



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2014 – 04:37 AM GMT

PDB ID : 3K5C
Title : Human BACE-1 complex with NB-216
Authors : Rondeau, J.-M.
Deposited on : 2009-10-07
Resolution : 2.12 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

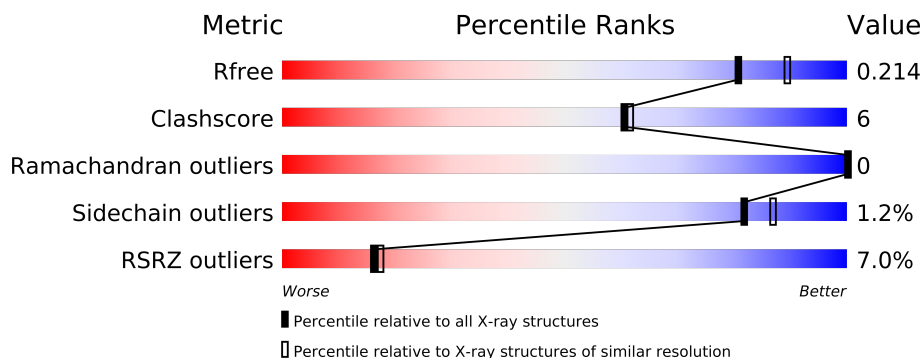
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.12 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}	66092	3409 (2.14-2.10)
Clashscore	79885	4090 (2.14-2.10)
Ramachandran outliers	78287	4048 (2.14-2.10)
Sidechain outliers	78261	4049 (2.14-2.10)
RSRZ outliers	66119	3410 (2.14-2.10)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	402	
1	B	402	
1	C	402	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9519 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

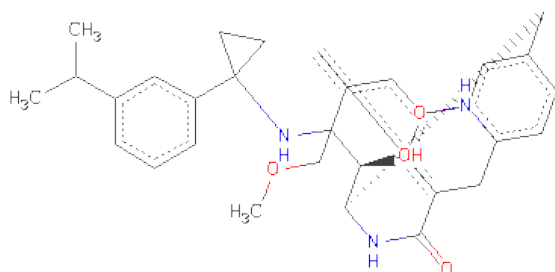
- Molecule 1 is a protein called Beta-secretase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	377	Total	C	N	O	S	0	0	0
			2966	1898	493	561	14			
1	B	377	Total	C	N	O	S	0	0	0
			2966	1898	493	561	14			
1	C	381	Total	C	N	O	S	0	0	0
			2993	1917	497	565	14			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	33P	GLY	-	expression tag	UNP P56817
A	34P	PRO	-	expression tag	UNP P56817
B	33P	GLY	-	expression tag	UNP P56817
B	34P	PRO	-	expression tag	UNP P56817
C	33P	GLY	-	expression tag	UNP P56817
C	34P	PRO	-	expression tag	UNP P56817

- Molecule 2 is (4S)-4-[(1R)-1-HYDROXY-2-({1-[3-(1-METHYLETHYL)PHENYL]CYCLOPROPYL}AMINO)ETHYL]-19-(METHOXYMETHYL)-11-OXA-3,16-DIAZATRICYCLO[15.3.1.1^{6,10}]DOCOSA-1(21),6(22),7,9,17,19-HEXAEN-2-ONE (three-letter code: 0BI) (formula: C₃₅H₄₅N₃O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			42	35	3	4		
2	B	1	Total	C	N	O	0	0
			42	35	3	4		
2	C	1	Total	C	N	O	0	0
			42	35	3	4		

- Molecule 3 is water.

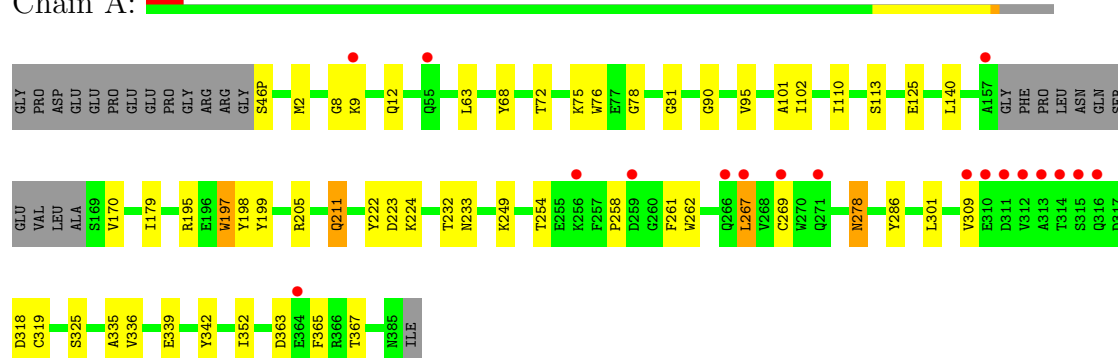
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	162	Total	O	0	0
			162	162		
3	B	135	Total	O	0	0
			135	135		
3	C	171	Total	O	0	0
			171	171		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

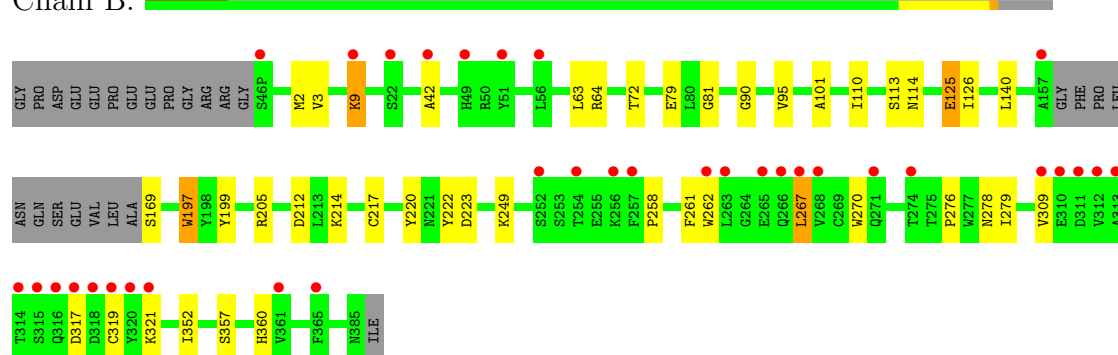
• Molecule 1: Beta-secretase 1

Chain A:



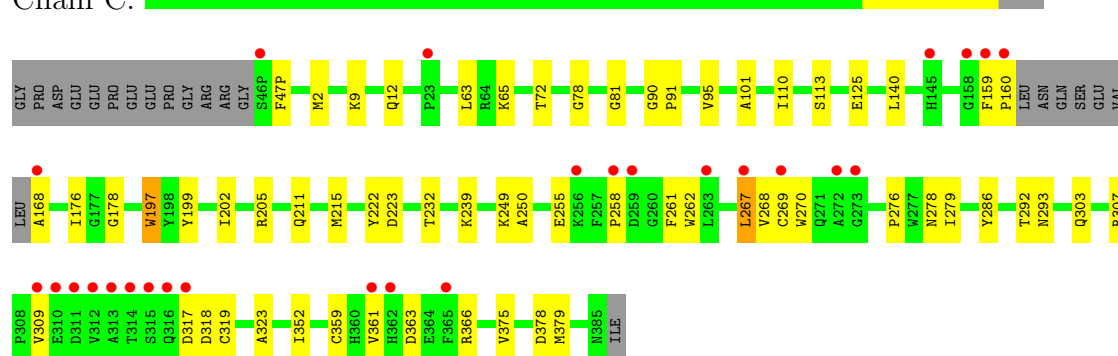
• Molecule 1: Beta-secretase 1

Chain B:



• Molecule 1: Beta-secretase 1

Chain C:



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	81.75Å 103.19Å 100.14Å 90.00° 104.44° 90.00°	Depositor
Resolution (Å)	70.55 – 2.12 70.55 – 2.12	Depositor EDS
% Data completeness (in resolution range)	99.6 (70.55-2.12) 99.8 (70.55-2.12)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.46 (at 2.12Å)	Xtriage
Refinement program	CNX 2005	Depositor
R, R_{free}	0.200 , 0.220 0.196 , 0.214	Depositor DCC
R_{free} test set	9103 reflections (11.10%)	DCC
Wilson B-factor (Å ²)	41.1	Xtriage
Anisotropy	0.221	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 44.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 91081 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9519	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.77% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 0BI

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	0.30	0/3041	0.47	0/4133
1	B	0.30	0/3041	0.47	0/4133
1	C	0.30	0/3070	0.48	0/4173
All	All	0.30	0/9152	0.48	0/12439

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2966	0	2875	36	0
1	B	2966	0	2875	32	0
1	C	2993	0	2899	43	0
2	A	42	0	45	1	0
2	B	42	0	45	1	0
2	C	42	0	45	2	0
3	A	162	0	0	2	0
3	B	135	0	0	0	0
3	C	171	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	9519	0	8784	106	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 6.

All (106) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:367:THR:H	1:C:211:GLN:HE22	1.19	0.85
1:A:267:LEU:HD23	1:A:267:LEU:H	1.48	0.74
1:B:267:LEU:HD23	1:B:267:LEU:H	1.55	0.72
1:C:255:GLU:OE1	1:C:270:TRP:NE1	2.26	0.69
1:C:267:LEU:H	1:C:267:LEU:HD23	1.58	0.69
1:B:9:LYS:HD2	1:B:9:LYS:N	2.10	0.66
1:A:46(P):SER:HB3	3:A:539:HOH:O	1.96	0.65
1:C:268:VAL:O	1:C:319:CYS:HA	1.97	0.63
1:C:9:LYS:HD3	1:C:168:ALA:HB3	1.79	0.63
1:A:367:THR:H	1:C:211:GLN:NE2	1.95	0.63
1:B:205:ARG:NH2	1:B:212:ASP:HB2	2.14	0.62
1:A:72:THR:HB	2:A:387:0BI:H37	1.81	0.62
1:A:267:LEU:HD12	1:A:319:CYS:HB3	1.85	0.59
1:A:233:ASN:HD22	1:A:325:SER:HG	1.51	0.57
1:C:267:LEU:HD13	1:C:309:VAL:CG2	2.36	0.56
1:B:2:MET:HG2	1:B:90:GLY:HA2	1.88	0.56
1:A:68:TYR:CD1	1:B:3:VAL:HG11	2.41	0.55
1:C:65:LYS:HD3	3:C:540:HOH:O	2.07	0.55
1:A:254:THR:HG23	1:C:278:ASN:OD1	2.07	0.55
1:C:269:CYS:HA	1:C:318:ASP:O	2.07	0.55
1:A:258:PRO:O	1:A:261:PHE:HB3	2.07	0.55
1:C:258:PRO:O	1:C:261:PHE:HB3	2.07	0.55
1:C:72:THR:HB	2:C:387:0BI:H37	1.88	0.54
1:A:335:ALA:O	1:A:339:GLU:HG3	2.07	0.54
1:B:9:LYS:NZ	1:B:114:ASN:HB2	2.23	0.53
1:A:2:MET:HG2	1:A:90:GLY:HA2	1.90	0.53
1:C:303:GLN:HB2	1:C:361:VAL:HG11	1.91	0.53
1:B:276:PRO:O	1:B:279:ILE:HG12	2.09	0.53
1:A:278:ASN:H	1:A:278:ASN:HD22	1.57	0.52
1:C:110:ILE:HB	1:C:113:SER:HB3	1.91	0.52
1:B:110:ILE:HB	1:B:113:SER:HB3	1.91	0.52
1:B:357:SER:O	1:B:360:HIS:HB3	2.10	0.52
1:C:159:PHE:HB3	1:C:160:PRO:HD2	1.92	0.52
1:A:110:ILE:HB	1:A:113:SER:HB3	1.92	0.52
1:B:258:PRO:O	1:B:261:PHE:HB3	2.09	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:267:LEU:HD13	1:B:309:VAL:CG2	2.41	0.51
1:B:270:TRP:O	1:B:317:ASP:HB3	2.10	0.51
1:C:2:MET:HG2	1:C:90:GLY:HA2	1.93	0.51
1:C:12:GLN:OE1	1:C:113:SER:HA	2.11	0.50
1:C:276:PRO:O	1:C:279:ILE:HG12	2.12	0.49
1:C:363:ASP:HB3	1:C:366:ARG:O	2.13	0.49
1:C:293:ASN:HA	1:C:375:VAL:HA	1.93	0.49
1:B:63:LEU:HG	1:B:81:GLY:HA2	1.95	0.48
1:B:267:LEU:HD13	1:B:309:VAL:HG21	1.95	0.48
1:C:125:GLU:HG2	1:C:197:TRP:HB3	1.95	0.48
1:A:125:GLU:OE2	1:A:195:ARG:NH2	2.44	0.48
1:B:199:TYR:HB3	1:B:352:ILE:HD11	1.94	0.48
1:C:267:LEU:HD13	1:C:309:VAL:HG21	1.96	0.48
1:A:199:TYR:HB3	1:A:352:ILE:HD11	1.95	0.48
1:C:249:LYS:HE2	1:C:262:TRP:CD1	2.49	0.47
1:B:72:THR:HB	2:B:387:OBI:H37	1.97	0.47
1:A:249:LYS:HE2	1:A:262:TRP:CD1	2.50	0.46
1:C:267:LEU:N	1:C:267:LEU:HD23	2.26	0.46
1:C:199:TYR:HB3	1:C:352:ILE:HD11	1.96	0.46
1:A:222:TYR:HA	1:A:223:ASP:HA	1.65	0.45
1:C:270:TRP:O	1:C:317:ASP:HB3	2.16	0.45
1:B:212:ASP:O	1:B:214:LYS:HG3	2.16	0.45
1:A:211:GLN:HB2	3:A:455:HOH:O	2.16	0.45
1:C:202:ILE:CD1	1:C:379:MET:HG3	2.46	0.45
1:C:63:LEU:HG	1:C:81:GLY:HA2	1.99	0.45
1:A:267:LEU:HD13	1:A:309:VAL:CG2	2.47	0.44
1:A:232:THR:O	1:A:336:VAL:HG13	2.17	0.44
1:A:211:GLN:HE21	1:A:211:GLN:HB2	1.65	0.44
1:C:359:CYS:SG	1:C:359:CYS:O	2.76	0.44
1:C:267:LEU:HD13	1:C:309:VAL:HG23	1.98	0.44
1:B:9:LYS:N	1:B:9:LYS:CD	2.81	0.44
1:C:202:ILE:HD11	1:C:379:MET:HG3	2.00	0.44
1:A:95:VAL:HG11	1:A:140:LEU:HA	2.00	0.44
1:C:205:ARG:HB3	1:C:286:TYR:HB2	2.00	0.43
1:C:307:ARG:HG3	1:C:323:ALA:HB2	1.99	0.43
1:B:249:LYS:HE2	1:B:262:TRP:CD1	2.52	0.43
1:C:215:MET:HE1	1:C:239:LYS:HG2	2.00	0.43
1:B:222:TYR:HA	1:B:223:ASP:HA	1.63	0.43
1:C:95:VAL:HG11	1:C:140:LEU:HA	2.00	0.43
1:B:125:GLU:HG3	1:B:125:GLU:O	2.17	0.43
1:C:78:GLY:HA3	1:C:101:ALA:O	2.18	0.43
1:A:9:LYS:HD2	1:A:12:GLN:OE1	2.19	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:301:LEU:HD13	1:A:363:ASP:HB2	2.02	0.42
1:B:267:LEU:HD23	1:B:267:LEU:N	2.28	0.42
1:C:65:LYS:HE2	1:C:65:LYS:HB2	1.95	0.42
1:C:292:THR:HG21	1:C:378:ASP:HB3	2.00	0.42
1:A:8:GLY:O	1:A:170:VAL:HG22	2.19	0.42
1:C:232:THR:OG1	2:C:387:OBI:H28	2.20	0.42
1:B:217:CYS:HA	1:B:220:TYR:CD1	2.55	0.42
1:A:269:CYS:HA	1:A:318:ASP:O	2.20	0.41
1:B:309:VAL:HG11	1:B:321:LYS:HG3	2.02	0.41
1:B:95:VAL:HG11	1:B:140:LEU:HA	2.02	0.41
1:A:205:ARG:HB3	1:A:286:TYR:HB2	2.02	0.41
1:C:91:PRO:HD3	1:C:176:ILE:HB	2.03	0.41
1:A:365:PHE:CG	1:C:250:ALA:HB1	2.56	0.41
1:B:126:ILE:HG23	1:B:197:TRP:HB2	2.03	0.41
1:B:42:ALA:CB	1:B:101:ALA:HB1	2.51	0.41
1:B:278:ASN:H	1:B:278:ASN:HD22	1.69	0.41
1:C:47(P):PHE:CZ	1:C:178:GLY:HA3	2.56	0.41
1:B:267:LEU:HD12	1:B:319:CYS:HB3	2.02	0.41
1:B:9:LYS:HZ1	1:B:114:ASN:HB2	1.84	0.41
1:A:179:ILE:HG23	1:A:342:TYR:HE2	1.86	0.41
1:C:222:TYR:HA	1:C:223:ASP:HA	1.63	0.41
1:A:63:LEU:HG	1:A:81:GLY:HA2	2.02	0.41
1:B:64:ARG:HA	1:B:79:GLU:OE2	2.20	0.41
1:A:78:GLY:HA3	1:A:101:ALA:O	2.21	0.41
1:A:197:TRP:N	1:A:197:TRP:CD1	2.89	0.41
1:A:8:GLY:C	1:A:170:VAL:HG22	2.42	0.40
1:A:76:TRP:HB2	1:A:102:ILE:HG23	2.03	0.40
1:B:197:TRP:CD1	1:B:197:TRP:N	2.89	0.40
1:A:198:TYR:CE2	1:A:224:LYS:HE3	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	373/402 (93%)	365 (98%)	8 (2%)	0	100	100
1	B	373/402 (93%)	361 (97%)	12 (3%)	0	100	100
1	C	377/402 (94%)	363 (96%)	14 (4%)	0	100	100
All	All	1123/1206 (93%)	1089 (97%)	34 (3%)	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	322/342 (94%)	317 (98%)	5 (2%)	75	79
1	B	322/342 (94%)	317 (98%)	5 (2%)	75	79
1	C	324/342 (95%)	322 (99%)	2 (1%)	92	95
All	All	968/1026 (94%)	956 (99%)	12 (1%)	82	87

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	75	LYS
1	A	197	TRP
1	A	211	GLN
1	A	267	LEU
1	A	278	ASN
1	B	9	LYS
1	B	125	GLU
1	B	169	SER
1	B	197	TRP
1	B	267	LEU
1	C	197	TRP
1	C	267	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	114	ASN
1	A	326	GLN
1	B	114	ASN
1	B	278	ASN
1	B	293	ASN
1	B	326	GLN
1	C	114	ASN
1	C	211	GLN
1	C	326	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

3 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	0BI	A	387	-	46,46,46	1.36	2 (4%)	64,64,64	1.10	7 (10%)
2	0BI	B	387	-	46,46,46	1.35	4 (8%)	64,64,64	1.08	7 (10%)
2	0BI	C	387	-	46,46,46	1.39	5 (10%)	64,64,64	1.10	6 (9%)

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	0BI	A	387	-	-	0/41/48/48	0/1/5/5
2	0BI	B	387	-	-	0/41/48/48	0/1/5/5
2	0BI	C	387	-	-	0/41/48/48	0/1/5/5

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	387	0BI	C65-C61	2.89	1.55	1.51
2	A	387	0BI	C65-C61	2.84	1.55	1.51
2	B	387	0BI	C65-C61	2.70	1.55	1.51
2	A	387	0BI	C62-C61	2.40	1.54	1.51
2	C	387	0BI	C62-C61	2.30	1.54	1.51
2	C	387	0BI	C72-C71	2.28	1.42	1.39
2	B	387	0BI	C62-C61	2.12	1.54	1.51
2	B	387	0BI	C69-C71	2.09	1.42	1.38
2	C	387	0BI	C56-C52	2.06	1.55	1.52
2	B	387	0BI	C72-C71	2.02	1.42	1.39
2	C	387	0BI	C37-C36	2.00	1.42	1.39

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	387	0BI	C62-C61-N59	-3.66	113.42	117.36
2	A	387	0BI	C62-C61-N59	-3.37	113.72	117.36
2	B	387	0BI	C62-C61-N59	-3.32	113.78	117.36
2	A	387	0BI	C19-O18-C11	2.85	125.20	117.94
2	C	387	0BI	C19-O18-C11	2.82	125.13	117.94
2	B	387	0BI	C62-C61-C68	-2.83	114.06	119.18
2	A	387	0BI	C62-C61-C68	-2.79	114.13	119.18
2	A	387	0BI	C56-N59-C61	-2.72	111.81	116.86
2	A	387	0BI	C36-C50-N1	2.67	121.34	116.89
2	C	387	0BI	C62-C61-C68	-2.62	114.44	119.18
2	B	387	0BI	C36-C50-N1	2.61	121.24	116.89
2	C	387	0BI	C36-C50-N1	2.55	121.15	116.89
2	B	387	0BI	C19-O18-C11	2.46	124.21	117.94
2	C	387	0BI	C56-N59-C61	-2.44	112.33	116.86
2	B	387	0BI	C56-N59-C61	-2.40	112.39	116.86
2	B	387	0BI	O51-C50-C36	-2.13	117.27	121.01
2	A	387	0BI	O51-C50-C36	-2.09	117.34	121.01

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	387	0BI	O54-C52-C56	-2.04	105.24	109.72
2	C	387	0BI	O51-C50-C36	-2.03	117.45	121.01
2	A	387	0BI	C3-N1-C50	2.02	127.29	122.92

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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 ⓘ

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	377/402 (93%)	0.58	18 (4%) 29 32	28, 44, 81, 116	0
1	B	377/402 (93%)	0.77	35 (9%) 9 9	27, 45, 85, 118	0
1	C	381/402 (94%)	0.66	27 (7%) 16 17	29, 44, 83, 116	0
All	All	1135/1206 (94%)	0.67	80 (7%) 16 17	27, 45, 84, 118	0

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	314	THR	13.8
1	B	312	VAL	12.8
1	C	312	VAL	12.0
1	A	312	VAL	9.3
1	C	158	GLY	9.0
1	B	313	ALA	7.8
1	A	314	THR	7.0
1	B	361	VAL	7.0
1	C	314	THR	7.0
1	C	315	SER	6.3
1	C	160	PRO	6.2
1	B	315	SER	5.9
1	B	256	LYS	5.8
1	C	168	ALA	5.7
1	C	310	GLU	5.5
1	C	313	ALA	5.5
1	A	311	ASP	5.3
1	C	159	PHE	5.3
1	B	310	GLU	5.1
1	A	157	ALA	4.8
1	A	316	GLN	4.7
1	B	271	GLN	4.6
1	B	311	ASP	4.5

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Mol	Chain	Res	Type	RSRZ
1	C	272	ALA	4.3
1	C	316	GLN	4.3
1	A	313	ALA	4.1
1	B	257	PHE	4.0
1	A	256	LYS	3.9
1	B	157	ALA	3.9
1	B	317	ASP	3.9
1	B	266	GLN	3.9
1	B	309	VAL	3.9
1	B	318	ASP	3.8
1	C	365	PHE	3.7
1	C	46(P)	SER	3.7
1	B	263	LEU	3.6
1	A	266	GLN	3.6
1	C	145	HIS	3.3
1	B	365	PHE	3.2
1	B	316	GLN	3.1
1	B	274	THR	2.9
1	B	265	GLU	2.9
1	C	309	VAL	2.9
1	C	267	LEU	2.9
1	B	268	VAL	2.8
1	B	319	CYS	2.8
1	B	320	TYR	2.8
1	B	46(P)	SER	2.8
1	C	311	ASP	2.8
1	A	9	LYS	2.7
1	B	267	LEU	2.7
1	B	254	THR	2.6
1	A	267	LEU	2.6
1	A	315	SER	2.5
1	B	262	TRP	2.5
1	B	9	LYS	2.5
1	C	269	CYS	2.5
1	C	256	LYS	2.4
1	B	51	TYR	2.4
1	B	252	SER	2.4
1	C	361	VAL	2.4
1	B	321	LYS	2.3
1	A	364	GLU	2.3
1	B	22	SER	2.2
1	A	271	GLN	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	49	HIS	2.2
1	A	310	GLU	2.2
1	A	55	GLN	2.2
1	C	317	ASP	2.1
1	A	259	ASP	2.1
1	C	258	PRO	2.1
1	A	269	CYS	2.1
1	B	56	LEU	2.1
1	C	259	ASP	2.1
1	C	273	GLY	2.1
1	B	42	ALA	2.0
1	C	362	HIS	2.0
1	A	309	VAL	2.0
1	C	23	PRO	2.0
1	C	263	LEU	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	0BI	A	387	42/42	0.16	-	30,36,44,59	0
2	0BI	C	387	42/42	0.14	-	31,36,44,59	0
2	0BI	B	387	42/42	0.16	-	31,37,43,58	0

6.5 Other polymers ⓘ

There are no such residues in this entry.