



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:34 PM UTC

PDB ID : 3ASB / pdb_00003asb
Title : Crystal structure of PLP-bound LL-diaminopimelate aminotransferase from *Chlamydia trachomatis*
Authors : Watanabe, N.; James, M.N.
Deposited on : 2010-12-10
Resolution : 2.70 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

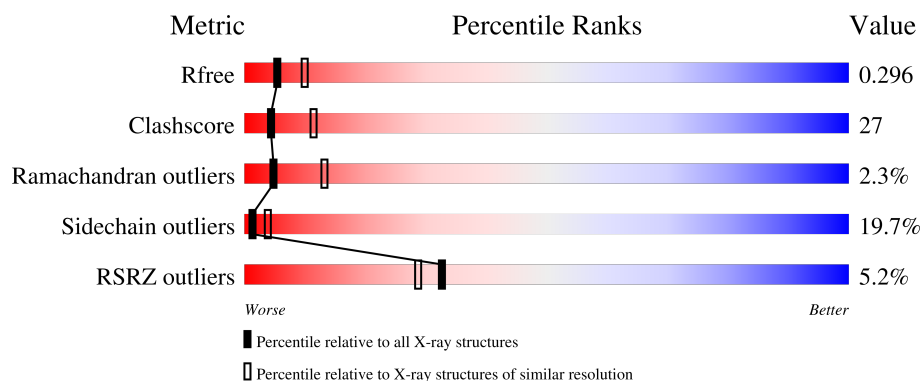
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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))
R_{free}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	400	5% 44% 35% 11% 9%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2926 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 LL-diaminopimelate aminotransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	364	Total	C	N	O	P	S	0	0	0
			2887	1871	480	528	1	7			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	395	HIS	-	expression tag	UNP O84395
A	396	HIS	-	expression tag	UNP O84395
A	397	HIS	-	expression tag	UNP O84395
A	398	HIS	-	expression tag	UNP O84395
A	399	HIS	-	expression tag	UNP O84395
A	400	HIS	-	expression tag	UNP O84395

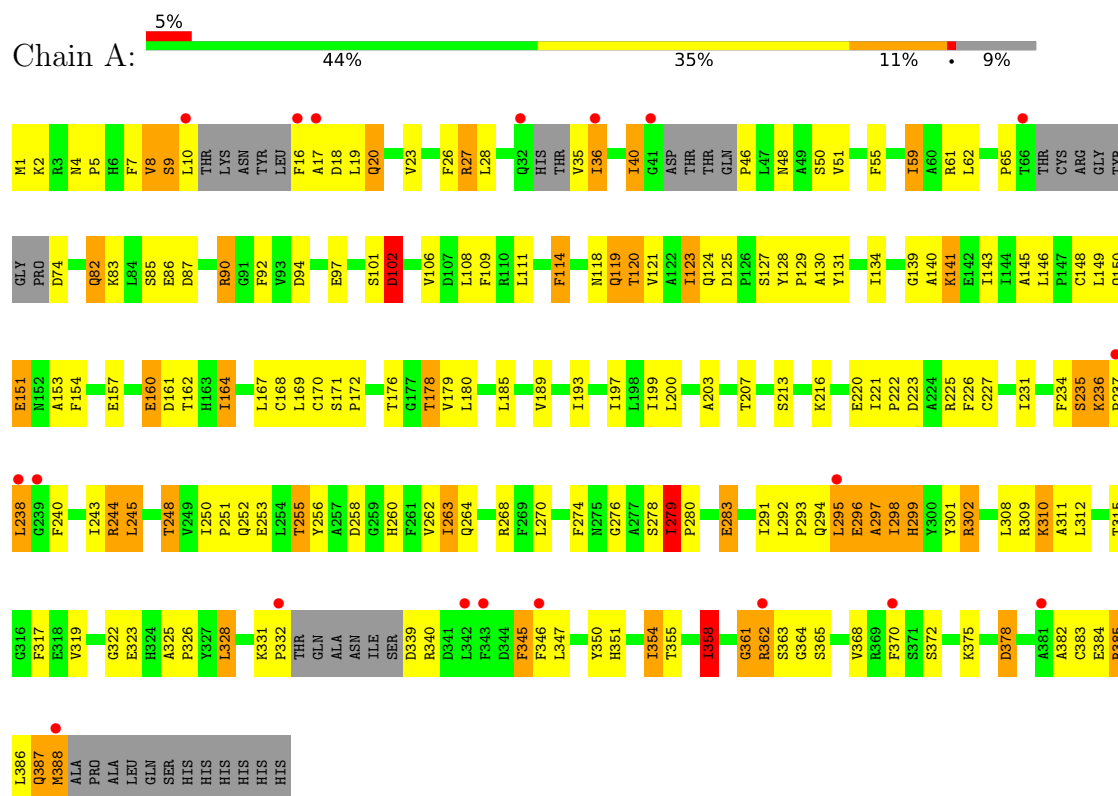
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	39	Total	O	0	0
			39	39		

3 Residue-property plots [i](#)

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 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.

- Molecule 1: LL-diaminopimelate aminotransferase



4 Data and refinement statistics

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	110.49Å 110.49Å 204.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.70 40.00 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.1 (40.00-2.70) 99.0 (40.00-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.70 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.5.0055	Depositor
R, R_{free}	0.222 , 0.287 0.223 , 0.296	Depositor DCC
R_{free} test set	900 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	73.1	Xtriage
Anisotropy	0.186	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 56.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2926	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.12% 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

Bond lengths and bond angles in the following residue types are not validated in this section: LLP

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	1.17	4/2934 (0.1%)	1.31	22/3975 (0.6%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	203	ALA	CA-CB	-7.37	1.43	1.53
1	A	170	CYS	N-CA	-6.90	1.38	1.46
1	A	221	ILE	CA-C	-5.22	1.49	1.52
1	A	86	GLU	N-CA	-5.11	1.40	1.46

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	361	GLY	N-CA-C	8.04	120.38	111.85
1	A	226	PHE	N-CA-C	-6.19	105.00	113.30
1	A	276	GLY	N-CA-C	6.01	120.46	112.65
1	A	102	ASP	N-CA-C	5.70	119.86	113.02
1	A	151	GLU	N-CA-C	-5.67	105.17	111.36
1	A	213	SER	N-CA-C	5.57	117.15	111.14
1	A	296	GLU	N-CA-C	-5.46	105.72	112.38
1	A	179	VAL	CB-CA-C	-5.41	103.08	110.77
1	A	36	ILE	N-CA-C	5.41	115.37	108.12
1	A	279	ILE	CA-C-N	-5.40	113.34	119.28
1	A	279	ILE	C-N-CA	-5.40	113.34	119.28
1	A	82	GLN	N-CA-C	5.39	116.84	110.97
1	A	227	CYS	N-CA-C	5.38	121.64	112.99
1	A	245	LEU	N-CA-C	5.33	116.73	107.93
1	A	120	THR	CB-CA-C	5.29	118.87	109.72
1	A	65	PRO	CA-C-N	5.19	131.05	121.70
1	A	65	PRO	C-N-CA	5.19	131.05	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	199	ILE	CB-CA-C	-5.19	103.10	110.83
1	A	119	GLN	CA-C-N	-5.17	115.87	123.00
1	A	119	GLN	C-N-CA	-5.17	115.87	123.00
1	A	250	ILE	CB-CA-C	-5.13	105.28	111.18
1	A	322	GLY	N-CA-C	5.05	122.25	115.59

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	2887	0	2834	153	0
2	A	39	0	0	3	0
All	All	2926	0	2834	153	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 27.

All (153) 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:358:ILE:H	1:A:358:ILE:HD13	1.17	1.10
1:A:17:ALA:HA	1:A:18:ASP:HB3	1.15	1.09
1:A:358:ILE:HD12	1:A:365:SER:HA	1.42	1.01
1:A:354:ILE:HG21	1:A:370:PHE:HE1	1.23	1.00
1:A:46:PRO:HD3	2:A:415:HOH:O	1.64	0.97
1:A:17:ALA:HA	1:A:18:ASP:CB	1.96	0.94
1:A:354:ILE:HG21	1:A:370:PHE:CE1	2.03	0.92
1:A:4:ASN:HB3	1:A:7:PHE:HB2	1.53	0.89
1:A:236:LLP:H5'1	1:A:236:LLP:HE3	1.53	0.89
1:A:1:MET:HE3	1:A:2:LYS:O	1.76	0.86
1:A:46:PRO:CD	2:A:415:HOH:O	2.20	0.86
1:A:123:ILE:HD12	1:A:124:GLN:H	1.38	0.85
1:A:358:ILE:H	1:A:358:ILE:CD1	1.83	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:GLN:O	1:A:23:VAL:HG12	1.78	0.84
1:A:354:ILE:HG12	1:A:368:VAL:HG21	1.58	0.84
1:A:345:PHE:C	1:A:345:PHE:CD1	2.56	0.83
1:A:358:ILE:HD13	1:A:358:ILE:N	1.95	0.82
1:A:235:SER:O	1:A:238:LEU:HD13	1.80	0.82
1:A:94:ASP:HB3	1:A:97:GLU:HG3	1.61	0.80
1:A:296:GLU:HA	1:A:299:HIS:HB2	1.64	0.80
1:A:311:ALA:HB1	1:A:386:LEU:HD12	1.64	0.79
1:A:17:ALA:CA	1:A:18:ASP:HB3	2.07	0.78
1:A:141:LYS:NZ	1:A:141:LYS:HB3	1.99	0.78
1:A:234:PHE:O	1:A:237:PRO:HD3	1.84	0.77
1:A:27:ARG:HH11	1:A:27:ARG:HG3	1.48	0.76
1:A:83:LYS:HE2	1:A:87:ASP:OD2	1.86	0.76
1:A:350:TYR:HD1	1:A:385:ARG:O	1.69	0.76
1:A:141:LYS:HB3	1:A:141:LYS:HZ3	1.53	0.73
1:A:236:LLP:HE3	1:A:236:LLP:C5'	2.19	0.71
1:A:27:ARG:NH2	1:A:35:VAL:O	2.24	0.71
1:A:118:ASN:HA	1:A:141:LYS:HG2	1.74	0.69
1:A:27:ARG:HG3	1:A:27:ARG:NH1	2.06	0.68
1:A:354:ILE:CG2	1:A:370:PHE:CE1	2.76	0.68
1:A:354:ILE:HG12	1:A:368:VAL:CG2	2.23	0.67
1:A:345:PHE:C	1:A:345:PHE:HD1	2.00	0.67
1:A:234:PHE:O	1:A:236:LLP:N	2.29	0.65
1:A:176:THR:OG1	1:A:178:THR:HB	1.96	0.65
1:A:48:ASN:HD22	1:A:51:VAL:H	1.45	0.65
1:A:292:LEU:HB3	1:A:293:PRO:HD3	1.79	0.65
1:A:295:LEU:HB3	1:A:297:ALA:HB3	1.79	0.65
1:A:354:ILE:C	1:A:354:ILE:CD1	2.71	0.64
1:A:240:PHE:HD2	1:A:243:ILE:HD12	1.63	0.63
1:A:296:GLU:C	1:A:298:ILE:H	2.06	0.63
1:A:5:PRO:C	1:A:7:PHE:H	2.06	0.63
1:A:97:GLU:OE2	1:A:252:GLN:HG2	2.01	0.61
1:A:384:GLU:O	1:A:386:LEU:N	2.28	0.61
1:A:350:TYR:CD1	1:A:385:ARG:O	2.54	0.61
1:A:279:ILE:HG12	1:A:280:PRO:HD3	1.84	0.60
1:A:387:GLN:H	1:A:388:MET:HE3	1.67	0.60
1:A:61:ARG:NH1	1:A:283:GLU:OE2	2.35	0.60
1:A:296:GLU:N	1:A:296:GLU:OE1	2.35	0.59
1:A:310:LYS:O	1:A:310:LYS:HD2	2.03	0.59
1:A:236:LLP:HA	1:A:238:LEU:HD22	1.85	0.59
1:A:118:ASN:HD22	1:A:141:LYS:HG2	1.69	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:164:ILE:HD11	1:A:197:ILE:CG2	2.34	0.57
1:A:207:THR:HG22	1:A:207:THR:O	2.03	0.57
1:A:164:ILE:HD11	1:A:197:ILE:HG21	1.87	0.56
1:A:296:GLU:C	1:A:298:ILE:N	2.64	0.56
1:A:311:ALA:O	1:A:315:THR:HG22	2.05	0.56
1:A:240:PHE:CD2	1:A:243:ILE:HD12	2.39	0.56
1:A:278:SER:OG	1:A:280:PRO:HD2	2.05	0.56
1:A:237:PRO:N	1:A:238:LEU:HA	2.20	0.56
1:A:384:GLU:C	1:A:386:LEU:H	2.14	0.55
1:A:354:ILE:C	1:A:354:ILE:HD12	2.31	0.55
1:A:185:LEU:O	1:A:189:VAL:HG23	2.07	0.55
1:A:234:PHE:CE2	1:A:245:LEU:HD23	2.43	0.54
1:A:361:GLY:O	1:A:364:GLY:N	2.37	0.53
1:A:118:ASN:HA	1:A:141:LYS:CG	2.39	0.53
1:A:354:ILE:HG22	1:A:370:PHE:CD1	2.44	0.53
1:A:5:PRO:C	1:A:7:PHE:N	2.68	0.52
1:A:171:SER:HA	1:A:172:PRO:C	2.34	0.52
1:A:234:PHE:C	1:A:237:PRO:HD3	2.35	0.52
1:A:378:ASP:N	1:A:378:ASP:OD1	2.36	0.52
1:A:48:ASN:ND2	1:A:50:SER:H	2.08	0.52
1:A:130:ALA:O	1:A:134:ILE:HG13	2.09	0.51
1:A:16:PHE:HA	1:A:18:ASP:HB3	1.91	0.51
1:A:123:ILE:HD13	1:A:168:CYS:HB2	1.92	0.51
1:A:123:ILE:HD12	1:A:124:GLN:N	2.18	0.51
1:A:351:HIS:HB2	1:A:385:ARG:HH11	1.76	0.51
1:A:109:PHE:HB2	1:A:134:ILE:HD13	1.93	0.51
1:A:141:LYS:NZ	1:A:141:LYS:CB	2.68	0.51
1:A:301:TYR:O	1:A:302:ARG:C	2.54	0.51
1:A:361:GLY:HA2	1:A:362:ARG:CZ	2.41	0.50
1:A:119:GLN:HA	1:A:119:GLN:OE1	2.11	0.50
1:A:207:THR:O	1:A:207:THR:CG2	2.60	0.49
1:A:354:ILE:CG2	1:A:370:PHE:CD1	2.94	0.49
1:A:118:ASN:O	1:A:141:LYS:HG3	2.12	0.49
1:A:5:PRO:O	1:A:8:VAL:HG22	2.12	0.49
1:A:345:PHE:HD1	1:A:346:PHE:N	2.10	0.49
1:A:124:GLN:HG2	1:A:148:CYS:SG	2.52	0.49
1:A:164:ILE:CG1	1:A:197:ILE:HD13	2.42	0.49
1:A:354:ILE:HG22	1:A:370:PHE:HD1	1.78	0.49
1:A:102:ASP:OD2	1:A:274:PHE:CZ	2.66	0.48
1:A:345:PHE:CD1	1:A:345:PHE:O	2.65	0.48
1:A:315:THR:HG23	1:A:317:PHE:CD1	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:ALA:HB1	1:A:19:LEU:H	1.78	0.48
1:A:90:ARG:CZ	1:A:216:LYS:HD2	2.44	0.48
1:A:59:ILE:HA	1:A:62:LEU:HD12	1.96	0.48
1:A:62:LEU:HD11	1:A:280:PRO:HG3	1.96	0.48
1:A:235:SER:O	1:A:238:LEU:CD1	2.56	0.47
1:A:193:ILE:CD1	1:A:223:ASP:HB3	2.45	0.47
1:A:309:ARG:HG3	1:A:319:VAL:HG11	1.97	0.47
1:A:92:PHE:CE2	1:A:220:GLU:HG2	2.50	0.47
1:A:102:ASP:HB2	1:A:244:ARG:HG3	1.97	0.47
1:A:339:ASP:HB3	1:A:358:ILE:HG13	1.97	0.46
1:A:9:SER:H	1:A:10:LEU:HD13	1.79	0.46
1:A:125:ASP:HB2	1:A:145:ALA:HB1	1.96	0.46
1:A:258:ASP:CG	1:A:260:HIS:HD1	2.24	0.46
1:A:222:PRO:O	1:A:223:ASP:HB2	2.15	0.46
1:A:153:ALA:O	1:A:154:PHE:HB2	2.16	0.46
1:A:237:PRO:HD2	1:A:240:PHE:CD1	2.51	0.46
1:A:111:LEU:O	1:A:114:PHE:HB3	2.16	0.45
1:A:255:THR:HG23	1:A:256:TYR:O	2.16	0.45
1:A:16:PHE:HA	1:A:17:ALA:HA	1.86	0.45
1:A:244:ARG:HD3	1:A:244:ARG:HA	1.68	0.45
1:A:308:LEU:HD23	1:A:383:CYS:SG	2.57	0.45
1:A:319:VAL:CG1	1:A:328:LEU:HD23	2.47	0.45
1:A:27:ARG:NE	1:A:35:VAL:HB	2.32	0.44
1:A:139:GLY:O	1:A:140:ALA:C	2.61	0.44
1:A:375:LYS:HB3	1:A:375:LYS:HE3	1.78	0.44
1:A:331:LYS:HA	1:A:332:PRO:HD3	1.73	0.44
1:A:128:TYR:CD2	1:A:236:LLP:H2'3	2.54	0.43
1:A:294:GLN:O	1:A:295:LEU:C	2.62	0.43
1:A:7:PHE:O	1:A:10:LEU:HD13	2.19	0.43
1:A:253:GLU:N	1:A:253:GLU:OE1	2.51	0.43
1:A:128:TYR:HA	1:A:129:PRO:HD3	1.74	0.43
1:A:123:ILE:HD11	1:A:131:TYR:CD2	2.54	0.43
1:A:5:PRO:O	1:A:7:PHE:N	2.52	0.42
1:A:26:PHE:CD1	1:A:26:PHE:C	2.96	0.42
1:A:234:PHE:CD2	1:A:245:LEU:HD23	2.54	0.42
1:A:319:VAL:HG13	1:A:328:LEU:HD23	2.01	0.42
1:A:101:SER:HB3	2:A:421:HOH:O	2.19	0.42
1:A:315:THR:HG23	1:A:317:PHE:HD1	1.84	0.42
1:A:4:ASN:HA	1:A:5:PRO:HD3	1.96	0.42
1:A:237:PRO:O	1:A:298:ILE:CG2	2.68	0.42
1:A:234:PHE:O	1:A:237:PRO:CD	2.63	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:270:LEU:HD23	1:A:270:LEU:HA	1.85	0.41
1:A:347:LEU:HD12	1:A:347:LEU:HA	1.78	0.41
1:A:262:VAL:O	1:A:263:ILE:C	2.63	0.41
1:A:27:ARG:HH11	1:A:27:ARG:CG	2.21	0.41
1:A:108:LEU:HD21	1:A:131:TYR:HD1	1.85	0.41
1:A:160:GLU:H	1:A:160:GLU:CD	2.27	0.41
1:A:225:ARG:HB3	1:A:251:PRO:HB3	2.03	0.41
1:A:231:ILE:HG22	1:A:248:THR:OG1	2.20	0.41
1:A:325:ALA:HB1	1:A:326:PRO:CD	2.50	0.41
1:A:164:ILE:HG12	1:A:197:ILE:HD13	2.01	0.41
1:A:36:ILE:HG21	1:A:382:ALA:HB2	2.03	0.41
1:A:240:PHE:O	1:A:243:ILE:HG13	2.21	0.41
1:A:291:ILE:C	1:A:291:ILE:HD12	2.46	0.40
1:A:354:ILE:HD12	1:A:354:ILE:O	2.21	0.40
1:A:94:ASP:CB	1:A:97:GLU:HG3	2.43	0.40
1:A:382:ALA:O	1:A:386:LEU:HG	2.21	0.40
1:A:51:VAL:HG11	1:A:243:ILE:HD11	2.04	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	351/400 (88%)	305 (87%)	38 (11%)	8 (2%)	5 13

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	235	SER
1	A	295	LEU
1	A	385	ARG
1	A	40	ILE

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Mol	Chain	Res	Type
1	A	297	ALA
1	A	114	PHE
1	A	8	VAL
1	A	358	ILE

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	304/336 (90%)	244 (80%)	60 (20%)	1 4

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	SER
1	A	20	GLN
1	A	27	ARG
1	A	28	LEU
1	A	40	ILE
1	A	55	PHE
1	A	59	ILE
1	A	74	ASP
1	A	82	GLN
1	A	85	SER
1	A	90	ARG
1	A	102	ASP
1	A	106	VAL
1	A	120	THR
1	A	121	VAL
1	A	123	ILE
1	A	127	SER
1	A	141	LYS
1	A	143	ILE
1	A	146	LEU
1	A	149	LEU
1	A	150	GLN

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Mol	Chain	Res	Type
1	A	151	GLU
1	A	157	GLU
1	A	160	GLU
1	A	161	ASP
1	A	162	THR
1	A	164	ILE
1	A	167	LEU
1	A	169	LEU
1	A	178	THR
1	A	180	LEU
1	A	200	LEU
1	A	238	LEU
1	A	244	ARG
1	A	248	THR
1	A	255	THR
1	A	263	ILE
1	A	264	GLN
1	A	268	ARG
1	A	279	ILE
1	A	283	GLU
1	A	298	ILE
1	A	299	HIS
1	A	302	ARG
1	A	310	LYS
1	A	312	LEU
1	A	323	GLU
1	A	328	LEU
1	A	340	ARG
1	A	345	PHE
1	A	354	ILE
1	A	355	THR
1	A	358	ILE
1	A	362	ARG
1	A	363	SER
1	A	372	SER
1	A	378	ASP
1	A	387	GLN
1	A	388	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	48	ASN
1	A	82	GLN
1	A	118	ASN
1	A	152	ASN
1	A	252	GLN
1	A	264	GLN
1	A	387	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
1	LLP	A	236	1	23,24,25	2.52	10 (43%)	25,32,34	1.85	4 (16%)

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
1	LLP	A	236	1	-	11/16/17/19	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	236	LLP	C4'-NZ	5.93	1.47	1.27
1	A	236	LLP	CE-NZ	-5.84	1.33	1.46
1	A	236	LLP	C4-C4'	3.84	1.54	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	236	LLP	CD-CE	-3.67	1.38	1.51
1	A	236	LLP	P-OP2	3.27	1.67	1.54
1	A	236	LLP	CD-CG	-2.54	1.39	1.51
1	A	236	LLP	P-OP3	2.52	1.64	1.54
1	A	236	LLP	P-OP4	2.37	1.67	1.60
1	A	236	LLP	OP4-C5'	-2.15	1.37	1.44
1	A	236	LLP	C2'-C2	2.00	1.53	1.50

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	236	LLP	OP4-C5'-C5	6.51	121.57	109.36
1	A	236	LLP	CD-CE-NZ	3.10	119.03	110.83
1	A	236	LLP	C4-C4'-NZ	-2.85	110.88	124.04
1	A	236	LLP	OP4-P-OP1	2.33	112.75	106.44

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	236	LLP	C4-C4'-NZ-CE
1	A	236	LLP	C5'-OP4-P-OP2
1	A	236	LLP	N-CA-CB-CG
1	A	236	LLP	CG-CD-CE-NZ
1	A	236	LLP	C3-C4-C4'-NZ
1	A	236	LLP	CA-CB-CG-CD
1	A	236	LLP	CD-CE-NZ-C4'
1	A	236	LLP	C-CA-CB-CG
1	A	236	LLP	C5'-OP4-P-OP1
1	A	236	LLP	C5-C4-C4'-NZ
1	A	236	LLP	CE-CD-CG-CB

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	236	LLP	5	0

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 ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2		OWAB(Å ²)	Q < 0.9
1	A	363/400 (90%)	0.51	19 (5%)	33 29	48, 75, 116, 123	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	16	PHE	4.9
1	A	238	LEU	4.6
1	A	36	ILE	4.1
1	A	17	ALA	4.0
1	A	10	LEU	3.6
1	A	388	MET	3.3
1	A	32	GLN	3.3
1	A	381	ALA	2.6
1	A	66	THR	2.5
1	A	295	LEU	2.5
1	A	41	GLY	2.5
1	A	239	GLY	2.4
1	A	370	PHE	2.3
1	A	346	PHE	2.3
1	A	342	LEU	2.2
1	A	237	PRO	2.2
1	A	362	ARG	2.1
1	A	343	PHE	2.1
1	A	332	PRO	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
1	LLP	A	236	24/25	0.86	0.13	96,113,124,125	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.