



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:51 AM UTC

PDB ID : 7YTH / pdb_00007yth
Title : Structure of OCPx1 from Nostoc flagelliforme CCNUN1
Authors : Yang, Y.W.; Liu, K.; Chen, S.Z.; Chen, M.; Qiu, B.S.
Deposited on : 2022-08-14
Resolution : 1.93 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
Xtrriage (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

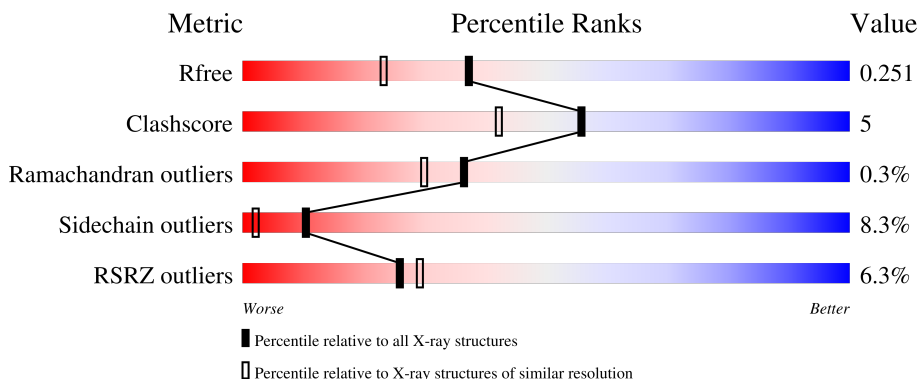
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.93 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	1452 (1.94-1.94)
Clashscore	190562	1494 (1.94-1.94)
Ramachandran outliers	187476	1479 (1.94-1.94)
Sidechain outliers	187428	1479 (1.94-1.94)
RSRZ outliers	180081	1453 (1.94-1.94)

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	326	 4% 81% 13% . .
1	B	326	 5% 83% 12% . .
1	C	326	 4% 82% 12% . .
1	D	326	 11% 80% 11% . 5%

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 10296 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 Ketosteroid isomerase-related protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	314	2433	1562	406	453	12	0	0	0
1	B	315	2437	1564	407	454	12	0	0	0
1	C	314	2439	1566	408	453	12	0	0	0
1	D	311	2409	1548	399	450	12	0	0	0

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	321	HIS	-	expression tag	UNP A0A2K8SJT8
A	322	HIS	-	expression tag	UNP A0A2K8SJT8
A	323	HIS	-	expression tag	UNP A0A2K8SJT8
A	324	HIS	-	expression tag	UNP A0A2K8SJT8
A	325	HIS	-	expression tag	UNP A0A2K8SJT8
A	326	HIS	-	expression tag	UNP A0A2K8SJT8
B	321	HIS	-	expression tag	UNP A0A2K8SJT8
B	322	HIS	-	expression tag	UNP A0A2K8SJT8
B	323	HIS	-	expression tag	UNP A0A2K8SJT8
B	324	HIS	-	expression tag	UNP A0A2K8SJT8
B	325	HIS	-	expression tag	UNP A0A2K8SJT8
B	326	HIS	-	expression tag	UNP A0A2K8SJT8
C	321	HIS	-	expression tag	UNP A0A2K8SJT8
C	322	HIS	-	expression tag	UNP A0A2K8SJT8
C	323	HIS	-	expression tag	UNP A0A2K8SJT8
C	324	HIS	-	expression tag	UNP A0A2K8SJT8
C	325	HIS	-	expression tag	UNP A0A2K8SJT8
C	326	HIS	-	expression tag	UNP A0A2K8SJT8
D	321	HIS	-	expression tag	UNP A0A2K8SJT8
D	322	HIS	-	expression tag	UNP A0A2K8SJT8
D	323	HIS	-	expression tag	UNP A0A2K8SJT8

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Chain	Residue	Modelled	Actual	Comment	Reference
D	324	HIS	-	expression tag	UNP A0A2K8SJT8
D	325	HIS	-	expression tag	UNP A0A2K8SJT8
D	326	HIS	-	expression tag	UNP A0A2K8SJT8

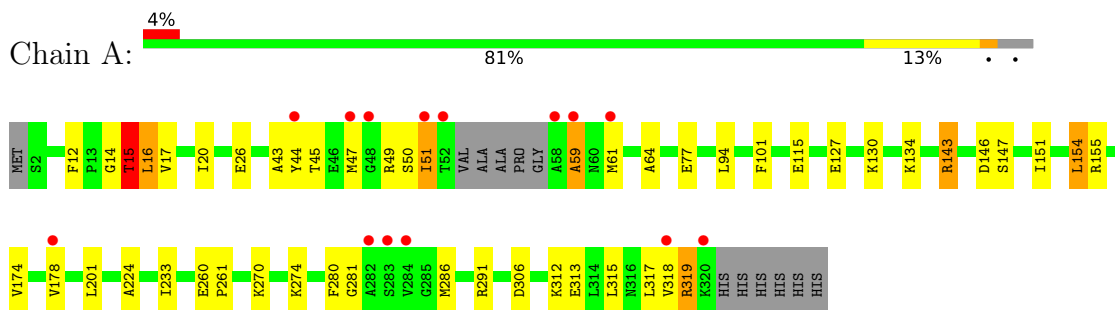
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	199	Total 199	O 199	0	0
2	B	107	Total 107	O 107	0	0
2	C	157	Total 157	O 157	0	0
2	D	115	Total 115	O 115	0	0

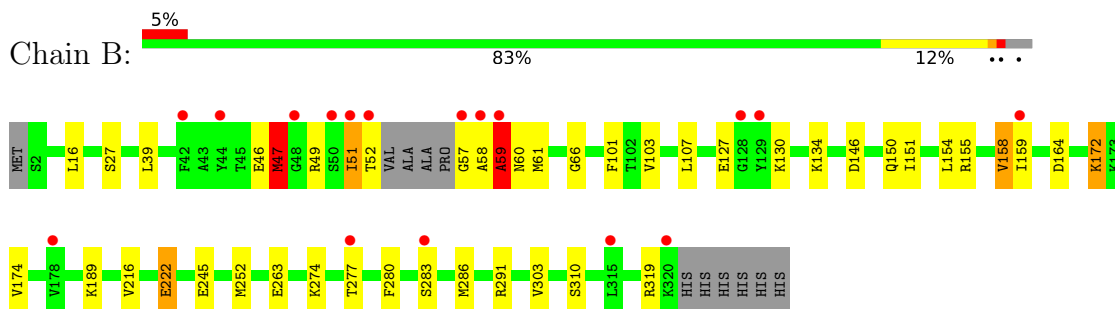
3 Residue-property plots [i](#)

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.

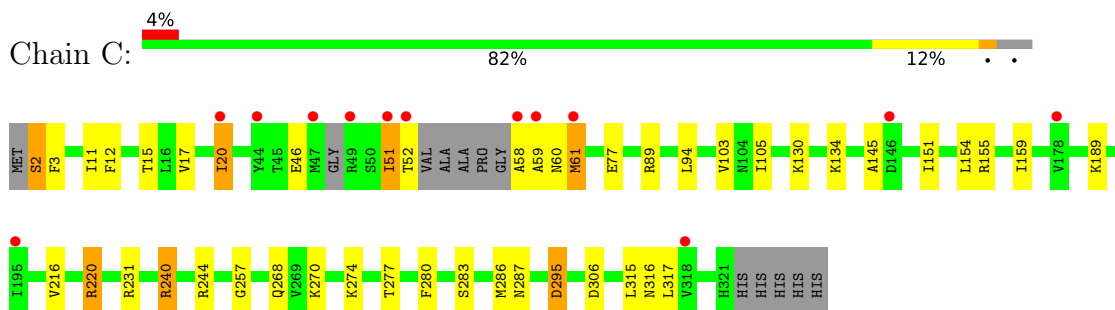
- Molecule 1: Ketosteroid isomerase-related protein



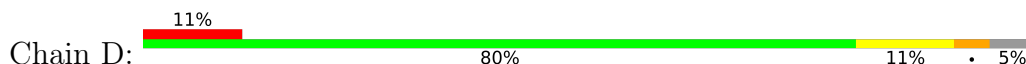
- Molecule 1: Ketosteroid isomerase-related protein

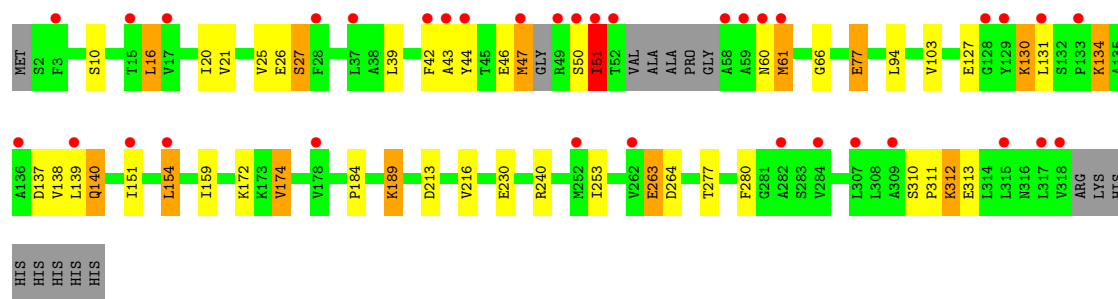


- Molecule 1: Ketosteroid isomerase-related protein



- Molecule 1: Ketosteroid isomerase-related protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	79.78Å 122.95Å 141.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	67.01 – 1.93 67.01 – 1.93	Depositor EDS
% Data completeness (in resolution range)	98.5 (67.01-1.93) 98.5 (67.01-1.93)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.65 (at 1.92Å)	Xtrriage
Refinement program	REFMAC 5.8.0257	Depositor
R, R_{free}	0.199 , 0.248 0.206 , 0.251	Depositor DCC
R_{free} test set	5103 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	38.8	Xtrriage
Anisotropy	0.080	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 38.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10296	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.14% 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 i

5.1 Standard geometry i

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	1.09	1/2487 (0.0%)	1.35	7/3383 (0.2%)
1	B	1.03	0/2491	1.39	3/3388 (0.1%)
1	C	1.07	2/2493 (0.1%)	1.37	4/3390 (0.1%)
1	D	1.07	0/2462	1.40	7/3350 (0.2%)
All	All	1.06	3/9933 (0.0%)	1.38	21/13511 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	261	PRO	C-O	-5.64	1.17	1.23
1	C	11	ILE	C-O	5.38	1.30	1.24
1	C	220	ARG	C-O	5.36	1.30	1.23

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	77	GLU	CB-CG-CD	6.62	123.86	112.60
1	D	264	ASP	CA-CB-CG	6.61	119.21	112.60
1	C	268	GLN	CB-CA-C	5.99	119.64	109.75
1	D	42	PHE	CB-CA-C	5.87	120.54	110.79
1	C	59	ALA	CA-C-N	5.84	128.11	120.28
1	C	59	ALA	C-N-CA	5.84	128.11	120.28
1	B	101	PHE	CA-CB-CG	5.70	119.50	113.80
1	B	59	ALA	CA-C-N	5.65	128.42	120.28
1	B	59	ALA	C-N-CA	5.65	128.42	120.28
1	A	146	ASP	CA-CB-CG	5.60	118.20	112.60
1	C	295	ASP	CA-CB-CG	5.46	118.06	112.60
1	D	189	LYS	CB-CA-C	5.45	119.16	109.72
1	A	101	PHE	CA-CB-CG	5.31	119.11	113.80
1	D	230	GLU	CA-C-N	5.31	127.77	120.39
1	D	230	GLU	C-N-CA	5.31	127.77	120.39
1	A	59	ALA	CA-C-N	5.21	127.78	120.28
1	A	59	ALA	C-N-CA	5.21	127.78	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	15	THR	CB-CA-C	5.19	119.37	110.81
1	A	201	LEU	CA-C-N	5.15	125.70	119.98
1	A	201	LEU	C-N-CA	5.15	125.70	119.98
1	D	127	GLU	N-CA-C	-5.06	102.56	110.10

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2433	0	2438	28	0
1	B	2437	0	2441	20	0
1	C	2439	0	2441	20	0
1	D	2409	0	2408	29	0
2	A	199	0	0	3	0
2	B	107	0	0	3	0
2	C	157	0	0	0	0
2	D	115	0	0	1	0
All	All	10296	0	9728	91	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 5.

All (91) 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:14:GLY:HA3	2:A:505:HOH:O	1.82	0.79
1:D:140:GLN:HE21	1:D:140:GLN:HA	1.53	0.74
1:C:220:ARG:HD2	1:C:295:ASP:OD2	1.89	0.71
1:D:103:VAL:HG13	1:D:159:ILE:HD11	1.73	0.70
1:D:312:LYS:N	1:D:312:LYS:HD2	2.07	0.69
1:C:103:VAL:HG13	1:C:159:ILE:HD11	1.78	0.66
1:A:155:ARG:NH2	2:A:401:HOH:O	2.24	0.65
1:C:151:ILE:HD12	1:C:280:PHE:CZ	2.33	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:51:ILE:HD12	1:D:51:ILE:HD12	1.79	0.64
1:D:26:GLU:OE1	1:D:26:GLU:HA	1.99	0.63
1:D:140:GLN:HE21	1:D:140:GLN:CA	2.10	0.63
1:A:43:ALA:HB1	1:A:154:LEU:HD21	1.81	0.62
1:A:51:ILE:HG13	1:B:51:ILE:HG21	1.81	0.62
1:D:47:MET:SD	1:D:154:LEU:HD23	2.40	0.61
1:A:15:THR:HB	1:A:270:LYS:HZ1	1.66	0.61
1:D:131:LEU:HB2	2:D:444:HOH:O	2.01	0.60
1:D:16:LEU:HB2	1:D:313:GLU:CD	2.26	0.60
1:B:274:LYS:HA	1:B:286:MET:O	2.01	0.59
1:A:12:PHE:O	1:A:15:THR:CG2	2.51	0.59
1:A:44:TYR:OH	1:A:115:GLU:OE2	2.21	0.58
1:C:240:ARG:HH21	1:C:244:ARG:NH2	2.03	0.56
1:B:103:VAL:HG13	1:B:159:ILE:HD11	1.87	0.55
1:C:316:ASN:C	1:C:317:LEU:HD22	2.31	0.55
1:D:134:LYS:O	1:D:137:ASP:HB2	2.07	0.54
1:C:240:ARG:NH2	1:C:244:ARG:NH1	2.55	0.54
1:A:17:VAL:HG12	1:A:306:ASP:HB3	1.91	0.53
1:A:47:MET:HE1	1:A:154:LEU:HD23	1.91	0.53
1:A:15:THR:HB	1:A:270:LYS:NZ	2.24	0.52
1:D:61:MET:HE1	1:D:174:VAL:HG11	1.91	0.52
1:A:51:ILE:HB	1:B:51:ILE:HB	1.92	0.52
1:A:12:PHE:O	1:A:15:THR:HG23	2.10	0.52
1:D:39:LEU:CD2	1:D:139:LEU:HB2	2.40	0.52
1:D:43:ALA:HB1	1:D:154:LEU:HD21	1.91	0.52
1:B:66:GLY:HA3	1:D:66:GLY:HA3	1.92	0.52
1:A:260:GLU:OE1	1:A:319:ARG:HA	2.10	0.51
1:A:49:ARG:H	1:A:49:ARG:HD2	1.75	0.50
1:D:310:SER:HB2	1:D:311:PRO:HD2	1.93	0.50
1:A:47:MET:HE3	1:A:151:ILE:CD1	2.42	0.50
1:C:94:LEU:C	1:C:94:LEU:HD13	2.37	0.50
1:A:281:GLY:HA3	1:B:51:ILE:HD12	1.93	0.49
1:B:252:MET:HE1	2:B:441:HOH:O	2.13	0.49
1:C:240:ARG:HH21	1:C:244:ARG:HH22	1.60	0.48
1:B:172:LYS:HE3	2:B:489:HOH:O	2.13	0.48
1:D:130:LYS:CE	1:D:130:LYS:HA	2.42	0.47
1:A:94:LEU:C	1:A:94:LEU:HD13	2.39	0.47
1:C:20:ILE:CD1	1:C:145:ALA:HB2	2.44	0.47
1:D:213:ASP:OD1	1:D:240:ARG:HD3	2.15	0.47
1:C:274:LYS:HE3	1:C:287:ASN:OD1	2.15	0.47
1:D:27:SER:HB2	1:D:138:VAL:HG22	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:47:MET:SD	1:D:154:LEU:CD2	3.03	0.46
1:A:47:MET:HE3	1:A:151:ILE:HD13	1.97	0.46
1:C:220:ARG:CD	1:C:295:ASP:OD2	2.61	0.46
1:D:263:GLU:H	1:D:263:GLU:HG3	1.46	0.46
1:D:21:VAL:O	1:D:25:VAL:HG23	2.16	0.46
1:A:16:LEU:HB2	1:A:313:GLU:CD	2.42	0.45
1:D:50:SER:C	1:D:51:ILE:HG12	2.38	0.45
1:A:12:PHE:O	1:A:15:THR:HG22	2.16	0.45
1:A:51:ILE:HD12	1:B:51:ILE:HG13	1.98	0.45
1:D:39:LEU:HD21	1:D:139:LEU:HB2	1.99	0.45
1:D:94:LEU:HD13	1:D:94:LEU:C	2.42	0.45
1:A:151:ILE:HD12	1:A:280:PHE:CZ	2.52	0.44
1:B:245:GLU:OE2	2:B:401:HOH:O	2.21	0.44
1:B:107:LEU:HD23	1:B:158:VAL:CG2	2.47	0.44
1:B:280:PHE:O	1:B:283:SER:HB2	2.17	0.44
1:D:151:ILE:HD12	1:D:280:PHE:CZ	2.53	0.44
1:C:61:MET:HE3	1:C:105:ILE:HD13	1.98	0.43
1:D:312:LYS:N	1:D:312:LYS:CD	2.79	0.43
1:B:291:ARG:O	1:B:303:VAL:HA	2.19	0.43
1:B:58:ALA:O	1:B:59:ALA:HB3	2.19	0.43
1:B:164:ASP:OD1	1:B:164:ASP:C	2.62	0.43
1:C:2:SER:HB2	1:C:3:PHE:H	1.72	0.42
1:A:147:SER:O	1:A:151:ILE:HG12	2.18	0.42
1:B:39:LEU:HD23	1:B:39:LEU:HA	1.86	0.42
1:C:58:ALA:C	1:C:60:ASN:H	2.27	0.42
1:A:274:LYS:HA	1:A:286:MET:O	2.19	0.42
1:D:16:LEU:HD22	1:D:313:GLU:OE1	2.20	0.42
1:D:130:LYS:HA	1:D:130:LYS:HD2	1.87	0.42
1:C:12:PHE:O	1:C:15:THR:HG22	2.20	0.41
1:C:274:LYS:HA	1:C:286:MET:O	2.19	0.41
1:C:17:VAL:HG12	1:C:306:ASP:HB3	2.02	0.41
1:B:146:ASP:O	1:B:150:GLN:HG3	2.20	0.41
1:C:240:ARG:NH2	1:C:244:ARG:HH12	2.19	0.41
1:B:47:MET:HA	1:B:49:ARG:NH1	2.36	0.41
1:B:52:THR:CG2	1:B:57:GLY:HA2	2.50	0.41
1:B:222:GLU:CD	1:B:222:GLU:H	2.29	0.41
1:A:15:THR:HG21	1:A:291:ARG:HD3	2.03	0.40
1:A:143:ARG:HD3	2:A:567:HOH:O	2.22	0.40
1:D:27:SER:CB	1:D:138:VAL:HG22	2.51	0.40
1:A:59:ALA:HB1	1:A:64:ALA:HB1	2.03	0.40
1:A:224:ALA:HA	1:A:233:ILE:O	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:257:GLY:HA2	1:C:270:LYS:O	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	310/326 (95%)	301 (97%)	9 (3%)	0	100	100
1	B	311/326 (95%)	300 (96%)	8 (3%)	3 (1%)	12	5
1	C	308/326 (94%)	297 (96%)	11 (4%)	0	100	100
1	D	305/326 (94%)	298 (98%)	6 (2%)	1 (0%)	36	30
All	All	1234/1304 (95%)	1196 (97%)	34 (3%)	4 (0%)	36	30

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	47	MET
1	B	59	ALA
1	B	46	GLU
1	D	51	ILE

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	260/269 (97%)	239 (92%)	21 (8%)	11	2
1	B	260/269 (97%)	238 (92%)	22 (8%)	10	2
1	C	261/269 (97%)	242 (93%)	19 (7%)	13	3
1	D	258/269 (96%)	234 (91%)	24 (9%)	8	1
All	All	1039/1076 (97%)	953 (92%)	86 (8%)	10	2

All (86) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	THR
1	A	16	LEU
1	A	20	ILE
1	A	26	GLU
1	A	45	THR
1	A	50	SER
1	A	51	ILE
1	A	61	MET
1	A	77	GLU
1	A	127	GLU
1	A	130	LYS
1	A	134	LYS
1	A	143	ARG
1	A	154	LEU
1	A	174	VAL
1	A	178	VAL
1	A	312	LYS
1	A	315	LEU
1	A	317	LEU
1	A	318	VAL
1	A	319	ARG
1	B	16	LEU
1	B	27	SER
1	B	47	MET
1	B	51	ILE
1	B	60	ASN
1	B	61	MET
1	B	127	GLU
1	B	130	LYS
1	B	134	LYS
1	B	151	ILE
1	B	154	LEU
1	B	155	ARG

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Mol	Chain	Res	Type
1	B	158	VAL
1	B	172	LYS
1	B	174	VAL
1	B	189	LYS
1	B	216	VAL
1	B	222	GLU
1	B	263	GLU
1	B	277	THR
1	B	310	SER
1	B	319	ARG
1	C	2	SER
1	C	20	ILE
1	C	46	GLU
1	C	51	ILE
1	C	52	THR
1	C	61	MET
1	C	77	GLU
1	C	89	ARG
1	C	130	LYS
1	C	134	LYS
1	C	154	LEU
1	C	155	ARG
1	C	189	LYS
1	C	216	VAL
1	C	231	ARG
1	C	240	ARG
1	C	277	THR
1	C	283	SER
1	C	315	LEU
1	D	10	SER
1	D	16	LEU
1	D	20	ILE
1	D	27	SER
1	D	44	TYR
1	D	46	GLU
1	D	47	MET
1	D	51	ILE
1	D	60	ASN
1	D	61	MET
1	D	77	GLU
1	D	130	LYS
1	D	134	LYS

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Mol	Chain	Res	Type
1	D	140	GLN
1	D	154	LEU
1	D	172	LYS
1	D	174	VAL
1	D	184	PRO
1	D	189	LYS
1	D	216	VAL
1	D	253	ILE
1	D	263	GLU
1	D	277	THR
1	D	312	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	GLN
1	A	36	GLN
1	A	73	GLN
1	A	206	ASN
1	A	248	GLN
1	B	60	ASN
1	B	73	GLN
1	B	166	ASN
1	C	73	GLN
1	C	297	GLN
1	D	36	GLN
1	D	140	GLN
1	D	297	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	314/326 (96%)	0.36	14 (4%) 38 43	28, 40, 67, 98	0
1	B	315/326 (96%)	0.65	17 (5%) 31 35	32, 49, 77, 114	0
1	C	314/326 (96%)	0.47	13 (4%) 41 46	29, 43, 72, 104	0
1	D	311/326 (95%)	0.81	35 (11%) 10 11	29, 50, 81, 117	0
All	All	1254/1304 (96%)	0.57	79 (6%) 26 29	28, 45, 75, 117	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	58	ALA	6.0
1	B	58	ALA	4.9
1	D	318	VAL	4.5
1	B	51	ILE	4.2
1	D	44	TYR	4.2
1	D	52	THR	4.2
1	C	58	ALA	4.1
1	B	52	THR	4.1
1	D	128	GLY	4.0
1	A	58	ALA	3.9
1	C	178	VAL	3.8
1	A	318	VAL	3.8
1	D	47	MET	3.6
1	D	131	LEU	3.5
1	B	59	ALA	3.4
1	B	283	SER	3.4
1	A	52	THR	3.4
1	B	178	VAL	3.4
1	B	42	PHE	3.4
1	B	44	TYR	3.3
1	C	52	THR	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	129	TYR	3.2
1	D	59	ALA	3.2
1	C	51	ILE	3.2
1	A	283	SER	3.2
1	A	51	ILE	3.1
1	D	51	ILE	3.1
1	D	139	LEU	3.0
1	D	49	ARG	2.9
1	A	48	GLY	2.9
1	D	28	PHE	2.9
1	D	151	ILE	2.8
1	D	17	VAL	2.8
1	A	59	ALA	2.8
1	D	3	PHE	2.8
1	D	133	PRO	2.8
1	D	43	ALA	2.7
1	D	178	VAL	2.7
1	D	262	VAL	2.7
1	A	47	MET	2.6
1	C	44	TYR	2.6
1	B	57	GLY	2.6
1	B	315	LEU	2.6
1	C	61	MET	2.4
1	D	15	THR	2.4
1	C	146	ASP	2.4
1	D	282	ALA	2.4
1	D	309	ALA	2.4
1	A	284	VAL	2.4
1	D	37	LEU	2.4
1	A	282	ALA	2.4
1	D	252	MET	2.4
1	A	320	LYS	2.4
1	D	61	MET	2.3
1	D	60	ASN	2.3
1	D	129	TYR	2.2
1	B	159	ILE	2.2
1	D	317	LEU	2.2
1	B	50	SER	2.2
1	A	44	TYR	2.2
1	C	20	ILE	2.2
1	C	47	MET	2.2
1	C	318	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	277	THR	2.2
1	D	50	SER	2.2
1	D	284	VAL	2.1
1	C	49	ARG	2.1
1	D	42	PHE	2.1
1	B	48	GLY	2.1
1	D	154	LEU	2.1
1	D	315	LEU	2.1
1	C	59	ALA	2.1
1	D	136	ALA	2.1
1	B	320	LYS	2.0
1	C	195	ILE	2.0
1	B	128	GLY	2.0
1	A	178	VAL	2.0
1	D	307	LEU	2.0
1	A	61	MET	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

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.