



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:28 AM UTC

PDB ID : 3SKF / pdb_00003skf
Title : Crystal structure of beta-site app-cleaving enzyme 1 (BACE-WT) complex with (2S)-2-((3S)-3-(acetylamino)-3-(butan-2-yl)-2-oxopyrrolidin-1-yl)-N-((2S,3R)-3-hydroxy-4-((3-methoxybenzyl)amino)-1-phenylbutan-2-yl)-4-phenylbutanamide
Authors : Muckelbauer, J.K.
Deposited on : 2011-06-22
Resolution : 3.00 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](#) 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

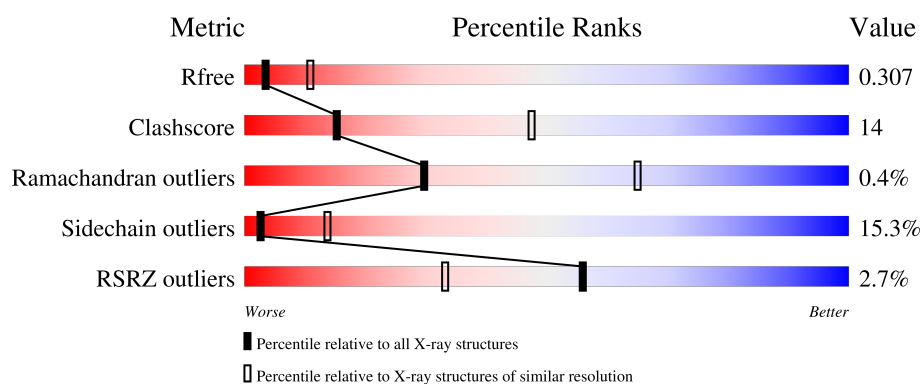
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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 $\leq 5\%$. 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	455	
1	B	455	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6361 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 Beta-secretase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	387	Total	C	N	O	S	0	0	0
			3038	1946	505	573	14			
1	B	387	Total	C	N	O	S	0	0	0
			3039	1947	505	573	14			

There are 28 discrepancies between the modelled and reference sequences:

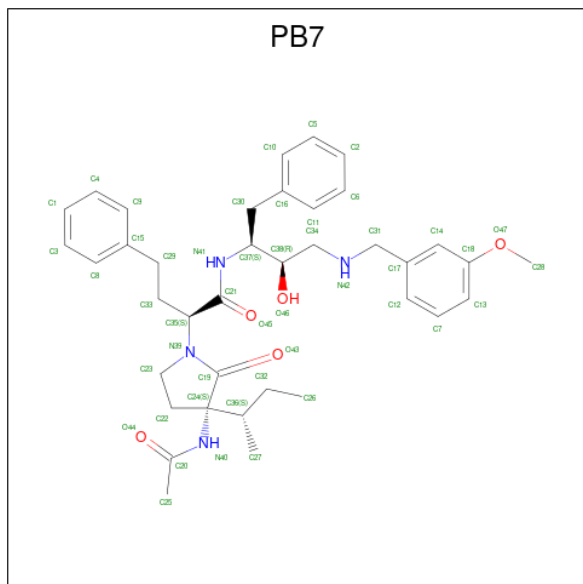
Chain	Residue	Modelled	Actual	Comment	Reference
A	-61	MET	-	expression tag	UNP P56817
A	-60	ALA	-	expression tag	UNP P56817
A	-59	SER	-	expression tag	UNP P56817
A	-58	MET	-	expression tag	UNP P56817
A	-57	THR	-	expression tag	UNP P56817
A	-56	GLY	-	expression tag	UNP P56817
A	-55	GLY	-	expression tag	UNP P56817
A	-54	GLN	-	expression tag	UNP P56817
A	-53	GLN	-	expression tag	UNP P56817
A	-52	MET	-	expression tag	UNP P56817
A	-51	GLY	-	expression tag	UNP P56817
A	-50	ARG	-	expression tag	UNP P56817
A	-49	GLY	-	expression tag	UNP P56817
A	-48	SER	-	expression tag	UNP P56817
B	-61	MET	-	expression tag	UNP P56817
B	-60	ALA	-	expression tag	UNP P56817
B	-59	SER	-	expression tag	UNP P56817
B	-58	MET	-	expression tag	UNP P56817
B	-57	THR	-	expression tag	UNP P56817
B	-56	GLY	-	expression tag	UNP P56817
B	-55	GLY	-	expression tag	UNP P56817
B	-54	GLN	-	expression tag	UNP P56817
B	-53	GLN	-	expression tag	UNP P56817
B	-52	MET	-	expression tag	UNP P56817
B	-51	GLY	-	expression tag	UNP P56817

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-50	ARG	-	expression tag	UNP P56817
B	-49	GLY	-	expression tag	UNP P56817
B	-48	SER	-	expression tag	UNP P56817

- Molecule 2 is (2S)-2-[(3S)-3-(acetylamino)-3-[(2S)-butan-2-yl]-2-oxopyrrolidin-1-yl]-N-[(2S,3R)-3-hydroxy-4-[(3-methoxybenzyl)amino]-1-phenylbutan-2-yl]-4-phenylbutanamide (CCD ID: PB7) (formula: C₃₈H₅₀N₄O₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			47	38	4	5		
2	B	1	Total	C	N	O	0	0
			47	38	4	5		

- Molecule 3 is IODIDE ION (CCD ID: IOD) (formula: I).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	I	0	0
			1	1		
3	B	1	Total	I	0	0
			1	1		

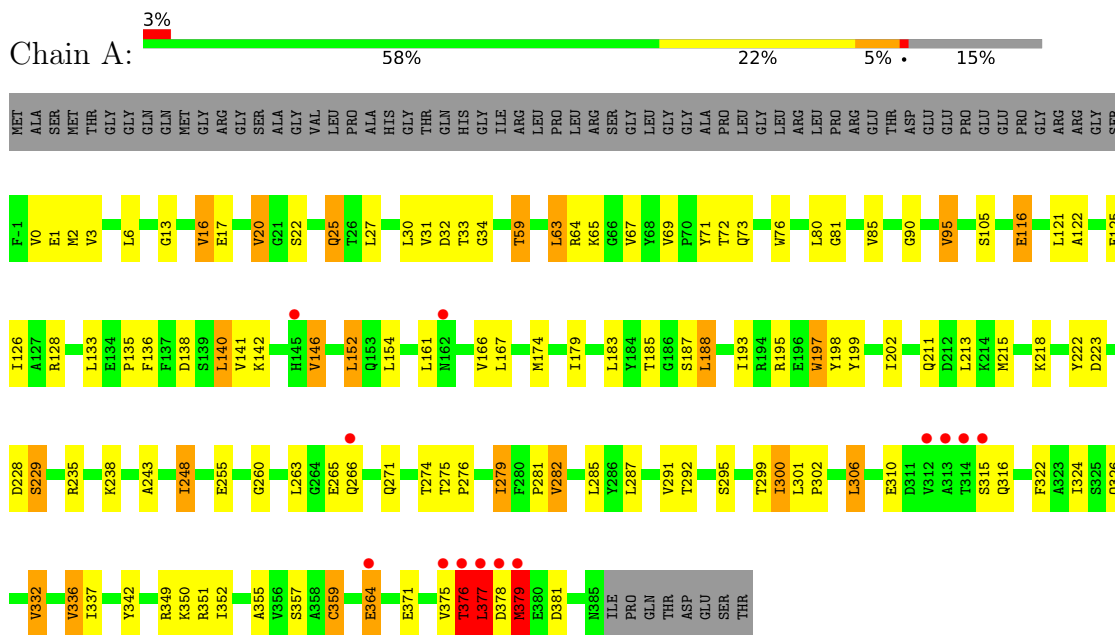
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	89	Total 89	O 89	0	0
4	B	99	Total 99	O 99	0	0

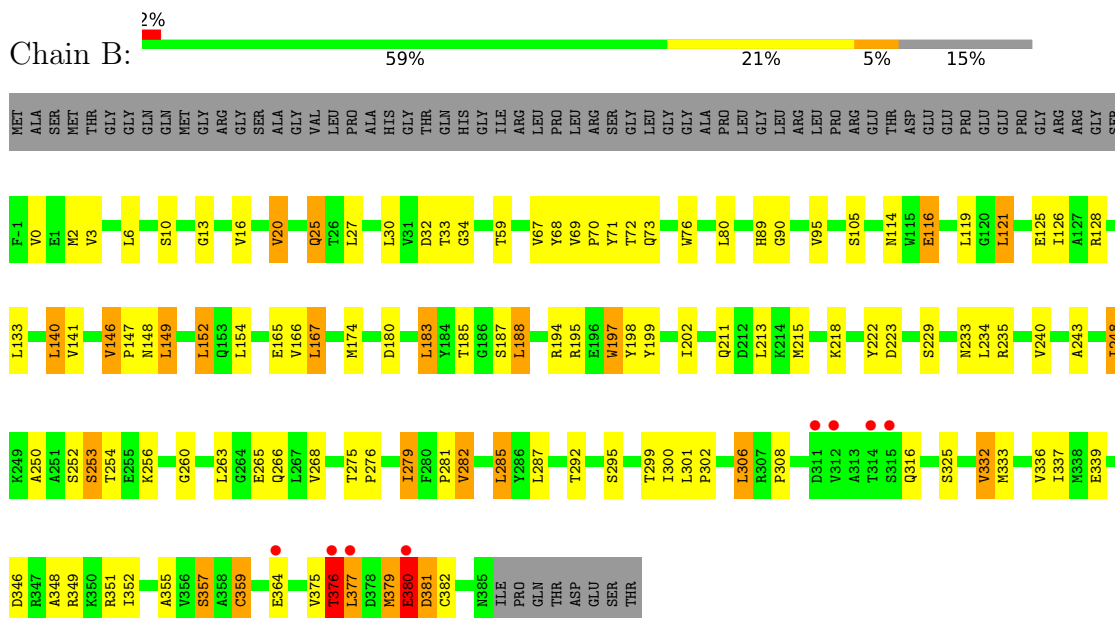
3 Residue-property plots

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 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: Beta-secretase 1



• Molecule 1: Beta-secretase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	90.29Å 130.26Å 85.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.06 – 3.00 34.06 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.4 (34.06-3.00) 99.4 (34.06-3.00)	Depositor EDS
R_{merge}	0.23	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.63 (at 3.00Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.252 , 0.313 0.252 , 0.307	Depositor DCC
R_{free} test set	1064 reflections (5.14%)	wwPDB-VP
Wilson B-factor (Å ²)	40.8	Xtriage
Anisotropy	0.069	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 41.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.010 for l,-k,h	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	6361	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 39.64 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.0775e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PB7, IOD

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	0.58	0/3116	0.88	5/4238 (0.1%)
1	B	0.57	0/3117	0.88	2/4239 (0.0%)
All	All	0.58	0/6233	0.88	7/8477 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	2

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	379	MET	N-CA-C	7.02	123.14	112.54
1	A	379	MET	CA-C-N	-6.35	111.43	122.36
1	A	379	MET	C-N-CA	-6.35	111.43	122.36
1	B	381	ASP	N-CA-C	-6.04	106.26	113.21
1	B	332	VAL	CB-CA-C	5.67	117.47	110.73
1	A	377	LEU	N-CA-C	5.39	116.91	109.15
1	A	315	SER	N-CA-C	5.07	116.88	110.33

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	379	MET	Peptide

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Mol	Chain	Res	Type	Group
1	B	380	GLU	Peptide

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	3038	0	2950	82	0
1	B	3039	0	2952	88	0
2	A	47	0	50	2	0
2	B	47	0	50	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	89	0	0	5	0
4	B	99	0	0	2	0
All	All	6361	0	6002	165	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 14.

All (165) 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:254:THR:HB	1:B:279:ILE:HD11	1.24	1.18
1:A:276:PRO:O	1:A:279:ILE:HG22	1.63	0.96
1:A:235:ARG:HB2	1:A:332:VAL:HG23	1.47	0.95
1:B:67:VAL:HB	1:B:80:LEU:HD21	1.50	0.94
1:B:276:PRO:O	1:B:279:ILE:HG22	1.73	0.88
1:B:235:ARG:HB2	1:B:332:VAL:HG12	1.59	0.84
1:A:67:VAL:HB	1:A:80:LEU:HD21	1.61	0.80
1:B:254:THR:HB	1:B:279:ILE:CD1	2.12	0.77
1:B:13:GLY:HA3	1:B:30:LEU:HD11	1.70	0.74
1:A:310:GLU:HG2	4:A:481:HOH:O	1.91	0.71
1:B:2:MET:HG2	1:B:90:GLY:HA2	1.72	0.70
1:A:2:MET:HG2	1:A:90:GLY:HA2	1.73	0.69
1:B:188:LEU:HD13	1:B:355:ALA:HB2	1.75	0.69
1:B:295:SER:CB	1:B:377:LEU:HD23	2.23	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:282:VAL:HG13	1:B:299:THR:HG23	1.77	0.66
1:A:34:GLY:O	2:A:394:PB7:H14	1.95	0.66
1:A:6:LEU:HD12	1:A:16:VAL:HG12	1.78	0.66
1:A:67:VAL:HB	1:A:80:LEU:CD2	2.25	0.66
1:A:379:MET:O	1:A:379:MET:CG	2.43	0.66
1:A:140:LEU:HD13	1:A:146:VAL:HG11	1.79	0.65
1:B:215:MET:HE1	1:B:240:VAL:HA	1.79	0.64
1:A:276:PRO:HB2	1:A:279:ILE:CG2	2.28	0.64
1:B:380:GLU:C	1:B:382:CYS:H	2.05	0.63
1:A:379:MET:O	1:A:379:MET:HG3	1.99	0.63
1:A:276:PRO:HB2	1:A:279:ILE:HG22	1.80	0.63
1:A:215:MET:HE3	1:A:243:ALA:HB3	1.81	0.62
1:A:22:SER:HB2	1:A:59:THR:OG1	2.01	0.61
1:B:33:THR:H	1:B:229:SER:HB3	1.66	0.60
1:B:197:TRP:H	1:B:197:TRP:HD1	1.47	0.60
1:A:33:THR:H	1:A:229:SER:CB	2.15	0.60
1:A:282:VAL:HG13	1:A:299:THR:HG23	1.84	0.60
1:A:235:ARG:HB2	1:A:332:VAL:CG2	2.29	0.60
1:B:197:TRP:H	1:B:197:TRP:CD1	2.19	0.59
1:A:188:LEU:HD13	1:A:355:ALA:HB2	1.85	0.59
1:B:69:VAL:HG22	1:B:128:ARG:HB2	1.85	0.59
1:A:300:ILE:HD13	1:A:337:ILE:HD12	1.84	0.58
1:B:33:THR:H	1:B:229:SER:CB	2.15	0.58
1:A:376:THR:OG1	1:B:364:GLU:CD	2.47	0.57
1:A:95:VAL:HG11	1:A:140:LEU:HA	1.86	0.57
1:B:295:SER:CB	1:B:377:LEU:CD2	2.83	0.56
1:B:6:LEU:HD11	1:B:174:MET:HB2	1.87	0.56
1:A:197:TRP:CG	1:A:198:TYR:H	2.24	0.56
1:A:69:VAL:HG22	1:A:128:ARG:HB2	1.88	0.56
1:A:125:GLU:OE2	1:A:195:ARG:NH2	2.38	0.55
1:B:285:LEU:HD23	1:B:333:MET:HE2	1.89	0.55
1:B:125:GLU:OE2	1:B:195:ARG:NH2	2.37	0.54
1:A:197:TRP:H	1:A:197:TRP:HD1	1.55	0.54
1:B:180:ASP:HB3	1:B:183:LEU:HD22	1.89	0.54
1:B:260:GLY:C	1:B:266:GLN:HB3	2.32	0.54
1:B:375:VAL:O	1:B:377:LEU:N	2.39	0.54
1:A:64:ARG:NE	4:A:420:HOH:O	2.40	0.54
1:B:148:ASN:HB3	1:B:348:ALA:HB2	1.91	0.53
1:A:179:ILE:HG23	1:A:342:TYR:HE2	1.74	0.53
1:A:375:VAL:O	1:A:377:LEU:N	2.41	0.53
1:B:197:TRP:CG	1:B:198:TYR:H	2.27	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:89:HIS:HD2	4:B:457:HOH:O	1.92	0.52
1:A:199:TYR:HB3	1:A:352:ILE:HD11	1.92	0.52
1:A:248:ILE:HG13	1:A:281:PRO:HG2	1.91	0.51
1:A:197:TRP:H	1:A:197:TRP:CD1	2.27	0.51
1:B:32:ASP:HA	1:B:229:SER:HB3	1.92	0.51
1:B:375:VAL:C	1:B:377:LEU:H	2.18	0.51
1:A:271:GLN:O	1:A:274:THR:HB	2.11	0.51
1:A:375:VAL:C	1:A:377:LEU:H	2.19	0.51
1:B:13:GLY:HA3	1:B:30:LEU:CD1	2.40	0.51
1:B:199:TYR:HB3	1:B:352:ILE:HD11	1.93	0.51
1:B:6:LEU:CD1	1:B:174:MET:HB2	2.41	0.51
1:B:67:VAL:HB	1:B:80:LEU:CD2	2.33	0.50
1:B:215:MET:HE3	1:B:243:ALA:HB3	1.94	0.50
1:B:295:SER:HB3	1:B:377:LEU:CD2	2.42	0.50
1:B:295:SER:HB2	1:B:377:LEU:HD23	1.94	0.50
1:A:33:THR:H	1:A:229:SER:HB3	1.75	0.50
1:B:34:GLY:O	2:B:394:PB7:H14	2.12	0.50
1:A:364:GLU:OE2	1:B:376:THR:N	2.40	0.49
1:A:33:THR:H	1:A:229:SER:HB2	1.76	0.49
1:A:6:LEU:CD1	1:A:16:VAL:HG12	2.41	0.49
1:B:27:LEU:HD22	1:B:116:GLU:HG3	1.94	0.49
1:A:13:GLY:HA3	1:A:30:LEU:CD1	2.42	0.49
1:A:279:ILE:HA	1:B:211:GLN:HG3	1.94	0.49
1:B:380:GLU:C	1:B:382:CYS:N	2.71	0.48
1:A:63:LEU:HD22	1:A:81:GLY:HA2	1.95	0.48
1:B:194:ARG:HG3	1:B:202:ILE:HD11	1.95	0.48
1:B:119:LEU:HG	1:B:121:LEU:HD22	1.95	0.48
1:A:13:GLY:HA3	1:A:30:LEU:HD11	1.94	0.48
2:A:394:PB7:H32A	2:A:394:PB7:O43	2.13	0.48
1:B:301:LEU:HB3	1:B:302:PRO:HD2	1.96	0.48
1:A:302:PRO:O	1:A:306:LEU:HB2	2.14	0.47
1:A:228:ASP:HB3	1:A:332:VAL:HG12	1.95	0.47
1:A:2:MET:CG	1:A:90:GLY:HA2	2.42	0.47
1:A:152:LEU:CD1	1:A:174:MET:HG3	2.45	0.47
1:A:260:GLY:C	1:A:266:GLN:HB3	2.40	0.47
1:B:215:MET:CE	1:B:243:ALA:HB3	2.44	0.47
1:A:135:PRO:HD2	1:A:138:ASP:OD2	2.15	0.47
1:A:364:GLU:CD	1:B:376:THR:OG1	2.58	0.46
1:B:71:TYR:HE2	1:B:76:TRP:HE1	1.62	0.46
1:B:222:TYR:HA	1:B:223:ASP:HA	1.69	0.46
1:B:300:ILE:HD13	1:B:337:ILE:HD12	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:248:ILE:HG13	1:A:281:PRO:CG	2.46	0.46
1:A:222:TYR:HA	1:A:223:ASP:HA	1.72	0.46
1:A:336:VAL:HG12	1:A:337:ILE:HD13	1.98	0.46
1:B:149:LEU:HD22	1:B:346:ASP:HA	1.98	0.46
1:B:248:ILE:HG13	1:B:281:PRO:CG	2.46	0.46
1:A:27:LEU:HD22	1:A:116:GLU:HG3	1.98	0.45
1:B:349:ARG:HH11	1:B:351:ARG:NH2	2.13	0.45
1:B:188:LEU:HD13	1:B:355:ALA:CB	2.43	0.45
1:A:32:ASP:HA	1:A:229:SER:HB3	1.97	0.45
1:A:6:LEU:HD11	1:A:174:MET:HB2	1.98	0.45
1:B:282:VAL:CG1	1:B:299:THR:HG23	2.46	0.45
1:B:379:MET:HG2	1:B:380:GLU:HB3	1.99	0.45
1:B:233:ASN:HD22	1:B:325:SER:HB3	1.81	0.44
1:A:6:LEU:CD1	1:A:174:MET:HB2	2.47	0.44
1:B:282:VAL:HG13	1:B:299:THR:CG2	2.47	0.44
1:A:65:LYS:HE2	4:A:473:HOH:O	2.17	0.44
1:B:10:SER:HB3	1:B:339:GLU:OE2	2.16	0.44
1:A:322:PHE:CE2	1:A:324:ILE:HB	2.53	0.44
1:B:67:VAL:HG12	1:B:80:LEU:HD22	1.99	0.44
1:B:287:LEU:HD21	1:B:333:MET:HE1	2.00	0.44
1:A:357:SER:C	1:A:359:CYS:H	2.25	0.43
1:B:235:ARG:HG3	1:B:332:VAL:CG1	2.48	0.43
1:B:302:PRO:O	1:B:306:LEU:HB2	2.18	0.43
1:A:218:LYS:HG3	1:A:381:ASP:O	2.17	0.43
1:B:69:VAL:HG22	1:B:128:ARG:CB	2.47	0.43
1:B:114:ASN:HD22	1:B:167:LEU:CD2	2.31	0.43
1:A:33:THR:HG23	1:A:229:SER:HB2	2.00	0.43
1:A:16:VAL:HG13	1:A:31:VAL:CG2	2.48	0.43
1:A:85:VAL:CG1	1:A:136:PHE:HE1	2.31	0.43
1:B:254:THR:CB	1:B:279:ILE:HD11	2.18	0.43
1:A:211:GLN:HG3	1:B:279:ILE:HA	2.01	0.43
1:A:85:VAL:HG12	1:A:95:VAL:O	2.19	0.42
1:A:275:THR:HA	1:A:276:PRO:HD3	1.85	0.42
1:A:349:ARG:HH11	1:A:351:ARG:NH2	2.17	0.42
1:B:68:TYR:CD2	1:B:70:PRO:HD3	2.54	0.42
1:A:85:VAL:HG11	1:A:136:PHE:HE1	1.83	0.42
1:B:20:VAL:O	1:B:25:GLN:HB2	2.19	0.42
1:B:379:MET:O	1:B:380:GLU:HG3	2.20	0.42
1:B:375:VAL:C	1:B:377:LEU:N	2.77	0.42
1:B:140:LEU:CD1	1:B:146:VAL:HG11	2.50	0.42
1:B:250:ALA:O	1:B:253:SER:OG	2.38	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:308:PRO:HD2	4:B:398:HOH:O	2.20	0.42
1:A:295:SER:HB2	1:A:379:MET:HE1	2.00	0.42
1:A:357:SER:C	1:A:359:CYS:N	2.75	0.42
1:B:256:LYS:HD3	1:B:256:LYS:HA	1.88	0.42
1:A:122:ALA:HA	1:A:199:TYR:CE2	2.55	0.42
1:A:193:ILE:HB	1:A:350:LYS:HG2	2.01	0.42
1:A:17:GLU:HB3	4:A:445:HOH:O	2.20	0.41
1:A:20:VAL:O	1:A:25:GLN:HB2	2.19	0.41
1:A:238:LYS:CG	1:A:326:GLN:NE2	2.84	0.41
1:A:16:VAL:HG13	1:A:31:VAL:HG23	2.03	0.41
1:A:71:TYR:HE2	1:A:76:TRP:HE1	1.67	0.41
1:A:238:LYS:HG3	1:A:326:GLN:NE2	2.35	0.41
1:B:146:VAL:HA	1:B:147:PRO:HD3	1.90	0.41
1:B:357:SER:C	1:B:359:CYS:H	2.28	0.41
1:A:371:GLU:HG2	4:A:458:HOH:O	2.21	0.41
1:B:152:LEU:CD1	1:B:174:MET:HG3	2.50	0.41
1:B:197:TRP:CD1	1:B:197:TRP:N	2.83	0.41
1:A:71:TYR:HE2	1:A:76:TRP:NE1	2.18	0.41
1:A:375:VAL:C	1:A:377:LEU:N	2.78	0.41
1:B:248:ILE:HG13	1:B:281:PRO:HG2	2.03	0.41
1:B:357:SER:C	1:B:359:CYS:N	2.77	0.41
1:B:218:LYS:HG3	1:B:381:ASP:O	2.22	0.40
1:B:275:THR:HA	1:B:276:PRO:HD3	1.88	0.40
1:A:376:THR:HG23	1:B:364:GLU:OE1	2.21	0.40
1:B:68:TYR:CE2	1:B:70:PRO:HD3	2.56	0.40
1:B:140:LEU:HD13	1:B:146:VAL:HG11	2.04	0.40
1:B:72:THR:OG1	2:B:394:PB7:H12	2.21	0.40
1:B:154:LEU:O	1:B:339:GLU:HA	2.21	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	385/455 (85%)	366 (95%)	18 (5%)	1 (0%)	36	70
1	B	385/455 (85%)	363 (94%)	20 (5%)	2 (0%)	24	60
All	All	770/910 (85%)	729 (95%)	38 (5%)	3 (0%)	30	65

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	376	THR
1	B	376	THR
1	B	380	GLU

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	329/381 (86%)	274 (83%)	55 (17%)	2	12
1	B	329/381 (86%)	283 (86%)	46 (14%)	3	16
All	All	658/762 (86%)	557 (85%)	101 (15%)	3	14

All (101) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	0	VAL
1	A	1	GLU
1	A	3	VAL
1	A	16	VAL
1	A	20	VAL
1	A	25	GLN
1	A	59	THR
1	A	63	LEU
1	A	72	THR
1	A	73	GLN
1	A	95	VAL
1	A	105	SER
1	A	116	GLU

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Mol	Chain	Res	Type
1	A	121	LEU
1	A	126	ILE
1	A	133	LEU
1	A	140	LEU
1	A	141	VAL
1	A	142	LYS
1	A	146	VAL
1	A	152	LEU
1	A	154	LEU
1	A	161	LEU
1	A	166	VAL
1	A	167	LEU
1	A	183	LEU
1	A	185	THR
1	A	187	SER
1	A	188	LEU
1	A	197	TRP
1	A	202	ILE
1	A	213	LEU
1	A	229	SER
1	A	248	ILE
1	A	255	GLU
1	A	263	LEU
1	A	265	GLU
1	A	279	ILE
1	A	282	VAL
1	A	285	LEU
1	A	287	LEU
1	A	291	VAL
1	A	292	THR
1	A	300	ILE
1	A	301	LEU
1	A	306	LEU
1	A	316	GLN
1	A	332	VAL
1	A	336	VAL
1	A	359	CYS
1	A	364	GLU
1	A	376	THR
1	A	377	LEU
1	A	378	ASP
1	A	379	MET

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Mol	Chain	Res	Type
1	B	0	VAL
1	B	3	VAL
1	B	16	VAL
1	B	20	VAL
1	B	25	GLN
1	B	59	THR
1	B	73	GLN
1	B	95	VAL
1	B	105	SER
1	B	116	GLU
1	B	121	LEU
1	B	126	ILE
1	B	133	LEU
1	B	140	LEU
1	B	141	VAL
1	B	146	VAL
1	B	149	LEU
1	B	152	LEU
1	B	165	GLU
1	B	166	VAL
1	B	167	LEU
1	B	183	LEU
1	B	185	THR
1	B	187	SER
1	B	188	LEU
1	B	197	TRP
1	B	213	LEU
1	B	234	LEU
1	B	248	ILE
1	B	252	SER
1	B	253	SER
1	B	263	LEU
1	B	265	GLU
1	B	268	VAL
1	B	279	ILE
1	B	282	VAL
1	B	285	LEU
1	B	292	THR
1	B	306	LEU
1	B	316	GLN
1	B	336	VAL
1	B	357	SER

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Mol	Chain	Res	Type
1	B	359	CYS
1	B	376	THR
1	B	377	LEU
1	B	380	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	GLN
1	A	98	ASN
1	A	153	GLN
1	A	233	ASN
1	A	326	GLN
1	B	89	HIS
1	B	211	GLN

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](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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
2	PB7	B	394	-	50,50,50	5.86	28 (56%)	57,68,68	1.27	6 (10%)
2	PB7	A	394	-	50,50,50	5.79	28 (56%)	57,68,68	1.11	4 (7%)

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	PB7	B	394	-	-	5/46/62/62	0/4/4/4
2	PB7	A	394	-	-	7/46/62/62	0/4/4/4

All (56) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	394	PB7	O43-C19	13.46	1.42	1.22
2	B	394	PB7	O43-C19	13.43	1.42	1.22
2	B	394	PB7	O45-C21	10.58	1.43	1.23
2	A	394	PB7	O45-C21	10.01	1.42	1.23
2	A	394	PB7	C19-N39	9.12	1.49	1.34
2	B	394	PB7	O44-C20	9.11	1.43	1.23
2	B	394	PB7	C7-C13	8.86	1.54	1.38
2	A	394	PB7	C7-C13	8.71	1.53	1.38
2	A	394	PB7	C5-C10	8.70	1.53	1.38
2	B	394	PB7	C20-N40	8.67	1.48	1.34
2	A	394	PB7	O44-C20	8.66	1.42	1.23
2	B	394	PB7	C7-C12	8.65	1.53	1.38
2	A	394	PB7	C7-C12	8.58	1.53	1.38
2	A	394	PB7	C20-N40	8.57	1.48	1.34
2	B	394	PB7	C14-C18	8.56	1.53	1.39
2	B	394	PB7	C4-C9	8.48	1.53	1.38
2	A	394	PB7	C14-C17	8.46	1.53	1.39
2	B	394	PB7	C14-C17	8.45	1.53	1.39
2	A	394	PB7	C14-C18	8.42	1.52	1.39
2	B	394	PB7	C19-N39	8.34	1.48	1.34
2	B	394	PB7	C13-C18	8.34	1.54	1.38
2	B	394	PB7	C3-C8	8.34	1.53	1.38
2	B	394	PB7	C5-C10	8.33	1.53	1.38
2	A	394	PB7	C6-C11	8.32	1.53	1.38
2	B	394	PB7	C6-C11	8.24	1.53	1.38
2	A	394	PB7	C3-C8	8.21	1.52	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	394	PB7	C4-C9	8.20	1.52	1.38
2	A	394	PB7	C13-C18	7.87	1.53	1.38
2	B	394	PB7	C21-N41	7.65	1.50	1.34
2	B	394	PB7	C8-C15	7.54	1.53	1.38
2	A	394	PB7	C10-C16	7.38	1.53	1.38
2	B	394	PB7	C11-C16	7.34	1.53	1.38
2	B	394	PB7	C12-C17	7.26	1.53	1.38
2	B	394	PB7	C10-C16	7.25	1.53	1.38
2	A	394	PB7	C12-C17	7.20	1.53	1.38
2	B	394	PB7	C9-C15	7.08	1.52	1.38
2	B	394	PB7	C4-C1	7.07	1.53	1.38
2	A	394	PB7	C11-C16	7.04	1.52	1.38
2	A	394	PB7	C8-C15	7.03	1.52	1.38
2	A	394	PB7	C21-N41	7.01	1.49	1.34
2	A	394	PB7	C6-C2	6.96	1.53	1.38
2	A	394	PB7	C9-C15	6.94	1.52	1.38
2	B	394	PB7	C3-C1	6.87	1.53	1.38
2	B	394	PB7	C6-C2	6.86	1.53	1.38
2	A	394	PB7	C5-C2	6.82	1.53	1.38
2	B	394	PB7	C5-C2	6.79	1.53	1.38
2	A	394	PB7	C4-C1	6.68	1.52	1.38
2	A	394	PB7	C3-C1	6.59	1.52	1.38
2	A	394	PB7	C35-N39	3.67	1.50	1.46
2	B	394	PB7	O47-C18	3.32	1.44	1.37
2	A	394	PB7	O47-C18	3.18	1.43	1.37
2	B	394	PB7	C24-N40	2.76	1.50	1.45
2	B	394	PB7	C35-N39	2.59	1.49	1.46
2	A	394	PB7	C24-N40	2.57	1.49	1.45
2	A	394	PB7	C37-N41	2.48	1.50	1.46
2	B	394	PB7	C37-N41	2.13	1.49	1.46

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	394	PB7	C23-N39-C19	-3.69	111.12	113.47
2	A	394	PB7	C27-C36-C32	-3.56	107.25	110.83
2	B	394	PB7	O44-C20-N40	3.45	127.01	122.23
2	B	394	PB7	C27-C36-C32	-3.41	107.40	110.83
2	A	394	PB7	C23-N39-C19	-3.03	111.54	113.47
2	B	394	PB7	C25-C20-N40	-2.82	112.43	115.94
2	A	394	PB7	C30-C37-C38	2.31	115.41	111.66
2	B	394	PB7	C24-N40-C20	2.16	125.44	121.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	394	PB7	C35-N39-C19	2.08	124.60	122.12
2	B	394	PB7	C22-C23-N39	-2.00	101.06	103.19

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	394	PB7	C22-C24-C36-C32
2	B	394	PB7	N42-C34-C38-O46
2	A	394	PB7	C13-C18-O47-C28
2	A	394	PB7	C14-C18-O47-C28
2	B	394	PB7	C17-C31-N42-C34
2	B	394	PB7	N42-C34-C38-C37
2	B	394	PB7	C21-C35-N39-C23
2	A	394	PB7	C19-C24-C36-C32
2	A	394	PB7	C33-C35-N39-C23
2	B	394	PB7	C33-C35-N39-C23
2	A	394	PB7	C10-C16-C30-C37
2	A	394	PB7	C11-C16-C30-C37

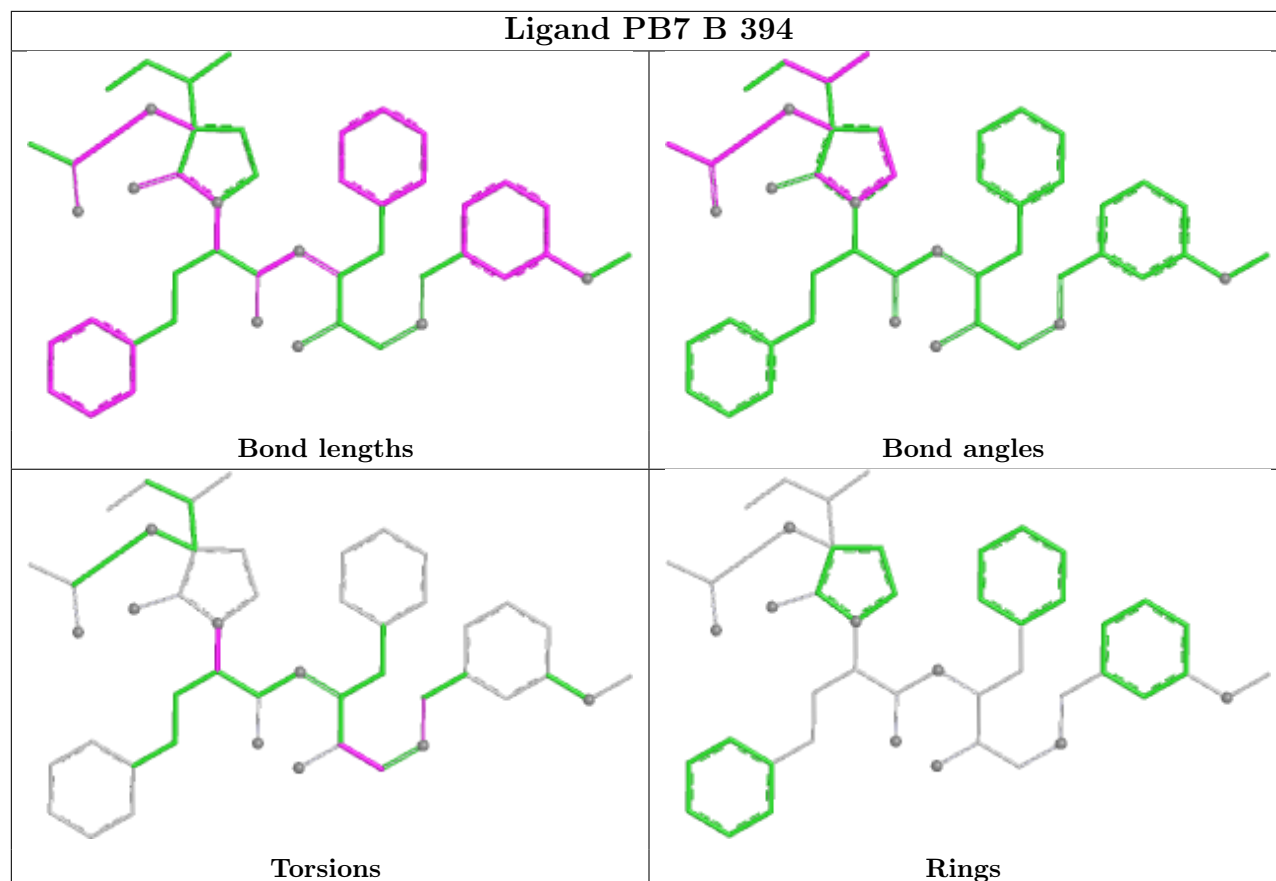
There are no ring outliers.

2 monomers are involved in 4 short contacts:

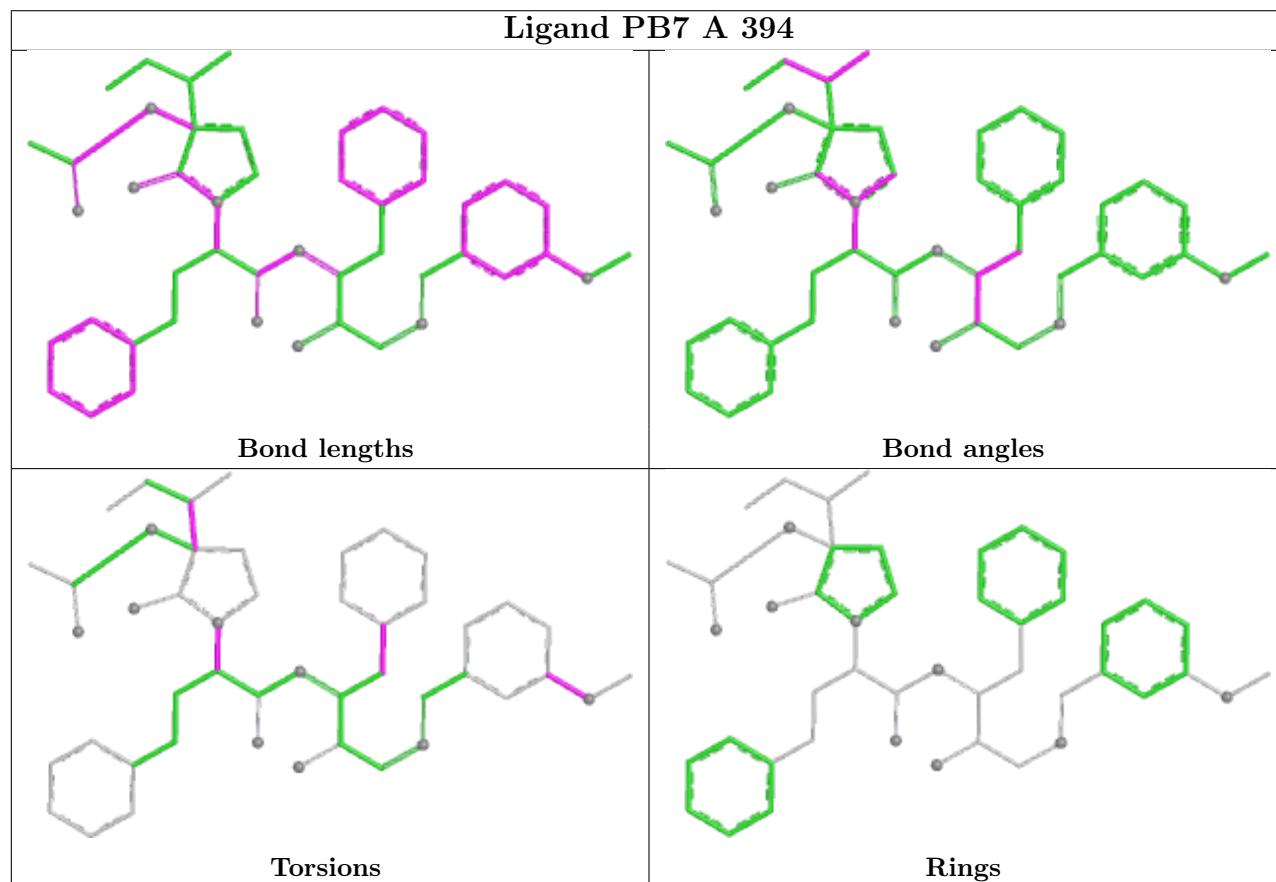
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	394	PB7	2	0
2	A	394	PB7	2	0

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.

Ligand PB7 B 394



Ligand PB7 A 394



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, 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	387/455 (85%)	0.18	13 (3%) 48 27	22, 32, 35, 36	0
1	B	387/455 (85%)	0.19	8 (2%) 63 40	22, 32, 35, 36	0
All	All	774/910 (85%)	0.19	21 (2%) 56 33	22, 32, 35, 36	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	314	THR	5.2
1	A	314	THR	3.8
1	A	312	VAL	3.6
1	B	315	SER	3.4
1	A	313	ALA	3.3
1	B	376	THR	3.0
1	A	379	MET	3.0
1	B	312	VAL	3.0
1	A	266	GLN	2.8
1	A	145	HIS	2.8
1	A	364	GLU	2.7
1	A	377	LEU	2.7
1	B	311	ASP	2.6
1	A	315	SER	2.6
1	A	378	ASP	2.6
1	B	364	GLU	2.4
1	A	375	VAL	2.4
1	B	377	LEU	2.3
1	A	376	THR	2.2
1	A	162	ASN	2.2
1	B	380	GLU	2.1

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](#)

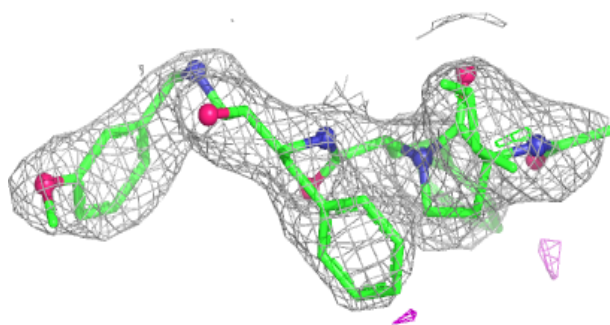
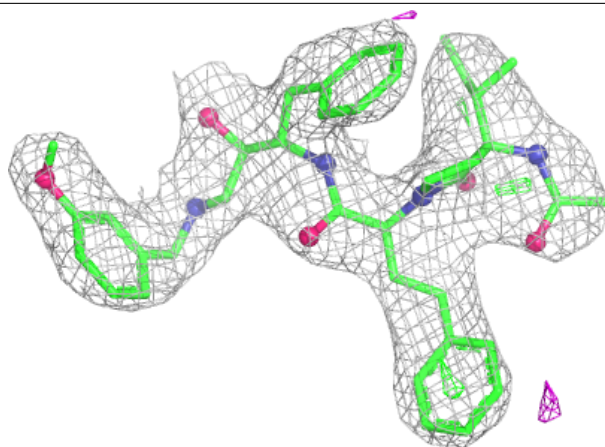
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
3	IOD	B	395	1/1	0.59	0.30	294,294,294,294	0
3	IOD	A	395	1/1	0.66	0.30	233,233,233,233	0
2	PB7	A	394	47/47	0.91	0.16	30,31,32,32	0
2	PB7	B	394	47/47	0.93	0.13	30,31,32,32	0

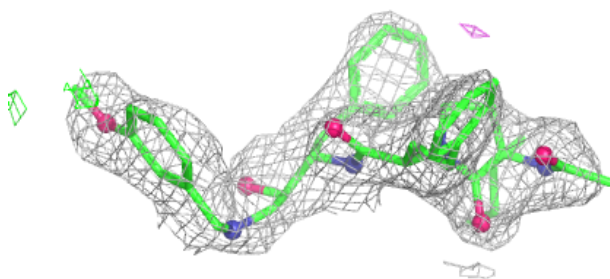
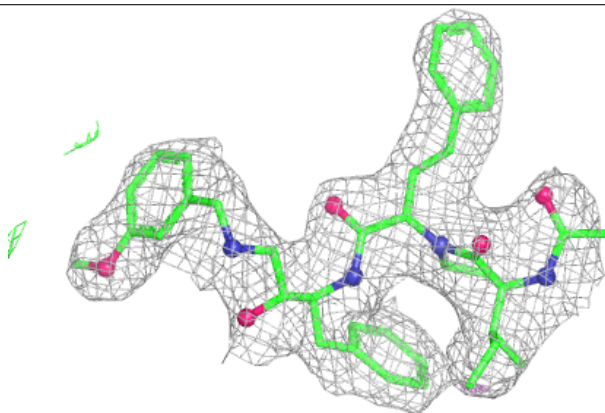
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around PB7 A 394:

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

**Electron density around PB7 B 394:**

$2mF_o-DF_c$ (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.