



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 01:39 AM UTC

PDB ID : 1XDK / pdb_00001xdk
Title : Crystal Structure of the RARbeta/RXRalpha Ligand Binding Domain Heterodimer in Complex with 9-cis Retinoic Acid and a Fragment of the TRAP220 Coactivator
Authors : Pogenberg, V.; Guichou, J.F.; Vivat-Hannah, V.; Kammerer, S.; Perez, E.; Germain, P.; De Lera, A.R.; Gronemeyer, H.; Royer, C.A.; Bourguet, W.
Deposited on : 2004-09-07
Resolution : 2.90 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 ⓘ 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:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

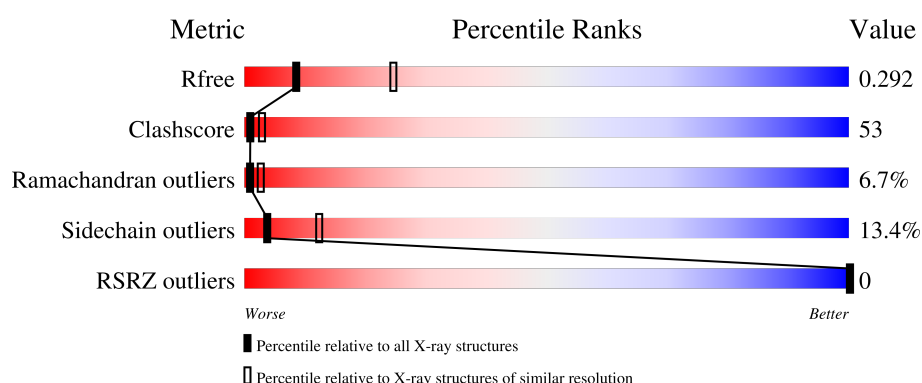
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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.



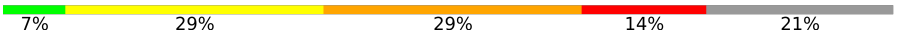
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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

Mol	Chain	Length	Quality of chain
1	A	238	
1	E	238	
2	B	303	
2	F	303	
3	C	14	

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Mol	Chain	Length	Quality of chain
3	D	14	 43% 29% 29%
3	G	14	 7% 29% 29% 14% 21%
3	H	14	 43% 29% 29%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7682 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 Retinoic acid receptor RXR-alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	217	Total	C	N	O	S	0	0	0
			1700	1091	290	309	10			
1	E	217	Total	C	N	O	S	0	0	0
			1700	1091	290	309	10			

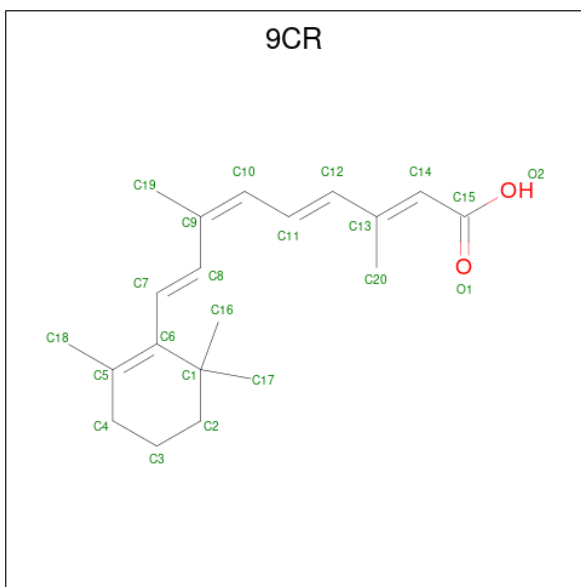
- Molecule 2 is a protein called Retinoic acid receptor, beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	245	Total	C	N	O	S	0	0	0
			1919	1220	325	360	14			
2	F	245	Total	C	N	O	S	0	0	0
			1919	1220	325	360	14			

- Molecule 3 is a protein called Thyroid Receptor Associated Protein 220.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	11	Total	C	N	O	S	0	0	0
			90	57	16	15	2			
3	D	10	Total	C	N	O	S	0	0	0
			82	53	14	13	2			
3	G	11	Total	C	N	O	S	0	0	0
			90	57	16	15	2			
3	H	10	Total	C	N	O	S	0	0	0
			82	53	14	13	2			

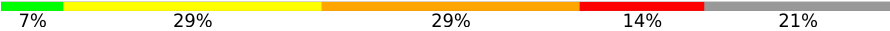
- Molecule 4 is (9cis)-retinoic acid (CCD ID: 9CR) (formula: C₂₀H₂₈O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			22	20	2		
4	B	1	Total	C	O	0	0
			22	20	2		
4	E	1	Total	C	O	0	0
			22	20	2		
4	F	1	Total	C	O	0	0
			22	20	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	4	Total	O	0	0
			4	4		
5	B	2	Total	O	0	0
			2	2		
5	E	4	Total	O	0	0
			4	4		
5	F	2	Total	O	0	0
			2	2		

Chain G:  7% 29% 29% 14% 21%



● Molecule 3: Thyroid Receptor Associated Protein 220

Chain H:  43% 29% 29%



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	115.70Å 115.70Å 247.20Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.90 30.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.6 (30.00-2.90) 99.5 (30.00-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.30 (at 2.90Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.253 , 0.296 0.253 , 0.292	Depositor DCC
R_{free} test set	1372 reflections (3.19%)	wwPDB-VP
Wilson B-factor (Å ²)	74.0	Xtriage
Anisotropy	0.605	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 87.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.480 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7682	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.34% 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: 9CR

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	0.91	0/1734	1.38	23/2345 (1.0%)
1	E	0.91	0/1734	1.38	23/2345 (1.0%)
2	B	0.81	2/1952 (0.1%)	1.26	19/2638 (0.7%)
2	F	0.82	1/1952 (0.1%)	1.26	20/2638 (0.8%)
3	C	0.75	0/91	1.50	2/121 (1.7%)
3	D	0.76	0/83	1.03	2/110 (1.8%)
3	G	0.76	0/91	1.50	2/121 (1.7%)
3	H	0.77	0/83	1.05	2/110 (1.8%)
All	All	0.86	3/7720 (0.0%)	1.32	93/10428 (0.9%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	393	MET	SD-CE	5.52	1.93	1.79
2	F	393	MET	SD-CE	5.49	1.93	1.79
2	B	395	ILE	CA-CB	5.15	1.59	1.53

The worst 5 of 93 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	240	ILE	N-CA-C	-10.73	100.21	111.58
2	F	296	GLY	CA-C-N	10.29	130.60	119.28
2	F	296	GLY	C-N-CA	10.29	130.60	119.28
2	B	296	GLY	CA-C-N	10.10	130.39	119.28
2	B	296	GLY	C-N-CA	10.10	130.39	119.28

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	1700	0	1728	204	0
1	E	1700	0	1728	207	0
2	B	1919	0	1979	205	0
2	F	1919	0	1979	202	0
3	C	90	0	93	18	0
3	D	82	0	87	14	0
3	G	90	0	93	19	0
3	H	82	0	87	12	0
4	A	22	0	27	3	0
4	B	22	0	27	3	0
4	E	22	0	27	2	0
4	F	22	0	27	3	0
5	A	4	0	0	0	0
5	B	2	0	0	0	0
5	E	4	0	0	1	0
5	F	2	0	0	0	0
All	All	7682	0	7882	821	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 53.

The worst 5 of 821 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)
2:B:343:LYS:HA	2:B:346:GLU:HG3	1.26	1.16
2:F:343:LYS:HA	2:F:346:GLU:HG3	1.26	1.14
2:F:315:ASP:HB3	2:F:359:ARG:HH12	1.09	1.13
2:F:357:ARG:HG3	2:F:357:ARG:HH11	1.13	1.09
1:A:237:VAL:HG23	1:A:238:GLU:H	1.14	1.07

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	213/238 (90%)	165 (78%)	33 (16%)	15 (7%)	1	2
1	E	213/238 (90%)	163 (76%)	34 (16%)	16 (8%)	1	2
2	B	243/303 (80%)	200 (82%)	34 (14%)	9 (4%)	2	11
2	F	243/303 (80%)	202 (83%)	32 (13%)	9 (4%)	2	11
3	C	9/14 (64%)	2 (22%)	2 (22%)	5 (56%)	0	0
3	D	8/14 (57%)	5 (62%)	1 (12%)	2 (25%)	0	0
3	G	9/14 (64%)	2 (22%)	2 (22%)	5 (56%)	0	0
3	H	8/14 (57%)	5 (62%)	1 (12%)	2 (25%)	0	0
All	All	946/1138 (83%)	744 (79%)	139 (15%)	63 (7%)	1	3

5 of 63 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	238	GLU
1	A	287	TRP
1	A	448	GLY
2	B	412	GLY
3	C	643	PRO

5.3.2 Protein sidechains

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	184/204 (90%)	163 (89%)	21 (11%)	5	18

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	184/204 (90%)	163 (89%)	21 (11%)	5	18
2	B	216/274 (79%)	185 (86%)	31 (14%)	3	11
2	F	216/274 (79%)	186 (86%)	30 (14%)	3	11
3	C	11/13 (85%)	8 (73%)	3 (27%)	0	1
3	D	10/13 (77%)	8 (80%)	2 (20%)	1	4
3	G	11/13 (85%)	8 (73%)	3 (27%)	0	1
3	H	10/13 (77%)	8 (80%)	2 (20%)	1	4
All	All	842/1008 (84%)	729 (87%)	113 (13%)	4	12

5 of 113 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	235	MET
3	G	651	ASP
1	E	369	LYS
3	G	644	MET
2	F	369	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 18 such sidechains are listed below:

Mol	Chain	Res	Type
2	F	250	GLN
3	H	647	ASN
3	G	641	ASN
3	D	647	ASN
2	F	197	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$
4	9CR	E	1500	-	22,22,22	2.19	7 (31%)	30,30,30	2.24	8 (26%)
4	9CR	F	1600	-	22,22,22	2.65	5 (22%)	30,30,30	2.43	8 (26%)
4	9CR	A	500	-	22,22,22	2.24	6 (27%)	30,30,30	2.38	9 (30%)
4	9CR	B	600	-	22,22,22	2.49	4 (18%)	30,30,30	2.55	10 (33%)

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
4	9CR	E	1500	-	-	4/15/32/32	0/1/1/1
4	9CR	F	1600	-	-	4/15/32/32	0/1/1/1
4	9CR	A	500	-	-	4/15/32/32	0/1/1/1
4	9CR	B	600	-	-	4/15/32/32	0/1/1/1

The worst 5 of 22 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	1600	9CR	C1-C6	8.94	1.65	1.53
4	B	600	9CR	C1-C6	8.14	1.64	1.53
4	A	500	9CR	C1-C6	7.12	1.62	1.53
4	E	1500	9CR	C1-C6	6.67	1.62	1.53
4	F	1600	9CR	C5-C6	5.56	1.43	1.34

The worst 5 of 35 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	1600	9CR	C18-C5-C6	7.41	132.57	124.48
4	B	600	9CR	C18-C5-C6	7.37	132.53	124.48
4	A	500	9CR	C18-C5-C6	6.27	131.33	124.48
4	E	1500	9CR	C18-C5-C6	6.03	131.07	124.48
4	F	1600	9CR	C15-C14-C13	5.79	137.02	128.40

There are no chirality outliers.

5 of 16 torsion outliers are listed below:

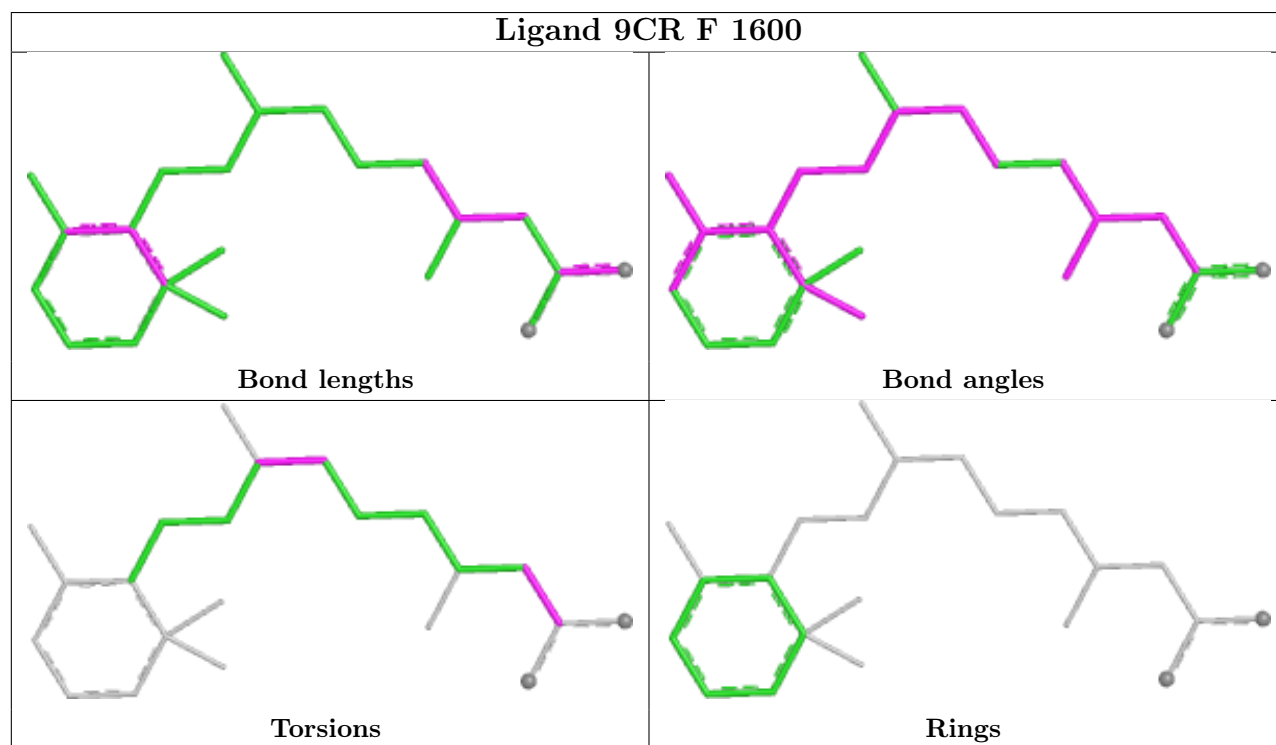
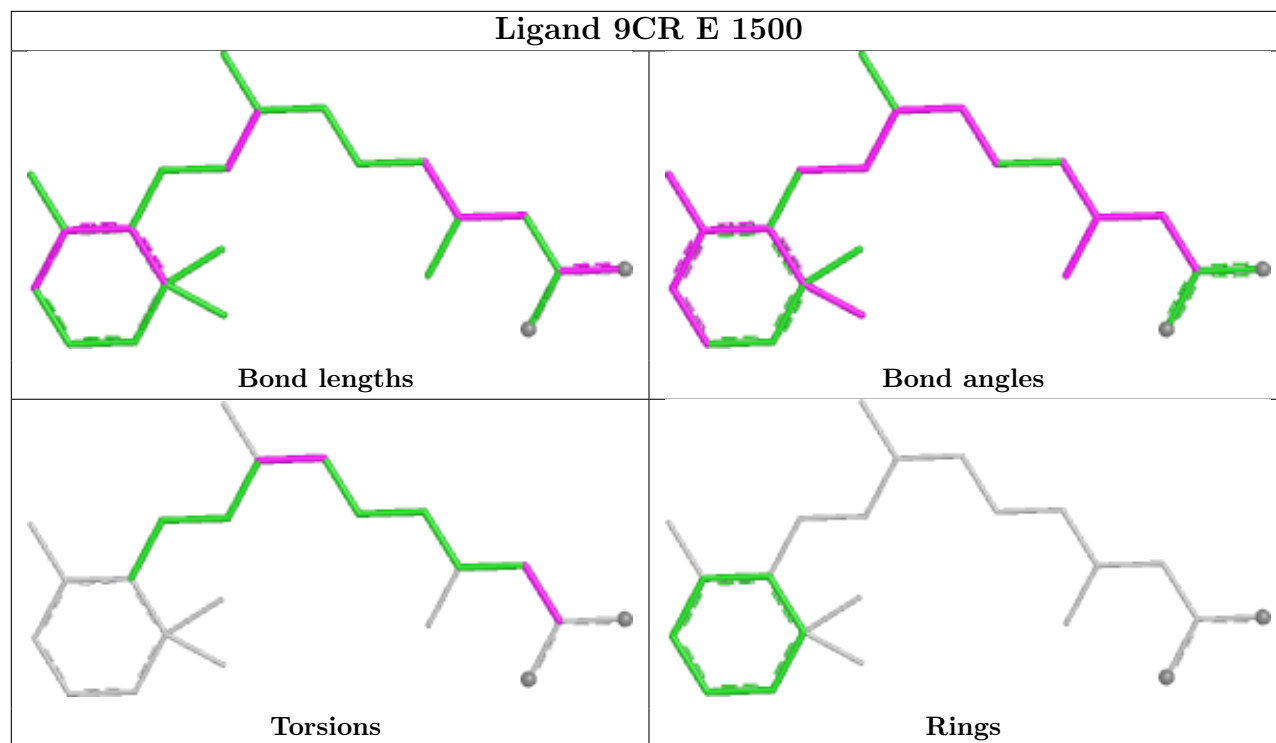
Mol	Chain	Res	Type	Atoms
4	B	600	9CR	C13-C14-C15-O1
4	F	1600	9CR	C13-C14-C15-O1
4	F	1600	9CR	C13-C14-C15-O2
4	B	600	9CR	C13-C14-C15-O2
4	E	1500	9CR	C13-C14-C15-O2

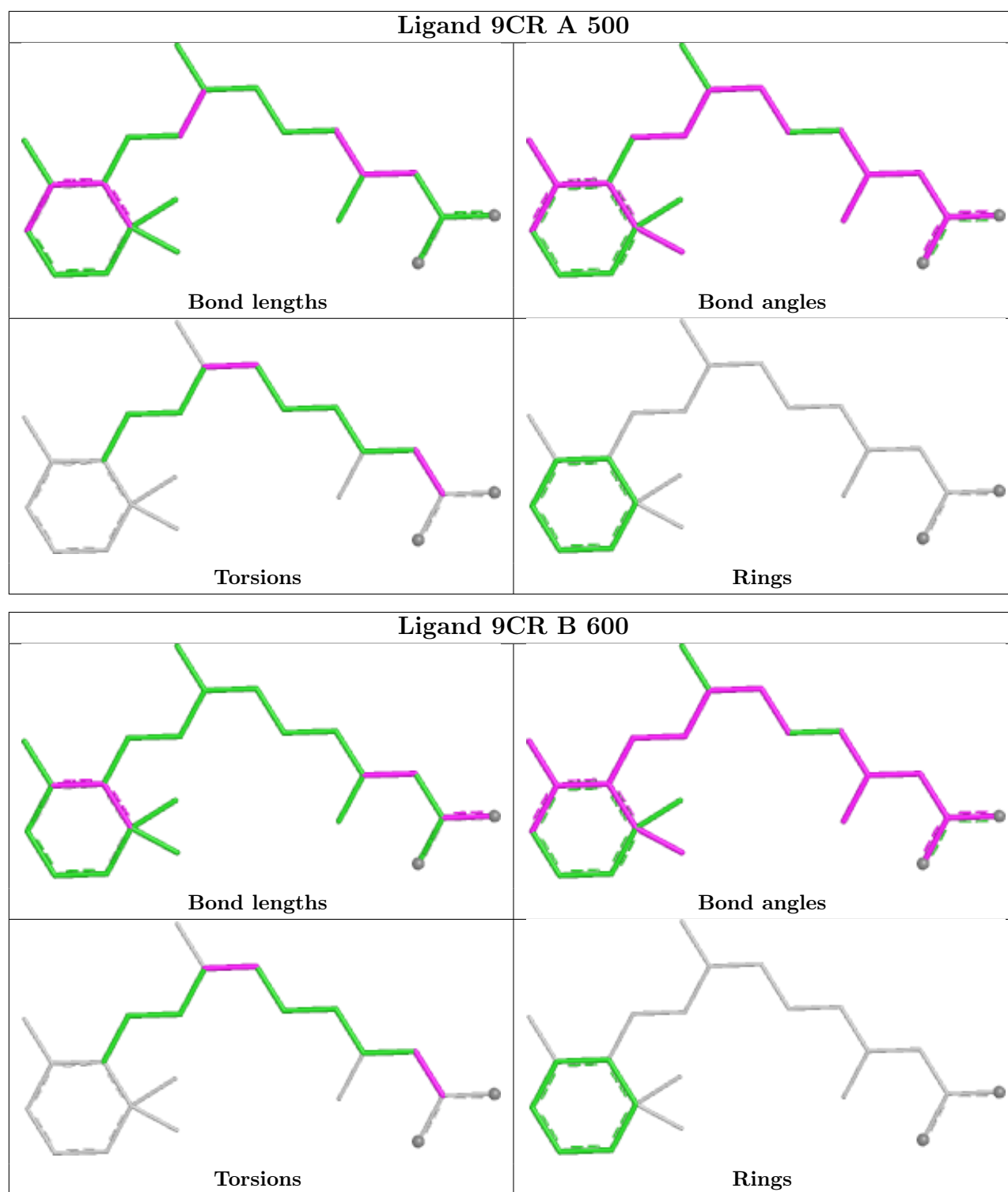
There are no ring outliers.

4 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	1500	9CR	2	0
4	F	1600	9CR	3	0
4	A	500	9CR	3	0
4	B	600	9CR	3	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 ⓘ

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	217/238 (91%)	-1.26	0 100 100	39, 71, 93, 104	0
1	E	217/238 (91%)	-1.27	0 100 100	40, 71, 92, 105	0
2	B	245/303 (80%)	-1.16	0 100 100	45, 75, 96, 104	0
2	F	245/303 (80%)	-1.16	0 100 100	44, 75, 96, 105	0
3	C	11/14 (78%)	-0.93	0 100 100	91, 96, 102, 103	0
3	D	10/14 (71%)	-1.12	0 100 100	91, 96, 101, 106	0
3	G	11/14 (78%)	-1.01	0 100 100	90, 97, 104, 104	0
3	H	10/14 (71%)	-0.97	0 100 100	92, 97, 102, 106	0
All	All	966/1138 (84%)	-1.20	0 100 100	39, 74, 97, 106	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

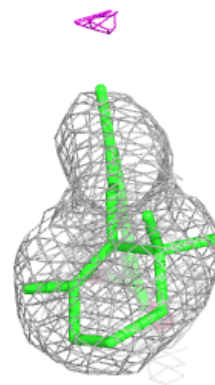
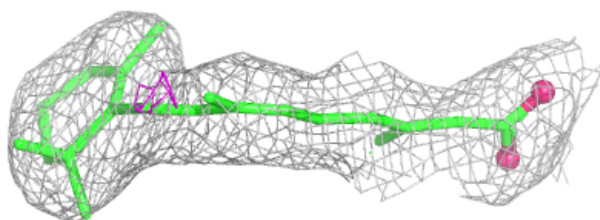
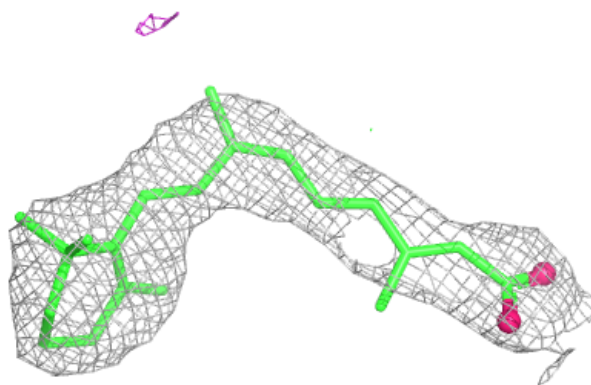
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	9CR	A	500	22/22	0.99	0.05	48,57,66,68	0
4	9CR	B	600	22/22	0.99	0.05	69,72,76,77	0
4	9CR	F	1600	22/22	0.99	0.05	69,72,75,77	0
4	9CR	E	1500	22/22	1.00	0.04	48,57,69,71	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.

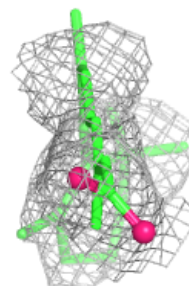
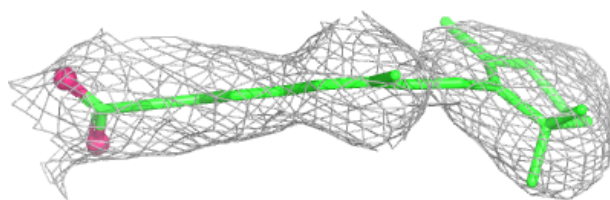
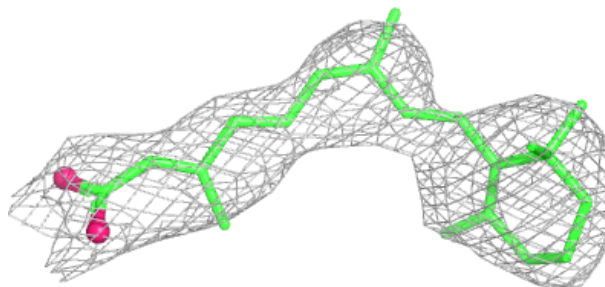
Electron density around 9CR A 500:

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

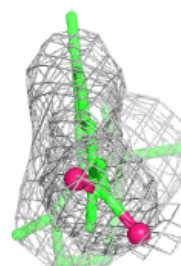
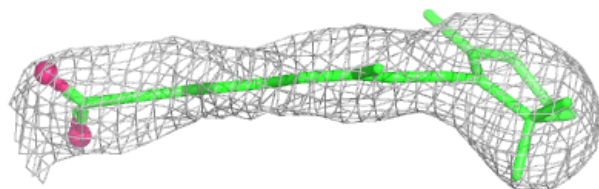
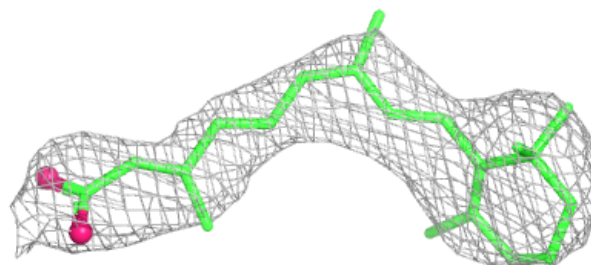


Electron density around 9CR B 600:

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

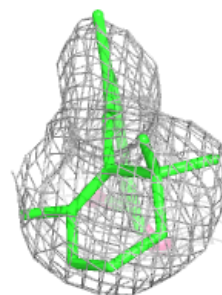
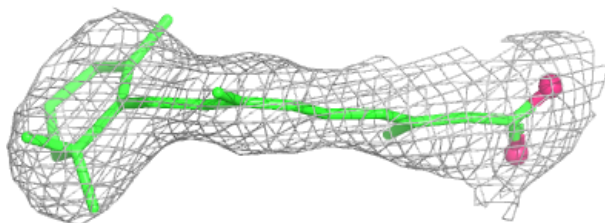
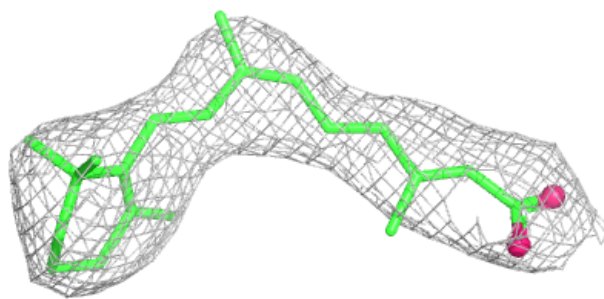
**Electron density around 9CR F 1600:**

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



Electron density around 9CR E 1500:

$2mF_o-DF_c$ (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.