



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 05:32 PM GMT

PDB ID : 1UBX
Title : STRUCTURE OF FARNESYL PYROPHOSPHATE SYNTHETASE
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Deposited on : 1996-10-14
Resolution : 2.50 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>

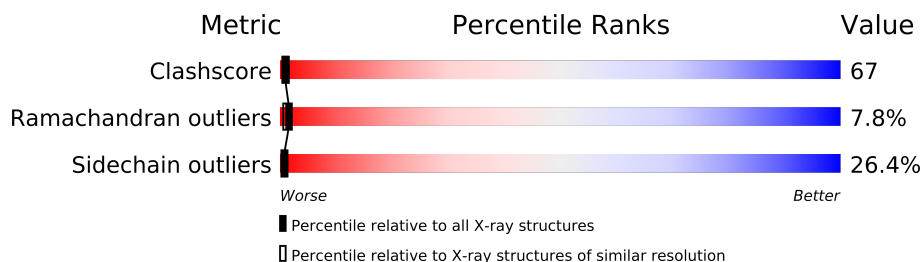
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
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 2.50 Å.

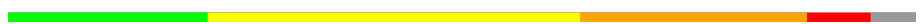
Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	79885	3562 (2.50-2.50)
Ramachandran outliers	78287	3480 (2.50-2.50)
Sidechain outliers	78261	3482 (2.50-2.50)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	367	

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 2874 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 FARNESYL DIPHOSPHATE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	348	Total	C	N	O	S	0	0	0
			2804	1786	477	527	14			

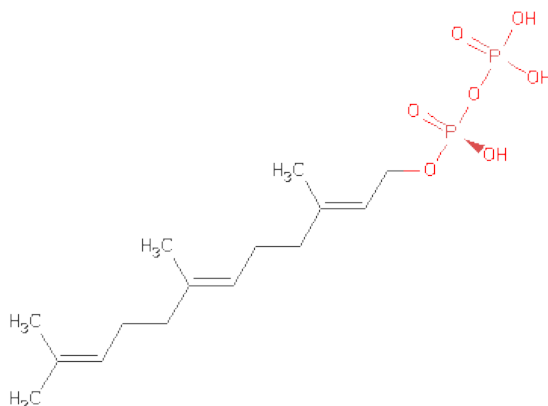
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	112	ALA	PHE	CONFLICT	UNP P08836
A	113	SER	PHE	CONFLICT	UNP P08836

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is FARNESYL DIPHOSPHATE (three-letter code: FPP) (formula: C₁₅H₂₈O₇P₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			24	15	7	2		

- Molecule 4 is water.

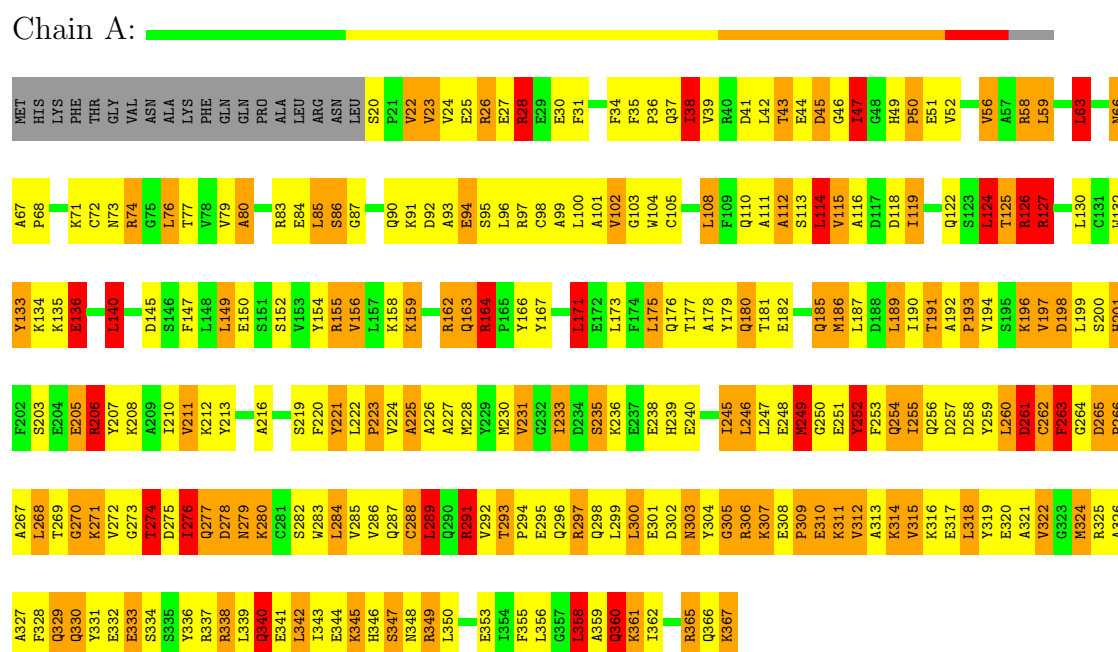
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	45	Total	O	0	0
			45	45		

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.

Note EDS was not executed.

• Molecule 1: FARNESYL DIPHOSPHATE SYNTHASE



4 Data and refinement statistics

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

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	88.70Å 88.70Å 276.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.50	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.50)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT V. 5-E	Depositor
R, R_{free}	(Not available) , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2874	wwPDB-VP
Average B, all atoms (Å ²)	41.0	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: MG, FPP

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	1.20	9/2859 (0.3%)	1.53	41/3856 (1.1%)

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	A	0	16

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	136	GLU	CG-CD	7.32	1.62	1.51
1	A	25	GLU	CG-CD	6.74	1.62	1.51
1	A	176	GLN	CG-CD	6.36	1.65	1.51
1	A	361	LYS	CD-CE	6.12	1.66	1.51
1	A	166	TYR	CD1-CE1	-5.66	1.30	1.39
1	A	156	VAL	CB-CG1	-5.47	1.41	1.52
1	A	205	GLU	CG-CD	5.22	1.59	1.51
1	A	263	PHE	CB-CG	-5.05	1.42	1.51
1	A	25	GLU	CB-CG	5.03	1.61	1.52

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	130	LEU	CB-CG-CD2	-9.77	94.39	111.00
1	A	76	LEU	CB-CG-CD1	-9.53	94.80	111.00
1	A	108	LEU	CB-CG-CD2	-9.41	95.00	111.00
1	A	140	LEU	CA-CB-CG	-9.20	94.13	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	171	LEU	CB-CG-CD1	-8.50	96.54	111.00
1	A	38	ILE	CB-CA-C	-7.98	95.65	111.60
1	A	206	ARG	NE-CZ-NH2	7.97	124.29	120.30
1	A	284	LEU	CB-CG-CD1	7.66	124.02	111.00
1	A	63	LEU	CA-CB-CG	7.40	132.32	115.30
1	A	276	ILE	CG1-CB-CG2	-7.27	95.41	111.40
1	A	246	LEU	CB-CG-CD2	7.23	123.30	111.00
1	A	127	ARG	NE-CZ-NH1	7.05	123.83	120.30
1	A	155	ARG	NE-CZ-NH2	-7.03	116.79	120.30
1	A	342	LEU	CA-CB-CG	6.92	131.21	115.30
1	A	164	ARG	NE-CZ-NH2	-6.90	116.85	120.30
1	A	231	VAL	CB-CA-C	-6.39	99.25	111.40
1	A	270	GLY	N-CA-C	-6.32	97.30	113.10
1	A	118	ASP	CB-CG-OD1	-6.27	112.65	118.30
1	A	288	CYS	CA-CB-SG	6.25	125.26	114.00
1	A	342	LEU	CB-CG-CD1	6.15	121.45	111.00
1	A	155	ARG	NE-CZ-NH1	5.97	123.29	120.30
1	A	263	PHE	CB-CG-CD1	-5.90	116.67	120.80
1	A	280	LYS	CD-CE-NZ	5.84	125.13	111.70
1	A	278	ASP	CB-CG-OD2	5.82	123.54	118.30
1	A	85	LEU	CB-CG-CD2	5.80	120.85	111.00
1	A	247	LEU	CB-CG-CD2	5.78	120.83	111.00
1	A	289	LEU	CA-CB-CG	-5.76	102.05	115.30
1	A	119	ILE	CB-CA-C	-5.75	100.11	111.60
1	A	263	PHE	N-CA-C	5.72	126.45	111.00
1	A	291	ARG	N-CA-C	-5.66	95.71	111.00
1	A	261	ASP	CB-CG-OD1	5.57	123.31	118.30
1	A	358	LEU	CB-CG-CD1	-5.49	101.66	111.00
1	A	66	ASN	N-CA-C	5.41	125.60	111.00
1	A	289	LEU	CB-CG-CD2	-5.34	101.92	111.00
1	A	338	ARG	NE-CZ-NH1	5.28	122.94	120.30
1	A	249	MET	CB-CG-SD	-5.24	96.69	112.40
1	A	124	LEU	CB-CG-CD1	-5.23	102.11	111.00
1	A	47	ILE	CB-CA-C	-5.20	101.21	111.60
1	A	206	ARG	NE-CZ-NH1	-5.17	117.71	120.30
1	A	114	LEU	CB-CG-CD1	-5.14	102.27	111.00
1	A	114	LEU	CB-CG-CD2	-5.05	102.42	111.00

There are no chirality outliers.

All (16) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	102	VAL	Mainchain
1	A	112	ALA	Mainchain
1	A	115	VAL	Mainchain
1	A	126	ARG	Mainchain
1	A	133	TYR	Sidechain
1	A	135	LYS	Mainchain
1	A	149	LEU	Mainchain
1	A	179	TYR	Mainchain
1	A	221	TYR	Sidechain
1	A	225	ALA	Mainchain
1	A	252	TYR	Sidechain
1	A	263	PHE	Sidechain
1	A	28	ARG	Mainchain
1	A	340	GLN	Mainchain
1	A	80	ALA	Mainchain
1	A	98	CYS	Mainchain

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	2804	0	2790	380	0
2	A	1	0	0	0	0
3	A	24	0	25	3	0
4	A	45	0	0	8	0
All	All	2874	0	2815	380	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 67.

All (380) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:299:LEU:HD13	1:A:318:LEU:HD22	1.24	1.14
1:A:298:GLN:HA	1:A:301:GLU:HG2	1.31	1.08
1:A:296:GLN:HB3	1:A:318:LEU:HD21	1.39	1.02
1:A:300:LEU:HD13	1:A:301:GLU:H	1.18	1.02

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:300:LEU:HD13	1:A:301:GLU:N	1.73	1.02
1:A:249:MET:CE	1:A:358:LEU:HB3	1.89	1.01
1:A:366:GLN:HG2	1:A:367:LYS:HE3	1.39	1.00
1:A:275:ASP:HB2	1:A:280:LYS:HZ2	1.29	0.96
1:A:80:ALA:HA	1:A:83:ARG:NH1	1.83	0.94
1:A:249:MET:HE2	1:A:358:LEU:HB3	1.48	0.93
1:A:293:THR:HG23	1:A:296:GLN:HE22	1.37	0.90
1:A:277:GLN:HB2	1:A:307:LYS:HB3	1.57	0.87
1:A:185:GLN:HA	1:A:185:GLN:HE21	1.39	0.87
1:A:332:GLU:OE1	1:A:365:ARG:HG2	1.74	0.87
1:A:191:THR:O	1:A:193:PRO:HD3	1.75	0.86
1:A:298:GLN:HA	1:A:301:GLU:CG	2.05	0.86
1:A:329:GLN:NE2	1:A:365:ARG:NE	2.25	0.85
1:A:205:GLU:HB2	4:A:438:HOH:O	1.76	0.84
1:A:80:ALA:HA	1:A:83:ARG:HH12	1.43	0.83
1:A:329:GLN:HE21	1:A:365:ARG:HE	1.26	0.82
1:A:220:PHE:CE2	1:A:250:GLY:HA2	2.16	0.81
1:A:328:PHE:HZ	1:A:365:ARG:HB2	1.46	0.79
1:A:111:ALA:O	1:A:115:VAL:HG23	1.83	0.79
1:A:312:VAL:O	1:A:315:VAL:HG23	1.83	0.78
1:A:177:THR:O	1:A:181:THR:HG23	1.82	0.78
1:A:249:MET:HE1	1:A:358:LEU:HB3	1.66	0.78
1:A:92:ASP:OD2	1:A:94:GLU:HB3	1.83	0.77
1:A:79:VAL:HG21	1:A:100:LEU:HD23	1.66	0.77
1:A:318:LEU:HA	1:A:321:ALA:HB3	1.66	0.77
1:A:277:GLN:HG3	1:A:278:ASP:N	1.98	0.76
1:A:298:GLN:CA	1:A:301:GLU:HG2	2.15	0.76
1:A:190:ILE:HG22	1:A:191:THR:N	2.01	0.76
1:A:285:VAL:HG22	1:A:319:TYR:OH	1.85	0.75
1:A:292:VAL:CG1	1:A:297:ARG:HB2	2.16	0.75
1:A:328:PHE:CZ	1:A:365:ARG:HB2	2.21	0.75
1:A:296:GLN:HB3	1:A:318:LEU:CD2	2.15	0.75
1:A:252:TYR:CE2	1:A:332:GLU:HG3	2.21	0.75
1:A:318:LEU:O	1:A:322:VAL:HG23	1.87	0.74
1:A:283:TRP:O	1:A:287:GLN:HG2	1.88	0.74
1:A:198:ASP:C	1:A:199:LEU:HD12	2.08	0.74
1:A:330:GLN:HA	1:A:333:GLU:HG2	1.70	0.74
1:A:22:VAL:O	1:A:26:ARG:HG2	1.88	0.73
1:A:317:GLU:O	1:A:321:ALA:N	2.20	0.73
1:A:255:ILE:HG22	1:A:256:GLN:N	2.04	0.73
1:A:38:ILE:O	1:A:41:ASP:HB2	1.88	0.73
1:A:248:GLU:OE1	1:A:338:ARG:HG2	1.88	0.72

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:275:ASP:HB2	1:A:280:LYS:NZ	2.02	0.72
1:A:264:GLY:HA3	1:A:367:LYS:HZ1	1.53	0.72
1:A:226:ALA:O	1:A:230:MET:HG3	1.89	0.72
1:A:314:LYS:O	1:A:317:GLU:HB3	1.90	0.72
1:A:245:ILE:HG13	1:A:246:LEU:N	2.04	0.72
1:A:35:PHE:HB3	1:A:36:PRO:HD3	1.72	0.71
1:A:329:GLN:NE2	1:A:365:ARG:HE	1.87	0.71
1:A:284:LEU:CD2	1:A:324:MET:HG3	2.20	0.70
1:A:276:ILE:N	1:A:280:LYS:HD3	2.06	0.70
1:A:282:SER:O	1:A:286:VAL:HG23	1.90	0.70
1:A:38:ILE:CG2	1:A:42:LEU:HD12	2.22	0.70
1:A:245:ILE:HA	1:A:342:LEU:HD23	1.74	0.69
1:A:274:THR:HG23	1:A:275:ASP:N	2.03	0.69
1:A:200:SER:OG	1:A:201:HIS:N	2.25	0.69
1:A:303:ASN:HD22	1:A:311:LYS:HG2	1.56	0.69
1:A:336:TYR:O	1:A:340:GLN:HG2	1.92	0.69
1:A:293:THR:HG23	1:A:296:GLN:NE2	2.07	0.69
1:A:211:VAL:HG12	1:A:212:LYS:N	2.08	0.69
1:A:295:GLU:HG2	1:A:296:GLN:OE1	1.92	0.69
1:A:245:ILE:HD11	1:A:355:PHE:HB3	1.75	0.69
1:A:278:ASP:C	1:A:280:LYS:H	1.97	0.68
1:A:41:ASP:HA	4:A:432:HOH:O	1.93	0.68
1:A:299:LEU:HD13	1:A:318:LEU:CD2	2.14	0.68
1:A:220:PHE:O	1:A:223:PRO:HD2	1.93	0.68
1:A:196:LYS:O	1:A:198:ASP:N	2.27	0.68
1:A:196:LYS:HG3	1:A:198:ASP:OD2	1.93	0.68
1:A:163:GLN:C	1:A:164:ARG:HG2	2.14	0.67
1:A:365:ARG:HH11	1:A:365:ARG:HG3	1.60	0.67
1:A:311:LYS:O	1:A:314:LYS:N	2.23	0.67
1:A:317:GLU:O	1:A:320:GLU:HB2	1.94	0.67
1:A:203:SER:O	1:A:206:ARG:N	2.27	0.67
1:A:356:LEU:O	1:A:359:ALA:HB3	1.95	0.66
1:A:28:ARG:HE	1:A:73:ASN:HD21	1.44	0.66
1:A:329:GLN:HA	1:A:329:GLN:OE1	1.91	0.66
1:A:303:ASN:HB3	1:A:311:LYS:CB	2.25	0.66
1:A:278:ASP:O	1:A:280:LYS:HG2	1.95	0.65
1:A:259:TYR:OH	1:A:325:ARG:HG2	1.96	0.65
1:A:71:LYS:HE3	4:A:431:HOH:O	1.97	0.65
1:A:341:GLU:HA	1:A:344:GLU:OE2	1.97	0.65
1:A:285:VAL:O	1:A:288:CYS:N	2.30	0.65
1:A:154:TYR:CZ	1:A:175:LEU:HD13	2.32	0.65
1:A:349:ARG:O	1:A:350:LEU:HD23	1.96	0.65

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:296:GLN:O	1:A:299:LEU:N	2.30	0.64
1:A:196:LYS:HE3	1:A:196:LYS:O	1.96	0.64
1:A:329:GLN:HE21	1:A:365:ARG:NE	1.90	0.64
1:A:338:ARG:CG	1:A:342:LEU:HD13	2.27	0.64
1:A:245:ILE:CD1	1:A:355:PHE:HB3	2.28	0.64
1:A:92:ASP:O	1:A:95:SER:N	2.31	0.64
1:A:304:TYR:O	1:A:306:ARG:N	2.31	0.64
1:A:312:VAL:HG12	1:A:313:ALA:N	2.11	0.64
1:A:191:THR:C	1:A:193:PRO:HD3	2.19	0.63
1:A:255:ILE:O	1:A:258:ASP:HB2	1.99	0.63
1:A:59:LEU:O	1:A:59:LEU:HD22	1.99	0.63
1:A:275:ASP:HB3	1:A:280:LYS:CD	2.29	0.63
1:A:125:THR:HG23	4:A:418:HOH:O	1.98	0.63
1:A:252:TYR:HE2	1:A:332:GLU:HG3	1.64	0.63
1:A:186:MET:HE2	1:A:186:MET:HA	1.81	0.63
1:A:173:LEU:O	1:A:177:THR:HG23	1.99	0.62
1:A:185:GLN:CA	1:A:185:GLN:HE21	2.12	0.62
1:A:263:PHE:CE2	1:A:324:MET:HB2	2.34	0.62
1:A:102:VAL:O	1:A:105:CYS:HB2	1.99	0.62
1:A:338:ARG:HD3	4:A:428:HOH:O	1.98	0.62
1:A:303:ASN:HB3	1:A:311:LYS:HB3	1.81	0.62
1:A:294:PRO:HD2	1:A:295:GLU:OE1	2.00	0.62
1:A:342:LEU:O	1:A:346:HIS:HB2	2.00	0.62
1:A:154:TYR:CE1	1:A:175:LEU:HD13	2.33	0.62
1:A:309:PRO:O	1:A:313:ALA:HB3	2.00	0.62
1:A:126:ARG:HG3	1:A:127:ARG:N	2.15	0.62
1:A:147:PHE:O	1:A:150:GLU:HB3	1.99	0.61
1:A:327:ALA:HB1	4:A:424:HOH:O	1.99	0.61
1:A:285:VAL:HG21	1:A:304:TYR:CZ	2.35	0.61
1:A:194:VAL:HA	1:A:197:VAL:HA	1.81	0.61
1:A:293:THR:CG2	1:A:296:GLN:HE22	2.12	0.61
1:A:318:LEU:O	1:A:321:ALA:N	2.33	0.61
1:A:303:ASN:HB3	1:A:311:LYS:CG	2.31	0.61
1:A:74:ARG:O	1:A:77:THR:HB	2.00	0.61
1:A:338:ARG:HG2	1:A:342:LEU:HD13	1.83	0.61
1:A:110:GLN:O	1:A:114:LEU:HG	2.00	0.61
1:A:275:ASP:CB	1:A:280:LYS:HZ2	2.08	0.60
1:A:228:MET:HE2	1:A:239:HIS:CE1	2.36	0.60
1:A:275:ASP:HB3	1:A:280:LYS:HD2	1.81	0.60
1:A:22:VAL:HG12	1:A:23:VAL:N	2.17	0.60
1:A:46:GLY:HA3	1:A:56:VAL:HG21	1.82	0.60
1:A:340:GLN:O	1:A:344:GLU:OE1	2.18	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:366:GLN:O	1:A:367:LYS:NZ	2.32	0.60
1:A:38:ILE:HG22	1:A:42:LEU:HD12	1.82	0.60
1:A:284:LEU:HD22	1:A:324:MET:HG3	1.83	0.59
1:A:296:GLN:OE1	1:A:296:GLN:N	2.21	0.59
1:A:154:TYR:CZ	1:A:175:LEU:CD1	2.85	0.59
1:A:345:LYS:O	1:A:345:LYS:HG2	2.01	0.59
1:A:278:ASP:O	1:A:280:LYS:N	2.33	0.59
1:A:277:GLN:OE1	1:A:307:LYS:HD2	2.03	0.59
1:A:284:LEU:HD22	1:A:324:MET:CG	2.33	0.59
1:A:264:GLY:CA	1:A:367:LYS:HZ1	2.16	0.58
1:A:86:SER:OG	1:A:91:LYS:HE3	2.03	0.58
1:A:274:THR:O	1:A:275:ASP:OD1	2.21	0.58
1:A:101:ALA:O	1:A:103:GLY:N	2.37	0.58
1:A:220:PHE:HE2	1:A:250:GLY:HA2	1.66	0.58
1:A:353:GLU:OE1	1:A:353:GLU:N	2.36	0.58
1:A:26:ARG:HH11	1:A:26:ARG:HG3	1.68	0.58
1:A:43:THR:O	1:A:47:ILE:HG12	2.04	0.57
1:A:35:PHE:O	1:A:39:VAL:HG23	2.04	0.57
1:A:259:TYR:CE2	1:A:263:PHE:HB3	2.40	0.57
1:A:31:PHE:CE1	1:A:104:TRP:CD1	2.92	0.57
1:A:266:PRO:O	1:A:267:ALA:HB3	2.05	0.57
1:A:104:TRP:O	1:A:108:LEU:HG	2.04	0.57
1:A:296:GLN:O	1:A:300:LEU:HD12	2.05	0.56
1:A:66:ASN:HD21	1:A:132:TRP:HB2	1.70	0.56
1:A:292:VAL:HG11	1:A:297:ARG:HB2	1.85	0.56
1:A:26:ARG:NH1	1:A:26:ARG:HG3	2.19	0.56
1:A:255:ILE:O	1:A:258:ASP:N	2.39	0.56
1:A:235:SER:OG	1:A:238:GLU:HB2	2.04	0.56
1:A:311:LYS:O	1:A:315:VAL:HG22	2.05	0.56
1:A:326:ALA:O	1:A:329:GLN:N	2.39	0.56
1:A:233:ILE:HD13	1:A:350:LEU:CD2	2.35	0.56
1:A:279:ASN:OD1	1:A:304:TYR:HD2	1.89	0.56
1:A:366:GLN:C	1:A:367:LYS:HG3	2.25	0.56
1:A:155:ARG:HG2	1:A:155:ARG:HH11	1.71	0.56
1:A:31:PHE:HD1	1:A:100:LEU:HD22	1.71	0.55
1:A:265:ASP:O	1:A:268:LEU:HB2	2.07	0.55
1:A:71:LYS:O	1:A:72:CYS:HB2	2.06	0.55
1:A:270:GLY:CA	1:A:271:LYS:HE3	2.36	0.55
1:A:278:ASP:HB2	1:A:280:LYS:HD2	1.88	0.55
1:A:35:PHE:HB3	1:A:36:PRO:CD	2.36	0.55
1:A:275:ASP:CB	1:A:280:LYS:NZ	2.68	0.55
1:A:318:LEU:O	1:A:322:VAL:N	2.40	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:275:ASP:OD2	1:A:278:ASP:OD2	2.26	0.54
1:A:318:LEU:HD12	1:A:318:LEU:C	2.28	0.54
1:A:224:VAL:HG21	1:A:246:LEU:HD11	1.90	0.54
1:A:220:PHE:CD2	1:A:250:GLY:HA2	2.43	0.54
1:A:259:TYR:CZ	1:A:263:PHE:HB3	2.43	0.54
1:A:297:ARG:O	1:A:301:GLU:HB3	2.08	0.54
1:A:263:PHE:CE2	1:A:324:MET:CB	2.91	0.54
1:A:104:TRP:O	1:A:108:LEU:N	2.37	0.54
1:A:46:GLY:C	1:A:56:VAL:HG21	2.28	0.54
1:A:339:LEU:HD23	1:A:340:GLN:OE1	2.08	0.54
1:A:262:CYS:HB2	1:A:316:LYS:NZ	2.23	0.54
1:A:20:SER:C	1:A:22:VAL:H	2.10	0.54
1:A:248:GLU:HG3	1:A:338:ARG:CZ	2.37	0.54
1:A:332:GLU:CD	1:A:365:ARG:HG2	2.28	0.53
1:A:285:VAL:HG21	1:A:304:TYR:OH	2.08	0.53
1:A:349:ARG:O	1:A:349:ARG:HG2	2.08	0.53
1:A:293:THR:HG1	1:A:295:GLU:HG2	1.74	0.53
1:A:316:LYS:O	1:A:319:TYR:HB2	2.07	0.53
1:A:190:ILE:O	1:A:193:PRO:HG3	2.09	0.53
1:A:262:CYS:SG	1:A:263:PHE:CD1	3.01	0.53
1:A:287:GLN:HB3	1:A:324:MET:CE	2.38	0.53
1:A:277:GLN:HG3	1:A:278:ASP:OD1	2.09	0.53
1:A:365:ARG:HH11	1:A:365:ARG:CG	2.22	0.52
1:A:113:SER:O	1:A:116:ALA:HB3	2.09	0.52
1:A:252:TYR:CZ	1:A:332:GLU:HG3	2.43	0.52
1:A:318:LEU:HD12	1:A:322:VAL:CG2	2.40	0.52
1:A:185:GLN:HA	1:A:185:GLN:NE2	2.19	0.51
1:A:35:PHE:CE2	1:A:39:VAL:HG21	2.44	0.51
1:A:259:TYR:O	1:A:261:ASP:N	2.43	0.51
1:A:259:TYR:CE2	1:A:325:ARG:HG2	2.45	0.51
1:A:328:PHE:HZ	1:A:365:ARG:CB	2.19	0.51
1:A:248:GLU:O	1:A:251:GLU:N	2.43	0.51
1:A:298:GLN:HA	1:A:301:GLU:OE2	2.10	0.51
1:A:127:ARG:N	4:A:417:HOH:O	2.27	0.51
1:A:111:ALA:O	1:A:112:ALA:C	2.48	0.51
1:A:287:GLN:OE1	1:A:291:ARG:NH1	2.43	0.51
1:A:181:THR:HA	1:A:213:TYR:O	2.10	0.51
1:A:46:GLY:CA	1:A:56:VAL:HG21	2.41	0.51
1:A:277:GLN:C	1:A:277:GLN:HE21	2.12	0.51
1:A:233:ILE:HD13	1:A:350:LEU:HD21	1.93	0.51
1:A:293:THR:OG1	1:A:295:GLU:HG2	2.11	0.51
1:A:304:TYR:CD2	1:A:305:GLY:N	2.78	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:252:TYR:OH	1:A:332:GLU:HG3	2.11	0.51
1:A:34:PHE:CD2	1:A:100:LEU:HB3	2.46	0.50
1:A:59:LEU:HD22	1:A:63:LEU:HD22	1.92	0.50
1:A:259:TYR:CZ	1:A:325:ARG:HG2	2.46	0.50
1:A:298:GLN:HA	1:A:301:GLU:CD	2.32	0.50
1:A:333:GLU:HG3	1:A:334:SER:N	2.26	0.50
1:A:366:GLN:CG	1:A:367:LYS:HE3	2.28	0.50
1:A:27:GLU:O	1:A:30:GLU:N	2.45	0.50
1:A:222:LEU:N	1:A:223:PRO:CD	2.75	0.50
1:A:262:CYS:SG	1:A:263:PHE:N	2.84	0.50
1:A:203:SER:O	1:A:206:ARG:HB3	2.12	0.49
1:A:220:PHE:CD2	1:A:250:GLY:CA	2.95	0.49
1:A:259:TYR:C	1:A:261:ASP:H	2.15	0.49
1:A:252:TYR:CD1	1:A:252:TYR:C	2.86	0.49
1:A:318:LEU:HA	1:A:321:ALA:CB	2.41	0.49
1:A:77:THR:HG22	1:A:223:PRO:HB2	1.94	0.49
1:A:97:ARG:O	1:A:100:LEU:HB2	2.13	0.49
1:A:119:ILE:O	1:A:122:GLN:NE2	2.45	0.49
1:A:116:ALA:HB3	3:A:401:FPP:H102	1.94	0.49
1:A:299:LEU:HD22	1:A:318:LEU:HD23	1.95	0.49
1:A:191:THR:HG22	1:A:192:ALA:N	2.28	0.49
1:A:292:VAL:HG13	1:A:293:THR:N	2.27	0.48
1:A:125:THR:HG23	1:A:126:ARG:N	2.28	0.48
1:A:296:GLN:CB	1:A:318:LEU:HD21	2.29	0.48
1:A:122:GLN:HA	1:A:133:TYR:OH	2.12	0.48
1:A:293:THR:CB	1:A:294:PRO:CD	2.91	0.48
1:A:248:GLU:OE1	1:A:342:LEU:HD22	2.13	0.48
1:A:96:LEU:O	1:A:99:ALA:HB3	2.14	0.48
1:A:79:VAL:HG21	1:A:100:LEU:CD2	2.42	0.48
1:A:338:ARG:HG3	1:A:342:LEU:HD13	1.94	0.48
1:A:293:THR:OG1	1:A:295:GLU:OE1	2.32	0.48
1:A:263:PHE:CD2	1:A:325:ARG:HG3	2.49	0.48
1:A:178:ALA:O	1:A:182:GLU:HG3	2.13	0.48
1:A:296:GLN:O	1:A:299:LEU:HB3	2.14	0.47
1:A:35:PHE:N	1:A:36:PRO:HD2	2.28	0.47
1:A:27:GLU:OE2	1:A:83:ARG:NH1	2.47	0.47
1:A:262:CYS:HB2	1:A:316:LYS:HZ2	1.80	0.47
1:A:84:GLU:OE1	1:A:353:GLU:HB2	2.14	0.47
1:A:297:ARG:NH1	1:A:301:GLU:OE1	2.48	0.47
1:A:285:VAL:CG2	1:A:304:TYR:CE1	2.98	0.47
1:A:365:ARG:NH1	1:A:365:ARG:CG	2.78	0.47
1:A:207:TYR:O	1:A:210:ILE:HB	2.15	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:20:SER:O	1:A:24:VAL:HG12	2.15	0.47
1:A:59:LEU:HD13	1:A:63:LEU:HD22	1.97	0.47
1:A:292:VAL:CG1	1:A:293:THR:N	2.76	0.47
1:A:49:HIS:C	1:A:51:GLU:H	2.17	0.47
1:A:270:GLY:HA3	1:A:271:LYS:HE3	1.97	0.47
1:A:119:ILE:O	1:A:119:ILE:HG22	2.15	0.47
1:A:45:ASP:OD1	1:A:45:ASP:N	2.47	0.47
1:A:274:THR:HG23	1:A:275:ASP:HA	1.96	0.46
1:A:266:PRO:N	1:A:366:GLN:HE22	2.11	0.46
1:A:275:ASP:O	1:A:276:ILE:HG13	2.16	0.46
1:A:311:LYS:O	1:A:312:VAL:C	2.53	0.46
1:A:189:LEU:N	1:A:189:LEU:HD13	2.28	0.46
1:A:191:THR:CG2	1:A:192:ALA:N	2.78	0.46
1:A:74:ARG:HH21	1:A:219:SER:CB	2.28	0.46
1:A:196:LYS:HD2	1:A:196:LYS:HA	1.77	0.46
1:A:311:LYS:O	1:A:314:LYS:HB3	2.16	0.46
1:A:259:TYR:HD1	1:A:260:LEU:HD23	1.80	0.46
1:A:329:GLN:NE2	1:A:365:ARG:CZ	2.79	0.46
1:A:182:GLU:HG2	3:A:401:FPP:C12	2.45	0.46
1:A:329:GLN:C	1:A:331:TYR:H	2.19	0.46
1:A:206:ARG:HG3	1:A:207:TYR:N	2.29	0.46
1:A:340:GLN:O	1:A:343:ILE:N	2.41	0.46
1:A:49:HIS:HB3	1:A:52:VAL:HG22	1.97	0.46
1:A:304:TYR:CG	1:A:305:GLY:N	2.84	0.45
1:A:264:GLY:CA	1:A:367:LYS:NZ	2.79	0.45
1:A:111:ALA:C	1:A:115:VAL:HG23	2.34	0.45
1:A:276:ILE:HG21	1:A:315:VAL:HG21	1.98	0.45
1:A:171:LEU:HD12	1:A:171:LEU:HA	1.49	0.45
1:A:310:GLU:O	1:A:314:LYS:HB3	2.15	0.45
1:A:224:VAL:O	1:A:227:ALA:HB3	2.16	0.45
1:A:158:LYS:O	1:A:162:ARG:HB2	2.17	0.45
1:A:289:LEU:O	1:A:289:LEU:HD23	2.15	0.45
1:A:287:GLN:HA	1:A:291:ARG:NH1	2.32	0.45
1:A:339:LEU:O	1:A:340:GLN:C	2.54	0.45
1:A:275:ASP:HB2	1:A:280:LYS:CE	2.47	0.44
1:A:211:VAL:CG1	1:A:212:LYS:N	2.78	0.44
1:A:63:LEU:HD13	1:A:149:LEU:CD2	2.48	0.44
1:A:101:ALA:C	1:A:103:GLY:N	2.69	0.44
1:A:49:HIS:CD2	1:A:50:PRO:HD2	2.51	0.44
1:A:308:GLU:HA	1:A:309:PRO:HD2	1.44	0.44
1:A:86:SER:HA	1:A:90:GLN:OE1	2.17	0.44
1:A:67:ALA:N	1:A:68:PRO:CD	2.79	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:360:GLN:HG2	1:A:360:GLN:O	2.17	0.44
1:A:327:ALA:O	1:A:331:TYR:N	2.51	0.44
1:A:329:GLN:O	1:A:333:GLU:HB3	2.18	0.44
1:A:303:ASN:OD1	1:A:303:ASN:N	2.50	0.44
1:A:189:LEU:HD12	1:A:189:LEU:HA	1.33	0.44
1:A:124:LEU:HD13	1:A:124:LEU:HA	1.68	0.44
1:A:198:ASP:HB3	1:A:200:SER:OG	2.18	0.44
1:A:343:ILE:O	1:A:347:SER:HB2	2.18	0.44
1:A:140:LEU:HA	1:A:140:LEU:HD12	0.89	0.44
1:A:285:VAL:O	1:A:288:CYS:HB3	2.18	0.43
1:A:186:MET:HA	1:A:186:MET:CE	2.46	0.43
1:A:279:ASN:OD1	1:A:304:TYR:CD2	2.69	0.43
1:A:208:LYS:HA	1:A:211:VAL:HB	1.99	0.43
1:A:134:LYS:HA	1:A:134:LYS:HD3	1.65	0.43
1:A:318:LEU:HD12	1:A:322:VAL:HG22	2.00	0.43
1:A:154:TYR:OH	1:A:175:LEU:HD12	2.18	0.43
1:A:275:ASP:C	1:A:280:LYS:HD3	2.38	0.43
1:A:303:ASN:HB3	1:A:311:LYS:HG2	2.01	0.43
1:A:228:MET:CE	1:A:239:HIS:CE1	3.01	0.43
1:A:187:LEU:HA	1:A:187:LEU:HD23	1.72	0.43
1:A:304:TYR:C	1:A:306:ARG:H	2.21	0.43
1:A:259:TYR:CE2	1:A:263:PHE:CB	3.02	0.43
1:A:108:LEU:HD23	1:A:108:LEU:HA	1.66	0.43
1:A:149:LEU:O	1:A:152:SER:HB2	2.19	0.43
1:A:177:THR:HA	1:A:180:GLN:HG3	2.00	0.43
1:A:274:THR:HG23	1:A:275:ASP:CA	2.48	0.43
1:A:278:ASP:C	1:A:280:LYS:N	2.68	0.43
1:A:299:LEU:HD23	1:A:299:LEU:C	2.39	0.43
1:A:332:GLU:OE2	1:A:365:ARG:HB3	2.19	0.42
1:A:58:ARG:NH1	1:A:145:ASP:OD2	2.53	0.42
1:A:276:ILE:O	1:A:305:GLY:HA2	2.20	0.42
1:A:116:ALA:HB3	3:A:401:FPP:C10	2.49	0.42
1:A:318:LEU:O	1:A:321:ALA:HB3	2.19	0.42
1:A:56:VAL:O	1:A:59:LEU:HB3	2.19	0.42
1:A:300:LEU:C	1:A:302:ASP:N	2.72	0.42
1:A:266:PRO:O	1:A:267:ALA:CB	2.68	0.42
1:A:339:LEU:O	1:A:342:LEU:HB2	2.19	0.42
1:A:344:GLU:N	1:A:344:GLU:OE1	2.51	0.42
1:A:127:ARG:O	1:A:127:ARG:HG3	2.20	0.42
1:A:96:LEU:HA	1:A:96:LEU:HD23	1.71	0.42
1:A:276:ILE:HG21	1:A:276:ILE:HD13	1.45	0.42
1:A:186:MET:CA	1:A:186:MET:CE	2.96	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:262:CYS:HB2	1:A:316:LYS:HE3	2.01	0.42
1:A:292:VAL:HG13	1:A:297:ARG:HB2	1.99	0.42
1:A:312:VAL:C	1:A:314:LYS:H	2.23	0.42
1:A:331:TYR:O	1:A:334:SER:N	2.53	0.42
1:A:348:ASN:OD1	1:A:348:ASN:O	2.38	0.42
1:A:329:GLN:C	1:A:331:TYR:N	2.73	0.42
1:A:206:ARG:HB2	1:A:206:ARG:HE	1.66	0.42
1:A:273:GLY:O	1:A:274:THR:HB	2.20	0.41
1:A:275:ASP:CB	1:A:280:LYS:CE	2.97	0.41
1:A:318:LEU:O	1:A:322:VAL:CG2	2.65	0.41
1:A:191:THR:OG1	1:A:206:ARG:NH1	2.53	0.41
1:A:186:MET:HE3	1:A:186:MET:HB2	1.47	0.41
1:A:149:LEU:HD23	1:A:149:LEU:HA	1.89	0.41
1:A:342:LEU:O	1:A:346:HIS:CD2	2.74	0.41
1:A:101:ALA:O	1:A:102:VAL:C	2.57	0.41
1:A:254:GLN:HA	1:A:254:GLN:HE21	1.84	0.41
1:A:291:ARG:O	1:A:322:VAL:CG1	2.69	0.41
1:A:207:TYR:O	1:A:210:ILE:N	2.54	0.41
1:A:222:LEU:N	1:A:223:PRO:HD3	2.35	0.41
1:A:349:ARG:O	1:A:349:ARG:CG	2.68	0.41
1:A:87:GLY:O	1:A:90:GLN:N	2.53	0.41
1:A:49:HIS:HA	1:A:50:PRO:HD2	1.86	0.41
1:A:159:LYS:O	1:A:159:LYS:HG3	2.21	0.41
1:A:126:ARG:HG2	1:A:127:ARG:HG2	2.03	0.41
1:A:317:GLU:C	1:A:320:GLU:HB2	2.40	0.41
1:A:299:LEU:O	1:A:303:ASN:OD1	2.39	0.41
1:A:47:ILE:N	1:A:56:VAL:HG21	2.36	0.41
1:A:221:TYR:CE1	1:A:225:ALA:HB2	2.56	0.41
1:A:283:TRP:NE1	1:A:287:GLN:HG3	2.36	0.40
1:A:92:ASP:CG	1:A:95:SER:H	2.25	0.40
1:A:42:LEU:C	1:A:43:THR:HG23	2.41	0.40
1:A:248:GLU:HG3	1:A:338:ARG:NH1	2.36	0.40
1:A:336:TYR:CE1	1:A:340:GLN:OE1	2.74	0.40
1:A:253:PHE:C	1:A:253:PHE:CD1	2.94	0.40
1:A:26:ARG:H	1:A:26:ARG:HG2	1.68	0.40
1:A:276:ILE:HG13	1:A:316:LYS:NZ	2.37	0.40
1:A:212:LYS:O	1:A:216:ALA:HB3	2.22	0.40
1:A:348:ASN:ND2	4:A:447:HOH:O	2.50	0.40
1:A:156:VAL:O	1:A:156:VAL:HG12	2.21	0.40
1:A:85:LEU:HB3	1:A:231:VAL:HG11	2.03	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	346/367 (94%)	253 (73%)	66 (19%)	27 (8%)	1 1

All (27) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	127	ARG
1	A	193	PRO
1	A	198	ASP
1	A	276	ILE
1	A	279	ASN
1	A	309	PRO
1	A	349	ARG
1	A	136	GLU
1	A	162	ARG
1	A	197	VAL
1	A	260	LEU
1	A	265	ASP
1	A	305	GLY
1	A	311	LYS
1	A	360	GLN
1	A	47	ILE
1	A	167	TYR
1	A	223	PRO
1	A	233	ILE
1	A	266	PRO
1	A	274	THR
1	A	22	VAL
1	A	43	THR
1	A	289	LEU
1	A	50	PRO
1	A	93	ALA
1	A	345	LYS

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	299/315 (95%)	220 (74%)	79 (26%)	1 1

All (79) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	VAL
1	A	26	ARG
1	A	28	ARG
1	A	37	GLN
1	A	38	ILE
1	A	44	GLU
1	A	45	ASP
1	A	56	VAL
1	A	58	ARG
1	A	59	LEU
1	A	63	LEU
1	A	74	ARG
1	A	76	LEU
1	A	86	SER
1	A	94	GLU
1	A	114	LEU
1	A	124	LEU
1	A	125	THR
1	A	126	ARG
1	A	136	GLU
1	A	140	LEU
1	A	159	LYS
1	A	163	GLN
1	A	164	ARG
1	A	171	LEU
1	A	175	LEU
1	A	180	GLN
1	A	185	GLN
1	A	186	MET
1	A	189	LEU
1	A	191	THR

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Mol	Chain	Res	Type
1	A	196	LYS
1	A	201	HIS
1	A	206	ARG
1	A	211	VAL
1	A	235	SER
1	A	236	LYS
1	A	240	GLU
1	A	245	ILE
1	A	249	MET
1	A	252	TYR
1	A	254	GLN
1	A	255	ILE
1	A	257	ASP
1	A	261	ASP
1	A	262	CYS
1	A	268	LEU
1	A	269	THR
1	A	271	LYS
1	A	272	VAL
1	A	274	THR
1	A	277	GLN
1	A	289	LEU
1	A	291	ARG
1	A	293	THR
1	A	297	ARG
1	A	300	LEU
1	A	303	ASN
1	A	306	ARG
1	A	307	LYS
1	A	310	GLU
1	A	312	VAL
1	A	314	LYS
1	A	315	VAL
1	A	318	LEU
1	A	322	VAL
1	A	324	MET
1	A	329	GLN
1	A	330	GLN
1	A	333	GLU
1	A	337	ARG
1	A	340	GLN
1	A	347	SER

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Mol	Chain	Res	Type
1	A	358	LEU
1	A	360	GLN
1	A	361	LYS
1	A	362	ILE
1	A	365	ARG
1	A	367	LYS

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

Mol	Chain	Res	Type
1	A	66	ASN
1	A	73	ASN
1	A	185	GLN
1	A	239	HIS
1	A	254	GLN
1	A	329	GLN
1	A	330	GLN
1	A	346	HIS
1	A	348	ASN
1	A	360	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 ⓘ

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	$\# Z > 2$
3	FPP	A	401	-	23,23,23	1.40	3 (13%)	31,31,31	1.81	6 (19%)

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
3	FPP	A	401	-	-	0/25/25/25	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	401	FPP	PA-O3A	-3.18	1.54	1.59
3	A	401	FPP	PB-O3A	-2.88	1.55	1.60
3	A	401	FPP	C7-C8	2.13	1.37	1.32

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	FPP	PA-O3A-PB	-4.65	118.03	131.68
3	A	401	FPP	C4-C3-C5	3.23	120.30	115.39
3	A	401	FPP	O3A-PA-O1	3.23	117.83	103.41
3	A	401	FPP	O3B-PB-O3A	3.18	120.24	105.14
3	A	401	FPP	C10-C8-C9	3.03	120.00	115.39
3	A	401	FPP	O3B-PB-O1B	-2.80	101.30	110.44

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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.