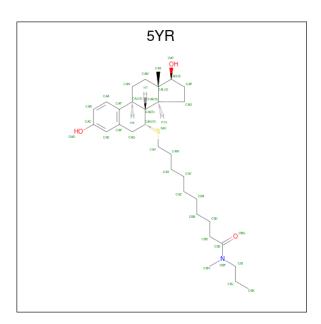
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	239	Total	С	N	О	S	0	2	0
1	A	. 239	1907	1222	325	344	16	U	2	
1	В	225	Total	С	N	О	S	0	0	0
1	Б	235	1862	1197	312	337	16	U	0	
1	С	240	Total	С	N	О	S	0	2	0
1		240	1919	1230	327	346	16	U	2	
1	D	237	Total	С	N	О	S	0	0	0
1	ע	231	1863	1197	312	338	16	U		

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	381	SER	CYS	engineered mutation	UNP P03372
A	417	SER	CYS	engineered mutation	UNP P03372
A	530	SER	CYS	engineered mutation	UNP P03372
A	536	SER	LEU	engineered mutation	UNP P03372
В	381	SER	CYS	engineered mutation	UNP P03372
В	417	SER	CYS	engineered mutation	UNP P03372
В	530	SER	CYS	engineered mutation	UNP P03372
В	536	SER	LEU	engineered mutation	UNP P03372
С	381	SER	CYS	engineered mutation	UNP P03372
С	417	SER	CYS	engineered mutation	UNP P03372
С	530	SER	CYS	engineered mutation	UNP P03372
С	536	SER	LEU	engineered mutation	UNP P03372
D	381	SER	CYS	engineered mutation	UNP P03372
D	417	SER	CYS	engineered mutation	UNP P03372
D	530	SER	CYS	engineered mutation	UNP P03372
D	536	SER	LEU	engineered mutation	UNP P03372

• Molecule 2 is 10-{[3,17beta-dihydroxyestra-1,3,5(10)-trien-7beta-yl]sulfanyl}-N-methyl-N-propyldecanamide (three-letter code: 5YR) (formula: $C_{32}H_{51}NO_3S$) (labeled as "Ligand of Interest" by depositor).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
9	Λ	1	Total	С	N	О	S	0	0
	Λ	1	37	32	1	3	1	0	
2	D	1	Total	С	N	О	S	0	0
	Ъ	1	37	32	1	3	1	0	
9	C	1	Total	С	N	О	S	0	0
	C	1	37	32	1	3	1	0	
9	D	1	Total	С	N	О	S	0	0
	D	1	37	32	1	3	1	0	

• Molecule 3 is water.

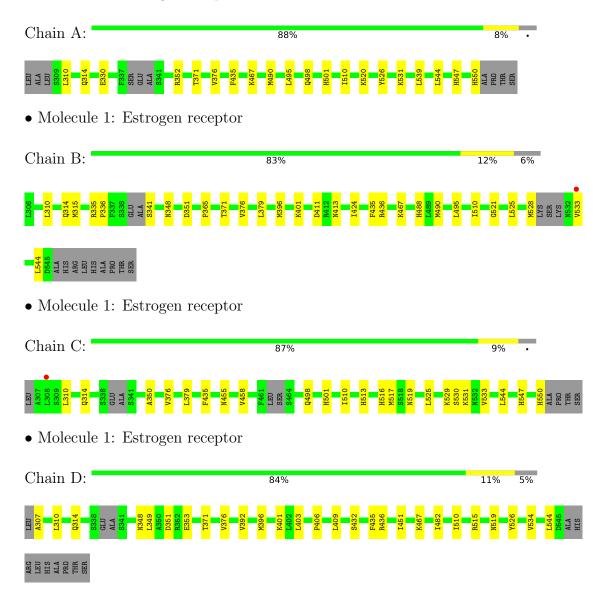
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	134	Total O 134 134	0	0
3	В	154	Total O 154 154	0	0
3	С	130	Total O 130 130	0	0
3	D	125	Total O 125 125	0	0



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: Estrogen receptor





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32	Depositor
Cell constants	58.16Å 58.16Å 274.39Å	Donositon
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.54 - 2.00	Depositor
rtesolution (A)	49.54 - 2.00	EDS
% Data completeness	94.5 (49.54-2.00)	Depositor
(in resolution range)	94.5 (49.54-2.00)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	2.62 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.182 , 0.225	Depositor
	0.183 , 0.224	DCC
R_{free} test set	3346 reflections $(5.06%)$	wwPDB-VP
Wilson B-factor (Å ²)	23.8	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.35, 31.1	EDS
L-test for twinning ²	$< L > = 0.51, < L^2> = 0.34$	Xtriage
	0.487 for -h,-k,l	
Estimated twinning fraction	0.488 for h,-h-k,-l	Xtriage
	0.487 for -k,-h,-l	
F_o, F_c correlation	0.96	EDS
Total number of atoms	8242	wwPDB-VP
Average B, all atoms (\mathring{A}^2)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.59% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



¹Intensities estimated from amplitudes.

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: 5YR

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		
IVIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.38	0/1944	0.54	0/2628	
1	В	0.40	0/1896	0.56	0/2563	
1	С	0.38	0/1955	0.56	0/2641	
1	D	0.38	0/1898	0.53	0/2568	
All	All	0.39	0/7693	0.55	0/10400	

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1907	0	1931	11	0
1	В	1862	0	1894	22	0
1	С	1919	0	1950	15	0
1	D	1863	0	1887	16	0
2	A	37	0	0	6	0
2	В	37	0	0	10	0
2	С	37	0	0	6	0
2	D	37	0	0	5	0
3	A	134	0	0	0	0

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\mathbf{N}	Iol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
	3	В	154	0	0	4	0
	3	С	130	0	0	0	0
	3	D	125	0	0	1	0
A	All	All	8242	0	7662	82	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 5.

The worst 5 of 82 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$egin{aligned} \operatorname{Clash} \ \operatorname{overlap}\ (\mathring{\mathbf{A}}) \end{aligned}$
2:A:601:5YR:CAJ	2:A:601:5YR:CAN	1.81	1.58
2:D:601:5YR:CAJ	2:D:601:5YR:CAN	1.81	1.58
2:C:601:5YR:CAJ	2:C:601:5YR:CAN	1.81	1.57
2:B:601:5YR:CAN	2:B:601:5YR:CAJ	1.81	1.55
1:B:533:VAL:HG13	2:B:601:5YR:CAX	1.82	1.08

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	Favoured Allowed		Perce	ntiles
1	A	$236/249 \ (95\%)$	230 (98%)	6 (2%)	0	100	100
1	В	229/249 (92%)	228 (100%)	1 (0%)	0	100	100
1	\mathbf{C}	235/249~(94%)	231 (98%)	4 (2%)	0	100	100
1	D	233/249 (94%)	230 (99%)	3 (1%)	0	100	100
All	All	933/996 (94%)	919 (98%)	14 (2%)	0	100	100

There are no Ramachandran outliers to report.



5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Rotameric Outliers		Percentiles		
1	A	213/224~(95%)	213 (100%)	0	100	100		
1	В	208/224 (93%)	207 (100%)	1 (0%)	88	92		
1	С	215/224~(96%)	215 (100%)	0	100	100		
1	D	207/224~(92%)	207 (100%)	0	100	100		
All	All	843/896 (94%)	842 (100%)	1 (0%)	93	95		

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

Mol	Chain	Res	Type	
1	В	528	MET	

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
1	В	519	ASN
1	С	519	ASN
1	D	519	ASN
1	D	513	HIS
1	В	513	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)

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



5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

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	Clasia	Res	Link	Bond lengths			Bond angles			
MIOI	Mol Type Chain			Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2	
2	5YR	В	601	-	38,40,40	5.69	18 (47%)	49,56,56	2.25	19 (38%)
2	5YR	D	601	-	38,40,40	5.78	18 (47%)	49,56,56	2.15	14 (28%)
2	5YR	A	601	-	38,40,40	5.72	17 (44%)	49,56,56	2.03	14 (28%)
2	5YR	С	601	-	38,40,40	5.74	18 (47%)	49,56,56	2.14	16 (32%)

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	5YR	В	601	-	-	7/20/63/63	0/4/4/4
2	5YR	D	601	-	-	7/20/63/63	0/4/4/4
2	5YR	A	601	-	-	5/20/63/63	0/4/4/4
2	5YR	С	601	-	-	6/20/63/63	0/4/4/4

The worst 5 of 71 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	Observed(A)	Ideal(Å)
2	A	601	5YR	CAN-CAJ	20.59	1.81	1.53
2	В	601	5YR	CAN-CAJ	20.47	1.81	1.53
2	D	601	5YR	CAN-CAJ	20.42	1.81	1.53
2	С	601	5YR	CAN-CAJ	20.27	1.81	1.53
2	В	601	5YR	CAF-CAJ	-16.23	1.28	1.52



The worst	5	of	63	bond	angle	outliers	are	listed	below:
110 WOID	$\overline{}$	O.	\circ	OIIG	WII SIC	Cathere	COL C	IID CCL	CIC III.

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^{o})$	$\operatorname{Ideal}({}^{o})$
2	D	601	5YR	CAM-CAN-CAJ	-6.25	103.91	112.33
2	С	601	5YR	CAM-CAN-CAJ	-5.90	104.38	112.33
2	A	601	5YR	CAM-CAN-CAJ	-5.51	104.90	112.33
2	В	601	5YR	CAF-CAJ-CAI	5.49	119.40	110.58
2	D	601	5YR	CAF-CAJ-CAI	5.35	119.17	110.58

There are no chirality outliers.

5 of 25 torsion outliers are listed below:

Mol	Chain	Chain Res Type		Atoms		
2	В	601	5YR	CBJ-CBI-NBF-CBE		
2	В	601	5YR	CBJ-CBI-NBF-CBH		
2	D	601	5YR	CBA-CBB-CBC-CBD		
2	В	601	5YR	CAY-CAZ-CBA-CBB		
2	D	601	5YR	CBC-CBD-CBE-OBG		

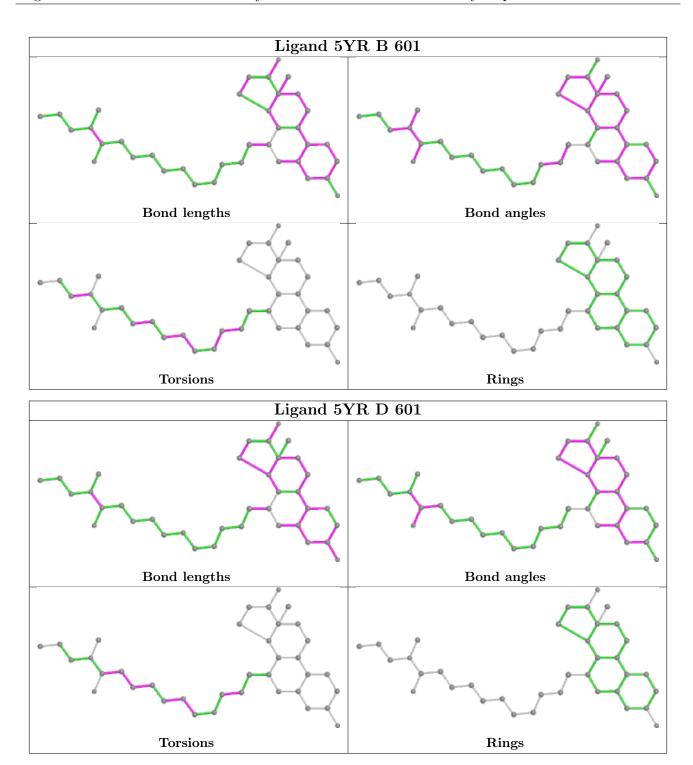
There are no ring outliers.

4 monomers are involved in 27 short contacts:

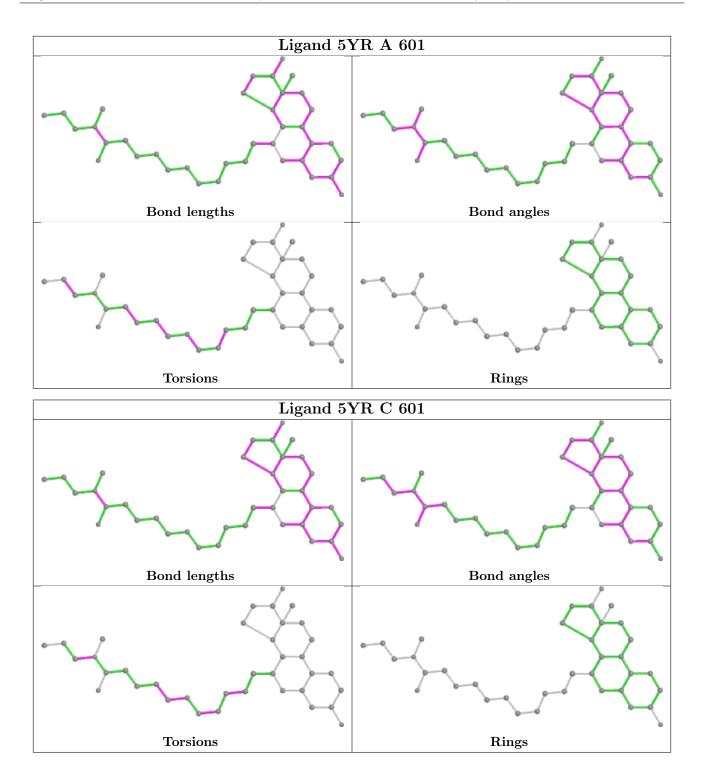
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	В	601	5YR	10	0
2	D	601	5YR	5	0
2	A	601	5YR	6	0
2	С	601	5YR	6	0

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.









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, 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	$\mathbf{OWAB}(\mathrm{\AA}^2)$	Q<0.9
1	A	239/249~(95%)	-0.44	0 100 100	13, 29, 56, 67	2 (0%)
1	В	235/249~(94%)	-0.46	1 (0%) 92 92	13, 29, 54, 67	1 (0%)
1	С	240/249 (96%)	-0.46	1 (0%) 92 92	13, 30, 55, 67	1 (0%)
1	D	237/249~(95%)	-0.45	0 100 100	13, 29, 56, 80	1 (0%)
All	All	951/996~(95%)	-0.45	2 (0%) 95 94	13, 29, 56, 80	5 (0%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	С	308	LEU	3.0
1	В	533	VAL	2.3

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 monosaccharides in this entry.

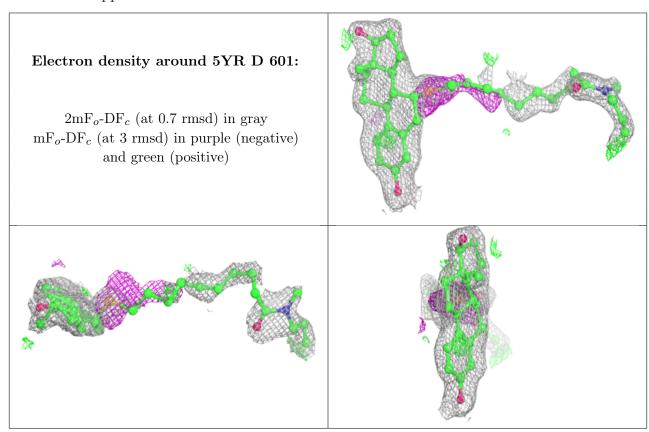
6.4 Ligands (i)

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, 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
2	5YR	D	601	37/37	0.86	0.19	17,32,67,69	0
2	5YR	С	601	37/37	0.87	0.19	18,33,62,66	0
2	5YR	A	601	37/37	0.88	0.20	16,30,58,61	0
2	5YR	В	601	37/37	0.89	0.20	19,31,72,74	0

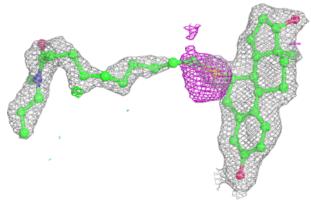
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

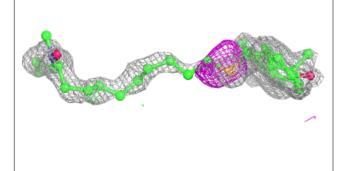


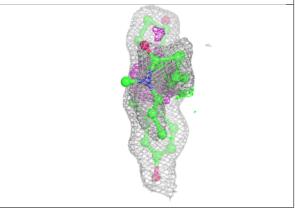


Electron density around 5YR C 601:

 $2 {\rm mF}_o\text{-}{\rm DF}_c$ (at 0.7 rmsd) in gray ${\rm mF}_o\text{-}{\rm DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)

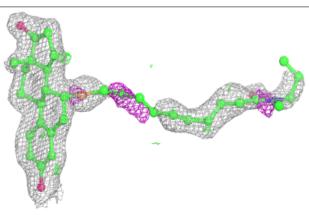


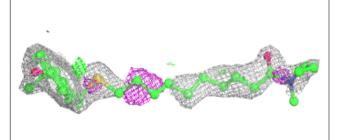


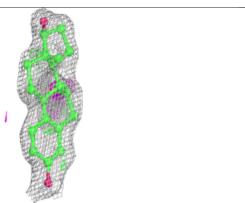


Electron density around 5YR A 601:

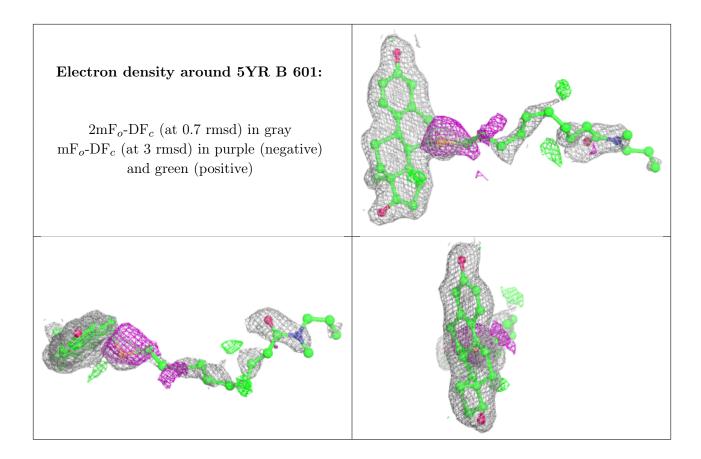
 $2 {\rm mF}_o\text{-}{\rm DF}_c$ (at 0.7 rmsd) in gray ${\rm mF}_o\text{-}{\rm DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)











6.5 Other polymers (i)

There are no such residues in this entry.

