



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 27, 2026 – 07:34 PM EDT

PDB ID : 4BKA / pdb_00004bka
Title : crystal structure of the human EphA4 ectodomain in complex with human ephrin A5
Authors : Seiradake, E.; Schaupp, A.; del Toro Ruiz, D.; Kaufmann, R.; Mitakidis, N.; Harlos, K.; Aricescu, A.R.; Klein, R.; Jones, E.Y.
Deposited on : 2013-04-23
Resolution : 5.30 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
Xtriage (Phenix) : 2.0
EDS : **FAILED**
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

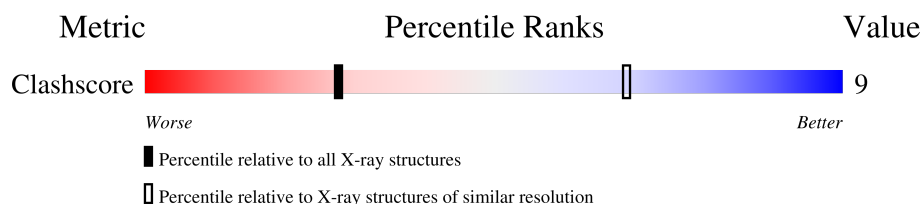
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 5.30 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	1098 (6.60-4.00)

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 $\leq 5\%$

Note EDS failed to run properly.

Mol	Chain	Length	Quality of chain
1	A	568	 73% 16% 11%
2	C	180	 67% 9% 24%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5087 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 EPHRIN TYPE-A RECEPTOR 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	507	3948	2461	682	779	26	0	0	0

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	expression tag	UNP P54764
A	-10	GLY	-	expression tag	UNP P54764
A	-9	ILE	-	expression tag	UNP P54764
A	-8	LEU	-	expression tag	UNP P54764
A	-7	PRO	-	expression tag	UNP P54764
A	-6	SER	-	expression tag	UNP P54764
A	-5	PRO	-	expression tag	UNP P54764
A	-4	GLY	-	expression tag	UNP P54764
A	-3	MET	-	expression tag	UNP P54764
A	-2	PRO	-	expression tag	UNP P54764
A	-1	ALA	-	expression tag	UNP P54764
A	0	LEU	-	expression tag	UNP P54764
A	1	LEU	-	expression tag	UNP P54764
A	2	SER	-	expression tag	UNP P54764
A	3	LEU	-	expression tag	UNP P54764
A	4	VAL	-	expression tag	UNP P54764
A	5	SER	-	expression tag	UNP P54764
A	6	LEU	-	expression tag	UNP P54764
A	7	LEU	-	expression tag	UNP P54764
A	8	SER	-	expression tag	UNP P54764
A	9	VAL	-	expression tag	UNP P54764
A	10	LEU	-	expression tag	UNP P54764
A	11	LEU	-	expression tag	UNP P54764
A	12	MET	-	expression tag	UNP P54764
A	13	GLY	-	expression tag	UNP P54764
A	14	CYS	-	expression tag	UNP P54764
A	15	VAL	-	expression tag	UNP P54764

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Chain	Residue	Modelled	Actual	Comment	Reference
A	16	ALA	-	expression tag	UNP P54764
A	17	GLU	-	expression tag	UNP P54764
A	18	THR	-	expression tag	UNP P54764
A	19	GLY	-	expression tag	UNP P54764
A	548	GLY	-	expression tag	UNP P54764
A	549	THR	-	expression tag	UNP P54764
A	550	LYS	-	expression tag	UNP P54764
A	551	HIS	-	expression tag	UNP P54764
A	552	HIS	-	expression tag	UNP P54764
A	553	HIS	-	expression tag	UNP P54764
A	554	HIS	-	expression tag	UNP P54764
A	555	HIS	-	expression tag	UNP P54764
A	556	HIS	-	expression tag	UNP P54764

- Molecule 2 is a protein called EPHRIN-A5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	137	Total	C	N	O	S	0	0	1
			1139	728	197	206	8			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	0	MET	-	expression tag	UNP P52803
C	1	GLY	-	expression tag	UNP P52803
C	2	ILE	-	expression tag	UNP P52803
C	3	LEU	-	expression tag	UNP P52803
C	4	PRO	-	expression tag	UNP P52803
C	5	SER	-	expression tag	UNP P52803
C	6	PRO	-	expression tag	UNP P52803
C	7	GLY	-	expression tag	UNP P52803
C	8	MET	-	expression tag	UNP P52803
C	9	PRO	-	expression tag	UNP P52803
C	10	ALA	-	expression tag	UNP P52803
C	11	LEU	-	expression tag	UNP P52803
C	12	LEU	-	expression tag	UNP P52803
C	13	SER	-	expression tag	UNP P52803
C	14	LEU	-	expression tag	UNP P52803
C	15	VAL	-	expression tag	UNP P52803
C	16	SER	-	expression tag	UNP P52803
C	17	LEU	-	expression tag	UNP P52803
C	18	LEU	-	expression tag	UNP P52803

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Chain	Residue	Modelled	Actual	Comment	Reference
C	19	SER	-	expression tag	UNP P52803
C	20	VAL	-	expression tag	UNP P52803
C	21	LEU	-	expression tag	UNP P52803
C	22	LEU	-	expression tag	UNP P52803
C	23	MET	-	expression tag	UNP P52803
C	24	GLY	-	expression tag	UNP P52803
C	25	CYS	-	expression tag	UNP P52803
C	26	VAL	-	expression tag	UNP P52803
C	27	ALA	-	expression tag	UNP P52803
C	28	GLU	-	expression tag	UNP P52803
C	29	THR	-	expression tag	UNP P52803
C	30	GLY	-	expression tag	UNP P52803
C	172	GLY	-	expression tag	UNP P52803
C	173	THR	-	expression tag	UNP P52803
C	174	LYS	-	expression tag	UNP P52803
C	175	HIS	-	expression tag	UNP P52803
C	176	HIS	-	expression tag	UNP P52803
C	177	HIS	-	expression tag	UNP P52803
C	178	HIS	-	expression tag	UNP P52803
C	179	HIS	-	expression tag	UNP P52803
C	180	HIS	-	expression tag	UNP P52803

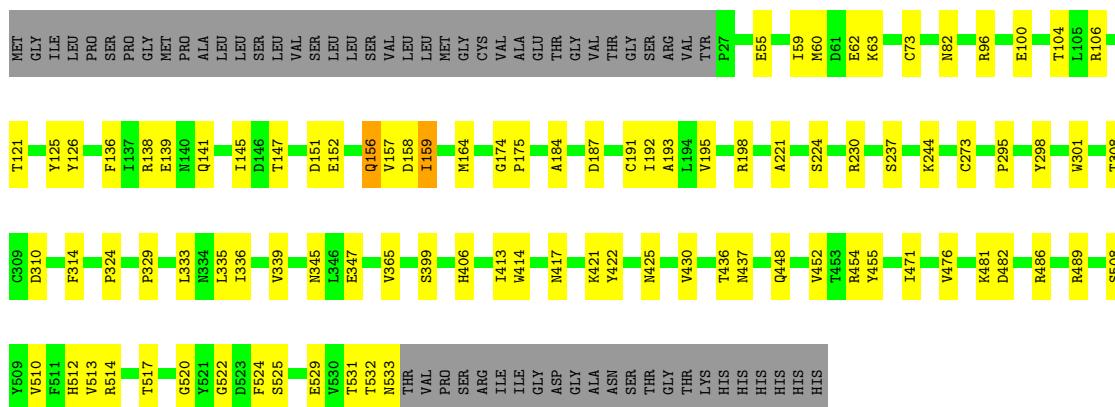
3 Residue-property plots

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 failed to run properly.

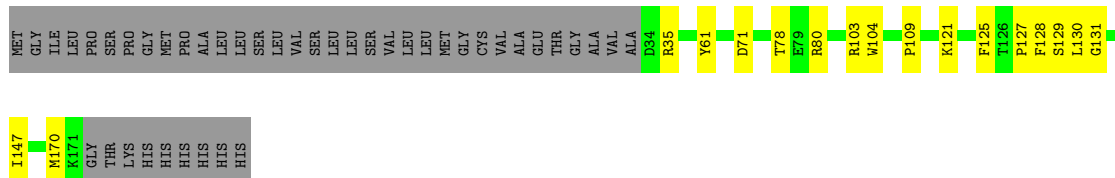
• Molecule 1: EPHRIN TYPE-A RECEPTOR 4

Chain A: 



• Molecule 2: EPHRIN-A5

Chain C: 



4 Data and refinement statistics

EDS failed to run properly - this section is therefore incomplete.

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	202.63Å 202.63Å 326.23Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	84.87 – 5.30	Depositor
% Data completeness (in resolution range)	96.3 (84.87-5.30)	Depositor
R_{merge}	0.74	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.67 (at 5.41Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.383 , 0.408	Depositor
Wilson B-factor (Å ²)	155.3	Xtriage
Anisotropy	0.356	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.31$, $\langle L^2 \rangle = 0.15$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5087	wwPDB-VP
Average B, all atoms (Å ²)	243.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.86% of the height of the origin peak. No significant pseudotranslation is detected.*

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

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	0.66	0/4032	1.15	4/5484 (0.1%)
2	C	0.63	1/1177 (0.1%)	1.03	0/1592
All	All	0.65	1/5209 (0.0%)	1.13	4/7076 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	170	MET	C-N	-5.73	1.25	1.33

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	333	LEU	N-CA-C	7.81	119.79	111.28
1	A	159	ILE	N-CA-C	5.23	116.00	110.72
1	A	156	GLN	N-CA-C	5.10	117.97	111.69
1	A	520	GLY	N-CA-C	5.09	117.45	110.58

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	3948	0	3801	86	104
2	C	1139	0	1053	61	8
All	All	5087	0	4854	89	104

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 9.

All (89) 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:195:VAL:HG22	2:C:128:PHE:CE1	1.46	1.46
1:A:195:VAL:CG2	2:C:128:PHE:HE1	1.31	1.39
1:A:164:MET:CE	2:C:130:LEU:HG	1.72	1.17
1:A:55:GLU:HG2	2:C:61:TYR:OH	1.46	1.13
1:A:55:GLU:OE1	2:C:61:TYR:OH	1.64	1.13
1:A:55:GLU:CG	2:C:61:TYR:OH	1.98	1.11
1:A:55:GLU:OE1	2:C:121:LYS:HD2	1.56	1.04
1:A:55:GLU:CD	2:C:61:TYR:HH	1.65	1.04
1:A:59:ILE:HG12	2:C:103:ARG:HH22	1.17	1.03
1:A:55:GLU:CD	2:C:61:TYR:OH	2.01	1.03
1:A:164:MET:HE2	2:C:130:LEU:CG	1.90	1.01
1:A:164:MET:HE2	2:C:130:LEU:HG	1.01	1.00
1:A:59:ILE:HG12	2:C:103:ARG:NH2	1.79	0.97
1:A:104:THR:HG21	2:C:129:SER:N	1.79	0.97
1:A:59:ILE:CG1	2:C:103:ARG:HH22	1.77	0.96
1:A:104:THR:HG21	2:C:129:SER:H	1.32	0.94
1:A:191:CYS:C	2:C:127:PRO:HB3	1.92	0.94
1:A:195:VAL:CG2	2:C:128:PHE:CE1	2.24	0.89
1:A:164:MET:SD	2:C:130:LEU:HB2	2.17	0.84
1:A:164:MET:CE	2:C:130:LEU:CG	2.53	0.80
1:A:55:GLU:OE1	2:C:121:LYS:CD	2.34	0.75
1:A:195:VAL:HG22	2:C:128:PHE:HE1	0.60	0.75
1:A:164:MET:CE	2:C:130:LEU:HB2	2.17	0.75
1:A:73:CYS:HB2	2:C:127:PRO:HG3	1.73	0.70
1:A:193:ALA:CB	2:C:128:PHE:HB2	2.21	0.70
1:A:59:ILE:HG23	2:C:103:ARG:HH12	1.57	0.69
1:A:164:MET:CE	2:C:130:LEU:CB	2.71	0.69
1:A:59:ILE:HG12	2:C:103:ARG:CZ	2.24	0.67
1:A:106:ARG:HH21	2:C:125:PHE:HB3	1.60	0.66
1:A:60:MET:CG	2:C:103:ARG:O	2.46	0.63
1:A:193:ALA:HB1	2:C:128:PHE:HB2	1.81	0.62
1:A:59:ILE:HG23	2:C:103:ARG:NH1	2.15	0.61
1:A:195:VAL:HG21	2:C:128:PHE:CE1	2.34	0.61
1:A:59:ILE:HG12	2:C:103:ARG:NH1	2.15	0.60
1:A:59:ILE:HG12	2:C:103:ARG:HH12	1.66	0.59
1:A:138:ARG:HB3	1:A:141:GLN:HB2	1.85	0.57
1:A:164:MET:HE1	2:C:130:LEU:HG	1.81	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:GLU:OE1	2:C:121:LYS:NZ	2.37	0.57
2:C:35:ARG:HG2	2:C:61:TYR:HB2	1.86	0.57
1:A:329:PRO:HD3	1:A:417:ASN:HB2	1.86	0.56
1:A:164:MET:HE1	2:C:130:LEU:CB	2.35	0.55
1:A:336:ILE:HB	1:A:347:GLU:HB3	1.87	0.55
1:A:471:ILE:HA	1:A:517:THR:HG22	1.87	0.55
1:A:100:GLU:HB3	1:A:198:ARG:HB2	1.88	0.55
1:A:106:ARG:HD3	1:A:157:VAL:HG23	1.90	0.54
1:A:59:ILE:HD13	2:C:131:GLY:HA2	1.89	0.54
1:A:125:TYR:HB2	1:A:184:ALA:HB3	1.90	0.53
2:C:80:ARG:HB3	2:C:147:ILE:HD12	1.90	0.53
1:A:104:THR:HB	2:C:128:PHE:HA	1.90	0.53
1:A:55:GLU:OE1	2:C:121:LYS:CE	2.58	0.51
1:A:298:TYR:HB2	1:A:310:ASP:HB3	1.93	0.51
1:A:193:ALA:CB	2:C:128:PHE:CB	2.88	0.50
1:A:513:VAL:H	1:A:525:SER:HB2	1.76	0.50
1:A:126:TYR:HB3	1:A:145:ILE:HD11	1.94	0.49
1:A:106:ARG:HB2	1:A:191:CYS:HB3	1.95	0.49
1:A:82:ASN:HB2	1:A:187:ASP:HB3	1.94	0.49
1:A:59:ILE:CG2	2:C:103:ARG:HH12	2.24	0.49
1:A:104:THR:HG21	2:C:129:SER:CA	2.43	0.47
1:A:406:HIS:H	1:A:437:ASN:HB2	1.79	0.47
1:A:104:THR:HG21	2:C:129:SER:CB	2.44	0.47
2:C:78:THR:HG21	2:C:109:PRO:HG3	1.97	0.47
1:A:191:CYS:SG	2:C:127:PRO:HG3	2.54	0.47
1:A:192:ILE:C	2:C:127:PRO:HB2	2.40	0.47
1:A:295:PRO:HD2	1:A:324:PRO:HB3	1.97	0.46
1:A:59:ILE:HG13	2:C:103:ARG:HH22	1.74	0.46
1:A:192:ILE:N	2:C:127:PRO:HB3	2.27	0.46
1:A:486:ARG:HD3	1:A:489:ARG:HH12	1.81	0.46
1:A:193:ALA:HB3	2:C:128:PHE:HA	1.98	0.45
1:A:339:VAL:HG21	1:A:436:THR:HA	1.98	0.45
1:A:514:ARG:HH11	1:A:522:GLY:H	1.65	0.45
1:A:60:MET:HG3	2:C:103:ARG:O	2.17	0.45
1:A:191:CYS:O	2:C:127:PRO:HB3	2.15	0.44
1:A:175:PRO:HB3	1:A:221:ALA:HB1	1.99	0.44
1:A:230:ARG:HG2	1:A:244:LYS:HB3	2.00	0.44
1:A:237:SER:HB2	1:A:273:CYS:SG	2.58	0.44
1:A:335:LEU:HD11	1:A:413:ILE:HD12	1.99	0.44
1:A:104:THR:CG2	2:C:129:SER:HB3	2.49	0.43
1:A:60:MET:HG3	2:C:104:TRP:HA	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:MET:HG2	2:C:103:ARG:O	2.17	0.42
1:A:121:THR:HB	1:A:147:THR:HG22	2.00	0.42
1:A:335:LEU:HD12	1:A:430:VAL:HG12	2.02	0.42
1:A:476:VAL:HG22	1:A:513:VAL:HG22	2.01	0.42
1:A:62:GLU:HA	1:A:63:LYS:HA	1.76	0.41
1:A:159:ILE:H	1:A:159:ILE:HG13	1.75	0.41
1:A:96:ARG:HG3	1:A:174:GLY:HA3	2.02	0.41
1:A:193:ALA:CB	2:C:128:PHE:CA	2.99	0.41
1:A:156:GLN:N	1:A:157:VAL:HA	2.36	0.41
1:A:164:MET:HE1	2:C:130:LEU:CG	2.41	0.40
1:A:55:GLU:OE1	2:C:61:TYR:CZ	2.66	0.40

All (104) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:301:TRP:CE2	1:A:414:TRP:CE3[5_675]	0.50	1.70
1:A:448:GLN:OE1	1:A:448:GLN:OE1[16_454]	0.67	1.53
1:A:481:LYS:CD	1:A:533:ASN:CG[16_454]	0.68	1.52
1:A:301:TRP:CZ2	1:A:414:TRP:CD2[5_675]	0.69	1.51
1:A:481:LYS:CE	1:A:533:ASN:OD1[16_454]	0.73	1.47
1:A:301:TRP:CZ2	1:A:414:TRP:CE2[5_675]	0.74	1.46
1:A:298:TYR:CD2	1:A:425:ASN:ND2[5_675]	0.76	1.44
1:A:152:GLU:OE2	1:A:512:HIS:NE2[6_555]	0.82	1.38
1:A:308:THR:OG1	1:A:421:LYS:O[5_675]	0.95	1.25
1:A:301:TRP:CE2	1:A:414:TRP:CD2[5_675]	0.97	1.23
1:A:454:ARG:NH1	1:A:482:ASP:OD1[16_454]	0.97	1.23
1:A:481:LYS:CD	1:A:533:ASN:ND2[16_454]	0.98	1.22
1:A:508:SER:OG	1:A:532:THR:O[16_454]	1.01	1.19
1:A:531:THR:O	1:A:531:THR:OG1[16_454]	1.01	1.19
1:A:301:TRP:NE1	1:A:414:TRP:CE3[5_675]	1.04	1.16
1:A:481:LYS:CE	1:A:533:ASN:CG[16_454]	1.06	1.14
1:A:158:ASP:OD2	1:A:455:TYR:OH[8_565]	1.11	1.09
1:A:301:TRP:CH2	1:A:414:TRP:CE2[5_675]	1.11	1.09
1:A:139:GLU:OE2	2:C:71:ASP:O[3_465]	1.13	1.07
1:A:531:THR:CG2	1:A:531:THR:CG2[16_454]	1.14	1.06
1:A:136:PHE:CE1	2:C:71:ASP:OD1[3_465]	1.15	1.05
1:A:531:THR:C	1:A:531:THR:OG1[16_454]	1.21	0.99
1:A:422:TYR:CZ	1:A:422:TYR:OH[5_675]	1.29	0.91
1:A:422:TYR:CE1	1:A:422:TYR:OH[5_675]	1.29	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:301:TRP:CD2	1:A:414:TRP:CE3[5_675]	1.30	0.90
1:A:422:TYR:CE2	1:A:422:TYR:CZ[5_675]	1.32	0.88
1:A:158:ASP:CG	1:A:455:TYR:OH[8_565]	1.37	0.83
1:A:301:TRP:CZ3	1:A:365:VAL:CG2[5_675]	1.37	0.83
1:A:152:GLU:OE2	1:A:512:HIS:CE1[6_555]	1.39	0.81
1:A:531:THR:CB	1:A:531:THR:CB[16_454]	1.39	0.81
1:A:481:LYS:CG	1:A:533:ASN:ND2[16_454]	1.41	0.79
1:A:531:THR:CA	1:A:531:THR:CB[16_454]	1.42	0.78
1:A:422:TYR:CE2	1:A:422:TYR:OH[5_675]	1.43	0.77
1:A:531:THR:C	1:A:531:THR:CB[16_454]	1.43	0.77
1:A:422:TYR:CD1	1:A:422:TYR:OH[5_675]	1.44	0.76
1:A:448:GLN:NE2	1:A:448:GLN:NE2[16_454]	1.45	0.75
1:A:452:VAL:O	1:A:529:GLU:OE1[16_454]	1.47	0.73
1:A:481:LYS:CD	1:A:533:ASN:OD1[16_454]	1.47	0.73
1:A:301:TRP:CD2	1:A:414:TRP:CZ3[5_675]	1.49	0.71
1:A:301:TRP:CH2	1:A:414:TRP:NE1[5_675]	1.50	0.70
1:A:301:TRP:CE2	1:A:414:TRP:CZ3[5_675]	1.51	0.69
1:A:422:TYR:CD2	1:A:422:TYR:OH[5_675]	1.57	0.63
1:A:422:TYR:CE1	1:A:422:TYR:CZ[5_675]	1.57	0.63
1:A:224:SER:OG	1:A:345:ASN:OD1[6_555]	1.61	0.59
1:A:422:TYR:CG	1:A:422:TYR:OH[5_675]	1.61	0.59
1:A:298:TYR:CE2	1:A:425:ASN:ND2[5_675]	1.62	0.58
1:A:482:ASP:OD2	1:A:533:ASN:C[16_454]	1.62	0.58
1:A:448:GLN:CD	1:A:448:GLN:OE1[16_454]	1.65	0.55
1:A:152:GLU:CD	1:A:512:HIS:NE2[6_555]	1.66	0.54
1:A:301:TRP:CZ2	1:A:414:TRP:NE1[5_675]	1.67	0.53
1:A:139:GLU:CD	2:C:71:ASP:O[3_465]	1.68	0.52
1:A:301:TRP:CE3	1:A:365:VAL:CG2[5_675]	1.69	0.51
1:A:301:TRP:CZ2	1:A:414:TRP:CG[5_675]	1.69	0.51
1:A:481:LYS:NZ	1:A:533:ASN:CB[16_454]	1.69	0.51
1:A:139:GLU:OE2	2:C:71:ASP:C[3_465]	1.72	0.48
1:A:301:TRP:CD1	1:A:414:TRP:CE3[5_675]	1.74	0.46
1:A:308:THR:CB	1:A:421:LYS:O[5_675]	1.74	0.46
1:A:152:GLU:OE2	1:A:512:HIS:CD2[6_555]	1.77	0.43
1:A:151:ASP:OD1	1:A:524:PHE:CZ[6_555]	1.78	0.42
1:A:481:LYS:NZ	1:A:533:ASN:OD1[16_454]	1.79	0.41
1:A:314:PHE:CZ	1:A:314:PHE:CZ[5_675]	1.81	0.39
1:A:481:LYS:NZ	1:A:533:ASN:CG[16_454]	1.81	0.39
1:A:301:TRP:CH2	1:A:414:TRP:CD2[5_675]	1.82	0.38
1:A:314:PHE:CE1	1:A:314:PHE:CE1[5_675]	1.83	0.37
1:A:301:TRP:NE1	1:A:414:TRP:CZ3[5_675]	1.84	0.36

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:301:TRP:CG	1:A:414:TRP:CZ3[5_675]	1.85	0.35
1:A:422:TYR:CE2	1:A:422:TYR:CE2[5_675]	1.85	0.35
1:A:508:SER:OG	1:A:532:THR:C[16_454]	1.87	0.33
1:A:301:TRP:CG	1:A:414:TRP:CE3[5_675]	1.88	0.32
1:A:298:TYR:CG	1:A:425:ASN:ND2[5_675]	1.89	0.31
1:A:301:TRP:CZ2	1:A:414:TRP:CE3[5_675]	1.89	0.31
1:A:151:ASP:OD1	1:A:524:PHE:CE2[6_555]	1.90	0.30
1:A:136:PHE:CZ	2:C:71:ASP:OD1[3_465]	1.91	0.29
1:A:301:TRP:CE2	1:A:414:TRP:CE2[5_675]	1.93	0.27
1:A:454:ARG:NH1	1:A:482:ASP:CG[16_454]	1.93	0.27
1:A:224:SER:CB	1:A:399:SER:OG[6_555]	1.94	0.26
1:A:152:GLU:CD	1:A:512:HIS:CE1[6_555]	1.95	0.25
1:A:301:TRP:CZ2	1:A:414:TRP:CZ2[5_675]	1.95	0.25
1:A:136:PHE:CE1	2:C:71:ASP:CG[3_465]	1.96	0.24
1:A:139:GLU:OE1	2:C:71:ASP:O[3_465]	1.97	0.23
1:A:531:THR:CB	1:A:531:THR:CG2[16_454]	1.97	0.23
1:A:301:TRP:CH2	1:A:414:TRP:CZ2[5_675]	1.98	0.22
1:A:454:ARG:CZ	1:A:482:ASP:OD1[16_454]	1.98	0.22
1:A:481:LYS:CD	1:A:533:ASN:CB[16_454]	1.99	0.21
1:A:298:TYR:CD2	1:A:425:ASN:CG[5_675]	2.00	0.20
1:A:531:THR:O	1:A:531:THR:CB[16_454]	2.00	0.20
1:A:301:TRP:CD1	1:A:414:TRP:CZ3[5_675]	2.01	0.19
1:A:481:LYS:CE	1:A:533:ASN:ND2[16_454]	2.01	0.19
1:A:301:TRP:NE1	1:A:414:TRP:CD2[5_675]	2.03	0.17
1:A:301:TRP:CZ2	1:A:414:TRP:CD1[5_675]	2.04	0.16
1:A:452:VAL:O	1:A:529:GLU:CD[16_454]	2.04	0.16
1:A:139:GLU:OE2	2:C:71:ASP:CA[3_465]	2.07	0.13
1:A:314:PHE:CE1	1:A:314:PHE:CZ[5_675]	2.09	0.11
1:A:152:GLU:OE1	1:A:510:VAL:CG1[6_555]	2.11	0.09
1:A:448:GLN:CD	1:A:448:GLN:NE2[16_454]	2.11	0.09
1:A:481:LYS:CG	1:A:533:ASN:CG[16_454]	2.12	0.08
1:A:152:GLU:CG	1:A:512:HIS:CE1[6_555]	2.13	0.07
1:A:301:TRP:CD2	1:A:414:TRP:CD2[5_675]	2.13	0.07
1:A:448:GLN:CD	1:A:448:GLN:CD[16_454]	2.13	0.07
1:A:298:TYR:CE2	1:A:425:ASN:CG[5_675]	2.14	0.06
1:A:308:THR:OG1	1:A:421:LYS:C[5_675]	2.15	0.05
1:A:301:TRP:CZ3	1:A:365:VAL:CB[5_675]	2.17	0.03
1:A:158:ASP:CB	1:A:455:TYR:OH[8_565]	2.18	0.02
1:A:481:LYS:CE	1:A:533:ASN:CB[16_454]	2.18	0.02

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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](#)

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	C	1
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	148:PRO	C	150:ASP	N	3.28
1	A	202:LYS	C	203:LYS	N	3.25

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS failed to run properly - this section is therefore empty.

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

EDS failed to run properly - this section is therefore empty.

6.3 Carbohydrates

EDS failed to run properly - this section is therefore empty.

6.4 Ligands

EDS failed to run properly - this section is therefore empty.

6.5 Other polymers

EDS failed to run properly - this section is therefore empty.