



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 1, 2026 – 08:15 PM UTC

PDB ID : 1K7L / pdb_00001k7l
Title : The 2.5 Angstrom resolution crystal structure of the human PPARalpha ligand binding domain bound with GW409544 and a co-activator peptide.
Authors : Xu, H.E.; Lambert, M.H.; Montana, V.G.; Moore, L.B.; Collins, J.L.; Oplinger, J.A.; Kliewer, S.A.; Gampe Jr., R.T.; McKee, D.D.; Moore, J.T.; Willson, T.M.
Deposited on : 2001-10-19
Resolution : 2.50 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

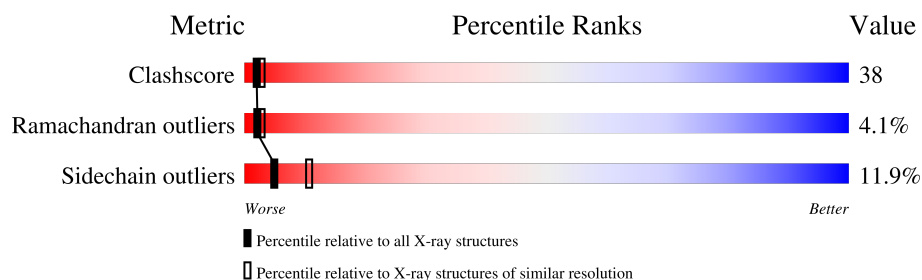
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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))
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	288	42% 41% 9% • 7%
1	C	288	50% 32% 10% • 7%
1	E	288	50% 32% 10% • 7%
1	G	288	43% 36% 13% • 7%
2	B	21	24% 38% 38%
2	D	21	24% 24% 14% 38%
2	F	21	19% 38% 5% 38%
2	H	21	14% 43% 5% 38%

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
3	544	A	469	X	-	-	-
3	544	C	470	X	-	-	-
3	544	E	501	X	-	-	-
3	544	G	601	X	-	-	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9308 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 Peroxisome proliferator activated receptor alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	267	Total	C	N	O	S	0	0	0
			2118	1359	355	386	18			
1	C	267	Total	C	N	O	S	0	0	0
			2118	1359	355	386	18			
1	E	267	Total	C	N	O	S	0	0	0
			2118	1359	355	386	18			
1	G	267	Total	C	N	O	S	0	0	0
			2118	1359	355	386	18			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	181	MET	-	expression tag	UNP Q07869
A	182	LYS	-	expression tag	UNP Q07869
A	183	LYS	-	expression tag	UNP Q07869
A	184	GLY	-	expression tag	UNP Q07869
A	185	HIS	-	expression tag	UNP Q07869
A	186	HIS	-	expression tag	UNP Q07869
A	187	HIS	-	expression tag	UNP Q07869
A	188	HIS	-	expression tag	UNP Q07869
A	189	HIS	-	expression tag	UNP Q07869
A	190	HIS	-	expression tag	UNP Q07869
A	191	GLY	-	expression tag	UNP Q07869
C	181	MET	-	expression tag	UNP Q07869
C	182	LYS	-	expression tag	UNP Q07869
C	183	LYS	-	expression tag	UNP Q07869
C	184	GLY	-	expression tag	UNP Q07869
C	185	HIS	-	expression tag	UNP Q07869
C	186	HIS	-	expression tag	UNP Q07869
C	187	HIS	-	expression tag	UNP Q07869
C	188	HIS	-	expression tag	UNP Q07869
C	189	HIS	-	expression tag	UNP Q07869
C	190	HIS	-	expression tag	UNP Q07869

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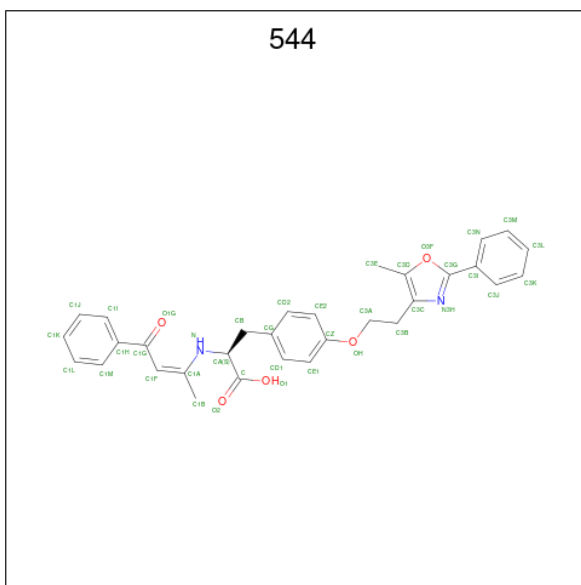
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Chain	Residue	Modelled	Actual	Comment	Reference
C	191	GLY	-	expression tag	UNP Q07869
E	181	MET	-	expression tag	UNP Q07869
E	182	LYS	-	expression tag	UNP Q07869
E	183	LYS	-	expression tag	UNP Q07869
E	184	GLY	-	expression tag	UNP Q07869
E	185	HIS	-	expression tag	UNP Q07869
E	186	HIS	-	expression tag	UNP Q07869
E	187	HIS	-	expression tag	UNP Q07869
E	188	HIS	-	expression tag	UNP Q07869
E	189	HIS	-	expression tag	UNP Q07869
E	190	HIS	-	expression tag	UNP Q07869
E	191	GLY	-	expression tag	UNP Q07869
G	181	MET	-	expression tag	UNP Q07869
G	182	LYS	-	expression tag	UNP Q07869
G	183	LYS	-	expression tag	UNP Q07869
G	184	GLY	-	expression tag	UNP Q07869
G	185	HIS	-	expression tag	UNP Q07869
G	186	HIS	-	expression tag	UNP Q07869
G	187	HIS	-	expression tag	UNP Q07869
G	188	HIS	-	expression tag	UNP Q07869
G	189	HIS	-	expression tag	UNP Q07869
G	190	HIS	-	expression tag	UNP Q07869
G	191	GLY	-	expression tag	UNP Q07869

- Molecule 2 is a protein called steroid receptor coactivator.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	13	Total	C	N	O	0	0	0
			110	68	24	18			
2	D	13	Total	C	N	O	0	0	0
			110	68	24	18			
2	F	13	Total	C	N	O	0	0	0
			110	68	24	18			
2	H	13	Total	C	N	O	0	0	0
			110	68	24	18			

- Molecule 3 is 2-(1-METHYL-3-OXO-3-PHENYL-PROPYLAMINO)-3-{4-[2-(5-METHYL-2-PHENYL-OXAZOL-4-YL)-ETHOXY]-PHENYL}-PROPIONIC ACID (CCD ID: 544) (formula: C₃₁H₃₀N₂O₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total 38	C 31	N 2	O 5	0	0
3	C	1	Total 38	C 31	N 2	O 5	0	0
3	E	1	Total 38	C 31	N 2	O 5	0	0
3	G	1	Total 38	C 31	N 2	O 5	0	0

- Molecule 4 is YTTRIUM (III) ION (CCD ID: YT3) (formula: Y).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	C	1	Total Y 1 1	0	0
4	E	1	Total Y 1 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	46	Total O 46 46	0	0
5	C	76	Total O 76 76	0	0
5	D	7	Total O 7 7	0	0

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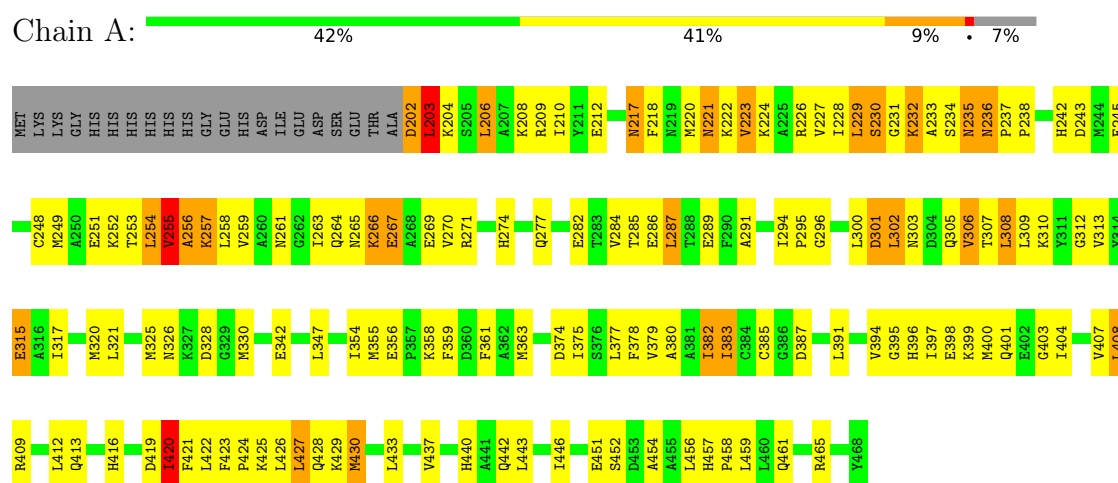
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	E	51	Total 51	O 51	0	0
5	F	1	Total 1	O 1	0	0
5	G	57	Total 57	O 57	0	0
5	H	4	Total 4	O 4	0	0

3 Residue-property plots

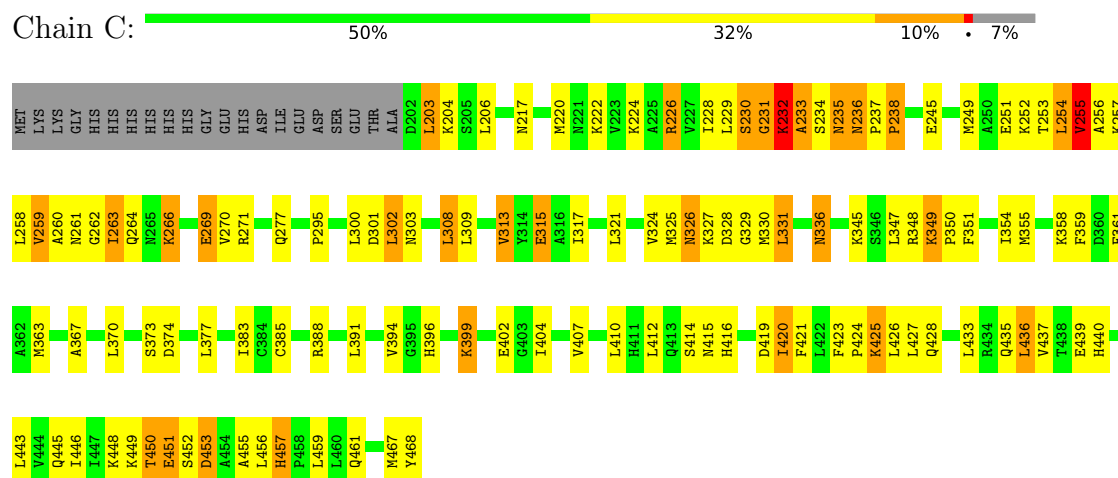
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

- Molecule 1: Peroxisome proliferator activated receptor alpha

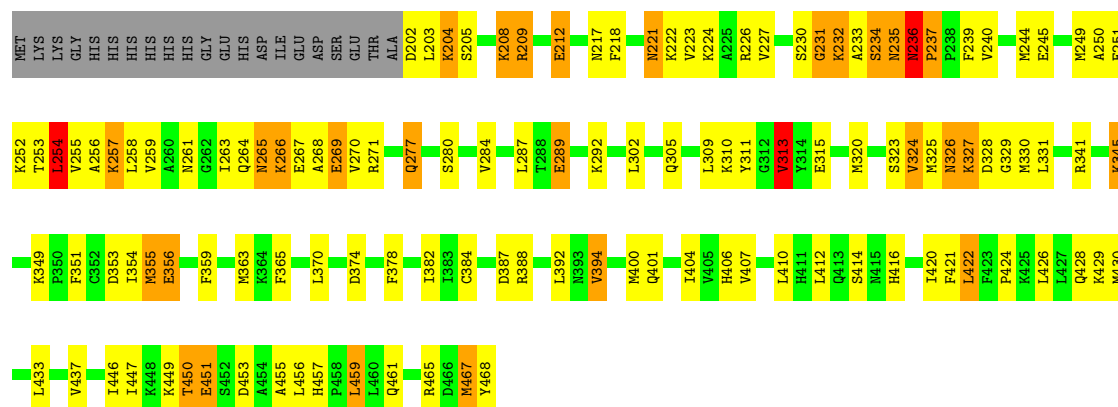


- Molecule 1: Peroxisome proliferator activated receptor alpha



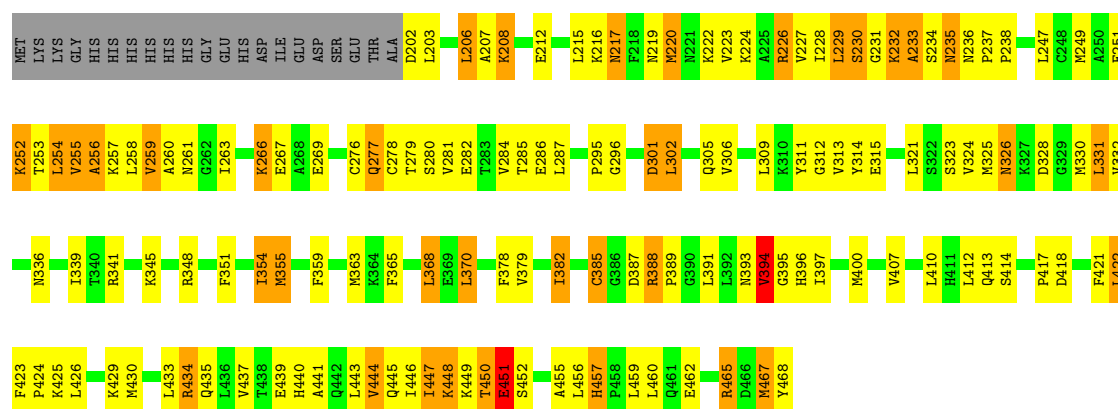
- Molecule 1: Peroxisome proliferator activated receptor alpha





- Molecule 1: Peroxisome proliferator activated receptor alpha

Chain G: 43% 36% 13% 7%



- Molecule 2: steroid receptor coactivator

Chain B: 24% 38% 38%



- Molecule 2: steroid receptor coactivator

Chain D: 24% 24% 14% 38%

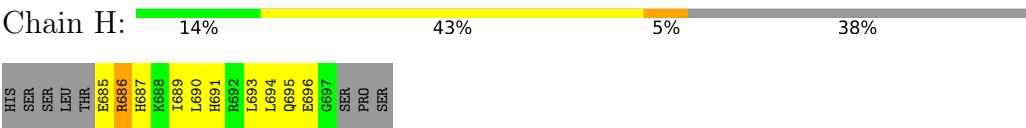


- Molecule 2: steroid receptor coactivator

Chain F: 19% 38% 5% 38%



● Molecule 2: steroid receptor coactivator



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	95.56Å 121.60Å 121.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.99 – 2.50	Depositor
% Data completeness (in resolution range)	98.3 (19.99-2.50)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
Refinement program	CNX 2000.1	Depositor
R, R_{free}	0.247 , 0.284	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9308	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

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: 544, YT3

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	0.48	0/2156	1.02	8/2906 (0.3%)
1	C	0.54	0/2156	1.03	11/2906 (0.4%)
1	E	0.49	0/2156	1.03	15/2906 (0.5%)
1	G	0.51	0/2156	1.03	11/2906 (0.4%)
2	B	0.35	0/111	0.74	0/147
2	D	0.49	0/111	1.02	0/147
2	F	0.45	0/111	0.81	0/147
2	H	0.35	0/111	0.60	0/147
All	All	0.50	0/9068	1.02	45/12212 (0.4%)

There are no bond length outliers.

The worst 5 of 45 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	256	ALA	N-CA-C	-14.58	95.07	112.86
1	E	204	LYS	N-CA-C	-12.72	98.40	114.56
1	G	256	ALA	N-CA-C	-12.70	95.50	112.26
1	A	354	ILE	N-CA-C	9.04	119.03	110.53
1	C	266	LYS	N-CA-C	-8.35	97.04	110.32

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	2118	0	2161	191	0
1	C	2118	0	2161	182	0
1	E	2118	0	2161	133	0
1	G	2118	0	2161	177	0
2	B	110	0	108	14	0
2	D	110	0	108	11	0
2	F	110	0	108	12	0
2	H	110	0	108	16	0
3	A	38	0	28	1	0
3	C	38	0	28	3	0
3	E	38	0	29	2	0
3	G	38	0	29	2	0
4	C	1	0	0	0	0
4	E	1	0	0	0	0
5	A	46	0	0	0	0
5	C	76	0	0	10	0
5	D	7	0	0	1	0
5	E	51	0	0	2	0
5	F	1	0	0	0	0
5	G	57	0	0	3	0
5	H	4	0	0	1	0
All	All	9308	0	9190	696	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 38.

The worst 5 of 696 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:E:253:THR:HG23	1:E:254:LEU:H	1.09	1.09
1:A:252:LYS:HG3	1:A:259:VAL:HG21	1.39	1.04
1:E:251:GLU:HA	1:E:255:VAL:HB	1.41	1.02
1:C:420:ILE:HG13	1:C:421:PHE:H	1.23	1.00
1:E:202:ASP:O	1:E:204:LYS:N	1.93	1.00

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	265/288 (92%)	233 (88%)	23 (9%)	9 (3%)	3	4
1	C	265/288 (92%)	233 (88%)	18 (7%)	14 (5%)	1	1
1	E	265/288 (92%)	240 (91%)	16 (6%)	9 (3%)	3	4
1	G	265/288 (92%)	236 (89%)	18 (7%)	11 (4%)	2	2
2	B	11/21 (52%)	8 (73%)	3 (27%)	0	100	100
2	D	11/21 (52%)	8 (73%)	3 (27%)	0	100	100
2	F	11/21 (52%)	9 (82%)	2 (18%)	0	100	100
2	H	11/21 (52%)	9 (82%)	0	2 (18%)	0	0
All	All	1104/1236 (89%)	976 (88%)	83 (8%)	45 (4%)	2	3

5 of 45 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	230	SER
1	A	232	LYS
1	A	236	ASN
1	A	420	ILE
1	C	232	LYS

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	234/252 (93%)	204 (87%)	30 (13%)	4	9

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	234/252 (93%)	214 (92%)	20 (8%)	10	22
1	E	234/252 (93%)	208 (89%)	26 (11%)	6	12
1	G	234/252 (93%)	199 (85%)	35 (15%)	3	6
2	B	11/20 (55%)	10 (91%)	1 (9%)	9	19
2	D	11/20 (55%)	8 (73%)	3 (27%)	0	1
2	F	11/20 (55%)	9 (82%)	2 (18%)	2	3
2	H	11/20 (55%)	11 (100%)	0	100	100
All	All	980/1088 (90%)	863 (88%)	117 (12%)	5	11

5 of 117 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	221	ASN
1	G	422	LEU
1	E	387	ASP
1	G	413	GLN
1	G	302	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 55 such sidechains are listed below:

Mol	Chain	Res	Type
2	D	691	HIS
1	E	299	ASN
2	H	695	GLN
1	G	305	GLN
2	D	695	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	544	G	601	-	40,41,41	2.81	20 (50%)	51,55,55	2.04	8 (15%)
3	544	C	470	-	40,41,41	2.75	20 (50%)	51,55,55	2.05	8 (15%)
3	544	A	469	-	40,41,41	2.74	23 (57%)	51,55,55	2.19	9 (17%)
3	544	E	501	-	40,41,41	2.75	20 (50%)	51,55,55	2.00	7 (13%)

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	544	G	601	-	1/1/4/6	1/30/30/30	0/4/4/4
3	544	C	470	-	1/1/4/6	1/30/30/30	0/4/4/4
3	544	A	469	-	1/1/4/6	2/30/30/30	0/4/4/4
3	544	E	501	-	1/1/4/6	1/30/30/30	0/4/4/4

The worst 5 of 83 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	G	601	544	C3G-N3H	7.71	1.41	1.30
3	C	470	544	C3G-N3H	7.62	1.40	1.30
3	A	469	544	C3G-N3H	7.51	1.40	1.30
3	E	501	544	C3G-N3H	7.40	1.40	1.30
3	C	470	544	C1A-N	6.25	1.42	1.33

The worst 5 of 32 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	469	544	C3E-C3D-C3C	-7.88	128.77	135.22
3	E	501	544	C3E-C3D-C3C	-6.90	129.58	135.22
3	G	601	544	C3E-C3D-C3C	-6.89	129.58	135.22
3	C	470	544	C3E-C3D-C3C	-6.88	129.59	135.22
3	A	469	544	C1F-C1A-N	-5.80	116.74	121.29

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	469	544	C1A
3	C	470	544	C1A
3	E	501	544	C1A
3	G	601	544	C1A

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	470	544	C-CA-N-C1A
3	E	501	544	OH-C3A-C3B-C3C
3	G	601	544	OH-C3A-C3B-C3C
3	A	469	544	C-CA-N-C1A
3	A	469	544	OH-C3A-C3B-C3C

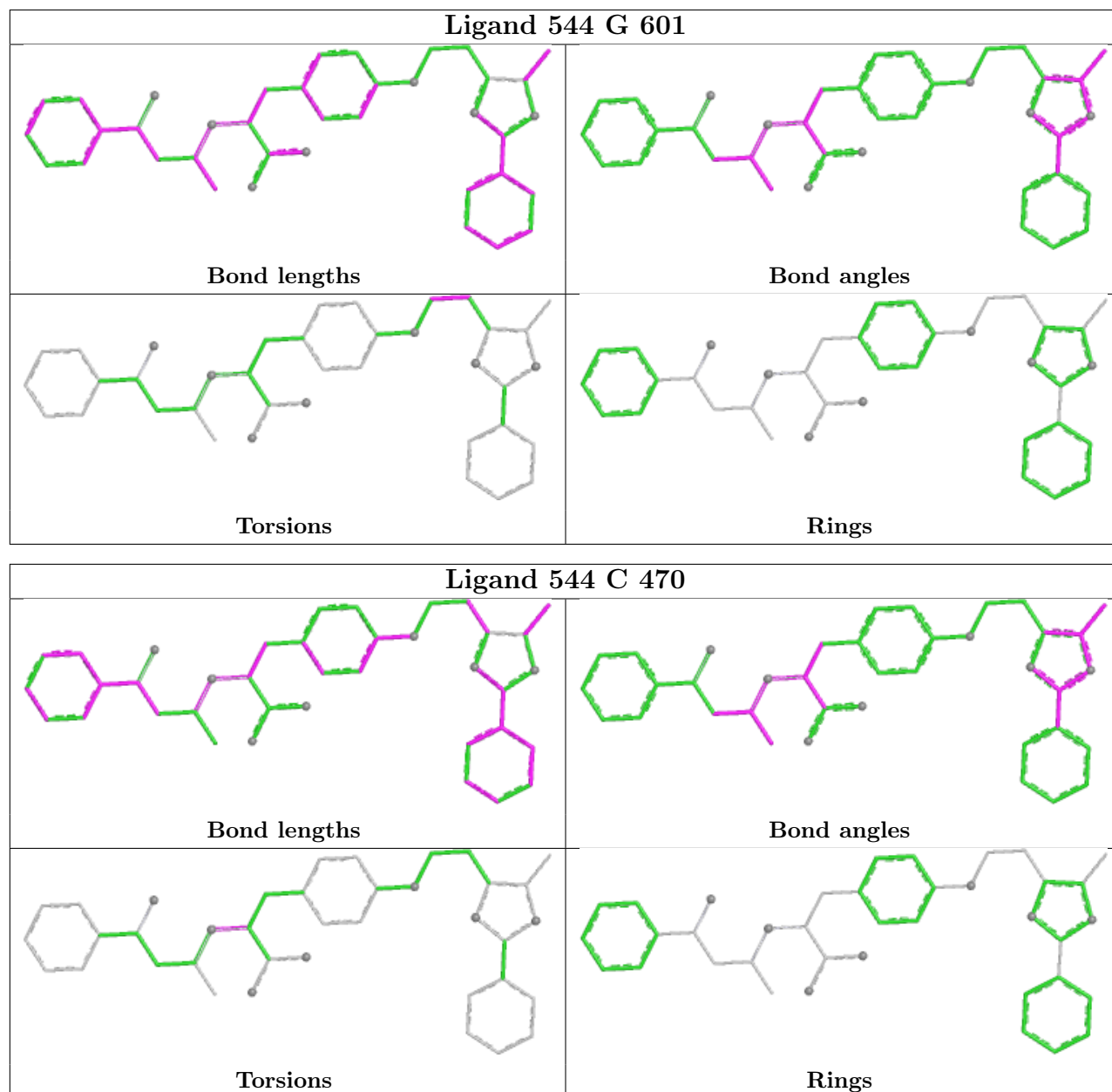
There are no ring outliers.

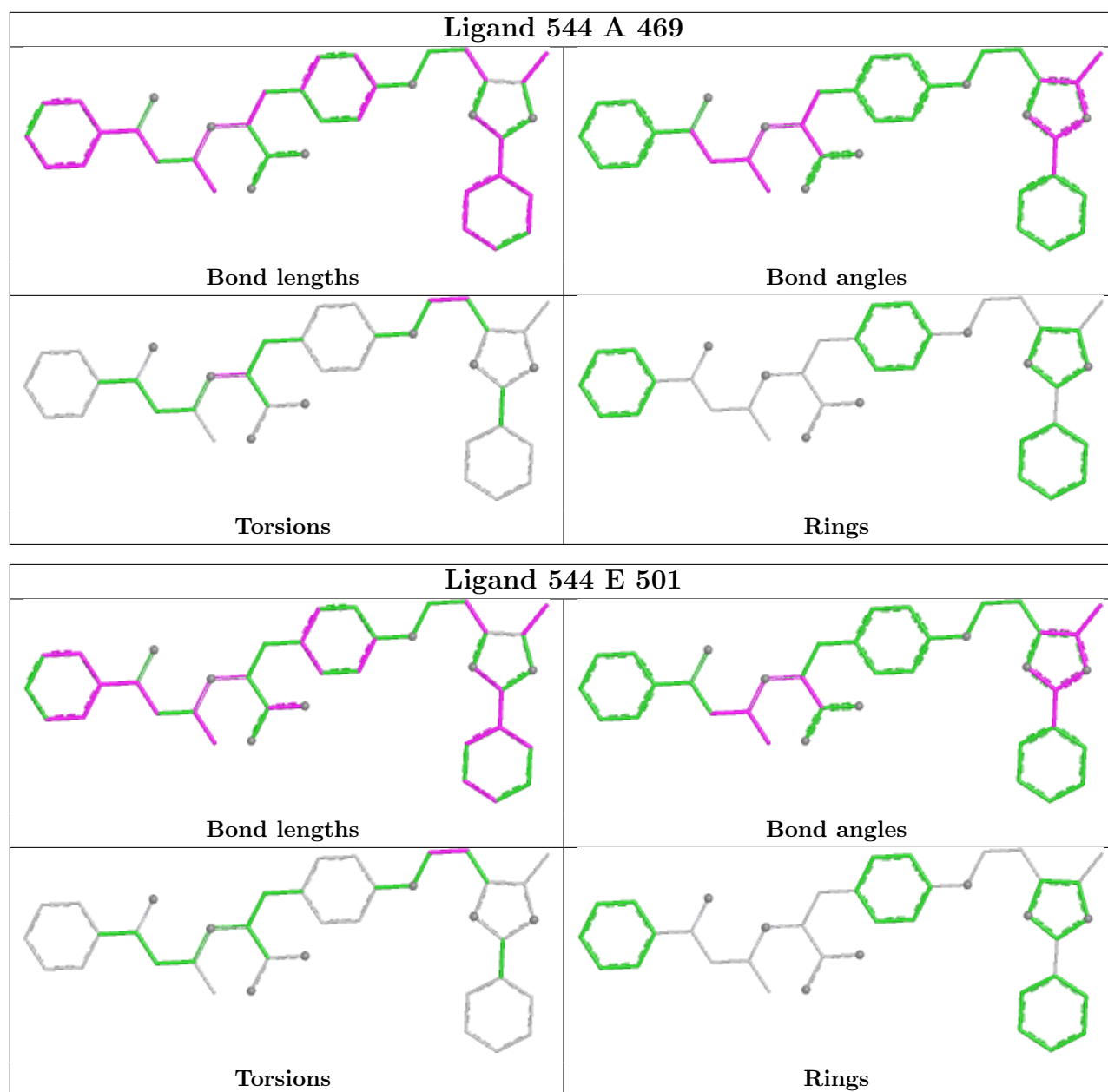
4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	G	601	544	2	0
3	C	470	544	3	0
3	A	469	544	1	0
3	E	501	544	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and

any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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](#)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section is therefore empty.

6.4 Ligands [i](#)

EDS was not executed - this section is therefore empty.

6.5 Other polymers [i](#)

EDS was not executed - this section is therefore empty.