



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:51 AM UTC

PDB ID : 1FPU / pdb_00001fpu
Title : CRYSTAL STRUCTURE OF ABL KINASE DOMAIN IN COMPLEX WITH
A SMALL MOLECULE INHIBITOR
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Deposited on : 2000-08-31
Resolution : 2.40 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)
Xtrriage (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

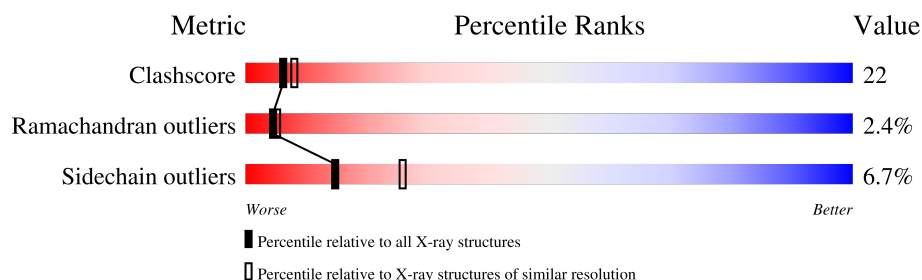
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)

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	293	
1	B	293	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4451 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.

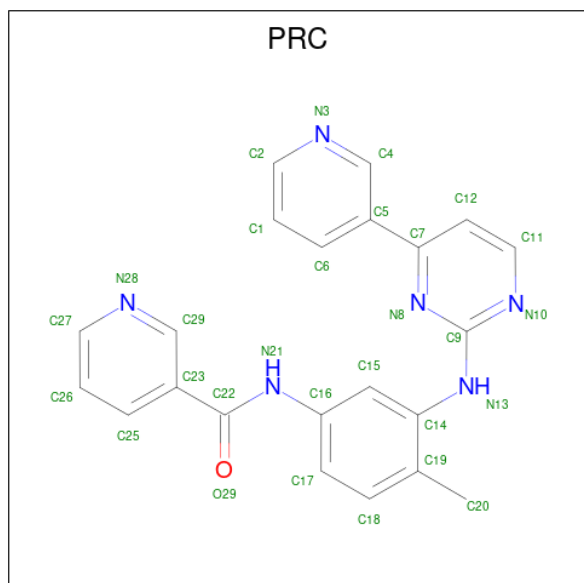
- Molecule 1 is a protein called PROTO-ONCOGENE TYROSINE-PROTEIN KINASE ABL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	275	Total	C	N	O	S	0	0	0
			2148	1389	352	391	16			
1	B	270	Total	C	N	O	S	0	0	0
			2146	1387	350	393	16			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	223	GLY	-	cloning artifact	UNP P00520
A	224	ALA	-	cloning artifact	UNP P00520
A	225	MET	-	cloning artifact	UNP P00520
A	226	ASP	-	cloning artifact	UNP P00520
A	227	PRO	-	cloning artifact	UNP P00520
A	228	SER	-	cloning artifact	UNP P00520

- Molecule 2 is N-[4-METHYL-3-[[4-(3-PYRIDINYL)-2-PYRIMIDINYL]AMINO]PHENYL]-3-PYRIDINECARBOXAMIDE (CCD ID: PRC) (formula: C₂₂H₁₈N₆O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			29	22	6	1		
2	B	1	Total	C	N	O	0	0
			29	22	6	1		

- Molecule 3 is water.

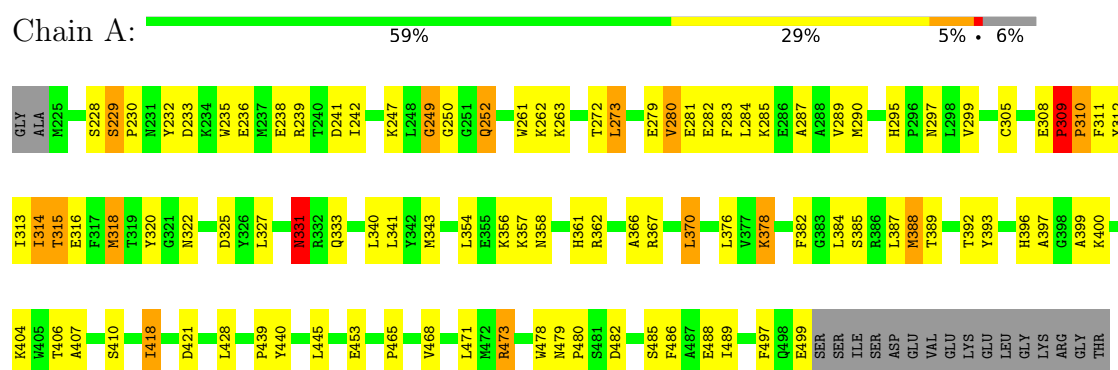
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	67	Total	O	0	0
			67	67		
3	B	32	Total	O	0	0
			32	32		

3 Residue-property plots

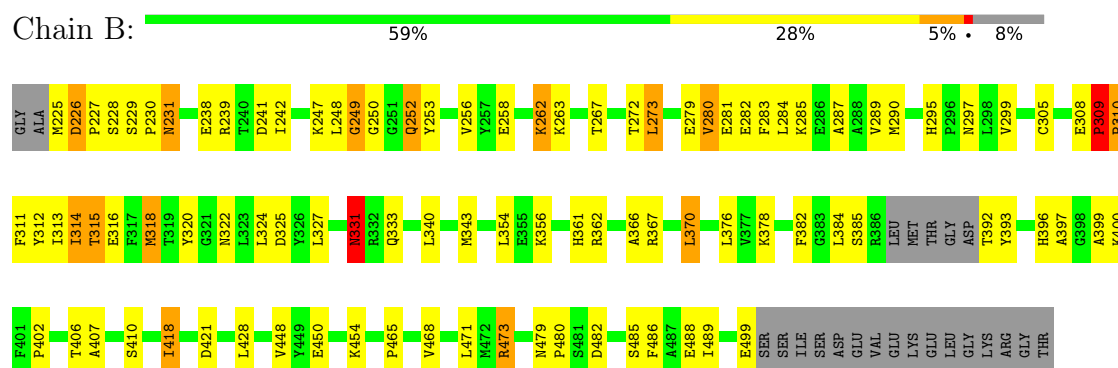
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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



• Molecule 1: PROTO-ONCOGENE TYROSINE-PROTEIN KINASE ABL



4 Data and refinement statistics

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

Property	Value	Source
Space group	F 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	110.70Å 146.10Å 152.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	99.00 – 2.40	Depositor
% Data completeness (in resolution range)	97.8 (99.00-2.40)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.239 , 0.264	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4451	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PRC

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	0/2206	0.97	11/2999 (0.4%)
1	B	0.45	0/2204	0.96	11/2993 (0.4%)
All	All	0.47	0/4410	0.97	22/5992 (0.4%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	230	PRO	N-CA-CB	7.55	110.36	103.34
1	B	320	TYR	N-CA-C	7.43	122.06	112.92
1	A	320	TYR	N-CA-C	6.88	121.39	112.92
1	A	389	THR	N-CA-C	-6.79	104.57	112.92
1	A	309	PRO	N-CA-C	6.53	118.67	110.70
1	B	309	PRO	N-CA-C	6.51	118.65	110.70
1	A	249	GLY	N-CA-C	6.39	121.70	114.67
1	B	249	GLY	N-CA-C	6.09	121.53	114.16
1	B	448	VAL	N-CA-C	5.82	116.01	110.42
1	A	421	ASP	N-CA-C	-5.74	105.00	112.23
1	B	421	ASP	N-CA-C	-5.46	104.97	111.69
1	A	331	ASN	N-CA-C	-5.45	100.22	108.67
1	A	392	THR	N-CA-C	5.44	116.89	111.07
1	B	407	ALA	CA-C-N	5.36	124.81	119.24
1	B	407	ALA	C-N-CA	5.36	124.81	119.24
1	A	407	ALA	CA-C-N	5.32	124.77	119.24
1	A	407	ALA	C-N-CA	5.32	124.77	119.24
1	A	232	TYR	N-CA-C	-5.21	102.68	110.23
1	B	226	ASP	CA-C-N	5.20	126.34	119.84
1	B	226	ASP	C-N-CA	5.20	126.34	119.84
1	B	267	THR	N-CA-C	-5.10	102.14	110.20
1	B	331	ASN	N-CA-C	-5.01	100.34	108.41

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	2148	0	2020	88	0
1	B	2146	0	2046	101	0
2	A	29	0	18	1	0
2	B	29	0	18	1	0
3	A	67	0	0	1	0
3	B	32	0	0	3	0
All	All	4451	0	4102	189	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 22.

All (189) 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:241:ASP:HA	1:B:262:LYS:HE3	1.38	1.05
1:B:262:LYS:HZ3	1:B:263:LYS:HG3	1.22	1.00
1:B:309:PRO:HB2	1:B:310:PRO:HD3	1.52	0.91
1:B:262:LYS:NZ	1:B:263:LYS:HG3	1.87	0.89
1:B:309:PRO:HB2	1:B:310:PRO:CD	2.03	0.89
1:A:309:PRO:HB2	1:A:310:PRO:HD3	1.53	0.87
1:B:262:LYS:HD3	1:B:263:LYS:N	1.89	0.87
1:A:309:PRO:HB2	1:A:310:PRO:CD	2.06	0.85
1:A:295:HIS:HD2	1:A:297:ASN:H	1.21	0.84
1:B:250:GLY:C	1:B:252:GLN:H	1.83	0.83
1:B:295:HIS:HD2	1:B:297:ASN:H	1.25	0.83
1:A:250:GLY:C	1:A:252:GLN:H	1.84	0.81
1:A:295:HIS:CD2	1:A:297:ASN:H	1.99	0.77
1:B:295:HIS:CD2	1:B:297:ASN:H	2.03	0.77
1:B:241:ASP:CG	1:B:262:LYS:HZ1	1.94	0.75
1:B:262:LYS:HD3	1:B:263:LYS:H	1.51	0.73
1:B:249:GLY:HA2	1:B:252:GLN:HG2	1.71	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:290:MET:HE1	1:A:313:ILE:HD13	1.71	0.71
1:B:262:LYS:CD	1:B:263:LYS:H	2.04	0.70
1:B:305:CYS:HB2	1:B:312:TYR:HB2	1.73	0.70
1:A:249:GLY:HA2	1:A:252:GLN:HG2	1.72	0.70
1:B:272:THR:HG22	1:B:273:LEU:H	1.57	0.69
1:A:331:ASN:HD22	1:A:331:ASN:C	2.00	0.69
1:B:272:THR:HG22	1:B:273:LEU:N	2.08	0.69
1:A:305:CYS:HB2	1:A:312:TYR:HB2	1.74	0.69
1:A:279:GLU:O	1:A:282:GLU:N	2.26	0.69
1:A:272:THR:HG22	1:A:273:LEU:N	2.07	0.68
1:B:241:ASP:CA	1:B:262:LYS:HE3	2.18	0.68
1:B:331:ASN:HD22	1:B:331:ASN:C	2.00	0.67
1:A:315:THR:HG22	1:A:316:GLU:O	1.94	0.67
1:B:279:GLU:O	1:B:282:GLU:N	2.27	0.67
1:A:272:THR:HG22	1:A:273:LEU:H	1.59	0.67
1:A:250:GLY:C	1:A:252:GLN:N	2.53	0.66
1:A:331:ASN:ND2	1:A:333:GLN:H	1.93	0.66
1:A:318:MET:HA	1:A:318:MET:HE2	1.77	0.66
1:B:290:MET:HE1	1:B:313:ILE:HD13	1.78	0.66
1:B:331:ASN:ND2	1:B:333:GLN:H	1.94	0.65
1:B:314:ILE:HD12	1:B:314:ILE:N	2.12	0.64
1:A:309:PRO:CB	1:A:310:PRO:HD3	2.28	0.64
1:B:392:THR:O	1:B:402:PRO:HA	1.98	0.64
1:B:309:PRO:CB	1:B:310:PRO:HD3	2.26	0.64
1:B:318:MET:HA	1:B:318:MET:HE2	1.80	0.63
1:B:418:ILE:HD11	1:B:486:PHE:HD2	1.64	0.63
1:B:309:PRO:CB	1:B:310:PRO:CD	2.76	0.62
1:A:241:ASP:OD2	1:A:263:LYS:HE3	2.00	0.62
1:A:331:ASN:HD21	1:A:333:GLN:HB2	1.64	0.62
1:A:314:ILE:HD12	1:A:314:ILE:N	2.15	0.61
1:B:262:LYS:CD	1:B:263:LYS:N	2.63	0.61
1:B:340:LEU:HA	1:B:343:MET:HE2	1.82	0.61
1:A:340:LEU:HA	1:A:343:MET:HE2	1.82	0.61
1:A:309:PRO:CB	1:A:310:PRO:CD	2.79	0.60
1:B:485:SER:OG	1:B:488:GLU:HG2	2.01	0.60
1:B:227:PRO:HG2	3:B:542:HOH:O	2.00	0.60
1:A:418:ILE:HD11	1:A:486:PHE:HD2	1.67	0.59
1:B:315:THR:HG22	1:B:316:GLU:O	2.02	0.59
1:A:239:ARG:HH12	1:A:310:PRO:HD2	1.67	0.58
1:B:250:GLY:C	1:B:252:GLN:N	2.53	0.58
1:B:331:ASN:HD21	1:B:333:GLN:HB2	1.68	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:239:ARG:HH12	1:B:310:PRO:HD2	1.69	0.57
1:A:331:ASN:HD22	1:A:333:GLN:H	1.52	0.57
1:B:393:TYR:HB3	1:B:400:LYS:HB3	1.87	0.56
1:A:485:SER:OG	1:A:488:GLU:HG2	2.04	0.56
1:A:297:ASN:HA	1:A:378:LYS:HG3	1.87	0.56
1:B:282:GLU:HA	1:B:282:GLU:OE1	2.05	0.55
1:A:282:GLU:OE1	1:A:282:GLU:HA	2.07	0.55
1:B:253:TYR:CE2	1:B:382:PHE:CD1	2.95	0.55
1:B:285:LYS:O	1:B:289:VAL:HG23	2.07	0.55
1:A:315:THR:HG21	2:A:1:PRC:H13N	1.71	0.55
1:A:295:HIS:HD2	1:A:297:ASN:N	1.99	0.55
1:B:465:PRO:HG2	1:B:468:VAL:CG2	2.37	0.55
1:A:465:PRO:HG2	1:A:468:VAL:CG2	2.37	0.55
1:A:249:GLY:CA	1:A:252:GLN:HG2	2.38	0.54
1:A:393:TYR:HB3	1:A:400:LYS:HB3	1.89	0.54
1:B:249:GLY:CA	1:B:252:GLN:HG2	2.36	0.54
1:B:290:MET:HE1	1:B:313:ILE:HG21	1.89	0.53
1:B:279:GLU:O	1:B:280:VAL:C	2.51	0.53
1:A:279:GLU:O	1:A:280:VAL:C	2.51	0.53
1:A:479:ASN:HD22	1:A:480:PRO:HD2	1.73	0.53
1:A:285:LYS:O	1:A:289:VAL:HG23	2.09	0.53
1:B:385:SER:HB2	1:B:400:LYS:HB2	1.91	0.53
1:B:230:PRO:O	1:B:231:ASN:CG	2.52	0.52
1:B:250:GLY:O	1:B:252:GLN:N	2.42	0.52
1:B:299:VAL:HG12	3:B:543:HOH:O	2.08	0.52
1:B:309:PRO:O	1:B:311:PHE:N	2.42	0.52
1:B:450:GLU:OE2	1:B:454:LYS:NZ	2.41	0.52
1:A:250:GLY:O	1:A:252:GLN:N	2.42	0.52
1:A:309:PRO:O	1:A:311:PHE:CD2	2.63	0.52
1:B:249:GLY:C	1:B:252:GLN:HG2	2.35	0.52
1:B:473:ARG:HH11	1:B:473:ARG:HG3	1.75	0.52
1:B:262:LYS:HZ3	1:B:263:LYS:CG	2.10	0.52
1:A:249:GLY:C	1:A:252:GLN:HG2	2.35	0.52
1:A:385:SER:HB2	1:A:400:LYS:HB2	1.91	0.52
1:B:331:ASN:HD22	1:B:333:GLN:H	1.57	0.52
1:A:387:LEU:O	1:A:388:MET:C	2.53	0.51
1:B:309:PRO:O	1:B:311:PHE:CD2	2.62	0.51
1:B:479:ASN:HD22	1:B:480:PRO:HD2	1.75	0.51
1:B:272:THR:CG2	1:B:273:LEU:H	2.23	0.51
1:A:309:PRO:O	1:A:311:PHE:N	2.43	0.51
1:A:397:ALA:C	1:A:399:ALA:H	2.18	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:397:ALA:C	1:B:399:ALA:H	2.18	0.51
1:B:322:ASN:HD21	1:B:325:ASP:H	1.59	0.50
1:A:309:PRO:O	1:A:310:PRO:C	2.55	0.50
1:A:331:ASN:C	1:A:331:ASN:ND2	2.68	0.50
1:B:418:ILE:HD11	1:B:486:PHE:CD2	2.45	0.50
1:B:331:ASN:C	1:B:331:ASN:ND2	2.67	0.50
1:B:253:TYR:OH	1:B:322:ASN:HB3	2.12	0.50
1:A:382:PHE:HB2	1:A:384:LEU:HG	1.94	0.50
1:A:473:ARG:HH11	1:A:473:ARG:HG3	1.76	0.49
1:B:272:THR:CG2	1:B:273:LEU:N	2.75	0.49
1:B:382:PHE:HB2	1:B:384:LEU:HG	1.93	0.49
1:B:241:ASP:OD2	1:B:263:LYS:HE3	2.12	0.49
1:A:272:THR:CG2	1:A:273:LEU:H	2.24	0.49
1:B:479:ASN:HB3	1:B:482:ASP:OD2	2.13	0.49
1:B:314:ILE:HD12	1:B:314:ILE:H	1.78	0.49
1:A:361:HIS:O	1:A:362:ARG:HB2	2.13	0.48
1:B:309:PRO:O	1:B:310:PRO:C	2.55	0.48
1:B:238:GLU:O	1:B:241:ASP:HB2	2.12	0.48
1:A:239:ARG:HH22	1:A:308:GLU:C	2.22	0.48
1:A:322:ASN:HD21	1:A:325:ASP:H	1.62	0.48
1:B:322:ASN:ND2	1:B:325:ASP:H	2.11	0.48
1:A:290:MET:HE1	1:A:313:ILE:HG21	1.96	0.48
1:A:327:LEU:HD21	1:A:343:MET:HE1	1.96	0.47
1:A:479:ASN:HB3	1:A:482:ASP:OD2	2.14	0.47
1:A:322:ASN:ND2	1:A:325:ASP:H	2.12	0.47
1:B:279:GLU:O	1:B:281:GLU:N	2.48	0.47
1:B:361:HIS:O	1:B:362:ARG:HB2	2.13	0.47
1:A:239:ARG:HH12	1:A:310:PRO:CD	2.29	0.46
1:B:290:MET:CE	1:B:313:ILE:HD13	2.46	0.46
1:B:287:ALA:HA	1:B:290:MET:HE3	1.97	0.46
1:A:238:GLU:O	1:A:241:ASP:HB2	2.15	0.46
1:A:396:HIS:O	1:A:396:HIS:ND1	2.48	0.46
1:B:249:GLY:HA2	1:B:252:GLN:CG	2.43	0.46
1:B:256:VAL:HG11	2:B:2:PRC:C11	2.45	0.46
1:B:295:HIS:HD2	1:B:297:ASN:N	2.03	0.46
1:A:370:LEU:HD13	1:A:382:PHE:HZ	1.80	0.46
1:B:471:LEU:HD21	1:B:489:ILE:HG23	1.98	0.46
1:A:356:LYS:HB2	1:A:356:LYS:HE3	1.78	0.45
1:A:272:THR:CG2	1:A:273:LEU:N	2.75	0.45
1:B:242:ILE:CD1	1:B:314:ILE:HG12	2.47	0.45
1:A:471:LEU:HD21	1:A:489:ILE:HG23	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:283:PHE:CE2	1:B:311:PHE:HB3	2.52	0.45
1:B:327:LEU:HD21	1:B:343:MET:HE1	1.98	0.45
1:B:228:SER:O	1:B:229:SER:C	2.59	0.45
1:A:233:ASP:HB3	1:A:236:GLU:HG2	1.98	0.45
1:A:242:ILE:CD1	1:A:314:ILE:HG12	2.47	0.45
1:A:290:MET:CE	1:A:313:ILE:HD13	2.45	0.44
1:A:299:VAL:HG22	1:A:315:THR:HG23	1.98	0.44
1:B:396:HIS:O	1:B:396:HIS:ND1	2.47	0.44
1:B:239:ARG:HH22	1:B:308:GLU:C	2.24	0.44
1:B:370:LEU:HD13	1:B:382:PHE:HZ	1.81	0.44
1:B:273:LEU:HD11	1:B:279:GLU:CB	2.47	0.44
1:B:239:ARG:HH12	1:B:310:PRO:CD	2.30	0.44
1:A:406:THR:CG2	1:A:410:SER:HB2	2.48	0.44
1:A:357:LYS:O	1:A:358:ASN:HB2	2.18	0.44
1:A:249:GLY:HA2	1:A:252:GLN:CG	2.45	0.43
1:A:418:ILE:HD11	1:A:486:PHE:CD2	2.49	0.43
1:B:473:ARG:HG3	1:B:473:ARG:NH1	2.33	0.43
1:B:406:THR:CG2	1:B:410:SER:HB2	2.49	0.43
1:A:279:GLU:O	1:A:281:GLU:N	2.52	0.43
1:B:397:ALA:C	1:B:399:ALA:N	2.76	0.42
1:B:253:TYR:HE2	1:B:382:PHE:CD1	2.36	0.42
1:A:453:GLU:HG3	1:A:478:TRP:CH2	2.55	0.42
1:A:261:TRP:O	1:A:262:LYS:C	2.62	0.42
1:A:485:SER:HG	1:A:488:GLU:HG2	1.84	0.42
1:B:299:VAL:HG22	1:B:315:THR:HG23	2.00	0.42
1:A:233:ASP:OD1	1:A:235:TRP:HB2	2.20	0.42
1:A:341:LEU:HD21	1:A:497:PHE:HA	2.02	0.42
1:B:262:LYS:HD2	1:B:263:LYS:H	1.81	0.42
1:A:249:GLY:HA2	1:A:252:GLN:NE2	2.34	0.42
1:B:241:ASP:OD2	1:B:262:LYS:NZ	2.49	0.42
1:B:249:GLY:HA2	1:B:252:GLN:NE2	2.35	0.41
1:A:366:ALA:N	1:A:428:LEU:HD13	2.34	0.41
1:A:473:ARG:HG3	1:A:473:ARG:NH1	2.35	0.41
1:A:479:ASN:HD22	1:A:480:PRO:CD	2.34	0.41
1:B:315:THR:CG2	1:B:316:GLU:N	2.83	0.41
1:B:322:ASN:ND2	1:B:324:LEU:HB3	2.36	0.41
1:B:356:LYS:HE3	1:B:356:LYS:HB2	1.78	0.41
1:A:314:ILE:HD12	1:A:314:ILE:H	1.85	0.41
1:A:367:ARG:HD3	3:A:544:HOH:O	2.21	0.41
1:A:404:LYS:HE3	1:A:445:LEU:HD23	2.02	0.41
1:A:439:PRO:O	1:A:440:TYR:C	2.63	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:248:LEU:HD21	1:B:258:GLU:HB2	2.02	0.41
1:B:340:LEU:HD23	1:B:343:MET:CE	2.51	0.41
1:A:233:ASP:OD1	1:A:235:TRP:N	2.52	0.41
1:A:283:PHE:CE2	1:A:311:PHE:HB3	2.56	0.40
1:B:366:ALA:N	1:B:428:LEU:HD13	2.35	0.40
1:A:228:SER:O	1:A:229:SER:CB	2.70	0.40
1:A:287:ALA:HA	1:A:290:MET:HE3	2.02	0.40
1:B:367:ARG:HD3	3:B:526:HOH:O	2.22	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	273/293 (93%)	244 (89%)	22 (8%)	7 (3%)	4	4
1	B	266/293 (91%)	236 (89%)	24 (9%)	6 (2%)	5	6
All	All	539/586 (92%)	480 (89%)	46 (8%)	13 (2%)	4	5

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	252	GLN
1	A	280	VAL
1	A	309	PRO
1	B	231	ASN
1	B	252	GLN
1	B	280	VAL
1	B	309	PRO
1	A	310	PRO
1	B	247	LYS
1	B	310	PRO

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Mol	Chain	Res	Type
1	A	247	LYS
1	A	388	MET
1	A	229	SER

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	214/255 (84%)	201 (94%)	13 (6%)	17	30
1	B	221/255 (87%)	205 (93%)	16 (7%)	13	23
All	All	435/510 (85%)	406 (93%)	29 (7%)	15	26

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	273	LEU
1	A	284	LEU
1	A	314	ILE
1	A	315	THR
1	A	318	MET
1	A	331	ASN
1	A	354	LEU
1	A	370	LEU
1	A	376	LEU
1	A	378	LYS
1	A	418	ILE
1	A	473	ARG
1	A	499	GLU
1	B	225	MET
1	B	226	ASP
1	B	262	LYS
1	B	273	LEU
1	B	284	LEU
1	B	314	ILE
1	B	315	THR

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Mol	Chain	Res	Type
1	B	318	MET
1	B	331	ASN
1	B	354	LEU
1	B	370	LEU
1	B	376	LEU
1	B	378	LYS
1	B	418	ILE
1	B	473	ARG
1	B	499	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	246	HIS
1	A	295	HIS
1	A	297	ASN
1	A	322	ASN
1	A	331	ASN
1	A	479	ASN
1	A	490	HIS
1	B	246	HIS
1	B	295	HIS
1	B	297	ASN
1	B	322	ASN
1	B	331	ASN
1	B	479	ASN

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

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	PRC	B	2	-	32,32,32	1.50	6 (18%)	43,43,43	2.82	14 (32%)
2	PRC	A	1	-	32,32,32	1.41	5 (15%)	43,43,43	2.74	12 (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	PRC	B	2	-	-	1/16/16/16	0/4/4/4
2	PRC	A	1	-	-	2/16/16/16	0/4/4/4

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2	PRC	C15-C16	3.07	1.44	1.39
2	A	1	PRC	C15-C14	2.79	1.43	1.39
2	A	1	PRC	C5-C7	-2.68	1.44	1.49
2	B	2	PRC	C15-C14	2.57	1.43	1.39
2	B	2	PRC	C7-N8	2.53	1.38	1.34
2	B	2	PRC	C1-C6	2.51	1.43	1.38
2	A	1	PRC	C25-C23	2.44	1.43	1.39
2	A	1	PRC	C15-C16	2.44	1.43	1.39
2	B	2	PRC	C9-N10	2.33	1.37	1.34
2	B	2	PRC	C25-C23	2.29	1.42	1.39
2	A	1	PRC	C9-N10	2.17	1.37	1.34

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	PRC	C7-N8-C9	8.98	123.72	116.81
2	A	1	PRC	N10-C9-N8	-8.44	118.26	126.42
2	B	2	PRC	N10-C9-N8	-8.44	118.26	126.42
2	B	2	PRC	C7-N8-C9	8.29	123.19	116.81
2	B	2	PRC	C11-N10-C9	8.00	122.11	115.42
2	A	1	PRC	C11-N10-C9	7.84	121.97	115.42
2	B	2	PRC	C20-C19-C14	5.25	127.17	121.23
2	A	1	PRC	C20-C19-C14	4.40	126.22	121.23
2	A	1	PRC	C12-C11-N10	-3.70	119.43	123.97
2	B	2	PRC	C12-C11-N10	-3.59	119.58	123.97
2	B	2	PRC	C27-N28-C29	3.29	122.61	116.85
2	B	2	PRC	C2-N3-C4	3.27	122.57	116.85
2	A	1	PRC	C27-N28-C29	3.19	122.44	116.85
2	B	2	PRC	C12-C7-N8	-3.14	117.83	121.97
2	A	1	PRC	C2-N3-C4	3.12	122.31	116.85
2	B	2	PRC	C20-C19-C18	-3.00	114.50	120.28
2	B	2	PRC	C19-C14-N13	2.91	123.72	118.68
2	B	2	PRC	C5-C7-N8	2.84	120.14	116.04
2	A	1	PRC	C12-C7-N8	-2.78	118.30	121.97
2	A	1	PRC	C20-C19-C18	-2.44	115.58	120.28
2	B	2	PRC	O29-C22-C23	-2.39	116.17	120.90
2	B	2	PRC	C15-C14-N13	-2.28	116.44	121.12
2	B	2	PRC	C23-C29-N28	-2.23	120.19	123.50
2	A	1	PRC	C19-C14-N13	2.15	122.40	118.68
2	A	1	PRC	C23-C29-N28	-2.02	120.49	123.50
2	A	1	PRC	O29-C22-C23	-2.00	116.93	120.90

There are no chirality outliers.

All (3) torsion outliers are listed below:

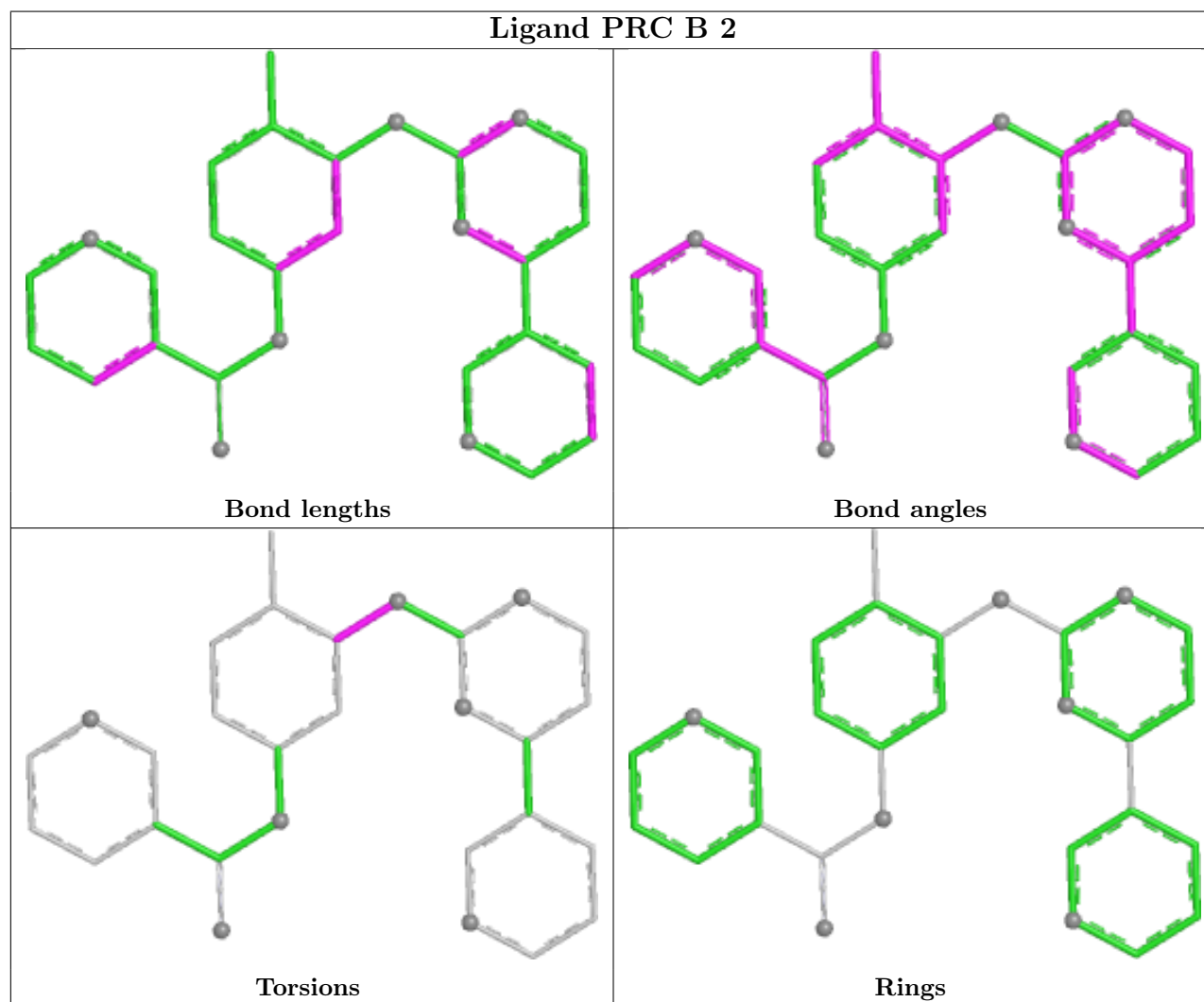
Mol	Chain	Res	Type	Atoms
2	A	1	PRC	C15-C14-N13-C9
2	A	1	PRC	C19-C14-N13-C9
2	B	2	PRC	C15-C14-N13-C9

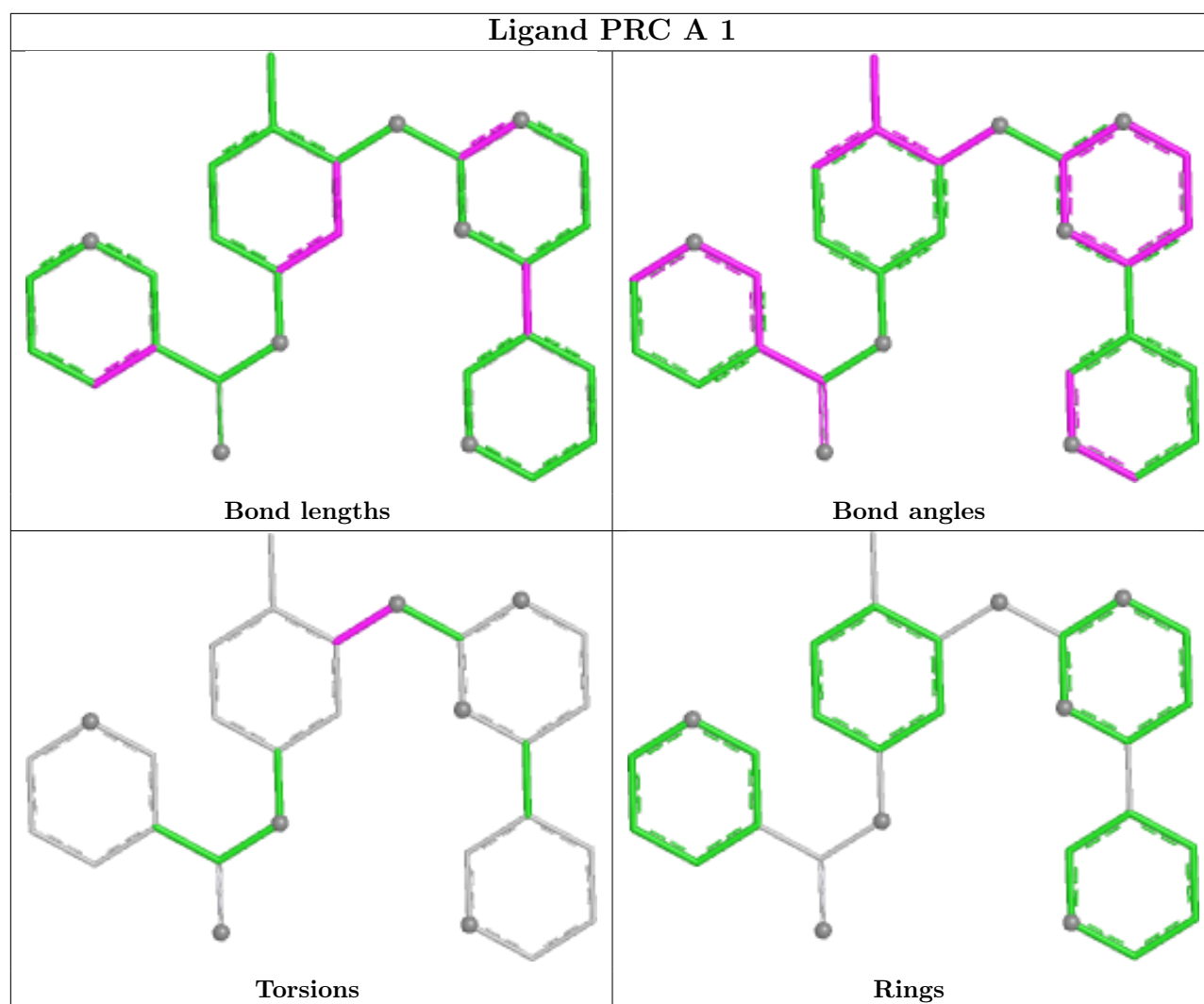
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2	PRC	1	0
2	A	1	PRC	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 [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

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.