



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:47 AM UTC

PDB ID : 1XN2 / pdb_00001xn2
Title : New substrate binding pockets for beta-secretase.
Authors : Turner III, R.T.; Hong, L.; Koelsch, G.; Ghosh, A.K.; Tang, J.
Deposited on : 2004-10-04
Resolution : 1.90 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
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

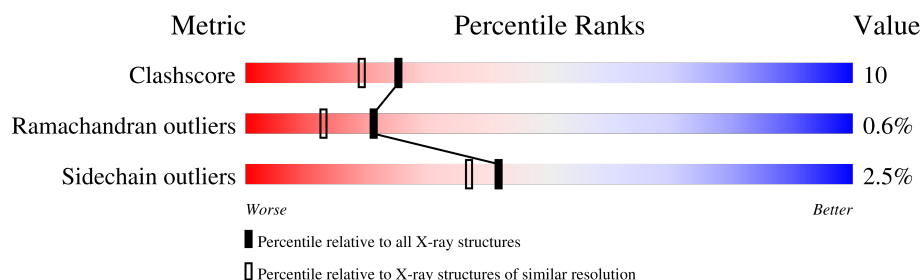
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)

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	389	80% 16% .
1	B	389	84% 14% ..
1	C	389	82% 15% .
1	D	389	82% 16% .
2	E	12	58% 17% 8% 17%
2	F	12	42% 25% 17% 17%
2	G	12	42% 33% 8% 17%
2	H	12	33% 8% 33% 8% 17%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13323 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	389	Total	C	N	O	S	0	0	0
			3053	1953	507	579	14			
1	B	389	Total	C	N	O	S	0	0	0
			3053	1953	507	579	14			
1	C	389	Total	C	N	O	S	0	0	0
			3053	1953	507	579	14			
1	D	389	Total	C	N	O	S	0	0	0
			3053	1953	507	579	14			

- Molecule 2 is a protein called OM03-4.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	10	Total	C	N	O	0	0	0
			97	66	13	18			
2	F	10	Total	C	N	O	0	0	0
			97	66	13	18			
2	G	10	Total	C	N	O	0	0	0
			97	66	13	18			
2	H	10	Total	C	N	O	0	0	0
			97	66	13	18			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	169	Total	O	0	0
			169	169		
3	B	194	Total	O	0	0
			194	194		
3	C	178	Total	O	0	0
			178	178		
3	D	174	Total	O	0	0
			174	174		

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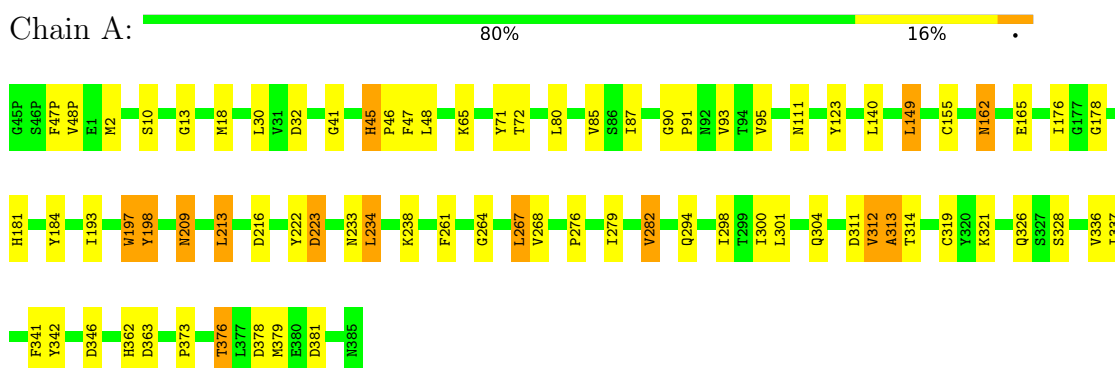
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	2	Total	O	0	0
			2	2		
3	F	2	Total	O	0	0
			2	2		
3	G	1	Total	O	0	0
			1	1		
3	H	3	Total	O	0	0
			3	3		

3 Residue-property plots

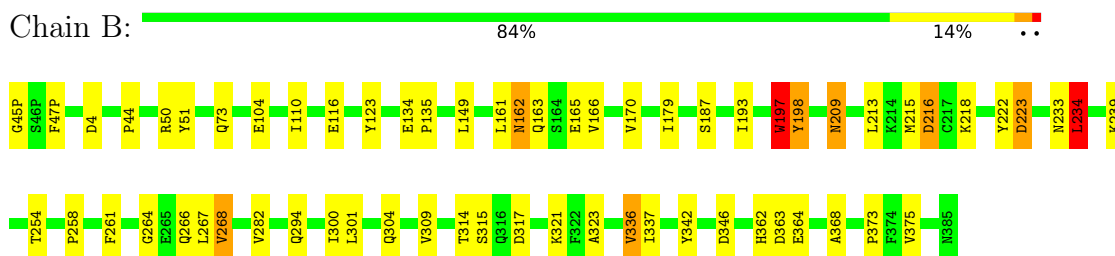
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

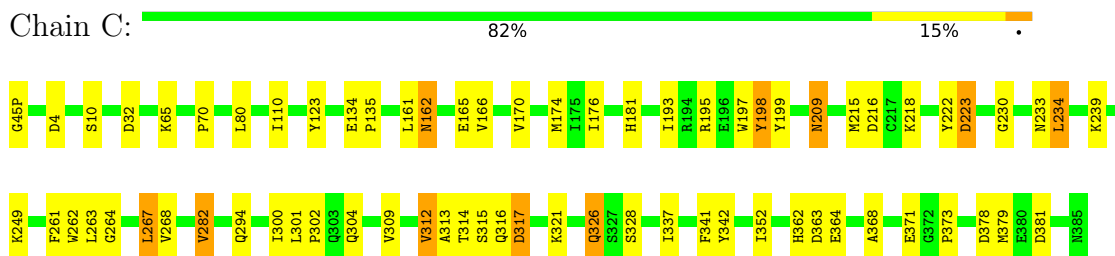
- Molecule 1: Beta-secretase 1



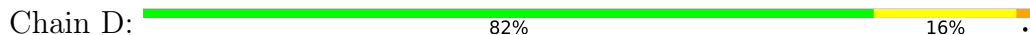
- Molecule 1: Beta-secretase 1

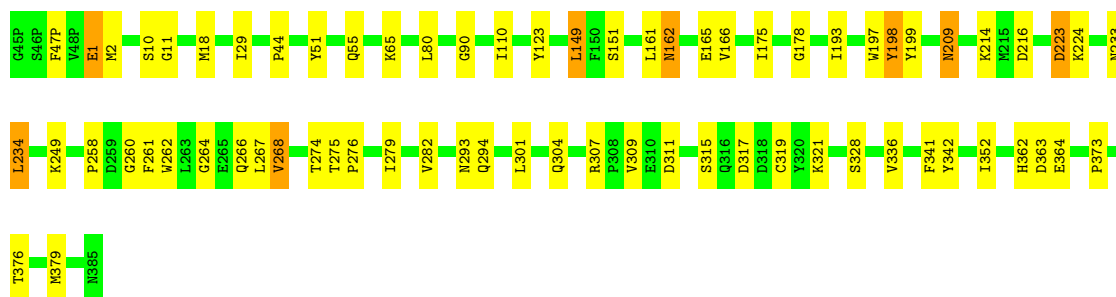


- Molecule 1: Beta-secretase 1



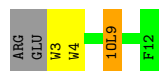
- Molecule 1: Beta-secretase 1





- Molecule 2: OM03-4

Chain E: 58% 17% 8% 17%



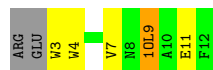
- Molecule 2: OM03-4

Chain F: 42% 25% 17% 17%



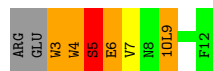
- Molecule 2: OM03-4

Chain G: 42% 33% 8% 17%



- Molecule 2: OM03-4

Chain H: 33% 8% 33% 8% 17%



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	86.22Å 130.78Å 88.74Å 90.00° 97.71° 90.00°	Depositor
Resolution (Å)	30.00 – 1.90	Depositor
% Data completeness (in resolution range)	(Not available) (30.00-1.90)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.193 , 0.220	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	13323	wwPDB-VP
Average B, all atoms (Å ²)	26.0	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: 1OL

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	0.63	0/3131	1.07	21/4256 (0.5%)
1	B	0.65	0/3131	1.05	12/4256 (0.3%)
1	C	0.61	0/3131	1.03	14/4256 (0.3%)
1	D	0.60	0/3131	1.02	13/4256 (0.3%)
2	E	0.50	0/87	0.82	0/116
2	F	0.60	0/87	0.69	0/116
2	G	0.56	0/87	0.77	0/116
2	H	0.55	0/87	1.30	1/116 (0.9%)
All	All	0.62	0/12872	1.04	61/17488 (0.3%)

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
2	E	0	2
2	F	0	2
2	G	0	2
2	H	0	2
All	All	0	8

There are no bond length outliers.

All (61) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	198	TYR	N-CA-C	-9.24	96.03	110.36
1	C	234	LEU	N-CA-C	-9.10	94.49	108.96
1	D	198	TYR	N-CA-C	-9.05	96.33	110.36
1	A	198	TYR	N-CA-C	-8.90	96.56	110.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	234	LEU	N-CA-C	-8.72	95.09	108.96
1	A	234	LEU	N-CA-C	-8.48	95.48	108.96
1	D	234	LEU	N-CA-C	-8.38	95.64	108.96
1	A	233	ASN	N-CA-C	7.97	120.81	110.53
1	B	123	TYR	N-CA-C	7.43	120.58	110.55
1	A	342	TYR	N-CA-C	-7.29	97.36	109.46
1	B	342	TYR	N-CA-C	-7.23	97.47	108.96
1	A	123	TYR	N-CA-C	7.21	120.29	110.55
1	C	342	TYR	N-CA-C	-7.09	97.69	109.46
1	B	198	TYR	N-CA-C	-7.07	99.41	110.36
1	C	282	VAL	N-CA-C	-6.89	102.54	110.05
1	B	233	ASN	N-CA-C	6.44	119.93	110.59
1	C	233	ASN	N-CA-C	6.35	119.80	110.59
1	D	341	PHE	N-CA-C	6.25	119.20	109.52
1	D	342	TYR	N-CA-C	-6.19	99.12	108.96
1	A	312	VAL	N-CA-C	-6.16	103.35	111.05
1	B	193	ILE	N-CA-C	-6.16	99.43	108.54
1	B	268	VAL	N-CA-C	-6.14	99.31	108.46
1	C	328	SER	N-CA-C	-6.13	105.08	113.30
1	A	41	GLY	N-CA-C	-6.10	104.22	111.36
1	C	193	ILE	N-CA-C	-6.09	98.91	108.23
1	D	193	ILE	N-CA-C	-6.08	99.55	108.54
1	D	216	ASP	N-CA-C	-6.07	101.06	110.10
1	A	193	ILE	N-CA-C	-6.05	99.59	108.54
1	D	10	SER	N-CA-C	5.95	117.76	111.28
1	A	328	SER	N-CA-C	-5.77	104.80	112.94
2	H	5	SER	N-CA-C	5.73	123.01	110.80
1	D	268	VAL	N-CA-C	-5.68	100.00	108.46
1	C	341	PHE	N-CA-C	5.68	118.32	109.52
1	A	10	SER	N-CA-C	5.64	117.43	111.28
1	A	341	PHE	N-CA-C	5.61	118.21	109.52
1	D	233	ASN	N-CA-C	5.60	118.71	110.59
1	C	216	ASP	N-CA-C	-5.58	101.78	110.10
1	C	302	PRO	N-CA-C	-5.51	106.68	113.84
1	A	216	ASP	N-CA-C	-5.49	101.92	110.10
1	C	123	TYR	N-CA-C	5.46	118.60	110.52
1	A	45	HIS	CA-C-N	5.39	125.72	119.47
1	A	45	HIS	C-N-CA	5.39	125.72	119.47
1	D	328	SER	N-CA-C	-5.38	106.08	113.30
1	A	213	LEU	N-CA-C	-5.32	105.65	111.82
1	B	213	LEU	N-CA-C	-5.28	105.70	111.82
1	B	216	ASP	N-CA-C	-5.25	102.28	110.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	312	VAL	N-CA-C	5.24	115.51	110.74
1	B	197	TRP	N-CA-C	-5.20	99.73	110.80
1	B	317	ASP	N-CA-C	5.15	117.50	110.24
1	B	368	ALA	N-CA-C	-5.13	103.76	110.53
1	C	10	SER	N-CA-C	5.11	117.51	111.33
1	D	223	ASP	N-CA-C	-5.10	99.94	110.80
1	D	123	TYR	N-CA-C	5.09	118.06	110.52
1	D	214	LYS	N-CA-C	5.06	118.61	111.52
1	A	282	VAL	N-CA-C	-5.06	102.42	109.45
1	A	313	ALA	N-CA-C	-5.04	102.45	109.96
1	A	155	CYS	N-CA-C	5.03	116.89	107.99
1	A	298	ILE	N-CA-C	-5.01	101.26	108.48
1	A	87	ILE	CA-C-N	5.01	125.03	119.32
1	A	87	ILE	C-N-CA	5.01	125.03	119.32
1	C	368	ALA	N-CA-C	-5.00	103.92	110.53

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	E	9	1OL	Peptide,Mainchain
2	F	9	1OL	Peptide,Mainchain
2	G	9	1OL	Peptide,Mainchain
2	H	9	1OL	Peptide,Mainchain

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3053	0	2959	60	0
1	B	3053	0	2959	62	0
1	C	3053	0	2959	65	0
1	D	3053	0	2959	54	0
2	E	97	0	74	5	0
2	F	97	0	74	10	0
2	G	97	0	74	8	0
2	H	97	0	74	18	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	169	0	0	1	0
3	B	194	0	0	0	0
3	C	178	0	0	4	0
3	D	174	0	0	1	0
3	E	2	0	0	0	0
3	F	2	0	0	1	0
3	G	1	0	0	0	0
3	H	3	0	0	0	0
All	All	13323	0	12132	239	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 10.

All (239) 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:A:312:VAL:HG22	1:C:313:ALA:HB3	1.12	1.07
2:H:4:TRP:O	2:H:5:SER:HB3	1.57	1.00
1:C:321:LYS:NZ	2:G:4:TRP:HE1	1.61	0.99
1:B:321:LYS:NZ	2:F:4:TRP:HE1	1.62	0.96
1:B:321:LYS:HZ2	2:F:4:TRP:HE1	1.17	0.93
1:A:46:PRO:HG3	1:C:263:LEU:HD12	1.52	0.92
1:A:321:LYS:HE2	3:A:553:HOH:O	1.73	0.88
1:C:321:LYS:HZ2	2:G:4:TRP:HE1	1.19	0.87
1:A:312:VAL:HG22	1:C:313:ALA:CB	2.05	0.83
1:B:336:VAL:HG12	1:B:337:ILE:HD13	1.62	0.82
1:C:4:ASP:CG	1:C:170:VAL:HG11	2.04	0.82
1:B:4:ASP:CB	1:B:170:VAL:HG11	2.12	0.80
1:C:301:LEU:H	1:C:304:GLN:HE21	1.27	0.80
1:B:162:ASN:C	1:B:162:ASN:HD22	1.91	0.79
1:C:312:VAL:HG21	2:E:3:TRP:N	1.97	0.78
1:A:162:ASN:C	1:A:162:ASN:HD22	1.90	0.78
1:D:267:LEU:HD22	1:D:309:VAL:HG21	1.65	0.77
1:C:321:LYS:NZ	2:G:4:TRP:NE1	2.30	0.77
1:D:1:GLU:H	1:D:1:GLU:CD	1.91	0.77
1:B:264:GLY:O	1:B:321:LYS:HE3	1.83	0.77
1:D:264:GLY:O	1:D:321:LYS:HE3	1.85	0.76
1:B:4:ASP:CG	1:B:170:VAL:HG11	2.11	0.75
2:H:3:TRP:HA	2:H:3:TRP:CE3	2.22	0.75
1:D:321:LYS:NZ	2:H:4:TRP:HE1	1.85	0.74
1:C:301:LEU:H	1:C:304:GLN:NE2	1.86	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:162:ASN:O	1:B:166:VAL:HG23	1.86	0.74
1:B:267:LEU:CD2	1:B:309:VAL:HG21	2.18	0.73
1:B:301:LEU:H	1:B:304:GLN:HE21	1.35	0.73
1:C:249:LYS:HE2	1:C:262:TRP:CD1	2.23	0.73
1:A:162:ASN:ND2	1:A:165:GLU:H	1.86	0.72
1:A:301:LEU:H	1:A:304:GLN:HE21	1.38	0.72
1:B:162:ASN:ND2	1:B:165:GLU:H	1.87	0.71
1:B:267:LEU:HD22	1:B:309:VAL:HG21	1.73	0.70
2:H:3:TRP:HA	2:H:3:TRP:HE3	1.54	0.70
1:A:264:GLY:O	1:A:321:LYS:HE3	1.92	0.70
1:C:321:LYS:HZ3	2:G:4:TRP:NE1	1.89	0.69
1:D:311:ASP:HB2	1:D:319:CYS:SG	2.34	0.68
1:B:45(P):GLY:HA2	1:B:179:ILE:O	1.93	0.68
1:B:261:PHE:CD1	1:B:268:VAL:HG23	2.29	0.68
1:A:238:LYS:HG3	1:A:326:GLN:NE2	2.09	0.68
1:A:267:LEU:H	1:A:267:LEU:CD2	2.07	0.67
1:C:162:ASN:ND2	1:C:165:GLU:H	1.92	0.67
1:C:312:VAL:HG11	2:E:3:TRP:N	2.09	0.67
1:C:267:LEU:HD23	1:C:309:VAL:HG21	1.75	0.67
1:D:162:ASN:ND2	1:D:165:GLU:H	1.92	0.66
1:B:301:LEU:H	1:B:304:GLN:NE2	1.93	0.66
1:D:267:LEU:CD2	1:D:309:VAL:HG21	2.26	0.66
1:B:321:LYS:HZ2	2:F:4:TRP:NE1	1.91	0.66
1:A:238:LYS:HG3	1:A:326:GLN:HE22	1.61	0.65
1:C:261:PHE:CD1	1:C:268:VAL:HG23	2.32	0.65
1:C:162:ASN:C	1:C:162:ASN:HD22	2.04	0.64
1:A:301:LEU:H	1:A:304:GLN:NE2	1.95	0.64
1:A:300:ILE:HD13	1:A:337:ILE:CD1	2.28	0.63
1:B:321:LYS:NZ	2:F:4:TRP:NE1	2.42	0.62
1:C:218:LYS:HE3	1:C:381:ASP:O	1.98	0.62
1:D:301:LEU:H	1:D:304:GLN:HE21	1.47	0.62
1:A:300:ILE:HD13	1:A:337:ILE:HD12	1.81	0.62
1:B:110:ILE:HD11	2:F:7:VAL:HG12	1.81	0.62
1:C:378:ASP:HB3	1:C:381:ASP:OD2	2.01	0.61
1:B:267:LEU:HD13	1:B:309:VAL:HG23	1.83	0.61
1:A:378:ASP:O	1:A:381:ASP:HB2	2.02	0.60
1:D:301:LEU:H	1:D:304:GLN:NE2	2.00	0.60
1:D:249:LYS:HE2	1:D:262:TRP:CD1	2.37	0.60
1:C:321:LYS:HZ3	2:G:4:TRP:CD1	2.21	0.59
1:B:209:ASN:ND2	1:B:282:VAL:H	2.00	0.59
1:C:161:LEU:HB2	1:C:166:VAL:HG22	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:321:LYS:HZ3	2:H:4:TRP:HE1	1.49	0.59
1:B:300:ILE:HD13	1:B:337:ILE:HD12	1.84	0.59
1:A:45:HIS:ND1	1:A:46:PRO:HD2	2.17	0.59
1:D:162:ASN:ND2	1:D:165:GLU:HB2	2.18	0.59
1:D:294:GLN:HG3	1:D:373:PRO:HB2	1.84	0.59
1:C:267:LEU:CD2	1:C:309:VAL:HG21	2.33	0.59
1:A:362:HIS:HD2	1:A:363:ASP:O	1.86	0.58
1:B:209:ASN:HD22	1:B:282:VAL:HG22	1.68	0.58
1:D:65:LYS:HE3	1:D:80:LEU:HD12	1.85	0.58
1:C:4:ASP:OD2	1:C:170:VAL:HG11	2.03	0.58
1:D:11:GLY:C	2:H:5:SER:HB2	2.29	0.58
1:D:149:LEU:C	1:D:149:LEU:HD23	2.29	0.58
2:H:4:TRP:O	2:H:5:SER:CB	2.40	0.58
1:C:313:ALA:C	1:C:315:SER:H	2.12	0.57
1:C:162:ASN:O	1:C:166:VAL:HG23	2.05	0.57
1:C:264:GLY:O	1:C:321:LYS:HE3	2.04	0.57
1:A:267:LEU:H	1:A:267:LEU:HD23	1.69	0.56
1:D:321:LYS:HZ2	2:H:4:TRP:NE1	2.03	0.56
1:D:198:TYR:CE2	1:D:224:LYS:HE3	2.40	0.56
1:D:11:GLY:O	2:H:5:SER:HB2	2.05	0.56
1:B:267:LEU:HD13	1:B:309:VAL:CG2	2.35	0.56
1:B:321:LYS:HZ3	2:F:4:TRP:HE1	1.50	0.56
1:A:209:ASN:HD22	1:A:282:VAL:HG22	1.69	0.56
1:A:312:VAL:H	1:C:313:ALA:HB2	1.69	0.56
1:B:197:TRP:CG	1:B:198:TYR:H	2.24	0.56
1:D:362:HIS:HD2	1:D:363:ASP:O	1.88	0.56
1:C:215:MET:HE1	1:C:239:LYS:HB3	1.87	0.55
1:A:162:ASN:C	1:A:162:ASN:ND2	2.63	0.55
1:D:2:MET:HG2	1:D:90:GLY:HA2	1.88	0.55
1:A:149:LEU:HB2	1:A:346:ASP:HA	1.89	0.55
1:B:44:PRO:HD3	1:B:51:TYR:CZ	2.42	0.55
1:D:321:LYS:NZ	2:H:4:TRP:NE1	2.55	0.54
1:B:149:LEU:HB2	1:B:346:ASP:HA	1.89	0.54
1:D:293:ASN:HA	1:D:376:THR:O	2.07	0.54
1:C:294:GLN:HG3	1:C:373:PRO:HB2	1.90	0.54
1:C:312:VAL:HG23	1:C:313:ALA:H	1.73	0.53
1:D:321:LYS:HZ2	2:H:4:TRP:HE1	1.56	0.53
1:B:161:LEU:HB2	1:B:166:VAL:HG22	1.89	0.53
1:B:234:LEU:HG	1:B:337:ILE:HD11	1.91	0.53
1:C:294:GLN:O	1:C:379:MET:HE1	2.09	0.53
1:D:162:ASN:C	1:D:162:ASN:HD22	2.17	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2:MET:HG2	1:A:90:GLY:HA2	1.90	0.52
1:A:197:TRP:CD1	1:A:198:TYR:H	2.28	0.52
1:A:261:PHE:CD1	1:A:268:VAL:HG23	2.44	0.52
1:B:162:ASN:HD21	1:B:165:GLU:H	1.55	0.52
1:B:258:PRO:HG3	1:B:266:GLN:HE21	1.73	0.52
1:B:362:HIS:HD2	1:B:363:ASP:O	1.92	0.52
1:C:199:TYR:HB3	1:C:352:ILE:HD11	1.91	0.52
1:A:197:TRP:CG	1:A:198:TYR:H	2.28	0.52
1:B:254:THR:HG23	3:C:479:HOH:O	2.08	0.51
1:D:199:TYR:HB3	1:D:352:ILE:HD11	1.92	0.51
2:G:3:TRP:O	2:G:4:TRP:HB3	2.10	0.51
1:B:294:GLN:HG3	1:B:373:PRO:HB2	1.92	0.51
1:C:197:TRP:CG	1:C:198:TYR:H	2.29	0.50
1:D:162:ASN:HD22	1:D:165:GLU:HB2	1.76	0.50
1:D:364:GLU:O	1:D:364:GLU:HG2	2.11	0.50
1:B:4:ASP:HB3	1:B:170:VAL:HG11	1.90	0.50
1:C:362:HIS:HD2	1:C:363:ASP:O	1.94	0.50
1:C:65:LYS:HE3	1:C:80:LEU:CD1	2.42	0.50
1:D:294:GLN:O	1:D:379:MET:HE1	2.11	0.50
1:D:307:ARG:NH1	2:H:5:SER:OG	2.45	0.50
1:B:364:GLU:O	1:B:364:GLU:HG2	2.11	0.50
1:D:65:LYS:HE3	1:D:80:LEU:CD1	2.42	0.50
1:B:73:GLN:HG2	3:F:719:HOH:O	2.11	0.49
1:C:326:GLN:HG2	3:C:463:HOH:O	2.11	0.49
1:A:321:LYS:HZ2	2:E:4:TRP:CD1	2.31	0.49
1:C:32:ASP:OD1	1:C:230:GLY:HA3	2.13	0.49
2:H:5:SER:OG	2:H:6:GLU:N	2.44	0.48
1:B:162:ASN:HD21	1:B:165:GLU:HG3	1.79	0.48
1:C:110:ILE:HD11	2:G:7:VAL:HG12	1.94	0.48
1:C:312:VAL:CG2	2:E:3:TRP:N	2.73	0.48
1:D:55:GLN:HB2	3:D:469:HOH:O	2.12	0.48
2:F:4:TRP:CD1	2:F:4:TRP:C	2.91	0.48
1:A:45:HIS:CG	1:A:46:PRO:HD2	2.48	0.48
1:C:209:ASN:ND2	1:C:282:VAL:H	2.11	0.48
1:A:47(P):PHE:CZ	1:A:178:GLY:HA3	2.48	0.48
1:B:4:ASP:CA	1:B:170:VAL:HG11	2.44	0.48
1:B:300:ILE:HD13	1:B:337:ILE:CD1	2.44	0.48
1:C:166:VAL:HG12	3:C:505:HOH:O	2.14	0.48
1:A:65:LYS:HE3	1:A:80:LEU:HD12	1.94	0.47
1:C:65:LYS:HE3	1:C:80:LEU:HD12	1.96	0.47
1:A:71:TYR:O	1:A:72:THR:C	2.56	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:209:ASN:ND2	1:D:282:VAL:H	2.12	0.47
1:A:95:VAL:HG11	1:A:140:LEU:HD12	1.97	0.47
1:B:73:GLN:HB3	2:F:8:ASN:HB2	1.96	0.47
1:C:222:TYR:O	1:C:223:ASP:CB	2.63	0.47
1:B:323:ALA:HB1	1:B:336:VAL:HG21	1.96	0.47
1:A:181:HIS:HA	1:A:184:TYR:CE1	2.50	0.47
1:D:11:GLY:CA	2:H:5:SER:HB2	2.45	0.47
1:B:50:ARG:O	1:B:116:GLU:HG2	2.15	0.47
1:D:44:PRO:HD3	1:D:51:TYR:CZ	2.51	0.46
1:A:47(P):PHE:CD1	1:A:47(P):PHE:N	2.83	0.46
1:B:216:ASP:OD2	1:B:218:LYS:HB2	2.15	0.46
1:B:261:PHE:CD1	1:B:268:VAL:CG2	2.98	0.46
1:A:311:ASP:HB2	1:A:319:CYS:SG	2.56	0.46
1:C:161:LEU:CB	1:C:166:VAL:HG22	2.44	0.46
1:D:258:PRO:HB2	1:D:266:GLN:OE1	2.16	0.46
1:B:161:LEU:CB	1:B:166:VAL:HG22	2.45	0.46
1:A:162:ASN:HD22	1:A:165:GLU:H	1.63	0.46
1:C:197:TRP:CD1	1:C:198:TYR:H	2.33	0.46
1:A:18:MET:HE2	1:A:85:VAL:CG1	2.46	0.46
1:A:311:ASP:CG	1:A:313:ALA:HB3	2.40	0.46
1:D:261:PHE:CD1	1:D:268:VAL:HG23	2.51	0.46
1:B:47(P):PHE:N	1:B:47(P):PHE:CD1	2.84	0.45
1:C:209:ASN:HD22	1:C:282:VAL:HG22	1.80	0.45
1:D:162:ASN:ND2	1:D:162:ASN:C	2.73	0.45
1:B:197:TRP:CD1	1:B:198:TYR:H	2.35	0.45
1:A:209:ASN:ND2	1:A:282:VAL:H	2.14	0.45
1:B:375:VAL:H	1:C:364:GLU:CD	2.24	0.45
1:D:197:TRP:CG	1:D:198:TYR:H	2.35	0.45
1:D:276:PRO:O	1:D:279:ILE:HG12	2.17	0.45
1:D:110:ILE:HD11	2:H:7:VAL:HG12	1.99	0.44
1:D:209:ASN:HD22	1:D:209:ASN:HA	1.65	0.44
1:A:45:HIS:CE1	1:A:46:PRO:HD2	2.52	0.44
1:A:222:TYR:O	1:A:223:ASP:CB	2.65	0.44
1:A:47:PHE:CE2	1:A:111:ASN:HB2	2.53	0.44
1:A:313:ALA:H	1:C:313:ALA:HB1	1.83	0.44
1:A:276:PRO:O	1:A:279:ILE:HG12	2.17	0.44
1:D:161:LEU:HB2	1:D:166:VAL:HG22	1.99	0.44
1:B:163:GLN:HA	1:B:163:GLN:NE2	2.32	0.44
1:D:260:GLY:C	1:D:266:GLN:HG2	2.43	0.44
1:C:316:GLN:O	1:C:317:ASP:C	2.60	0.44
1:C:162:ASN:ND2	1:C:162:ASN:C	2.72	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:18:MET:SD	1:D:29:ILE:HG13	2.58	0.44
1:A:312:VAL:H	1:C:313:ALA:CB	2.31	0.43
1:C:174:MET:HE3	1:C:176:ILE:HD11	2.00	0.43
1:C:371:GLU:HG2	3:C:556:HOH:O	2.18	0.43
1:B:4:ASP:OD2	1:B:170:VAL:HG11	2.17	0.43
1:C:209:ASN:HD22	1:C:209:ASN:HA	1.62	0.43
1:A:93:VAL:HG21	1:A:140:LEU:HD11	1.99	0.43
1:C:313:ALA:O	1:C:315:SER:N	2.52	0.43
1:A:376:THR:HB	1:A:379:MET:SD	2.59	0.43
1:B:222:TYR:O	1:B:223:ASP:CB	2.67	0.43
1:C:70:PRO:HG2	2:G:11:GLU:HB2	2.01	0.43
1:A:162:ASN:HD21	1:A:165:GLU:H	1.61	0.43
1:A:267:LEU:HD23	1:A:267:LEU:N	2.33	0.43
1:B:162:ASN:C	1:B:162:ASN:ND2	2.63	0.42
1:A:213:LEU:HD12	1:A:213:LEU:HA	1.80	0.42
1:A:30:LEU:HD23	1:A:32:ASP:HB2	2.02	0.42
1:C:162:ASN:HD22	1:C:165:GLU:H	1.65	0.42
1:C:300:ILE:HD13	1:C:337:ILE:HD13	2.01	0.42
1:D:274:THR:O	1:D:275:THR:C	2.62	0.42
1:B:323:ALA:CB	1:B:336:VAL:HG21	2.50	0.42
1:B:209:ASN:HD22	1:B:209:ASN:HA	1.64	0.42
1:B:267:LEU:HD22	1:B:309:VAL:CG2	2.46	0.42
1:C:45(P):GLY:N	1:C:181:HIS:CE1	2.87	0.42
1:B:110:ILE:CD1	2:F:7:VAL:HG12	2.48	0.42
1:C:313:ALA:C	1:C:315:SER:N	2.76	0.42
1:A:91:PRO:HD3	1:A:176:ILE:HB	2.01	0.42
1:A:294:GLN:HG3	1:A:373:PRO:HB2	2.02	0.42
1:A:312:VAL:HG23	1:C:315:SER:HB2	2.01	0.42
1:D:11:GLY:HA3	2:H:5:SER:HB2	2.01	0.42
1:A:46:PRO:CG	1:C:263:LEU:HD12	2.37	0.42
1:A:222:TYR:O	1:A:223:ASP:HB3	2.20	0.42
1:D:307:ARG:NH2	2:H:5:SER:OG	2.53	0.41
1:A:149:LEU:HD23	1:A:149:LEU:O	2.20	0.41
1:A:13:GLY:HA3	1:A:30:LEU:CD1	2.50	0.41
1:B:134:GLU:HA	1:B:135:PRO:HD3	1.86	0.41
1:B:321:LYS:HD2	2:F:4:TRP:CE2	2.55	0.41
1:D:162:ASN:ND2	1:D:165:GLU:N	2.66	0.41
1:B:215:MET:HE1	1:B:239:LYS:HB3	2.02	0.41
1:C:312:VAL:CB	2:E:3:TRP:N	2.83	0.41
1:C:195:ARG:HB3	1:C:197:TRP:CD1	2.56	0.41
1:A:65:LYS:HE3	1:A:80:LEU:CD1	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:47(P):PHE:CZ	1:D:178:GLY:HA3	2.56	0.41
1:B:267:LEU:CD1	1:B:309:VAL:HG21	2.50	0.41
1:D:151:SER:OG	1:D:175:ILE:HB	2.21	0.41
1:B:336:VAL:HG12	1:B:337:ILE:CD1	2.41	0.41
1:D:315:SER:C	1:D:317:ASP:H	2.28	0.40
1:D:1:GLU:CD	1:D:1:GLU:N	2.69	0.40
1:A:311:ASP:OD1	1:A:313:ALA:HB3	2.21	0.40
1:C:134:GLU:HA	1:C:135:PRO:HD3	1.98	0.40
1:D:11:GLY:HA3	2:H:5:SER:CB	2.51	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	387/389 (100%)	374 (97%)	12 (3%)	1 (0%)	36	29
1	B	387/389 (100%)	374 (97%)	12 (3%)	1 (0%)	36	29
1	C	387/389 (100%)	374 (97%)	10 (3%)	3 (1%)	16	8
1	D	387/389 (100%)	375 (97%)	11 (3%)	1 (0%)	36	29
2	E	7/12 (58%)	5 (71%)	2 (29%)	0	100	100
2	F	7/12 (58%)	6 (86%)	1 (14%)	0	100	100
2	G	7/12 (58%)	5 (71%)	2 (29%)	0	100	100
2	H	7/12 (58%)	3 (43%)	1 (14%)	3 (43%)	0	0
All	All	1576/1604 (98%)	1516 (96%)	51 (3%)	9 (1%)	21	13

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	223	ASP

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Mol	Chain	Res	Type
1	B	223	ASP
1	C	223	ASP
1	D	223	ASP
2	H	5	SER
1	C	314	THR
1	C	317	ASP
2	H	4	TRP
2	H	6	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	331/331 (100%)	320 (97%)	11 (3%)	33	26
1	B	331/331 (100%)	322 (97%)	9 (3%)	39	34
1	C	331/331 (100%)	326 (98%)	5 (2%)	57	56
1	D	331/331 (100%)	325 (98%)	6 (2%)	51	50
2	E	8/10 (80%)	8 (100%)	0	100	100
2	F	8/10 (80%)	6 (75%)	2 (25%)	0	0
2	G	8/10 (80%)	8 (100%)	0	100	100
2	H	8/10 (80%)	7 (88%)	1 (12%)	4	1
All	All	1356/1364 (99%)	1322 (98%)	34 (2%)	42	37

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48(P)	VAL
1	A	48	LEU
1	A	149	LEU
1	A	162	ASN
1	A	197	TRP
1	A	209	ASN
1	A	234	LEU

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Mol	Chain	Res	Type
1	A	267	LEU
1	A	314	THR
1	A	336	VAL
1	A	376	THR
1	B	104	GLU
1	B	162	ASN
1	B	187	SER
1	B	197	TRP
1	B	209	ASN
1	B	234	LEU
1	B	314	THR
1	B	315	SER
1	B	336	VAL
1	C	162	ASN
1	C	209	ASN
1	C	234	LEU
1	C	267	LEU
1	C	326	GLN
1	D	1	GLU
1	D	149	LEU
1	D	162	ASN
1	D	209	ASN
1	D	234	LEU
1	D	336	VAL
2	F	4	TRP
2	F	6	GLU
2	H	3	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (42) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	ASN
1	A	89	HIS
1	A	111	ASN
1	A	162	ASN
1	A	209	ASN
1	A	293	ASN
1	A	304	GLN
1	A	326	GLN
1	A	362	HIS
1	B	28	ASN
1	B	111	ASN

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Mol	Chain	Res	Type
1	B	145	HIS
1	B	162	ASN
1	B	181	HIS
1	B	209	ASN
1	B	266	GLN
1	B	304	GLN
1	B	362	HIS
1	C	28	ASN
1	C	55	GLN
1	C	89	HIS
1	C	111	ASN
1	C	162	ASN
1	C	181	HIS
1	C	209	ASN
1	C	294	GLN
1	C	304	GLN
1	C	362	HIS
1	C	385	ASN
1	D	12	GLN
1	D	28	ASN
1	D	49	HIS
1	D	53	GLN
1	D	73	GLN
1	D	89	HIS
1	D	111	ASN
1	D	162	ASN
1	D	209	ASN
1	D	293	ASN
1	D	294	GLN
1	D	304	GLN
1	D	362	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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	1OL	H	9	2	11,12,13	0.80	0	9,15,17	0.87	1 (11%)
2	1OL	F	9	2	11,12,13	1.09	1 (9%)	9,15,17	1.17	2 (22%)
2	1OL	G	9	2	11,12,13	0.71	0	9,15,17	0.83	1 (11%)
2	1OL	E	9	2	11,12,13	0.56	0	9,15,17	0.98	1 (11%)

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	1OL	H	9	2	-	1/14/14/16	-
2	1OL	F	9	2	-	3/14/14/16	-
2	1OL	G	9	2	-	1/14/14/16	-
2	1OL	E	9	2	-	3/14/14/16	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	9	1OL	C6-CA	3.16	1.56	1.53

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	9	1OL	O-C-C8	-2.32	120.15	125.33
2	F	9	1OL	C3-C4-CA	-2.28	110.86	115.76
2	H	9	1OL	O-C-C8	-2.25	120.31	125.33
2	E	9	1OL	O-C-C8	-2.23	120.34	125.33
2	G	9	1OL	O-C-C8	-2.12	120.61	125.33

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	9	1OL	O-C-C8-C7
2	F	9	1OL	O-C-C8-C7
2	G	9	1OL	O-C-C8-C7
2	H	9	1OL	O-C-C8-C7
2	E	9	1OL	C6-C7-C8-C9
2	E	9	1OL	C6-C7-C8-C
2	F	9	1OL	C6-C7-C8-C9
2	F	9	1OL	C6-C7-C8-C

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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.