



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:34 AM UTC

PDB ID : 1FPS / pdb_00001fps
Title : CRYSTAL STRUCTURE OF RECOMBINANT FARNESYL DIPHOSPHATE SYNTHASE AT 2.6 ANGSTROMS RESOLUTION
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Deposited on : 1994-06-30
Resolution : 2.60 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)	:	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

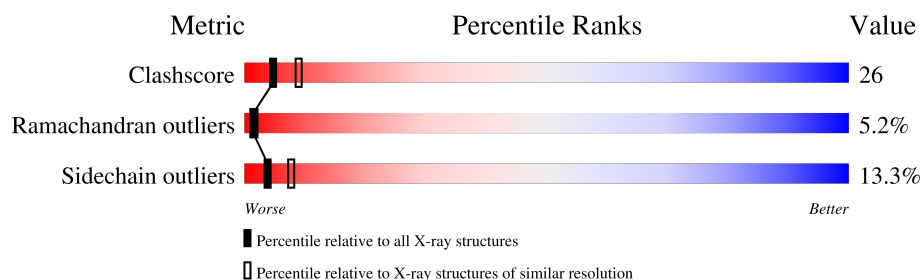
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)

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	348	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2894 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 FARNESYL DIPHOSPHATE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	348	Total	C	N	O	S	0	0	0
			2815	1798	477	526	14			

- Molecule 2 is water.

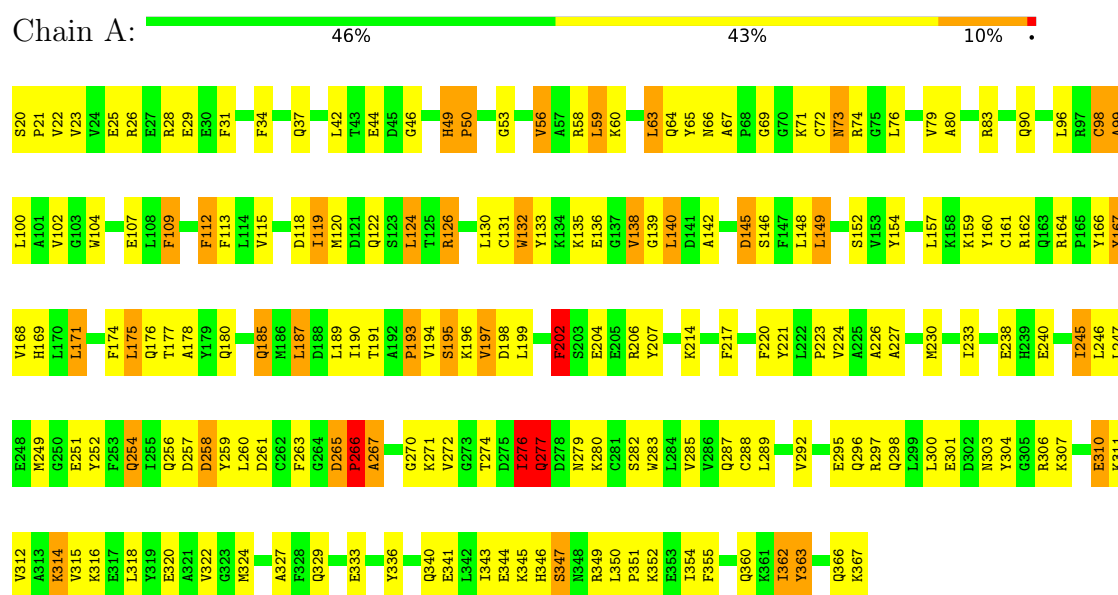
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	79	Total	O	0	0
			79	79		

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: FARNESYL DIPHOSPHATE SYNTHASE



4 Data and refinement statistics

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

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	88.60Å 88.60Å 276.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.60	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.60)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.194 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2894	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

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	0.71	0/2872	1.17	24/3873 (0.6%)

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	A	0	1

There are no bond length outliers.

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	276	ILE	N-CA-C	-8.83	104.55	113.47
1	A	263	PHE	N-CA-C	8.36	122.83	113.21
1	A	146	SER	N-CA-C	-8.34	101.22	111.40
1	A	49	HIS	CA-C-N	7.64	129.39	119.84
1	A	49	HIS	C-N-CA	7.64	129.39	119.84
1	A	66	ASN	N-CA-C	7.16	122.14	113.41
1	A	99	ALA	N-CA-C	-6.96	103.62	111.14
1	A	362	ILE	N-CA-C	6.35	117.45	111.48
1	A	109	PHE	N-CA-C	-5.72	105.13	111.36
1	A	161	CYS	N-CA-C	5.71	121.00	112.94
1	A	271	LYS	N-CA-C	-5.69	105.46	113.20
1	A	67	ALA	N-CA-C	5.65	119.86	112.17
1	A	277	GLN	N-CA-C	-5.64	105.23	111.71
1	A	160	TYR	N-CA-C	5.56	120.18	112.90
1	A	207	TYR	N-CA-C	-5.55	105.14	111.14
1	A	98	CYS	N-CA-C	-5.49	106.08	112.89
1	A	187	LEU	N-CA-C	-5.43	105.36	111.28
1	A	145	ASP	N-CA-C	-5.42	106.67	113.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	53	GLY	N-CA-C	5.41	119.43	112.83
1	A	119	ILE	CB-CA-C	-5.26	105.15	111.88
1	A	148	LEU	N-CA-C	5.24	116.99	111.28
1	A	21	PRO	N-CA-C	-5.17	106.34	113.53
1	A	258	ASP	N-CA-C	-5.17	106.30	113.18
1	A	42	LEU	N-CA-C	5.14	118.81	112.54

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	167	TYR	Sidechain

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	2815	0	2798	144	0
2	A	79	0	0	16	0
All	All	2894	0	2798	144	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 26.

All (144) 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:347:SER:HB2	1:A:352:LYS:HG2	1.47	0.95
1:A:329:GLN:HE22	1:A:333:GLU:HB2	1.41	0.86
1:A:288:CYS:SG	1:A:300:LEU:HD11	2.16	0.85
1:A:254:GLN:HE21	1:A:254:GLN:HA	1.42	0.82
1:A:177:THR:HA	1:A:180:GLN:HE21	1.48	0.78
1:A:185:GLN:HG3	2:A:539:HOH:O	1.84	0.77
1:A:59:LEU:HD22	1:A:63:LEU:HD22	1.66	0.77
1:A:318:LEU:O	1:A:322:VAL:HG23	1.84	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:199:LEU:HD23	1:A:289:LEU:HD21	1.67	0.75
1:A:245:ILE:HD11	1:A:355:PHE:CD1	2.22	0.74
1:A:220:PHE:O	1:A:223:PRO:HD2	1.87	0.73
1:A:327:ALA:HB2	2:A:506:HOH:O	1.87	0.73
1:A:329:GLN:NE2	1:A:333:GLU:HB2	2.02	0.73
1:A:132:TRP:HZ3	1:A:145:ASP:OD2	1.72	0.72
1:A:224:VAL:HG21	1:A:246:LEU:HD11	1.73	0.71
1:A:133:TYR:HB2	1:A:139:GLY:O	1.90	0.71
1:A:226:ALA:O	1:A:230:MET:HG3	1.91	0.71
1:A:311:LYS:O	1:A:315:VAL:HG23	1.90	0.70
1:A:120:MET:SD	2:A:415:HOH:O	2.49	0.70
1:A:60:LYS:HE2	1:A:64:GLN:OE1	1.92	0.69
1:A:245:ILE:HG21	2:A:461:HOH:O	1.93	0.68
1:A:303:ASN:HD22	1:A:314:LYS:HB3	1.60	0.67
1:A:238:GLU:HG3	1:A:349:ARG:HD3	1.76	0.66
1:A:276:ILE:HD12	1:A:316:LYS:HZ1	1.59	0.66
1:A:119:ILE:HD11	1:A:142:ALA:HB3	1.78	0.66
1:A:164:ARG:HG2	1:A:164:ARG:HH11	1.61	0.65
1:A:277:GLN:CD	1:A:307:LYS:HB2	2.23	0.63
1:A:245:ILE:HD11	1:A:355:PHE:CG	2.34	0.62
1:A:246:LEU:HD23	1:A:249:MET:HE3	1.81	0.62
1:A:133:TYR:HD1	2:A:416:HOH:O	1.81	0.62
1:A:260:LEU:HD21	1:A:366:GLN:H	1.64	0.62
1:A:347:SER:CB	1:A:352:LYS:HG2	2.24	0.62
1:A:20:SER:OG	1:A:23:VAL:HG23	1.99	0.61
1:A:154:TYR:CZ	1:A:175:LEU:HD13	2.34	0.61
1:A:276:ILE:HD12	1:A:316:LYS:NZ	2.16	0.61
1:A:277:GLN:OE1	1:A:307:LYS:HB2	2.01	0.60
1:A:221:TYR:HA	1:A:246:LEU:HD13	1.82	0.60
1:A:261:ASP:HA	2:A:544:HOH:O	2.01	0.60
1:A:71:LYS:O	1:A:72:CYS:HB2	2.03	0.59
1:A:197:VAL:HG13	1:A:279:ASN:ND2	2.18	0.59
1:A:65:TYR:CD1	1:A:132:TRP:CD1	2.90	0.58
1:A:280:LYS:HE3	1:A:282:SER:OG	2.05	0.57
1:A:34:PHE:O	1:A:37:GLN:HB3	2.05	0.57
1:A:28:ARG:HG2	1:A:76:LEU:HD11	1.87	0.57
1:A:157:LEU:HD11	1:A:174:PHE:CD2	2.39	0.56
1:A:191:THR:O	1:A:193:PRO:HD3	2.06	0.56
1:A:277:GLN:HE22	1:A:306:ARG:HD3	1.71	0.56
1:A:65:TYR:CD1	1:A:132:TRP:HD1	2.24	0.55
1:A:133:TYR:HA	1:A:138:VAL:HG12	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:300:LEU:HG	1:A:318:LEU:HD21	1.89	0.54
1:A:25:GLU:O	1:A:29:GLU:HG3	2.07	0.54
1:A:79:VAL:HG21	1:A:100:LEU:HD23	1.89	0.54
1:A:346:HIS:HB2	2:A:461:HOH:O	2.08	0.54
1:A:191:THR:OG1	1:A:206:ARG:NH1	2.41	0.54
1:A:336:TYR:O	1:A:340:GLN:HG2	2.08	0.54
1:A:197:VAL:HG13	1:A:279:ASN:CG	2.34	0.53
1:A:270:GLY:C	1:A:272:VAL:H	2.16	0.53
1:A:58:ARG:HH11	1:A:58:ARG:HG2	1.74	0.53
1:A:202:PHE:N	1:A:202:PHE:HD1	2.07	0.53
1:A:49:HIS:CD2	1:A:50:PRO:HD2	2.44	0.53
1:A:177:THR:HA	1:A:180:GLN:NE2	2.19	0.53
1:A:202:PHE:N	1:A:202:PHE:CD1	2.77	0.52
1:A:109:PHE:HZ	1:A:178:ALA:HA	1.75	0.52
1:A:345:LYS:O	1:A:345:LYS:HG2	2.08	0.52
1:A:164:ARG:HG2	1:A:164:ARG:NH1	2.25	0.51
1:A:132:TRP:CZ3	1:A:145:ASP:OD2	2.60	0.51
1:A:300:LEU:O	1:A:304:TYR:HB3	2.11	0.51
1:A:136:GLU:H	1:A:136:GLU:CD	2.18	0.51
1:A:65:TYR:CE1	1:A:132:TRP:HD1	2.29	0.50
1:A:149:LEU:O	1:A:152:SER:HB2	2.11	0.50
1:A:265:ASP:H	1:A:266:PRO:HD2	1.75	0.50
1:A:336:TYR:HB2	1:A:363:TYR:CD2	2.46	0.50
1:A:115:VAL:HA	2:A:473:HOH:O	2.11	0.49
1:A:28:ARG:NH2	1:A:71:LYS:O	2.45	0.49
1:A:191:THR:C	1:A:193:PRO:HD3	2.37	0.49
1:A:196:LYS:O	1:A:197:VAL:HB	2.13	0.49
1:A:349:ARG:HG3	2:A:463:HOH:O	2.12	0.49
1:A:132:TRP:O	1:A:138:VAL:HG11	2.12	0.49
1:A:276:ILE:HG21	1:A:312:VAL:HG13	1.94	0.49
1:A:194:VAL:O	1:A:195:SER:C	2.55	0.49
1:A:300:LEU:HG	1:A:318:LEU:CD2	2.43	0.49
1:A:120:MET:CE	1:A:189:LEU:HD21	2.43	0.48
1:A:220:PHE:C	1:A:223:PRO:HD2	2.37	0.48
1:A:190:ILE:HG22	1:A:191:THR:N	2.28	0.48
1:A:276:ILE:HD11	1:A:315:VAL:HG21	1.95	0.48
1:A:351:PRO:HG2	1:A:354:ILE:CD1	2.43	0.48
1:A:34:PHE:CE1	1:A:37:GLN:HG2	2.48	0.48
1:A:164:ARG:HD3	2:A:485:HOH:O	2.12	0.48
1:A:265:ASP:HB3	1:A:266:PRO:HD3	1.95	0.48
1:A:214:LYS:HD3	2:A:541:HOH:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:TYR:O	1:A:169:HIS:HB2	2.14	0.47
1:A:306:ARG:HD3	1:A:307:LYS:H	1.79	0.47
1:A:22:VAL:HG12	1:A:26:ARG:NH2	2.30	0.47
1:A:217:PHE:CE1	1:A:247:LEU:HD11	2.49	0.47
1:A:351:PRO:HG2	1:A:354:ILE:HD11	1.97	0.47
1:A:135:LYS:HB3	1:A:135:LYS:NZ	2.30	0.47
1:A:112:PHE:CZ	1:A:113:PHE:CE2	3.03	0.46
1:A:46:GLY:CA	1:A:56:VAL:HG11	2.46	0.46
1:A:297:ARG:HD2	1:A:301:GLU:OE2	2.15	0.46
1:A:265:ASP:O	1:A:267:ALA:N	2.48	0.46
1:A:224:VAL:CG2	1:A:246:LEU:HD11	2.45	0.46
1:A:46:GLY:HA3	1:A:56:VAL:HG11	1.96	0.46
1:A:252:TYR:CE1	1:A:362:ILE:HG22	2.51	0.45
1:A:138:VAL:HG12	1:A:139:GLY:N	2.31	0.45
1:A:295:GLU:HG2	1:A:296:GLN:H	1.81	0.45
1:A:276:ILE:O	1:A:276:ILE:HG12	2.15	0.45
1:A:64:GLN:HB3	2:A:469:HOH:O	2.17	0.44
1:A:171:LEU:O	1:A:175:LEU:HD22	2.17	0.44
1:A:98:CYS:O	1:A:102:VAL:HG23	2.17	0.44
1:A:28:ARG:HG2	1:A:76:LEU:CD1	2.47	0.44
1:A:120:MET:HE3	1:A:189:LEU:HD11	1.99	0.44
1:A:245:ILE:HD13	2:A:461:HOH:O	2.17	0.44
1:A:194:VAL:HG13	1:A:195:SER:N	2.33	0.43
1:A:124:LEU:O	1:A:130:LEU:HD12	2.18	0.43
1:A:259:TYR:C	1:A:261:ASP:H	2.26	0.43
1:A:130:LEU:O	1:A:131:CYS:C	2.62	0.43
1:A:224:VAL:O	1:A:227:ALA:N	2.52	0.43
1:A:256:GLN:C	1:A:258:ASP:H	2.27	0.43
1:A:162:ARG:HD3	2:A:532:HOH:O	2.18	0.43
1:A:140:LEU:HD12	1:A:140:LEU:HA	1.50	0.43
1:A:254:GLN:HA	1:A:254:GLN:NE2	2.20	0.43
1:A:343:ILE:HA	2:A:461:HOH:O	2.18	0.43
1:A:136:GLU:C	1:A:138:VAL:H	2.27	0.42
1:A:202:PHE:HD2	1:A:285:VAL:HG11	1.84	0.42
1:A:96:LEU:O	1:A:99:ALA:HB3	2.19	0.42
1:A:80:ALA:HA	1:A:83:ARG:NH1	2.34	0.42
1:A:196:LYS:HA	1:A:196:LYS:HD2	1.83	0.41
1:A:31:PHE:CE1	1:A:104:TRP:CD1	3.08	0.41
1:A:221:TYR:HA	1:A:246:LEU:CD1	2.48	0.41
1:A:72:CYS:O	1:A:73:ASN:C	2.62	0.41
1:A:295:GLU:HG2	1:A:296:GLN:N	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:ILE:HD11	1:A:142:ALA:CB	2.47	0.41
1:A:259:TYR:HD2	1:A:260:LEU:HD23	1.85	0.41
1:A:72:CYS:O	1:A:74:ARG:N	2.53	0.41
1:A:318:LEU:HD12	1:A:322:VAL:HG23	2.02	0.41
1:A:118:ASP:HB2	2:A:473:HOH:O	2.20	0.41
1:A:119:ILE:O	1:A:122:GLN:NE2	2.53	0.41
1:A:28:ARG:HA	1:A:76:LEU:HD21	2.02	0.41
1:A:28:ARG:NH2	1:A:71:LYS:HB3	2.36	0.41
1:A:336:TYR:HB2	1:A:363:TYR:HD2	1.86	0.41
1:A:118:ASP:OD1	1:A:126:ARG:NH1	2.54	0.40
1:A:283:TRP:CE2	1:A:287:GLN:HG3	2.55	0.40
1:A:193:PRO:O	1:A:197:VAL:HA	2.22	0.40
1:A:303:ASN:HD22	1:A:314:LYS:CB	2.31	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	346/348 (99%)	286 (83%)	42 (12%)	18 (5%)	1 2

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	138	VAL
1	A	195	SER
1	A	266	PRO
1	A	274	THR
1	A	44	GLU
1	A	198	ASP
1	A	202	PHE
1	A	310	GLU

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Mol	Chain	Res	Type
1	A	69	GLY
1	A	265	ASP
1	A	73	ASN
1	A	167	TYR
1	A	50	PRO
1	A	233	ILE
1	A	257	ASP
1	A	363	TYR
1	A	193	PRO
1	A	267	ALA

5.3.2 Protein sidechains [i](#)

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	300/300 (100%)	260 (87%)	40 (13%)	4 8

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	VAL
1	A	59	LEU
1	A	63	LEU
1	A	90	GLN
1	A	107	GLU
1	A	112	PHE
1	A	124	LEU
1	A	126	ARG
1	A	132	TRP
1	A	140	LEU
1	A	149	LEU
1	A	159	LYS
1	A	168	VAL
1	A	171	LEU
1	A	175	LEU
1	A	176	GLN

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Mol	Chain	Res	Type
1	A	185	GLN
1	A	187	LEU
1	A	197	VAL
1	A	202	PHE
1	A	204	GLU
1	A	240	GLU
1	A	245	ILE
1	A	251	GLU
1	A	254	GLN
1	A	266	PRO
1	A	276	ILE
1	A	277	GLN
1	A	292	VAL
1	A	298	GLN
1	A	310	GLU
1	A	314	LYS
1	A	320	GLU
1	A	324	MET
1	A	341	GLU
1	A	344	GLU
1	A	347	SER
1	A	350	LEU
1	A	360	GLN
1	A	367	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	GLN
1	A	66	ASN
1	A	122	GLN
1	A	180	GLN
1	A	254	GLN
1	A	290	GLN
1	A	298	GLN
1	A	303	ASN
1	A	329	GLN
1	A	366	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](#)

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.