



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:09 PM UTC

PDB ID : 2GYX / pdb_00002gyx
Title : Structure of aspartate semialdehyde dehydrogenase (ASADH) from *Streptococcus pneumoniae*
Authors : Faehnle, C.R.; Le Coq, J.; Liu, X.; Viola, R.E.
Deposited on : 2006-05-10
Resolution : 2.10 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
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

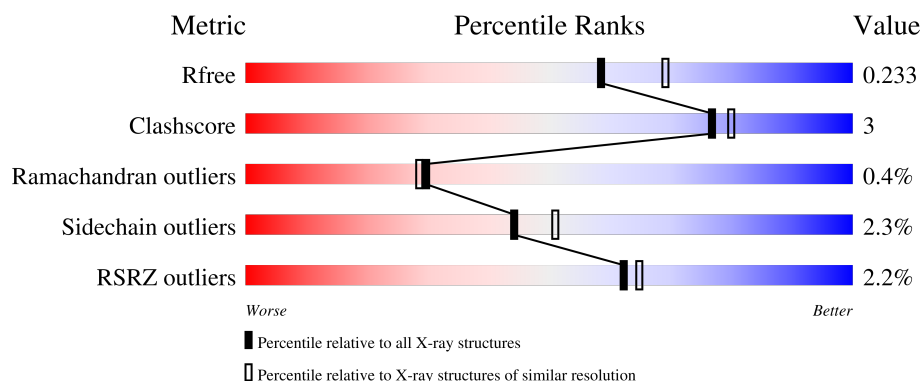
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	366	0% 90% 6% .
1	B	366	0% 85% 8% 7%
1	C	366	2% 88% 8% .
1	D	366	5% 84% 9% 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11565 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 Aspartate beta-semialdehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	352	Total	C	N	O	S	0	0	0
			2684	1699	450	522	13			
1	B	339	Total	C	N	O	S	0	0	0
			2594	1640	434	507	13			
1	C	350	Total	C	N	O	S	0	0	0
			2673	1693	448	519	13			
1	D	342	Total	C	N	O	S	0	0	0
			2611	1650	439	509	13			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	359	LEU	-	cloning artifact	UNP Q8DQ00
A	360	GLU	-	cloning artifact	UNP Q8DQ00
A	361	HIS	-	expression tag	UNP Q8DQ00
A	362	HIS	-	expression tag	UNP Q8DQ00
A	363	HIS	-	expression tag	UNP Q8DQ00
A	364	HIS	-	expression tag	UNP Q8DQ00
A	365	HIS	-	expression tag	UNP Q8DQ00
A	366	HIS	-	expression tag	UNP Q8DQ00
B	359	LEU	-	cloning artifact	UNP Q8DQ00
B	360	GLU	-	cloning artifact	UNP Q8DQ00
B	361	HIS	-	expression tag	UNP Q8DQ00
B	362	HIS	-	expression tag	UNP Q8DQ00
B	363	HIS	-	expression tag	UNP Q8DQ00
B	364	HIS	-	expression tag	UNP Q8DQ00
B	365	HIS	-	expression tag	UNP Q8DQ00
B	366	HIS	-	expression tag	UNP Q8DQ00
C	359	LEU	-	cloning artifact	UNP Q8DQ00
C	360	GLU	-	cloning artifact	UNP Q8DQ00
C	361	HIS	-	expression tag	UNP Q8DQ00
C	362	HIS	-	expression tag	UNP Q8DQ00
C	363	HIS	-	expression tag	UNP Q8DQ00

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Chain	Residue	Modelled	Actual	Comment	Reference
C	364	HIS	-	expression tag	UNP Q8DQ00
C	365	HIS	-	expression tag	UNP Q8DQ00
C	366	HIS	-	expression tag	UNP Q8DQ00
D	359	LEU	-	cloning artifact	UNP Q8DQ00
D	360	GLU	-	cloning artifact	UNP Q8DQ00
D	361	HIS	-	expression tag	UNP Q8DQ00
D	362	HIS	-	expression tag	UNP Q8DQ00
D	363	HIS	-	expression tag	UNP Q8DQ00
D	364	HIS	-	expression tag	UNP Q8DQ00
D	365	HIS	-	expression tag	UNP Q8DQ00
D	366	HIS	-	expression tag	UNP Q8DQ00

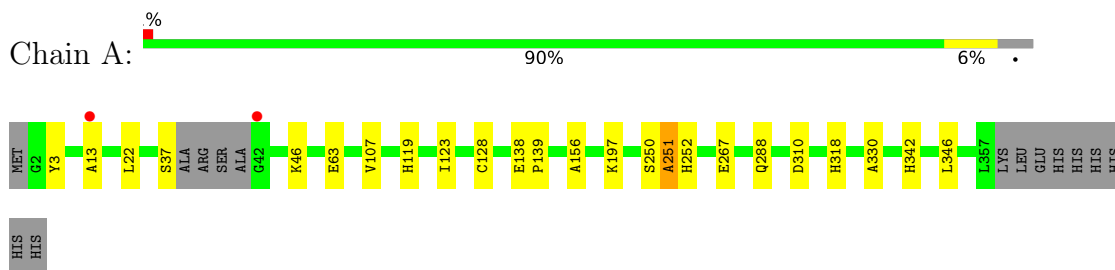
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	296	Total	O	0	0
			296	296		
2	B	232	Total	O	0	0
			232	232		
2	C	242	Total	O	0	0
			242	242		
2	D	233	Total	O	0	0
			233	233		

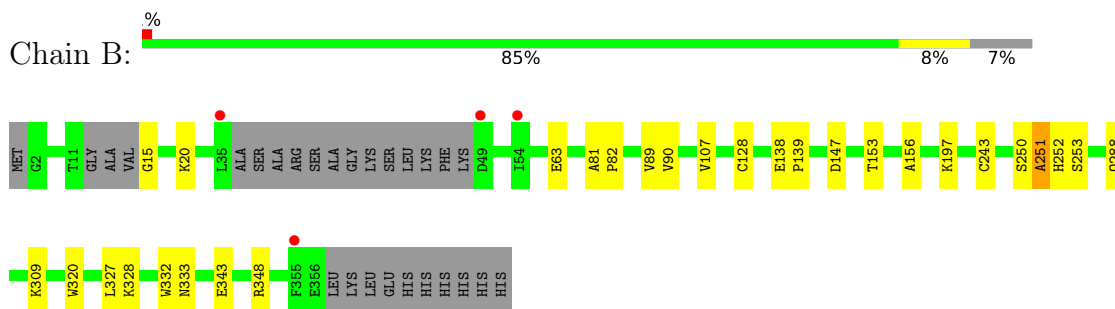
3 Residue-property plots [i](#)

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.

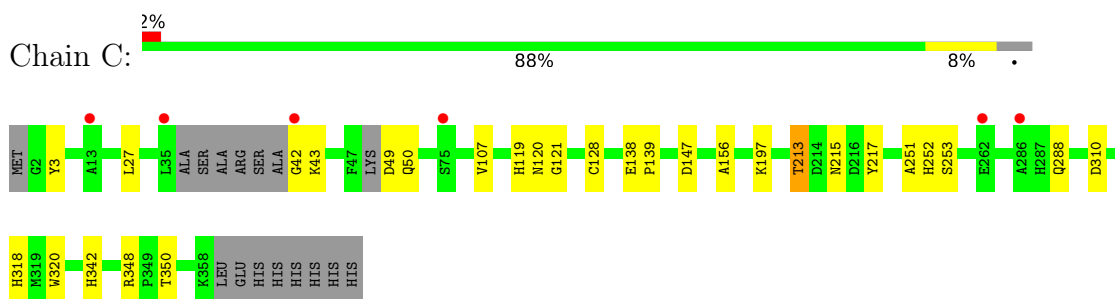
- Molecule 1: Aspartate beta-semialdehyde dehydrogenase



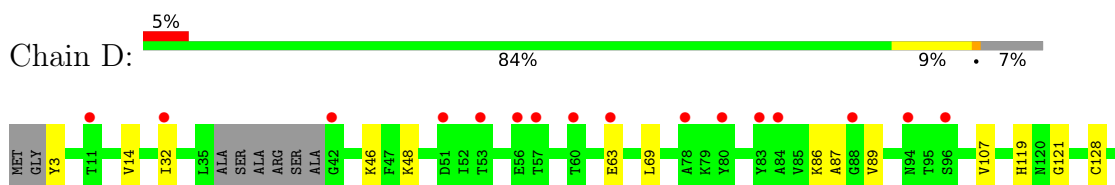
- Molecule 1: Aspartate beta-semialdehyde dehydrogenase

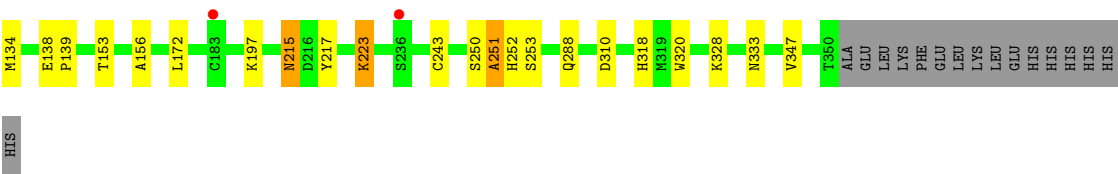


- Molecule 1: Aspartate beta-semialdehyde dehydrogenase



- Molecule 1: Aspartate beta-semialdehyde dehydrogenase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.48Å 78.69Å 243.41Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.11 – 2.10 29.11 – 2.10	Depositor EDS
% Data completeness (in resolution range)	94.6 (29.11-2.10) 94.5 (29.11-2.10)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.75 (at 2.10Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.178 , 0.221 0.189 , 0.233	Depositor DCC
R_{free} test set	3948 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	27.8	Xtriage
Anisotropy	0.246	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 56.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11565	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.56% 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	0.54	0/2732	0.79	2/3711 (0.1%)
1	B	0.50	0/2640	0.78	1/3589 (0.0%)
1	C	0.52	0/2720	0.79	1/3694 (0.0%)
1	D	0.58	4/2658 (0.2%)	0.80	1/3613 (0.0%)
All	All	0.54	4/10750 (0.0%)	0.79	5/14607 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	87	ALA	C-O	7.05	1.33	1.24
1	D	87	ALA	C-N	6.35	1.41	1.33
1	D	89	VAL	C-O	5.84	1.32	1.24
1	D	89	VAL	C-N	5.13	1.40	1.33

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	107	VAL	N-CA-C	5.79	116.22	108.11
1	B	107	VAL	N-CA-C	5.64	116.24	108.12
1	D	107	VAL	N-CA-C	5.53	116.08	108.12
1	A	13	ALA	N-CA-C	-5.24	105.75	111.82
1	A	107	VAL	N-CA-C	5.21	115.41	108.11

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	2684	0	2685	13	0
1	B	2594	0	2580	14	0
1	C	2673	0	2673	19	0
1	D	2611	0	2605	20	0
2	A	296	0	0	1	0
2	B	232	0	0	1	0
2	C	242	0	0	4	0
2	D	233	0	0	1	0
All	All	11565	0	10543	61	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 3.

All (61) 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:C:288:GLN:HE22	1:D:197:LYS:H	1.36	0.73
1:A:197:LYS:H	1:B:288:GLN:HE22	1.38	0.71
1:A:288:GLN:HE22	1:B:197:LYS:H	1.40	0.70
1:B:156:ALA:H	1:B:252:HIS:HD2	1.40	0.69
1:C:197:LYS:H	1:D:288:GLN:HE22	1.40	0.69
1:D:3:TYR:N	2:D:585:HOH:O	2.27	0.67
1:D:156:ALA:H	1:D:252:HIS:HD2	1.42	0.66
1:C:156:ALA:H	1:C:252:HIS:HD2	1.43	0.65
1:A:156:ALA:H	1:A:252:HIS:HD2	1.46	0.61
1:C:350:THR:HG23	2:C:537:HOH:O	2.04	0.58
1:D:128:CYS:SG	1:D:252:HIS:HE1	2.27	0.57
1:D:156:ALA:H	1:D:252:HIS:CD2	2.24	0.56
1:D:119:HIS:HD2	1:D:121:GLY:H	1.53	0.56
1:B:15:GLY:N	2:B:590:HOH:O	2.39	0.55
1:D:328:LYS:NZ	1:D:333:ASN:HD22	2.04	0.55
1:C:3:TYR:OH	1:C:342:HIS:HD2	1.90	0.55
1:C:49:ASP:N	2:C:599:HOH:O	2.40	0.54
1:A:310:ASP:OD1	1:A:318:HIS:HE1	1.91	0.53
1:C:119:HIS:HD2	1:C:121:GLY:H	1.57	0.53
1:C:128:CYS:SG	1:C:252:HIS:HE1	2.31	0.52
1:D:138:GLU:N	1:D:139:PRO:HD2	2.25	0.52
1:B:156:ALA:H	1:B:252:HIS:CD2	2.24	0.51
1:C:120:ASN:C	1:C:350:THR:HG21	2.35	0.51
1:D:310:ASP:OD1	1:D:318:HIS:HE1	1.93	0.51
1:C:156:ALA:H	1:C:252:HIS:CD2	2.26	0.51
1:C:138:GLU:N	1:C:139:PRO:HD2	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:310:ASP:OD1	1:C:318:HIS:HE1	1.92	0.50
1:B:138:GLU:N	1:B:139:PRO:HD2	2.26	0.50
1:D:215:ASN:HD21	1:D:217:TYR:HB2	1.77	0.50
1:A:138:GLU:N	1:A:139:PRO:HD2	2.27	0.49
1:B:128:CYS:SG	1:B:252:HIS:HE1	2.36	0.48
1:B:328:LYS:NZ	1:B:333:ASN:HD22	2.12	0.47
1:C:350:THR:HG22	2:C:410:HOH:O	2.13	0.47
1:D:215:ASN:C	1:D:215:ASN:HD22	2.22	0.47
1:D:215:ASN:ND2	1:D:217:TYR:H	2.12	0.46
1:A:123:ILE:HD12	1:A:123:ILE:N	2.31	0.46
1:D:253:SER:HB3	1:D:320:TRP:CZ2	2.51	0.45
1:C:213:THR:HB	1:C:217:TYR:O	2.16	0.45
1:D:328:LYS:HZ2	1:D:333:ASN:HD22	1.66	0.44
1:A:342:HIS:HE1	1:C:342:HIS:O	2.02	0.43
1:B:81:ALA:N	1:B:82:PRO:CD	2.82	0.42
1:A:128:CYS:SG	1:A:252:HIS:HE1	2.42	0.42
1:C:128:CYS:SG	1:C:252:HIS:CE1	3.12	0.42
1:A:119:HIS:HE1	1:A:346:LEU:O	2.02	0.42
1:B:250:SER:O	1:B:251:ALA:HB3	2.19	0.42
1:D:134:MET:HA	1:D:134:MET:HE2	2.02	0.42
1:C:42:GLY:N	2:C:592:HOH:O	2.53	0.42
1:D:250:SER:O	1:D:251:ALA:HB3	2.20	0.41
1:B:253:SER:HB3	1:B:320:TRP:CZ2	2.55	0.41
1:C:213:THR:HG22	1:C:215:ASN:OD1	2.20	0.41
1:D:3:TYR:CE1	1:D:347:VAL:HG11	2.55	0.41
1:A:3:TYR:OH	1:A:342:HIS:HD2	2.04	0.41
1:A:156:ALA:H	1:A:252:HIS:CD2	2.34	0.41
1:B:327:LEU:C	1:B:332:TRP:HB3	2.46	0.41
1:C:253:SER:HB3	1:C:320:TRP:CZ2	2.55	0.41
1:B:89:VAL:HG12	1:B:90:VAL:N	2.36	0.41
1:B:153:THR:OG1	1:B:243:CYS:HA	2.21	0.40
1:A:37:SER:C	2:A:657:HOH:O	2.64	0.40
1:D:153:THR:OG1	1:D:243:CYS:HA	2.21	0.40
1:A:250:SER:O	1:A:251:ALA:HB3	2.21	0.40
1:D:223:LYS:HB2	1:D:223:LYS:HE2	2.00	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	348/366 (95%)	335 (96%)	11 (3%)	2 (1%)	21	18
1	B	333/366 (91%)	321 (96%)	11 (3%)	1 (0%)	36	36
1	C	344/366 (94%)	330 (96%)	13 (4%)	1 (0%)	36	36
1	D	338/366 (92%)	325 (96%)	12 (4%)	1 (0%)	36	36
All	All	1363/1464 (93%)	1311 (96%)	47 (3%)	5 (0%)	30	28

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	251	ALA
1	A	251	ALA
1	C	251	ALA
1	D	251	ALA
1	A	330	ALA

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	289/302 (96%)	285 (99%)	4 (1%)	59	67
1	B	280/302 (93%)	274 (98%)	6 (2%)	47	54
1	C	288/302 (95%)	282 (98%)	6 (2%)	47	54
1	D	282/302 (93%)	272 (96%)	10 (4%)	32	35
All	All	1139/1208 (94%)	1113 (98%)	26 (2%)	44	51

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LEU
1	A	46	LYS
1	A	63	GLU
1	A	267	GLU
1	B	20	LYS
1	B	63	GLU
1	B	147	ASP
1	B	309	LYS
1	B	343	GLU
1	B	348	ARG
1	C	27	LEU
1	C	43	LYS
1	C	50	GLN
1	C	147	ASP
1	C	213	THR
1	C	348	ARG
1	D	14	VAL
1	D	32	ILE
1	D	46	LYS
1	D	48	LYS
1	D	63	GLU
1	D	69	LEU
1	D	86	LYS
1	D	172	LEU
1	D	215	ASN
1	D	223	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	100	GLN
1	A	119	HIS
1	A	252	HIS
1	A	287	HIS
1	A	288	GLN
1	A	318	HIS
1	A	333	ASN
1	A	342	HIS
1	B	100	GLN
1	B	119	HIS
1	B	169	GLN

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Mol	Chain	Res	Type
1	B	252	HIS
1	B	288	GLN
1	B	318	HIS
1	B	333	ASN
1	B	342	HIS
1	C	50	GLN
1	C	119	HIS
1	C	155	GLN
1	C	169	GLN
1	C	177	ASN
1	C	186	HIS
1	C	198	HIS
1	C	252	HIS
1	C	288	GLN
1	C	318	HIS
1	C	333	ASN
1	C	342	HIS
1	D	119	HIS
1	D	177	ASN
1	D	198	HIS
1	D	215	ASN
1	D	252	HIS
1	D	287	HIS
1	D	288	GLN
1	D	318	HIS
1	D	333	ASN
1	D	342	HIS

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 oligosaccharides 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

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	352/366 (96%)	-0.03	2 (0%) 85 87	18, 26, 38, 45	0
1	B	339/366 (92%)	0.22	4 (1%) 76 78	20, 32, 54, 86	0
1	C	350/366 (95%)	0.11	6 (1%) 69 71	17, 29, 52, 78	0
1	D	342/366 (93%)	0.32	18 (5%) 32 34	17, 32, 73, 102	0
All	All	1383/1464 (94%)	0.15	30 (2%) 62 65	17, 29, 59, 102	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	35	LEU	4.1
1	D	88	GLY	3.9
1	C	42	GLY	3.8
1	D	42	GLY	3.7
1	B	49	ASP	3.4
1	D	60	THR	3.4
1	A	13	ALA	3.1
1	D	57	THR	2.8
1	D	83	TYR	2.8
1	D	94	ASN	2.7
1	D	53	THR	2.7
1	C	286	ALA	2.6
1	B	35	LEU	2.6
1	C	262	GLU	2.5
1	D	56	GLU	2.4
1	B	54	ILE	2.4
1	D	236	SER	2.4
1	C	13	ALA	2.4
1	D	183	CYS	2.3
1	D	78	ALA	2.2
1	D	80	TYR	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	11	THR	2.1
1	D	32	ILE	2.1
1	A	42	GLY	2.1
1	D	84	ALA	2.1
1	D	96	SER	2.1
1	B	355	PHE	2.1
1	D	63	GLU	2.0
1	C	75	SER	2.0
1	D	51	ASP	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.