Full wwPDB X-ray Structure Validation Report ⓘ

Sep 6, 2023 – 08:14 PM EDT

PDB ID	:	8SXW
Title	:	X-ray crystal structure of UDP- 2,3-diacetamido-2,3-dideoxy-glucuronic acid-2-epimerase from Thermus thermophilus strain HB27, D98N mutation, apo structure at pH 6
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Deposited on	:	2023-05-24
Resolution	:	1.80 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:

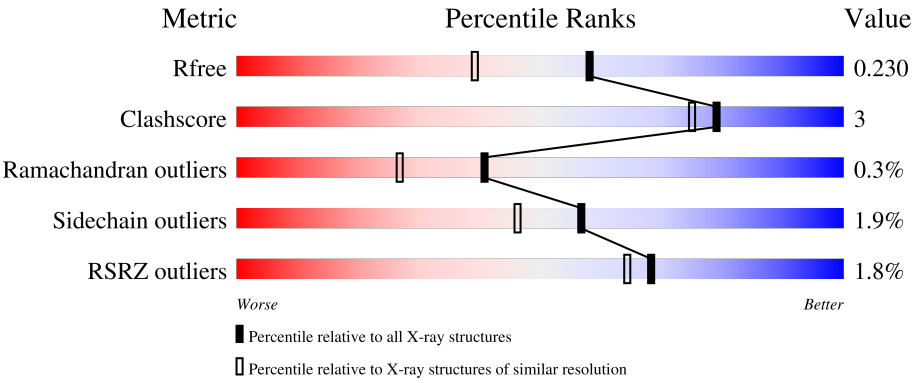
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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 <=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	367	90% 7%
1	B	367	90% 7%
1	C	367	5% 92% 6%
1	D	367	92% 6%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 12560 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 UDP-2,3-diacetamido-2,3-dideoxy-glucuronic acid-2-epimerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	361	Total	C	N	O	S	0	4	0
			2869	1822	514	524	9			
1	B	359	Total	C	N	O	S	0	6	0
			2846	1807	509	522	8			
1	C	363	Total	C	N	O	S	0	4	0
			2860	1818	508	525	9			
1	D	362	Total	C	N	O	S	0	5	0
			2871	1824	513	526	8			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q72KY0
A	0	HIS	-	expression tag	UNP Q72KY0
A	98	ASN	ASP	engineered mutation	UNP Q72KY0
B	-1	GLY	-	expression tag	UNP Q72KY0
B	0	HIS	-	expression tag	UNP Q72KY0
B	98	ASN	ASP	engineered mutation	UNP Q72KY0
C	-1	GLY	-	expression tag	UNP Q72KY0
C	0	HIS	-	expression tag	UNP Q72KY0
C	98	ASN	ASP	engineered mutation	UNP Q72KY0
D	-1	GLY	-	expression tag	UNP Q72KY0
D	0	HIS	-	expression tag	UNP Q72KY0
D	98	ASN	ASP	engineered mutation	UNP Q72KY0

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Cl	0	0
			1	1		
2	D	1	Total	Cl	0	0
			1	1		

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Na 1 1	0	0

- Molecule 4 is water.

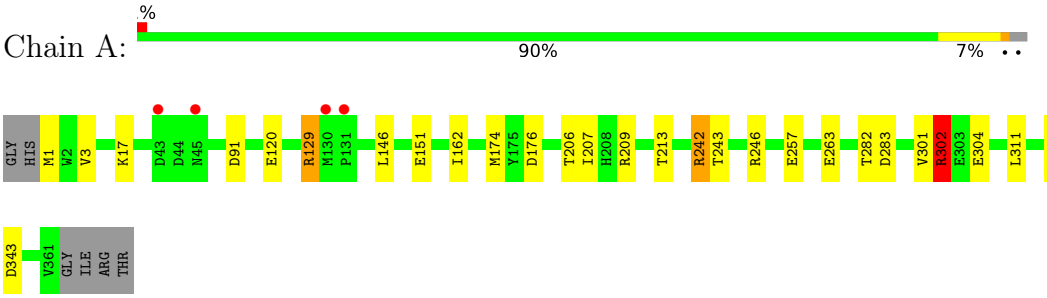
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	251	Total O 251 251	0	0
4	B	297	Total O 297 297	0	0
4	C	238	Total O 238 238	0	0
4	D	325	Total O 325 325	0	0

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.

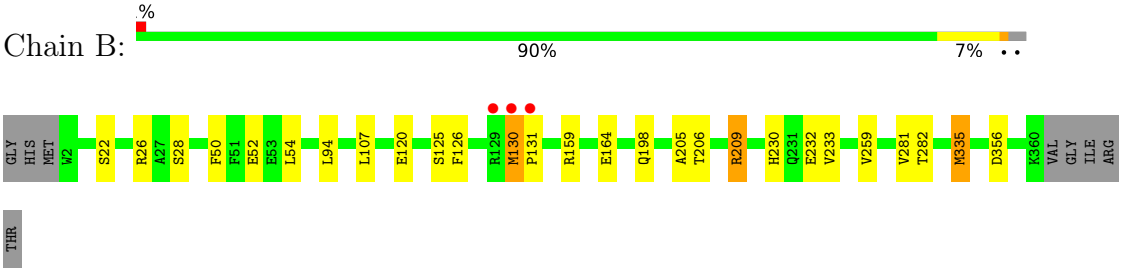
• Molecule 1: UDP-2,3-diacetamido-2,3-dideoxy-glucuronic acid-2-epimerase

Chain A:



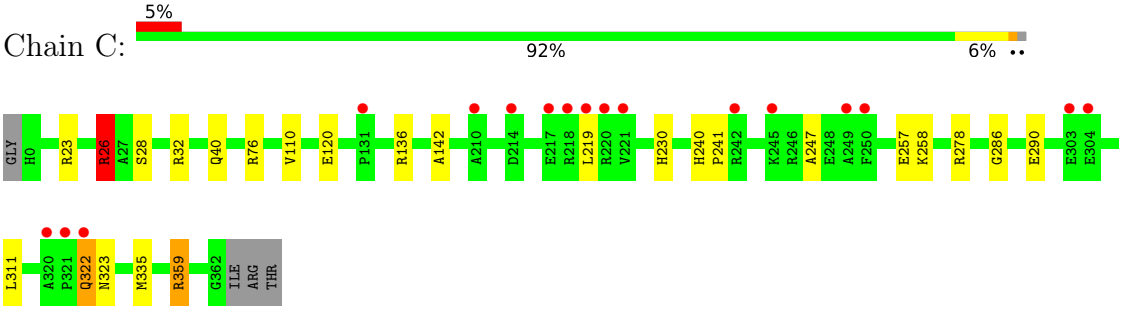
• Molecule 1: UDP-2,3-diacetamido-2,3-dideoxy-glucuronic acid-2-epimerase

Chain B:



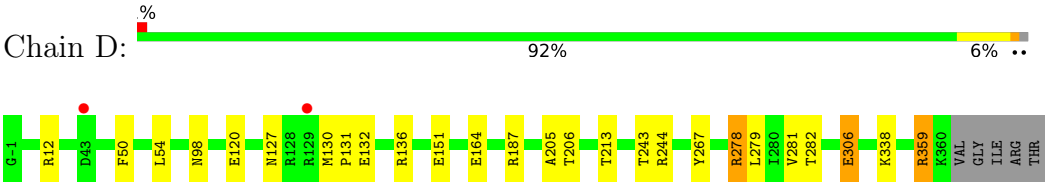
• Molecule 1: UDP-2,3-diacetamido-2,3-dideoxy-glucuronic acid-2-epimerase

Chain C:



• Molecule 1: UDP-2,3-diacetamido-2,3-dideoxy-glucuronic acid-2-epimerase

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	64.33Å 85.34Å 86.53Å 108.23° 110.06° 97.10°	Depositor
Resolution (Å)	41.72 – 1.80 41.68 – 1.80	Depositor EDS
% Data completeness (in resolution range)	94.1 (41.72-1.80) 94.1 (41.68-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$< I/\sigma(I) >$ [1]	1.72 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.8.0403	Depositor
R, R_{free}	0.178 , 0.224 0.187 , 0.230	Depositor DCC
R_{free} test set	7133 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å^2)	18.0	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 41.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	12560	wwPDB-VP
Average B, all atoms (Å^2)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.67% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: CL, NA

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	0.44	0/2935	0.72	1/3979 (0.0%)
1	B	0.46	0/2919	0.74	0/3956
1	C	0.39	0/2928	0.68	0/3973
1	D	0.43	0/2944	0.72	1/3994 (0.0%)
All	All	0.43	0/11726	0.72	2/15902 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	C	0	5
1	D	0	1
All	All	0	8

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	306	GLU	CB-CA-C	-5.29	99.82	110.40
1	A	302	ARG	NE-CZ-NH1	-5.00	117.80	120.30

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	302	ARG	Sidechain
1	B	209	ARG	Sidechain
1	C	136	ARG	Sidechain
1	C	26	ARG	Sidechain
1	C	32	ARG	Sidechain
1	C	359	ARG	Sidechain
1	C	76	ARG	Sidechain
1	D	278	ARG	Sidechain

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	2869	0	2898	16	0
1	B	2846	0	2878	13	0
1	C	2860	0	2883	15	0
1	D	2871	0	2889	20	0
2	B	1	0	0	0	0
2	D	1	0	0	0	0
3	B	1	0	0	0	0
4	A	251	0	0	3	0
4	B	297	0	0	4	1
4	C	238	0	0	4	1
4	D	325	0	0	11	2
All	All	12560	0	11548	64	2

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 3.

All (64) 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:D:278:ARG:HD3	4:D:554:HOH:O	1.82	0.79
1:B:232:GLU:OE2	4:B:501:HOH:O	2.01	0.77
1:B:209:ARG:HD2	4:B:521:HOH:O	1.83	0.77
1:C:290[B]:GLU:OE2	4:C:401:HOH:O	2.06	0.74
1:A:129:ARG:NH1	4:A:401:HOH:O	2.15	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:356:ASP:OD2	4:B:502:HOH:O	2.09	0.70
1:C:230:HIS:CD2	1:C:258:LYS:O	2.46	0.69
1:C:290[A]:GLU:OE1	4:C:402:HOH:O	2.10	0.68
1:B:164:GLU:O	4:B:503:HOH:O	2.11	0.67
1:A:242:ARG:HG2	1:A:243:THR:N	2.14	0.62
1:C:110[A]:VAL:HG11	1:C:142:ALA:O	2.01	0.60
1:D:54:LEU:HD21	1:D:267[B]:TYR:CD2	2.36	0.60
1:A:206:THR:O	1:A:207:ILE:HD12	2.03	0.58
1:C:26:ARG:NH2	4:C:405:HOH:O	2.36	0.58
1:D:267[A]:TYR:CD2	4:D:685:HOH:O	2.52	0.57
1:A:206:THR:C	1:A:207:ILE:HD12	2.26	0.56
1:A:336:ARG:NH2	4:A:403:HOH:O	2.21	0.55
1:D:267[A]:TYR:HD2	4:D:685:HOH:O	1.88	0.55
1:A:129:ARG:HH21	1:A:129:ARG:HG2	1.72	0.55
1:D:151:GLU:HG2	4:D:719:HOH:O	2.07	0.55
1:A:213:THR:O	1:A:246:ARG:HG2	2.07	0.54
1:D:338:LYS:NZ	4:D:501:HOH:O	2.25	0.54
1:A:146:LEU:HD13	1:A:162:ILE:HG21	1.91	0.53
1:B:130:MET:HG2	1:B:131:PRO:HD2	1.90	0.52
1:C:28:SER:OG	1:C:359:ARG:NH2	2.44	0.50
1:A:209:ARG:HD3	1:A:302:ARG:HD3	1.93	0.49
1:D:306:GLU:HG3	4:D:698:HOH:O	2.13	0.49
1:A:129:ARG:HG2	1:A:129:ARG:NH2	2.29	0.48
1:C:219:LEU:HD21	1:C:247:ALA:HB2	1.96	0.48
1:B:233:VAL:HG11	1:B:335:MET:SD	2.54	0.48
1:C:322:GLN:HG2	1:C:323:ASN:N	2.29	0.48
1:C:359:ARG:HD3	4:C:586:HOH:O	2.13	0.48
1:A:206:THR:O	1:A:282:THR:HA	2.14	0.47
1:A:301:VAL:O	1:A:302:ARG:HG2	2.15	0.46
1:C:278:ARG:HD2	1:C:335:MET:CG	2.45	0.46
1:C:23:ARG:HA	1:C:26:ARG:HD3	1.97	0.46
1:D:131:PRO:O	1:D:132:GLU:HB2	2.16	0.46
1:D:151:GLU:CG	4:D:719:HOH:O	2.63	0.45
1:D:359:ARG:HH21	1:D:359:ARG:HG3	1.82	0.45
1:B:126:PHE:HE1	1:B:159[B]:ARG:HH12	1.63	0.45
1:D:127:ASN:HB3	1:D:130:MET:HG3	1.99	0.45
1:D:130:MET:HB3	4:D:627:HOH:O	2.16	0.45
1:D:187:ARG:HD3	4:D:734:HOH:O	2.16	0.44
1:A:343:ASP:OD1	4:A:402:HOH:O	2.21	0.43
1:C:286:GLY:O	1:C:290[B]:GLU:HG2	2.19	0.43
1:A:17:LYS:HE3	1:A:174:MET:HG2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:164:GLU:HG2	4:D:743:HOH:O	2.19	0.43
1:D:98:ASN:OD1	1:D:136:ARG:HA	2.18	0.43
1:B:230:HIS:HB2	1:B:259:VAL:HG12	2.01	0.43
1:D:206:THR:O	1:D:282:THR:HA	2.19	0.43
1:B:206:THR:O	1:B:282:THR:HA	2.20	0.42
1:A:3:VAL:HG23	1:A:91:ASP:HB2	2.01	0.42
1:D:205:ALA:HA	1:D:281:VAL:O	2.19	0.42
1:A:283:ASP:HB2	1:A:302:ARG:HD3	2.02	0.42
1:C:40:GLN:HA	1:C:40:GLN:OE1	2.19	0.42
1:D:12:ARG:HG3	1:D:50:PHE:CE2	2.55	0.42
1:B:22[A]:SER:OG	1:B:26:ARG:NH1	2.53	0.42
1:D:244:ARG:NH1	4:D:508:HOH:O	2.35	0.41
1:D:213:THR:HG22	1:D:243:THR:HG23	2.02	0.41
1:B:205:ALA:HA	1:B:281:VAL:O	2.21	0.41
1:B:50:PHE:CE2	1:B:54:LEU:HD12	2.55	0.40
1:B:125:SER:O	1:B:126:PHE:HB2	2.22	0.40
1:C:240:HIS:CD2	1:C:241:PRO:HD2	2.56	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:537:HOH:O	4:D:597:HOH:O[1_565]	2.19	0.01
4:C:522:HOH:O	4:D:749:HOH:O[1_665]	2.19	0.01

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	363/367 (99%)	354 (98%)	8 (2%)	1 (0%)	41 27
1	B	363/367 (99%)	353 (97%)	9 (2%)	1 (0%)	41 27

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	C	365/367 (100%)	356 (98%)	8 (2%)	1 (0%)	41 27
1	D	365/367 (100%)	353 (97%)	11 (3%)	1 (0%)	41 27
All	All	1456/1468 (99%)	1416 (97%)	36 (2%)	4 (0%)	41 27

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	120	GLU
1	A	120	GLU
1	B	120	GLU
1	D	120	GLU

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	302/304 (99%)	291 (96%)	11 (4%)	35 20
1	B	300/304 (99%)	293 (98%)	7 (2%)	50 37
1	C	300/304 (99%)	296 (99%)	4 (1%)	69 62
1	D	301/304 (99%)	299 (99%)	2 (1%)	84 81
All	All	1203/1216 (99%)	1179 (98%)	24 (2%)	57 44

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	129	ARG
1	A	151[A]	GLU
1	A	151[B]	GLU
1	A	176	ASP
1	A	242	ARG
1	A	257	GLU
1	A	263	GLU

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Mol	Chain	Res	Type
1	A	304	GLU
1	A	311	LEU
1	A	322	GLN
1	B	28	SER
1	B	52	GLU
1	B	94	LEU
1	B	107	LEU
1	B	130	MET
1	B	198	GLN
1	B	335	MET
1	C	26	ARG
1	C	257	GLU
1	C	311	LEU
1	C	322	GLN
1	D	279	LEU
1	D	359	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	GLN
1	B	216	GLN
1	D	41	HIS
1	D	198	GLN

5.3.3 RNA

There are no RNA molecules 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 monosaccharides in this entry.

5.6 Ligand geometry

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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

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, 95^{th} 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	361/367 (98%)	-0.33	4 (1%) 80 78	12, 22, 52, 76	0
1	B	359/367 (97%)	-0.38	3 (0%) 86 84	13, 21, 43, 77	0
1	C	363/367 (98%)	-0.13	17 (4%) 31 25	10, 26, 76, 97	0
1	D	362/367 (98%)	-0.45	2 (0%) 89 87	10, 18, 44, 68	0
All	All	1445/1468 (98%)	-0.32	26 (1%) 68 64	10, 21, 57, 97	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	214	ASP	4.9
1	C	250	PHE	4.8
1	C	217	GLU	4.1
1	C	322	GLN	3.7
1	B	131	PRO	3.6
1	C	242	ARG	3.5
1	C	249	ALA	3.3
1	C	320	ALA	3.2
1	A	45	ASN	3.1
1	C	304	GLU	2.9
1	C	321	PRO	2.8
1	A	43[A]	ASP	2.7
1	D	129	ARG	2.7
1	C	245	LYS	2.6
1	B	130	MET	2.6
1	C	210	ALA	2.5
1	C	303	GLU	2.5
1	C	131	PRO	2.4
1	C	219	LEU	2.3
1	B	129	ARG	2.3
1	C	220	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	218	ARG	2.2
1	A	131	PRO	2.1
1	D	43	ASP	2.1
1	C	221	VAL	2.1
1	A	130	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 monosaccharides in this entry.

6.4 Ligands (i)

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, 95^{th} 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(Å^2)	Q<0.9
2	CL	B	401	1/1	0.92	0.08	41,41,41,41	0
3	NA	B	402	1/1	0.97	0.07	25,25,25,25	0
2	CL	D	401	1/1	0.99	0.04	22,22,22,22	0

6.5 Other polymers (i)

There are no such residues in this entry.