



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:22 AM UTC

PDB ID : 1C6Z / pdb_00001c6z
Title : ALTERNATE BINDING SITE FOR THE P1-P3 GROUP OF A CLASS OF POTENT HIV-1 PROTEASE INHIBITORS AS A RESULT OF CONCERTED STRUCTURAL CHANGE IN 80'S LOOP.
Authors : Munshi, S.
Deposited on : 1999-12-28
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 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

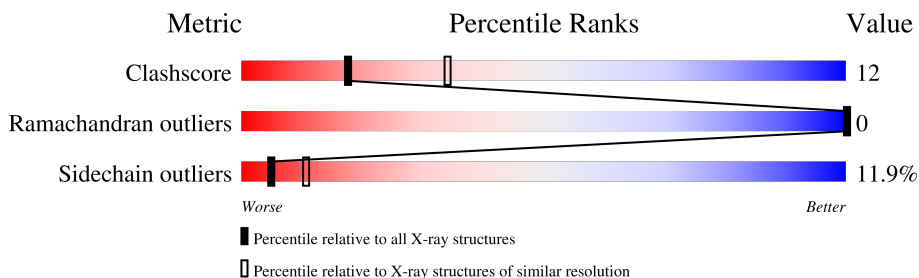
1 Overall quality at a glance

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	99	51% 37% 12%
1	B	99	55% 32% 12% .

2 Entry composition [i](#)

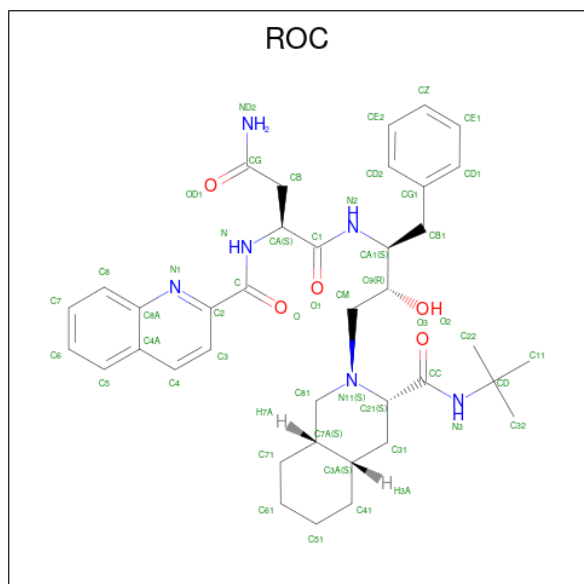
There are 3 unique types of molecules in this entry. The entry contains 1688 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PROTEIN (PROTEASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	99	Total	C	N	O	S	0	0	0
			757	488	129	136	4			
1	B	99	Total	C	N	O	S	0	0	0
			757	488	129	136	4			

- Molecule 2 is (2S)-N-[(2S,3R)-4-[(2S,3S,4aS,8aS)-3-(tert-butylcarbamoyl)-3,4,4a,5,6,7,8,8a-octahydro-1H-isoquinolin-2-yl]-3-hydroxy-1-phenyl-butan-2-yl]-2-(quinolin-2-ylcarbamoylamino)butanediamide (CCD ID: ROC) (formula: C₃₈H₅₀N₆O₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	1
			98	76	12	10		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	40	Total 40	O 40	0	0
3	B	36	Total 36	O 36	0	0

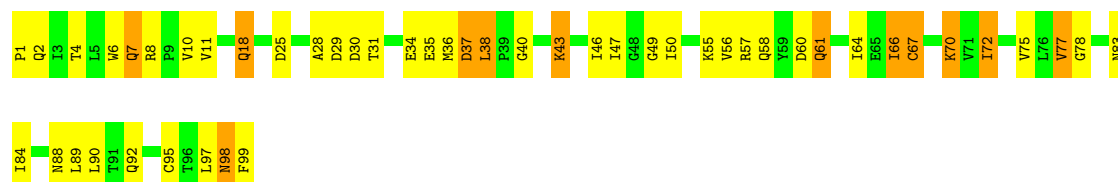
3 Residue-property plots [i](#)

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. 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.

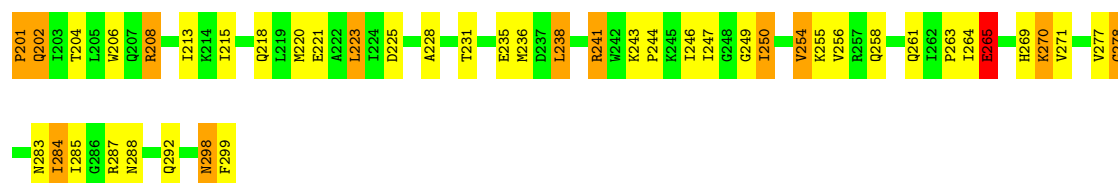
• Molecule 1: PROTEIN (PROTEASE)

Chain A: 



• Molecule 1: PROTEIN (PROTEASE)

Chain B: 



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	88.13Å 88.14Å 32.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.50	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.50)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.180 , 0.290	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1688	wwPDB-VP
Average B, all atoms (Å ²)	22.0	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: ROC

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	1.38	11/771 (1.4%)	1.56	19/1048 (1.8%)
1	B	1.34	11/771 (1.4%)	1.54	21/1048 (2.0%)
All	All	1.36	22/1542 (1.4%)	1.55	40/2096 (1.9%)

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	66	ILE	CA-C	-8.84	1.43	1.52
1	B	254	VAL	CA-C	8.44	1.63	1.52
1	A	30	ASP	CA-C	-7.92	1.43	1.52
1	A	25	ASP	CA-CB	7.28	1.63	1.53
1	B	241	ARG	CA-C	-7.18	1.43	1.52
1	B	292	GLN	CA-C	-6.94	1.43	1.52
1	B	206	TRP	CA-C	-6.89	1.43	1.52
1	B	265	GLU	CA-C	6.58	1.60	1.52
1	A	37	ASP	CA-CB	6.25	1.62	1.53
1	B	255	LYS	CA-C	-5.96	1.45	1.52
1	B	221	GLU	CA-C	-5.89	1.45	1.52
1	B	287	ARG	CA-C	5.76	1.60	1.52
1	B	244	PRO	CA-C	5.72	1.58	1.52
1	A	58	GLN	CA-C	-5.55	1.45	1.52
1	A	29	ASP	CA-C	-5.43	1.45	1.52
1	A	38	LEU	CA-C	5.35	1.59	1.52
1	A	6	TRP	CA-C	-5.32	1.46	1.52
1	A	90	LEU	CA-C	-5.27	1.45	1.52
1	A	61	GLN	CA-CB	5.15	1.61	1.53
1	B	238	LEU	CA-C	-5.15	1.45	1.52
1	B	284	ILE	CB-CG1	-5.12	1.43	1.53
1	A	46	ILE	CA-C	-5.00	1.46	1.52

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	7	GLN	OE1-CD-NE2	-14.17	108.43	122.60
1	B	218	GLN	OE1-CD-NE2	-11.56	111.04	122.60
1	A	1	PRO	CA-N-CD	-9.66	98.47	112.00
1	B	202	GLN	OE1-CD-NE2	-9.65	112.95	122.60
1	A	35	GLU	CA-CB-CG	9.30	132.70	114.10
1	A	61	GLN	OE1-CD-NE2	-9.04	113.56	122.60
1	A	83	ASN	OD1-CG-ND2	-8.11	114.49	122.60
1	A	61	GLN	CA-CB-CG	-7.51	99.08	114.10
1	B	201	PRO	CA-N-CD	-7.26	101.83	112.00
1	A	98	ASN	CB-CA-C	-6.97	95.87	109.68
1	A	7	GLN	N-CA-C	-6.90	99.99	110.14
1	A	58	GLN	OE1-CD-NE2	-6.83	115.77	122.60
1	A	43	LYS	CB-CA-C	6.79	119.82	109.56
1	B	218	GLN	CA-CB-CG	6.60	127.29	114.10
1	A	7	GLN	CG-CD-OE1	6.59	133.97	120.80
1	B	243	LYS	CB-CA-C	-6.58	99.62	109.56
1	B	298	ASN	OD1-CG-ND2	-6.56	116.04	122.60
1	B	261	GLN	OE1-CD-NE2	-6.23	116.37	122.60
1	B	265	GLU	CA-CB-CG	6.05	126.20	114.10
1	A	2	GLN	CA-CB-CG	-5.90	102.30	114.10
1	A	1	PRO	N-CA-CB	5.59	109.15	103.00
1	A	8	ARG	CA-C-O	-5.55	114.25	119.80
1	B	298	ASN	CB-CG-ND2	5.52	124.68	116.40
1	B	283	ASN	OD1-CG-ND2	-5.50	117.11	122.60
1	B	208	ARG	CA-C-N	5.49	126.70	119.84
1	B	208	ARG	C-N-CA	5.49	126.70	119.84
1	A	7	GLN	CA-CB-CG	5.47	125.03	114.10
1	B	241	ARG	CA-CB-CG	5.45	124.99	114.10
1	B	288	ASN	N-CA-C	-5.43	106.78	113.41
1	B	278	GLY	CA-C-N	5.39	126.58	119.84
1	B	278	GLY	C-N-CA	5.39	126.58	119.84
1	A	40	GLY	N-CA-C	5.23	117.36	112.08
1	B	255	LYS	CG-CD-CE	5.21	123.28	111.30
1	B	201	PRO	N-CA-CB	5.20	108.72	103.00
1	B	258	GLN	CA-CB-CG	-5.17	103.77	114.10
1	A	61	GLN	CG-CD-OE1	5.11	131.01	120.80
1	B	243	LYS	CB-CG-CD	5.10	123.03	111.30
1	A	67	CYS	CB-CA-C	-5.07	103.42	111.48
1	B	269	HIS	CB-CA-C	-5.05	101.78	110.22
1	A	18	GLN	OE1-CD-NE2	-5.00	117.59	122.60

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	757	0	797	23	4
1	B	757	0	795	23	1
2	B	98	0	98	2	0
3	A	40	0	0	4	3
3	B	36	0	0	5	4
All	All	1688	0	1690	41	7

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 12.

All (41) 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:B:256:VAL:HG12	1:B:278:GLY:HA3	1.42	0.96
1:A:92:GLN:HG3	3:A:333:HOH:O	1.67	0.91
1:A:50:ILE:HG22	1:B:249:GLY:HA3	1.61	0.83
1:A:66:ILE:HG22	1:A:67:CYS:SG	2.23	0.78
1:B:204:THR:O	3:B:324:HOH:O	2.10	0.68
1:B:265:GLU:HG2	1:B:270:LYS:HG2	1.77	0.66
1:A:70:LYS:HZ3	1:A:70:LYS:HB2	1.62	0.64
1:B:299:PHE:C	3:B:328:HOH:O	2.40	0.63
1:A:56:VAL:HG12	1:A:78:GLY:HA3	1.79	0.62
1:A:28:ALA:HB1	2:B:505[A]:ROC:H13	1.82	0.62
1:A:37:ASP:HB3	1:A:57:ARG:HH12	1.66	0.61
1:A:64:ILE:HD13	1:A:89:LEU:HD11	1.83	0.61
1:A:36:MET:O	1:A:77:VAL:HG11	1.99	0.61
1:A:31:THR:HG23	1:A:84:ILE:O	2.01	0.61
1:B:235:GLU:HA	1:B:277:VAL:CG1	2.30	0.60
1:B:236:MET:HE3	1:B:238:LEU:HD21	1.89	0.55
1:B:263:PRO:HD3	3:B:359:HOH:O	2.07	0.55
1:A:95:CYS:HA	1:B:299:PHE:HB3	1.92	0.51
1:B:264:ILE:HG23	1:B:271:VAL:HG23	1.92	0.51
1:A:72:ILE:O	3:A:333:HOH:O	2.19	0.51
1:B:231:THR:HG23	1:B:284:ILE:O	2.09	0.50
1:B:235:GLU:HA	1:B:277:VAL:HG13	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:ASP:HB3	1:A:61:GLN:HE21	1.78	0.49
1:A:55:LYS:NZ	3:A:339:HOH:O	2.44	0.48
1:A:98:ASN:HD21	1:B:202:GLN:HE21	1.61	0.46
1:B:208:ARG:HD3	3:B:325:HOH:O	2.15	0.46
1:B:223:LEU:HD23	1:B:225:ASP:HB2	1.97	0.46
1:A:99:PHE:CE1	1:B:201:PRO:HD2	2.52	0.44
1:B:213:ILE:HG21	1:B:285:ILE:HD11	2.00	0.43
1:B:256:VAL:HG12	1:B:278:GLY:CA	2.29	0.43
1:A:61:GLN:H	1:A:61:GLN:HG3	1.78	0.42
1:B:215:ILE:O	1:B:215:ILE:HG23	2.19	0.42
1:A:10:VAL:HG13	1:A:10:VAL:O	2.20	0.42
1:B:270:LYS:HG3	3:B:341:HOH:O	2.18	0.41
1:A:11:VAL:HG21	1:A:67:CYS:SG	2.60	0.41
1:A:70:LYS:HE3	3:A:329:HOH:O	2.21	0.41
1:B:228:ALA:HB1	2:B:505[B]:ROC:H321	2.02	0.41
1:B:238:LEU:HD23	1:B:238:LEU:HA	1.97	0.40
1:A:57:ARG:HD3	1:A:77:VAL:HG12	2.04	0.40
1:A:49:GLY:HA3	1:B:250:ILE:HG22	2.04	0.40
1:A:75:VAL:O	1:A:75:VAL:HG13	2.21	0.40

All (7) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:306:HOH:O	3:A:327:HOH:O[3_646]	0.39	1.81
3:A:339:HOH:O	3:A:351:HOH:O[3_656]	1.13	1.07
1:A:89:LEU:N	3:B:332:HOH:O[4_556]	1.41	0.79
1:A:89:LEU:CA	3:B:332:HOH:O[4_556]	1.51	0.69
1:B:241:ARG:NH1	3:A:333:HOH:O[4_456]	1.61	0.59
1:A:88:ASN:C	3:B:332:HOH:O[4_556]	1.89	0.31
1:A:89:LEU:CB	3:B:332:HOH:O[4_556]	2.15	0.05

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	97/99 (98%)	88 (91%)	9 (9%)	0	100	100
1	B	97/99 (98%)	88 (91%)	9 (9%)	0	100	100
All	All	194/198 (98%)	176 (91%)	18 (9%)	0	100	100

There are no Ramachandran outliers to report.

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	84/84 (100%)	73 (87%)	11 (13%)	4	8
1	B	84/84 (100%)	75 (89%)	9 (11%)	6	13
All	All	168/168 (100%)	148 (88%)	20 (12%)	5	11

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	THR
1	A	7	GLN
1	A	18	GLN
1	A	34	GLU
1	A	38	LEU
1	A	43	LYS
1	A	47	ILE
1	A	70	LYS
1	A	72	ILE
1	A	77	VAL
1	A	97	LEU
1	B	220	MET
1	B	223	LEU
1	B	246	ILE
1	B	247	ILE

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Mol	Chain	Res	Type
1	B	250	ILE
1	B	254	VAL
1	B	265	GLU
1	B	270	LYS
1	B	298	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	58	GLN
1	A	61	GLN
1	A	92	GLN
1	B	202	GLN
1	B	292	GLN
1	B	298	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	ROC	B	505[B]	-	53,53,53	1.40	6 (11%)	71,75,75	1.45	13 (18%)
2	ROC	B	505[A]	-	53,53,53	1.38	3 (5%)	71,75,75	1.43	12 (16%)

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	ROC	B	505[B]	-	-	3/41/64/64	0/5/5/5
2	ROC	B	505[A]	-	-	4/41/64/64	0/5/5/5

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	505[B]	ROC	O2-C9	-6.93	1.28	1.43
2	B	505[A]	ROC	O2-C9	-6.63	1.29	1.43
2	B	505[A]	ROC	CD2-CG1	2.41	1.43	1.38
2	B	505[B]	ROC	C2-C	-2.38	1.44	1.50
2	B	505[B]	ROC	CD2-CG1	2.27	1.43	1.38
2	B	505[A]	ROC	CD1-CG1	2.24	1.43	1.38
2	B	505[B]	ROC	CD1-CG1	2.13	1.43	1.38
2	B	505[B]	ROC	C81-N11	2.02	1.50	1.47
2	B	505[B]	ROC	CE1-CD1	2.01	1.42	1.38

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	505[B]	ROC	CM-N11-C81	-4.56	100.45	110.75
2	B	505[A]	ROC	CM-N11-C81	-4.15	101.39	110.75
2	B	505[B]	ROC	O2-C9-CA1	3.50	116.78	109.88
2	B	505[A]	ROC	C4A-C8A-N1	-3.44	117.51	122.25
2	B	505[B]	ROC	C4A-C8A-N1	-3.43	117.53	122.25
2	B	505[A]	ROC	C2-N1-C8A	3.26	122.49	117.56
2	B	505[B]	ROC	CD2-CG1-CD1	3.24	123.04	118.23
2	B	505[A]	ROC	O2-C9-CA1	3.12	116.02	109.88
2	B	505[B]	ROC	C2-N1-C8A	3.09	122.23	117.56
2	B	505[A]	ROC	CA-N-C	-2.60	115.33	121.56
2	B	505[B]	ROC	CE1-CD1-CG1	-2.49	117.10	120.61
2	B	505[B]	ROC	CE2-CD2-CG1	-2.48	117.12	120.61
2	B	505[A]	ROC	CD2-CG1-CD1	2.48	121.92	118.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	505[B]	ROC	C4-C3-C2	-2.43	116.26	118.84
2	B	505[A]	ROC	O3-CC-N3	2.37	128.10	123.15
2	B	505[B]	ROC	CB1-CG1-CD2	-2.35	116.53	120.90
2	B	505[B]	ROC	O3-CC-N3	2.33	128.02	123.15
2	B	505[A]	ROC	C4-C3-C2	-2.33	116.37	118.84
2	B	505[B]	ROC	C8-C8A-N1	2.27	122.00	118.69
2	B	505[A]	ROC	C8-C8A-N1	2.21	121.92	118.69
2	B	505[A]	ROC	O-C-C2	2.10	125.71	121.08
2	B	505[A]	ROC	CA-C1-N2	-2.07	112.22	116.63
2	B	505[B]	ROC	CB-CG-ND2	2.06	121.13	116.19
2	B	505[B]	ROC	CA1-N2-C1	-2.04	119.69	123.25
2	B	505[A]	ROC	O1-C1-N2	2.04	126.60	122.96

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	505[A]	ROC	O2-C9-CM-N11
2	B	505[B]	ROC	O2-C9-CM-N11
2	B	505[B]	ROC	CA-CB-CG-OD1
2	B	505[B]	ROC	CA-CB-CG-ND2
2	B	505[A]	ROC	O1-C1-CA-N
2	B	505[A]	ROC	N2-C1-CA-N
2	B	505[A]	ROC	CA-CB-CG-OD1

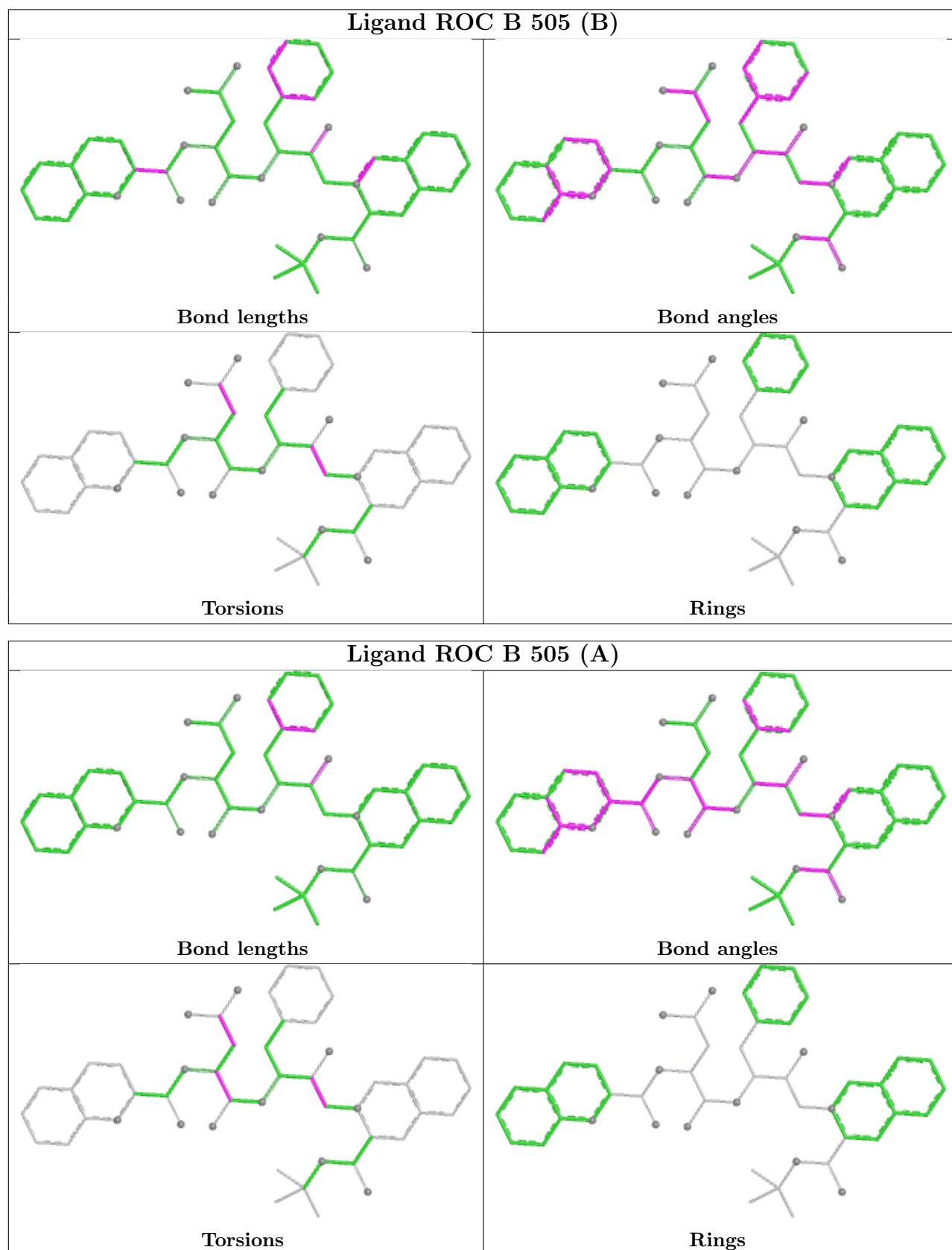
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	505[B]	ROC	1	0
2	B	505[A]	ROC	1	0

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

The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers

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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.