



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 12:46 PM GMT

PDB ID : 1HM4
Title : N219L PENTALENENE SYNTHASE
Authors : Seemann, M.; Paschall, C.M.; Christianson, D.W.; Cane, D.E.
Deposited on : 2000-12-04
Resolution : 3.47 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>

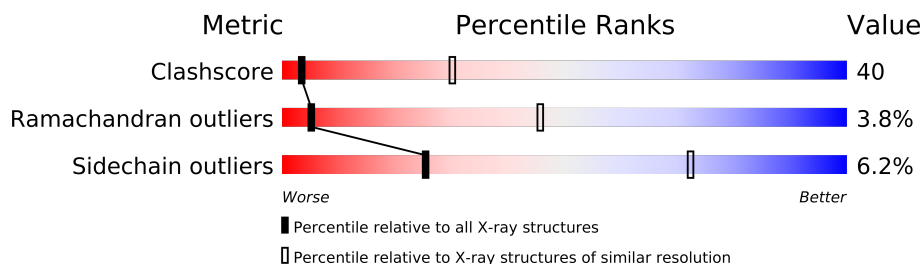
The following versions of software and data (see [references](#)) were used in the production of this report:

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 3.47 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	1001 (3.62-3.34)
Ramachandran outliers	78287	1001 (3.64-3.32)
Sidechain outliers	78261	1002 (3.64-3.32)

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	336	
1	B	336	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 4683 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 PENTALENENE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	289	Total	C	N	O	S	0	0	0
			2293	1433	418	431	11			
1	B	300	Total	C	N	O	S	0	0	0
			2390	1496	436	446	12			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	219	LEU	ASN	ENGINEERED	UNP Q55012
B	219	LEU	ASN	ENGINEERED	UNP Q55012

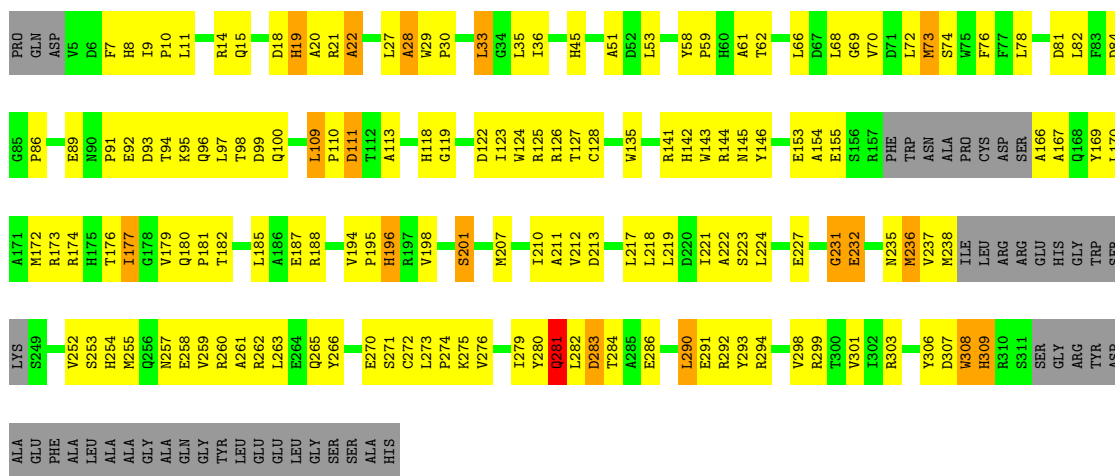
3 Residue-property plots

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.

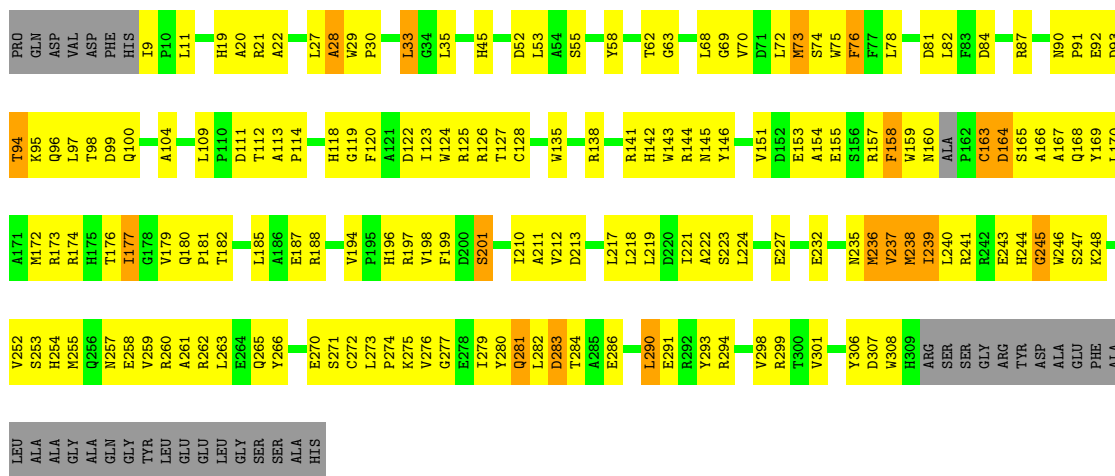
• Molecule 1: PENTALENENE SYNTHASE

Chain A:



• Molecule 1: PENTALENENE SYNTHASE

Chain B:



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, α , β , γ	182.00Å 182.00Å 56.40Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 3.47	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-3.47)	Depositor
R_{merge}	0.16	Depositor
R_{sym}	0.16	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.225 , 0.299	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4683	wwPDB-VP
Average B, all atoms (Å ²)	39.0	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	0.41	0/2346	0.62	0/3183
1	B	0.39	0/2449	0.61	0/3324
All	All	0.40	0/4795	0.61	0/6507

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2293	0	2189	159	0
1	B	2390	0	2282	203	0
All	All	4683	0	4471	362	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 40.

All (362) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:69:GLY:HA2	1:A:185:LEU:HD21	1.40	1.02
1:B:127:THR:HB	1:B:188:ARG:HD3	1.43	1.01

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:127:THR:HB	1:A:188:ARG:HD3	1.44	0.96
1:A:9:ILE:HA	1:A:263:LEU:HD13	1.59	0.82
1:A:109:LEU:N	1:A:109:LEU:HD12	1.95	0.82
1:B:160:ASN:HA	1:B:163:CYS:SG	2.21	0.81
1:B:69:GLY:HA2	1:B:185:LEU:HD21	1.59	0.81
1:A:109:LEU:H	1:A:109:LEU:HD12	1.46	0.81
1:B:277:GLY:HA2	1:B:282:LEU:HD23	1.61	0.80
1:B:244:HIS:HB2	1:B:246:TRP:CD1	2.17	0.79
1:A:258:GLU:O	1:A:262:ARG:HG2	1.84	0.78
1:A:283:ASP:HB3	1:A:286:GLU:HG3	1.65	0.77
1:B:53:LEU:HB2	1:B:308:TRP:HZ2	1.48	0.77
1:A:185:LEU:HD13	1:A:188:ARG:HD2	1.65	0.77
1:A:45:HIS:HE1	1:A:74:SER:OG	1.66	0.77
1:B:258:GLU:O	1:B:262:ARG:HG2	1.85	0.76
1:A:235:ASN:HD21	1:A:237:VAL:HG23	1.49	0.76
1:B:35:LEU:HD13	1:B:78:LEU:HD13	1.68	0.75
1:B:198:VAL:O	1:B:201:SER:HB3	1.87	0.75
1:B:185:LEU:HD13	1:B:188:ARG:HD2	1.68	0.75
1:B:223:SER:O	1:B:227:GLU:HG3	1.86	0.74
1:A:123:ILE:O	1:A:127:THR:HG23	1.88	0.74
1:A:194:VAL:HG12	1:A:195:PRO:HD2	1.66	0.74
1:A:33:LEU:HB3	1:A:35:LEU:HD13	1.68	0.74
1:B:194:VAL:HG21	1:B:293:TYR:CG	2.22	0.73
1:A:33:LEU:HD12	1:A:119:GLY:HA2	1.69	0.73
1:B:283:ASP:HB3	1:B:286:GLU:HG3	1.71	0.73
1:A:95:LYS:HA	1:A:98:THR:HB	1.71	0.72
1:A:223:SER:O	1:A:227:GLU:HG3	1.89	0.72
1:B:270:GLU:O	1:B:273:LEU:HB2	1.90	0.72
1:A:185:LEU:HD12	1:A:188:ARG:HB2	1.69	0.72
1:A:170:LEU:O	1:A:174:ARG:HB2	1.89	0.72
1:B:170:LEU:O	1:B:174:ARG:HB2	1.87	0.72
1:B:185:LEU:HD12	1:B:188:ARG:HB2	1.72	0.71
1:B:123:ILE:O	1:B:127:THR:HG23	1.90	0.71
1:A:270:GLU:O	1:A:273:LEU:HB2	1.90	0.71
1:B:157:ARG:HG3	1:B:158:PHE:N	2.06	0.71
1:B:95:LYS:HA	1:B:98:THR:HB	1.71	0.71
1:B:104:ALA:CB	1:B:109:LEU:HD21	2.20	0.71
1:A:33:LEU:HB3	1:A:35:LEU:CD1	2.21	0.70
1:B:261:ALA:O	1:B:265:GLN:HG3	1.91	0.70
1:B:213:ASP:O	1:B:217:LEU:HG	1.92	0.69
1:B:160:ASN:C	1:B:163:CYS:H	1.95	0.69
1:B:237:VAL:HG22	1:B:255:MET:HE1	1.73	0.69

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:280:TYR:HB2	1:B:282:LEU:HD21	1.74	0.69
1:A:45:HIS:CE1	1:A:74:SER:OG	2.46	0.68
1:A:144:ARG:HH11	1:A:144:ARG:HG3	1.57	0.68
1:B:141:ARG:O	1:B:144:ARG:HB2	1.93	0.68
1:B:235:ASN:O	1:B:238:MET:HG2	1.93	0.68
1:B:144:ARG:HH11	1:B:144:ARG:HG3	1.57	0.67
1:B:87:ARG:O	1:B:94:THR:HG22	1.95	0.67
1:A:213:ASP:O	1:A:217:LEU:HG	1.95	0.67
1:A:221:ILE:HD12	1:A:221:ILE:C	2.14	0.67
1:A:33:LEU:CB	1:A:35:LEU:HD13	2.24	0.67
1:B:235:ASN:O	1:B:237:VAL:N	2.28	0.67
1:B:122:ASP:O	1:B:126:ARG:HD3	1.95	0.66
1:B:104:ALA:HB2	1:B:109:LEU:HD21	1.78	0.66
1:A:298:VAL:O	1:A:301:VAL:HG12	1.95	0.66
1:A:218:LEU:O	1:A:221:ILE:HG13	1.96	0.66
1:A:261:ALA:O	1:A:265:GLN:HG3	1.95	0.65
1:A:271:SER:O	1:A:274:PRO:HD2	1.97	0.65
1:A:141:ARG:O	1:A:144:ARG:HB2	1.96	0.65
1:B:271:SER:O	1:B:274:PRO:HD2	1.97	0.64
1:B:235:ASN:HD22	1:B:237:VAL:HG23	1.63	0.64
1:B:158:PHE:H	1:B:158:PHE:HD2	1.45	0.64
1:B:255:MET:O	1:B:259:VAL:HG23	1.97	0.64
1:B:72:LEU:HA	1:B:123:ILE:HD13	1.80	0.64
1:A:185:LEU:CD1	1:A:188:ARG:HD2	2.27	0.64
1:A:53:LEU:HD23	1:A:73:MET:HG3	1.80	0.64
1:A:122:ASP:O	1:A:126:ARG:HD3	1.98	0.64
1:A:235:ASN:ND2	1:A:237:VAL:HG23	2.13	0.64
1:B:167:ALA:HA	1:B:170:LEU:HD13	1.79	0.63
1:B:81:ASP:HA	1:B:84:ASP:OD2	1.98	0.63
1:A:255:MET:O	1:A:259:VAL:HG23	1.97	0.63
1:B:277:GLY:O	1:B:282:LEU:HD23	1.97	0.63
1:B:163:CYS:O	1:B:164:ASP:HB2	1.98	0.63
1:A:135:TRP:HH2	1:A:187:GLU:HB2	1.64	0.63
1:B:58:TYR:OH	1:B:185:LEU:HD23	1.99	0.63
1:A:273:LEU:HD23	1:A:294:ARG:HD3	1.79	0.63
1:B:280:TYR:HB2	1:B:282:LEU:CD2	2.29	0.62
1:A:135:TRP:CH2	1:A:187:GLU:HB2	2.34	0.62
1:B:21:ARG:NH1	1:B:63:GLY:HA2	2.14	0.62
1:B:218:LEU:O	1:B:221:ILE:HG13	1.99	0.62
1:A:81:ASP:HA	1:A:84:ASP:OD2	1.98	0.62
1:B:254:HIS:HA	1:B:257:ASN:HB2	1.81	0.62
1:B:185:LEU:CD1	1:B:188:ARG:HD2	2.29	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:221:ILE:C	1:B:221:ILE:HD12	2.20	0.62
1:B:53:LEU:HB2	1:B:308:TRP:CZ2	2.32	0.62
1:B:53:LEU:HD23	1:B:73:MET:HG3	1.82	0.62
1:B:235:ASN:ND2	1:B:237:VAL:HG23	2.14	0.61
1:B:298:VAL:O	1:B:301:VAL:HG12	1.99	0.61
1:B:19:HIS:CG	1:B:20:ALA:N	2.67	0.61
1:B:306:TYR:O	1:B:308:TRP:N	2.31	0.61
1:B:237:VAL:HG22	1:B:255:MET:CE	2.29	0.60
1:A:290:LEU:HD23	1:A:290:LEU:C	2.21	0.60
1:A:19:HIS:HA	1:A:22:ALA:HB3	1.82	0.60
1:A:198:VAL:O	1:A:201:SER:HB3	2.01	0.60
1:B:224:LEU:HD22	1:B:237:VAL:HG21	1.82	0.60
1:A:290:LEU:HD23	1:A:291:GLU:N	2.16	0.60
1:A:68:LEU:HD21	1:A:127:THR:HG22	1.84	0.59
1:B:277:GLY:CA	1:B:282:LEU:HD23	2.31	0.59
1:B:198:VAL:HG11	1:B:290:LEU:HD12	1.84	0.59
1:A:167:ALA:HA	1:A:170:LEU:HD13	1.84	0.59
1:B:35:LEU:CD1	1:B:78:LEU:HD13	2.30	0.59
1:B:290:LEU:HD23	1:B:291:GLU:N	2.18	0.59
1:A:254:HIS:HA	1:A:257:ASN:HB2	1.83	0.59
1:A:91:PRO:HD3	1:A:154:ALA:O	2.03	0.59
1:B:170:LEU:N	1:B:170:LEU:HD12	2.18	0.59
1:B:104:ALA:HA	1:B:109:LEU:HD21	1.85	0.59
1:B:33:LEU:HD12	1:B:119:GLY:HA2	1.85	0.58
1:A:8:HIS:HD1	1:A:10:PRO:HD3	1.68	0.58
1:A:9:ILE:CD1	1:A:303:ARG:HA	2.34	0.58
1:B:104:ALA:CA	1:B:109:LEU:HD21	2.33	0.58
1:A:221:ILE:HD12	1:A:222:ALA:N	2.19	0.58
1:B:194:VAL:HG22	1:B:293:TYR:HB2	1.85	0.58
1:B:280:TYR:O	1:B:281:GLN:C	2.42	0.57
1:A:283:ASP:HB3	1:A:286:GLU:CG	2.33	0.57
1:B:290:LEU:HD23	1:B:290:LEU:C	2.25	0.57
1:A:18:ASP:O	1:A:20:ALA:N	2.37	0.57
1:A:180:GLN:HB3	1:A:181:PRO:HD3	1.85	0.57
1:B:275:LYS:O	1:B:279:ILE:HG13	2.05	0.57
1:B:165:SER:HB2	1:B:168:GLN:HE21	1.70	0.56
1:A:185:LEU:HA	1:A:188:ARG:CG	2.36	0.56
1:B:160:ASN:HA	1:B:163:CYS:HG	1.68	0.56
1:B:227:GLU:HB3	1:B:232:GLU:HB3	1.87	0.56
1:B:180:GLN:HB3	1:B:181:PRO:HD3	1.88	0.56
1:B:244:HIS:HB2	1:B:246:TRP:NE1	2.21	0.56
1:A:9:ILE:HG23	1:A:9:ILE:O	2.06	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:185:LEU:HA	1:A:188:ARG:HG3	1.87	0.55
1:B:235:ASN:O	1:B:236:MET:C	2.45	0.55
1:B:185:LEU:HA	1:B:188:ARG:CG	2.36	0.55
1:B:33:LEU:CD1	1:B:119:GLY:HA2	2.36	0.55
1:A:29:TRP:HB3	1:A:30:PRO:CD	2.36	0.55
1:B:96:GLN:O	1:B:100:GLN:HG3	2.07	0.55
1:B:164:ASP:OD1	1:B:172:MET:SD	2.65	0.54
1:A:170:LEU:HD12	1:A:170:LEU:N	2.22	0.54
1:B:194:VAL:HG21	1:B:293:TYR:CD1	2.43	0.54
1:B:283:ASP:HB3	1:B:286:GLU:CG	2.38	0.54
1:B:177:ILE:HD13	1:B:177:ILE:O	2.07	0.54
1:B:185:LEU:HA	1:B:188:ARG:HG3	1.90	0.54
1:A:15:GLN:NE2	1:A:19:HIS:CD2	2.75	0.54
1:B:236:MET:HG3	1:B:240:LEU:CD1	2.37	0.54
1:B:29:TRP:HB3	1:B:30:PRO:CD	2.38	0.54
1:A:283:ASP:CB	1:A:286:GLU:HG3	2.37	0.53
1:B:185:LEU:HA	1:B:188:ARG:CD	2.37	0.53
1:A:10:PRO:O	1:A:11:LEU:HD23	2.08	0.53
1:A:96:GLN:O	1:A:100:GLN:HG3	2.07	0.53
1:A:9:ILE:HD13	1:A:303:ARG:HA	1.91	0.53
1:A:109:LEU:CD1	1:A:109:LEU:N	2.68	0.53
1:B:53:LEU:HD23	1:B:73:MET:CG	2.38	0.53
1:B:221:ILE:HD12	1:B:222:ALA:N	2.23	0.53
1:B:33:LEU:HD12	1:B:119:GLY:CA	2.38	0.53
1:A:135:TRP:CH2	1:A:187:GLU:CB	2.91	0.53
1:A:166:ALA:O	1:A:169:TYR:HB3	2.09	0.53
1:A:185:LEU:HA	1:A:188:ARG:CD	2.39	0.53
1:B:237:VAL:O	1:B:241:ARG:HG3	2.09	0.53
1:A:35:LEU:HD23	1:A:78:LEU:HD13	1.90	0.53
1:B:157:ARG:CG	1:B:158:PHE:N	2.72	0.53
1:A:231:GLY:O	1:A:232:GLU:C	2.47	0.52
1:B:241:ARG:HG2	1:B:247:SER:O	2.09	0.52
1:B:45:HIS:HE1	1:B:74:SER:OG	1.92	0.52
1:A:82:LEU:HD23	1:A:97:LEU:HD21	1.90	0.52
1:A:51:ALA:HA	1:A:70:VAL:HG13	1.91	0.52
1:B:236:MET:HG3	1:B:240:LEU:HD11	1.91	0.52
1:A:280:TYR:O	1:A:282:LEU:N	2.43	0.52
1:A:113:ALA:HB3	1:A:118:HIS:NE2	2.25	0.52
1:A:275:LYS:O	1:A:279:ILE:HG13	2.10	0.52
1:B:166:ALA:O	1:B:169:TYR:HB3	2.10	0.52
1:A:15:GLN:NE2	1:A:19:HIS:NE2	2.58	0.52
1:A:135:TRP:CZ3	1:A:187:GLU:CD	2.83	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:144:ARG:NH1	1:B:144:ARG:HG3	2.24	0.51
1:A:196:HIS:N	1:A:196:HIS:ND1	2.59	0.51
1:B:11:LEU:HD22	1:B:299:ARG:CZ	2.41	0.51
1:B:273:LEU:N	1:B:274:PRO:HD2	2.26	0.51
1:B:243:GLU:OE1	1:B:244:HIS:NE2	2.44	0.51
1:A:153:GLU:C	1:A:155:GLU:H	2.13	0.51
1:A:177:ILE:HD13	1:A:177:ILE:O	2.09	0.51
1:A:7:PHE:HD2	1:A:306:TYR:CE1	2.28	0.51
1:B:160:ASN:O	1:B:163:CYS:N	2.32	0.51
1:B:45:HIS:CE1	1:B:74:SER:OG	2.64	0.51
1:B:177:ILE:HG23	1:B:179:VAL:HG23	1.93	0.50
1:B:163:CYS:O	1:B:164:ASP:CB	2.59	0.50
1:B:9:ILE:HG22	1:B:11:LEU:H	1.76	0.50
1:B:170:LEU:HB3	1:B:174:ARG:HD3	1.92	0.50
1:B:68:LEU:HD21	1:B:127:THR:HG22	1.92	0.50
1:A:273:LEU:N	1:A:274:PRO:HD2	2.26	0.50
1:A:187:GLU:OE2	1:A:194:VAL:HG23	2.12	0.50
1:A:53:LEU:HD23	1:A:73:MET:CG	2.40	0.50
1:B:124:TRP:O	1:B:125:ARG:C	2.48	0.50
1:A:254:HIS:O	1:A:258:GLU:HB2	2.11	0.49
1:B:170:LEU:HD12	1:B:170:LEU:H	1.77	0.49
1:B:273:LEU:HD23	1:B:294:ARG:HD3	1.94	0.49
1:B:219:LEU:C	1:B:219:LEU:HD13	2.32	0.49
1:B:254:HIS:O	1:B:258:GLU:HB2	2.12	0.49
1:B:166:ALA:HB3	1:B:243:GLU:OE2	2.12	0.49
1:B:170:LEU:H	1:B:170:LEU:CD1	2.25	0.49
1:B:142:HIS:HA	1:B:145:ASN:HD22	1.77	0.49
1:A:124:TRP:O	1:A:125:ARG:C	2.51	0.49
1:B:166:ALA:HB2	1:B:243:GLU:CD	2.33	0.49
1:A:306:TYR:C	1:A:308:TRP:H	2.16	0.49
1:B:153:GLU:C	1:B:155:GLU:H	2.15	0.49
1:B:52:ASP:O	1:B:55:SER:OG	2.31	0.49
1:B:262:ARG:HH11	1:B:262:ARG:HG3	1.77	0.49
1:B:244:HIS:O	1:B:245:GLY:C	2.51	0.49
1:A:272:CYS:O	1:A:276:VAL:HG23	2.13	0.49
1:B:272:CYS:O	1:B:276:VAL:HG23	2.12	0.49
1:A:69:GLY:CA	1:A:185:LEU:HD21	2.29	0.48
1:A:11:LEU:HD22	1:A:299:ARG:CZ	2.42	0.48
1:A:19:HIS:O	1:A:20:ALA:C	2.52	0.48
1:A:177:ILE:HG23	1:A:179:VAL:HG23	1.96	0.48
1:A:144:ARG:HG3	1:A:144:ARG:NH1	2.26	0.48
1:A:282:LEU:HD12	1:A:286:GLU:OE1	2.13	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:252:VAL:O	1:A:255:MET:HB2	2.14	0.48
1:A:219:LEU:HD13	1:A:219:LEU:C	2.34	0.48
1:A:170:LEU:H	1:A:170:LEU:CD1	2.26	0.47
1:A:135:TRP:CH2	1:A:187:GLU:CD	2.88	0.47
1:A:170:LEU:HB3	1:A:174:ARG:HD3	1.96	0.47
1:A:169:TYR:O	1:A:172:MET:HB2	2.13	0.47
1:B:68:LEU:HA	1:B:126:ARG:NH2	2.30	0.47
1:B:163:CYS:SG	1:B:164:ASP:N	2.88	0.47
1:B:169:TYR:O	1:B:172:MET:HB2	2.15	0.47
1:A:194:VAL:CG1	1:A:195:PRO:HD2	2.41	0.47
1:A:173:ARG:NH2	1:A:232:GLU:OE1	2.48	0.47
1:B:238:MET:O	1:B:239:ILE:C	2.52	0.47
1:B:169:TYR:CD2	1:B:239:ILE:HD12	2.50	0.47
1:A:170:LEU:HD12	1:A:170:LEU:H	1.80	0.47
1:A:14:ARG:HB3	1:A:59:PRO:HB3	1.96	0.47
1:A:58:TYR:OH	1:A:185:LEU:HD23	2.15	0.47
1:B:221:ILE:HG22	1:B:255:MET:HE1	1.97	0.47
1:B:213:ASP:HB3	1:B:262:ARG:NH1	2.30	0.47
1:B:252:VAL:O	1:B:255:MET:HB2	2.15	0.46
1:A:142:HIS:HA	1:A:145:ASN:HD22	1.80	0.46
1:B:138:ARG:HH21	1:B:196:HIS:HD2	1.63	0.46
1:B:263:LEU:O	1:B:266:TYR:HB3	2.15	0.46
1:B:19:HIS:O	1:B:22:ALA:HB3	2.15	0.46
1:B:138:ARG:HH21	1:B:196:HIS:CD2	2.33	0.46
1:B:127:THR:HG21	1:B:188:ARG:NH1	2.30	0.46
1:B:283:ASP:CB	1:B:286:GLU:HG3	2.43	0.46
1:A:273:LEU:CD2	1:A:294:ARG:HD3	2.46	0.46
1:B:238:MET:O	1:B:241:ARG:N	2.49	0.46
1:A:281:GLN:HB3	1:A:281:GLN:HE21	1.58	0.46
1:A:257:ASN:O	1:A:260:ARG:HB3	2.16	0.45
1:A:93:ASP:O	1:A:96:GLN:HB2	2.17	0.45
1:B:70:VAL:O	1:B:74:SER:HB2	2.16	0.45
1:A:292:ARG:HG3	1:A:292:ARG:HH11	1.82	0.45
1:A:9:ILE:CA	1:A:263:LEU:HD13	2.38	0.45
1:A:9:ILE:O	1:A:11:LEU:N	2.46	0.45
1:A:70:VAL:O	1:A:74:SER:HB2	2.17	0.45
1:B:273:LEU:O	1:B:276:VAL:HB	2.16	0.45
1:B:210:ILE:O	1:B:211:ALA:C	2.54	0.45
1:B:169:TYR:CD2	1:B:236:MET:HE1	2.52	0.45
1:B:166:ALA:HB2	1:B:243:GLU:OE1	2.16	0.45
1:B:27:LEU:O	1:B:28:ALA:C	2.55	0.45
1:B:166:ALA:H	1:B:239:ILE:HG22	1.81	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:151:VAL:O	1:B:154:ALA:N	2.50	0.45
1:A:8:HIS:ND1	1:A:10:PRO:HD3	2.32	0.45
1:A:93:ASP:HA	1:A:96:GLN:HE21	1.81	0.45
1:A:8:HIS:O	1:A:263:LEU:HD13	2.17	0.44
1:B:158:PHE:CD2	1:B:158:PHE:N	2.80	0.44
1:A:18:ASP:O	1:A:21:ARG:N	2.49	0.44
1:B:160:ASN:CA	1:B:163:CYS:H	2.30	0.44
1:A:170:LEU:HD11	1:A:236:MET:CE	2.47	0.44
1:B:93:ASP:O	1:B:96:GLN:HB2	2.17	0.44
1:A:173:ARG:HA	1:A:176:THR:OG1	2.17	0.44
1:A:143:TRP:O	1:A:146:TYR:HB3	2.18	0.44
1:A:262:ARG:HH11	1:A:262:ARG:HG3	1.82	0.44
1:A:210:ILE:O	1:A:211:ALA:C	2.55	0.44
1:B:135:TRP:CH2	1:B:187:GLU:HB2	2.52	0.44
1:B:90:ASN:HA	1:B:91:PRO:HD2	1.86	0.44
1:A:182:THR:O	1:A:185:LEU:HB3	2.17	0.44
1:B:75:TRP:CZ2	1:B:120:PHE:HB2	2.52	0.44
1:A:280:TYR:O	1:A:281:GLN:C	2.56	0.44
1:B:170:LEU:N	1:B:170:LEU:CD1	2.79	0.44
1:A:153:GLU:C	1:A:155:GLU:N	2.70	0.44
1:B:153:GLU:C	1:B:155:GLU:N	2.71	0.44
1:B:227:GLU:O	1:B:232:GLU:HB3	2.18	0.43
1:B:91:PRO:O	1:B:92:GLU:C	2.55	0.43
1:A:253:SER:C	1:A:255:MET:H	2.22	0.43
1:A:110:PRO:O	1:A:111:ASP:C	2.55	0.43
1:B:185:LEU:HA	1:B:188:ARG:HD2	2.01	0.43
1:A:27:LEU:O	1:A:28:ALA:C	2.55	0.43
1:B:143:TRP:O	1:B:146:TYR:HB3	2.17	0.43
1:B:283:ASP:CG	1:B:284:THR:N	2.72	0.43
1:A:91:PRO:O	1:A:92:GLU:C	2.56	0.43
1:A:27:LEU:O	1:A:30:PRO:HG2	2.19	0.43
1:B:127:THR:HB	1:B:188:ARG:CD	2.31	0.43
1:A:135:TRP:HH2	1:A:187:GLU:CB	2.30	0.43
1:A:86:PRO:O	1:A:89:GLU:HB2	2.19	0.43
1:B:221:ILE:HG22	1:B:255:MET:CE	2.48	0.43
1:B:213:ASP:HB3	1:B:262:ARG:HH12	1.84	0.43
1:B:194:VAL:CG2	1:B:293:TYR:HB2	2.49	0.43
1:B:111:ASP:HA	1:B:118:HIS:CE1	2.53	0.43
1:A:207:MET:HB3	1:A:207:MET:HE2	1.83	0.43
1:B:253:SER:C	1:B:255:MET:H	2.22	0.43
1:A:35:LEU:N	1:A:35:LEU:CD1	2.82	0.43
1:B:21:ARG:HH12	1:B:63:GLY:HA2	1.84	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:93:ASP:HA	1:B:96:GLN:HE21	1.84	0.42
1:B:69:GLY:CA	1:B:185:LEU:HD21	2.38	0.42
1:B:169:TYR:HD2	1:B:236:MET:HE1	1.84	0.42
1:A:135:TRP:HZ3	1:A:187:GLU:OE2	2.01	0.42
1:B:266:TYR:CE2	1:B:270:GLU:OE1	2.72	0.42
1:B:127:THR:CG2	1:B:188:ARG:HH11	2.32	0.42
1:B:276:VAL:HA	1:B:279:ILE:HD12	2.01	0.42
1:B:62:THR:O	1:B:62:THR:HG23	2.19	0.42
1:A:276:VAL:HA	1:A:279:ILE:HD12	2.01	0.42
1:A:170:LEU:N	1:A:170:LEU:CD1	2.82	0.42
1:A:179:VAL:O	1:A:182:THR:HB	2.20	0.42
1:A:8:HIS:O	1:A:263:LEU:CD1	2.67	0.42
1:B:173:ARG:HA	1:B:176:THR:OG1	2.19	0.42
1:B:273:LEU:HD12	1:B:273:LEU:HA	1.87	0.42
1:B:165:SER:CB	1:B:168:GLN:HE21	2.33	0.42
1:A:263:LEU:O	1:A:266:TYR:HB3	2.19	0.42
1:B:244:HIS:O	1:B:246:TRP:N	2.52	0.42
1:B:198:VAL:O	1:B:199:PHE:C	2.58	0.42
1:A:306:TYR:O	1:A:308:TRP:N	2.52	0.42
1:A:266:TYR:CE2	1:A:270:GLU:OE1	2.72	0.42
1:A:273:LEU:O	1:A:276:VAL:HB	2.19	0.42
1:B:160:ASN:HA	1:B:163:CYS:CB	2.49	0.42
1:A:14:ARG:HB2	1:A:59:PRO:HG2	2.02	0.42
1:B:238:MET:H	1:B:238:MET:HG2	1.44	0.42
1:B:235:ASN:HD22	1:B:237:VAL:CG2	2.31	0.42
1:A:194:VAL:HG22	1:A:293:TYR:CG	2.55	0.42
1:A:33:LEU:HD12	1:A:119:GLY:CA	2.46	0.42
1:A:35:LEU:HD12	1:A:35:LEU:N	2.35	0.42
1:B:194:VAL:CG2	1:B:293:TYR:CG	2.99	0.42
1:B:135:TRP:CZ3	1:B:187:GLU:OE1	2.73	0.42
1:B:82:LEU:HD23	1:B:97:LEU:HD21	2.01	0.42
1:A:224:LEU:HD22	1:A:237:VAL:HG21	2.02	0.42
1:A:30:PRO:HB2	1:A:36:ILE:HG12	2.00	0.42
1:B:69:GLY:HA2	1:B:185:LEU:CD2	2.41	0.41
1:B:146:TYR:HE1	1:B:176:THR:O	2.02	0.41
1:A:72:LEU:HA	1:A:123:ILE:HD13	2.00	0.41
1:B:27:LEU:O	1:B:30:PRO:HG2	2.19	0.41
1:A:111:ASP:O	1:A:113:ALA:N	2.53	0.41
1:B:113:ALA:HB1	1:B:114:PRO:HD2	2.03	0.41
1:B:169:TYR:OH	1:B:173:ARG:HD2	2.20	0.41
1:B:11:LEU:HD22	1:B:299:ARG:NE	2.35	0.41
1:A:282:LEU:HD11	1:A:286:GLU:HB3	2.01	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:62:THR:O	1:A:62:THR:HG23	2.20	0.41
1:B:72:LEU:HD11	1:B:76:PHE:CE1	2.56	0.41
1:B:58:TYR:CZ	1:B:185:LEU:HD23	2.55	0.41
1:B:127:THR:HG21	1:B:188:ARG:HH11	1.84	0.41
1:B:257:ASN:O	1:B:260:ARG:HB3	2.20	0.41
1:B:165:SER:N	1:B:168:GLN:NE2	2.68	0.41
1:B:236:MET:O	1:B:237:VAL:C	2.58	0.41
1:B:283:ASP:CG	1:B:284:THR:H	2.24	0.41
1:A:61:ALA:O	1:A:66:LEU:HB2	2.21	0.41
1:A:308:TRP:O	1:A:309:HIS:O	2.39	0.41
1:B:277:GLY:O	1:B:282:LEU:CD2	2.67	0.41
1:B:9:ILE:HG13	1:B:263:LEU:CD1	2.50	0.41
1:B:182:THR:O	1:B:185:LEU:HB3	2.21	0.41
1:A:135:TRP:CZ3	1:A:187:GLU:OE2	2.74	0.41
1:A:33:LEU:CD1	1:A:119:GLY:HA2	2.47	0.41
1:B:9:ILE:HG13	1:B:263:LEU:HD13	2.04	0.40
1:B:165:SER:HB2	1:B:168:GLN:NE2	2.34	0.40
1:A:169:TYR:OH	1:A:173:ARG:HD2	2.20	0.40
1:B:75:TRP:HB2	1:B:123:ILE:HD11	2.04	0.40
1:B:179:VAL:O	1:B:182:THR:HB	2.22	0.40
1:B:236:MET:HB2	1:B:236:MET:HE2	1.84	0.40
1:B:238:MET:O	1:B:240:LEU:N	2.54	0.40
1:A:283:ASP:CG	1:A:284:THR:N	2.75	0.40
1:B:141:ARG:HG2	1:B:145:ASN:HD21	1.87	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	283/336 (84%)	228 (81%)	44 (16%)	11 (4%)	5	45
1	B	296/336 (88%)	236 (80%)	49 (17%)	11 (4%)	5	47
All	All	579/672 (86%)	464 (80%)	93 (16%)	22 (4%)	5	46

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	281	GLN
1	A	309	HIS
1	B	164	ASP
1	B	236	MET
1	B	307	ASP
1	A	19	HIS
1	A	232	GLU
1	B	159	TRP
1	B	283	ASP
1	A	283	ASP
1	A	307	ASP
1	A	22	ALA
1	A	236	MET
1	B	237	VAL
1	B	239	ILE
1	B	281	GLN
1	A	28	ALA
1	A	231	GLY
1	B	28	ALA
1	B	245	GLY
1	B	212	VAL
1	A	212	VAL

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	235/272 (86%)	220 (94%)	15 (6%)	25	71
1	B	245/272 (90%)	230 (94%)	15 (6%)	26	73
All	All	480/544 (88%)	450 (94%)	30 (6%)	25	72

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	LEU
1	A	73	MET

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Mol	Chain	Res	Type
1	A	76	PHE
1	A	94	THR
1	A	99	ASP
1	A	109	LEU
1	A	111	ASP
1	A	128	CYS
1	A	177	ILE
1	A	196	HIS
1	A	201	SER
1	A	238	MET
1	A	281	GLN
1	A	290	LEU
1	A	308	TRP
1	B	33	LEU
1	B	73	MET
1	B	76	PHE
1	B	94	THR
1	B	99	ASP
1	B	112	THR
1	B	128	CYS
1	B	158	PHE
1	B	163	CYS
1	B	177	ILE
1	B	197	ARG
1	B	201	SER
1	B	238	MET
1	B	248	LYS
1	B	290	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
1	A	26	GLN
1	A	45	HIS
1	A	96	GLN
1	A	145	ASN
1	A	168	GLN
1	A	215	ASN
1	A	234	ASN
1	A	235	ASN
1	A	281	GLN

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Mol	Chain	Res	Type
1	B	15	GLN
1	B	26	GLN
1	B	45	HIS
1	B	96	GLN
1	B	145	ASN
1	B	160	ASN
1	B	168	GLN
1	B	209	GLN
1	B	215	ASN

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 ⓘ

There are no ligands in this entry.

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.