



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:49 AM UTC

PDB ID : 2QZ9 / pdb_00002qz9
Title : crystal structure of aspartate semialdehyde dehydrogenase II from vibrio cholerae
Authors : Viola, R.E.; Liu, X.; Ohren, J.F.; Faehnle, C.R.
Deposited on : 2007-08-16
Resolution : 2.20 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

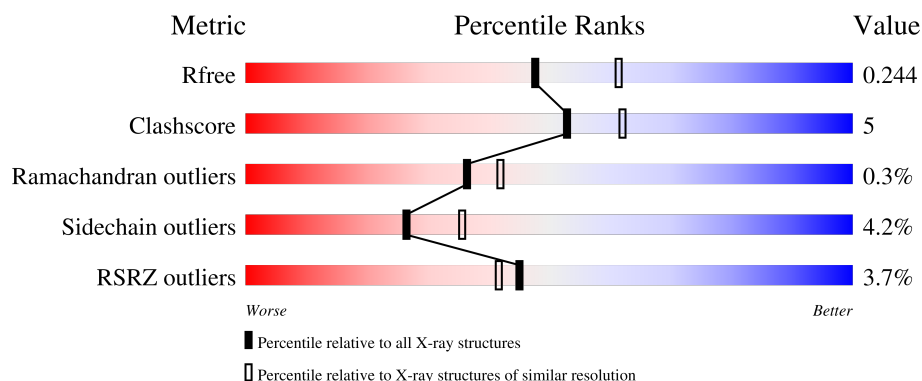
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	336	2% 84% 15%
1	B	336	8% 84% 15%
1	C	336	% 88% 11%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8266 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-semialdehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	336	Total	C	N	O	S	0	2	0
			2617	1655	447	504	11			
1	B	335	Total	C	N	O	S	0	2	0
			2611	1652	446	502	11			
1	C	336	Total	C	N	O	S	0	3	0
			2621	1657	447	506	11			

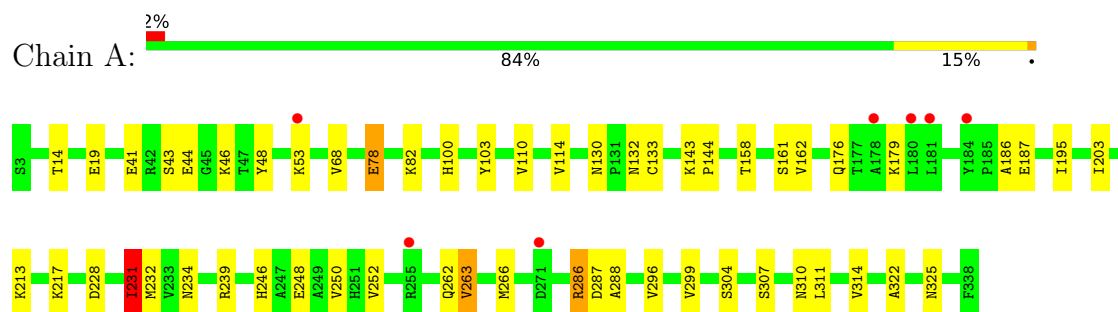
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	153	Total	O	0	0
			153	153		
2	B	103	Total	O	0	0
			103	103		
2	C	161	Total	O	0	0
			161	161		

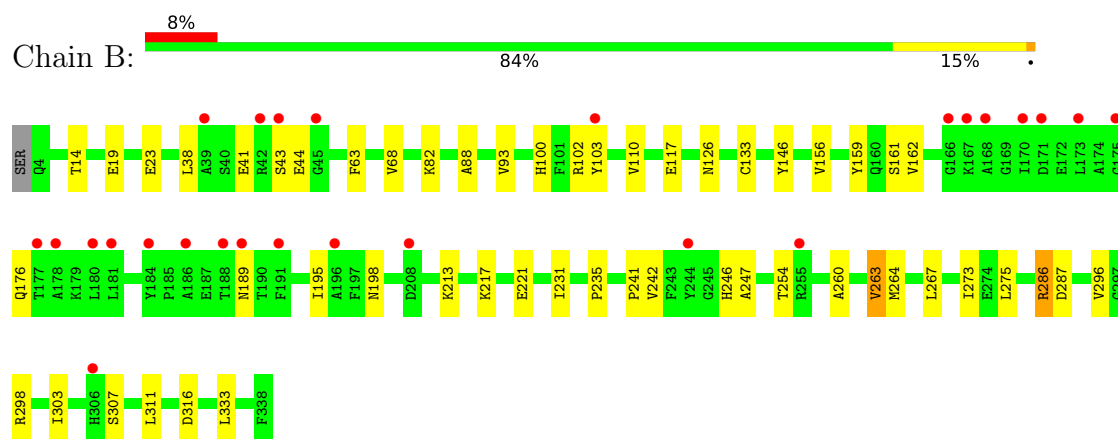
3 Residue-property plots [i](#)

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.

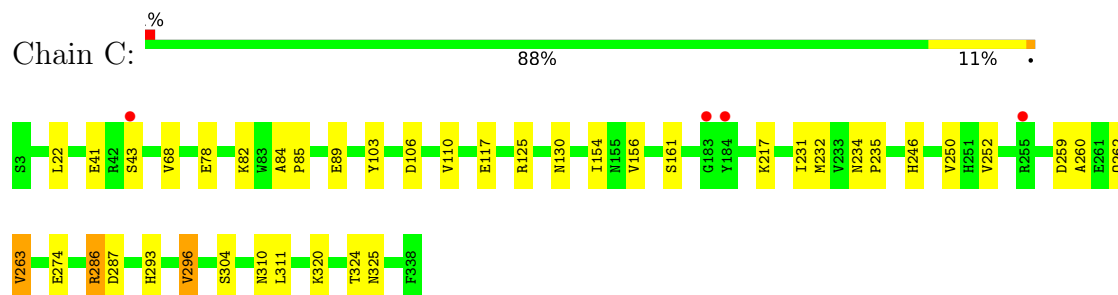
- Molecule 1: Aspartate-semialdehyde dehydrogenase



- Molecule 1: Aspartate-semialdehyde dehydrogenase



- Molecule 1: Aspartate-semialdehyde dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	119.33Å 85.66Å 116.10Å 90.00° 103.69° 90.00°	Depositor
Resolution (Å)	37.61 – 2.20 37.61 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.3 (37.61-2.20) 99.2 (37.61-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.06 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.198 , 0.250 0.195 , 0.244	Depositor DCC
R_{free} test set	2914 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	39.8	Xtriage
Anisotropy	0.013	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 38.7	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	0.96	EDS
Total number of atoms	8266	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.57% 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	0.62	0/2671	0.87	3/3633 (0.1%)
1	B	0.59	0/2665	0.83	2/3625 (0.1%)
1	C	0.61	0/2680	0.86	2/3645 (0.1%)
All	All	0.61	0/8016	0.85	7/10903 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	110	VAL	N-CA-C	6.54	117.27	108.11
1	B	100	HIS	N-CA-C	6.13	117.63	111.07
1	C	110	VAL	N-CA-C	5.96	116.45	108.11
1	A	110	VAL	N-CA-C	5.92	116.40	108.11
1	A	231	ILE	CB-CA-C	-5.25	104.93	111.08
1	C	296	VAL	N-CA-C	5.21	116.22	108.46
1	A	100	HIS	N-CA-C	5.01	117.12	111.11

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	2617	0	2538	33	0
1	B	2611	0	2533	27	0
1	C	2621	0	2540	26	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	153	0	0	3	0
2	B	103	0	0	2	0
2	C	161	0	0	2	0
All	All	8266	0	7611	77	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 (77) 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:234:ASN:HD21	1:C:310:ASN:HD21	1.00	0.93
1:B:103:TYR:OH	1:B:217:LYS:HG2	1.72	0.89
1:A:310:ASN:HD21	1:C:234:ASN:HD21	1.29	0.81
1:A:234:ASN:ND2	1:C:310:ASN:HD21	1.80	0.77
1:B:14:THR:HG21	1:B:43:SER:HB3	1.66	0.77
1:A:133:CYS:HB2	2:A:339:HOH:O	1.85	0.75
1:A:130:ASN:ND2	1:A:325:ASN:HD22	1.88	0.70
1:C:130:ASN:ND2	1:C:325:ASN:HD22	1.90	0.69
1:A:176:GLN:HE22	1:A:195:ILE:H	1.42	0.68
1:B:316:ASP:HB3	2:B:440:HOH:O	1.93	0.67
1:C:130:ASN:HD21	1:C:325:ASN:HD22	1.42	0.66
1:A:310:ASN:HD21	1:C:234:ASN:ND2	1.95	0.64
1:B:102:ARG:HD3	1:B:221:GLU:OE2	1.98	0.64
1:A:103:TYR:OH	1:A:217:LYS:HG2	1.97	0.63
1:A:234:ASN:HD21	1:C:310:ASN:ND2	1.84	0.63
1:A:310:ASN:ND2	1:C:234:ASN:HD21	1.97	0.61
1:A:232:MET:HB3	1:C:304:SER:HB2	1.83	0.61
1:C:78:GLU:H	1:C:78:GLU:CD	2.09	0.60
1:A:286:ARG:HD2	1:A:287:ASP:OD1	2.02	0.59
1:A:228:ASP:HB3	1:A:231:ILE:HD11	1.83	0.59
1:B:176:GLN:HE22	1:B:195:ILE:H	1.50	0.58
1:A:130:ASN:HD21	1:A:325:ASN:HD22	1.50	0.58
1:C:260:ALA:HA	1:C:263:VAL:CG1	2.34	0.57
1:B:254:THR:OG1	1:B:307:SER:HB3	2.05	0.57
1:A:304:SER:HB2	1:C:232:MET:HB3	1.87	0.56
1:A:161:SER:H	1:A:246:HIS:HD2	1.56	0.53
1:B:198:ASN:O	1:B:241:PRO:HD3	2.07	0.53
1:B:156:VAL:O	1:B:235:PRO:HA	2.09	0.53
1:C:286:ARG:HD2	1:C:287:ASP:OD1	2.10	0.52
1:B:286:ARG:HH11	1:B:287:ASP:CG	2.18	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:161:SER:H	1:B:246:HIS:HD2	1.58	0.50
1:A:133:CYS:SG	2:A:492:HOH:O	2.60	0.50
1:B:286:ARG:NH1	1:B:287:ASP:OD1	2.44	0.50
1:B:264:MET:O	1:B:275:LEU:HD22	2.11	0.50
1:B:267:LEU:HD22	1:B:273:ILE:HD13	1.93	0.50
1:C:103:TYR:OH	1:C:217:LYS:HG2	2.11	0.50
1:B:103:TYR:HH	1:B:217:LYS:HG2	1.75	0.50
1:A:161:SER:H	1:A:246:HIS:CD2	2.32	0.48
1:B:146:TYR:HE1	1:B:231:ILE:HD11	1.79	0.47
1:C:274:GLU:HB2	1:C:293:HIS:CD2	2.49	0.47
1:C:246:HIS:HE1	2:C:418:HOH:O	1.98	0.47
1:C:156:VAL:O	1:C:235:PRO:HA	2.15	0.47
1:B:260:ALA:HA	1:B:263:VAL:HG13	1.96	0.47
1:A:162:VAL:HA	1:A:239:ARG:HD2	1.97	0.46
1:B:275:LEU:HD12	1:B:296:VAL:HG23	1.97	0.46
1:C:161:SER:H	1:C:246:HIS:HD2	1.63	0.46
1:A:246:HIS:HE1	2:A:492:HOH:O	1.99	0.46
1:B:159:TYR:HB2	1:B:247:ALA:HB3	1.97	0.46
1:A:14:THR:HG21	1:A:43:SER:HB3	1.97	0.46
1:B:38:LEU:HD21	1:B:63:PHE:CD1	2.51	0.45
1:B:267:LEU:HD13	1:B:296:VAL:HG21	1.97	0.45
1:B:162:VAL:HG22	1:B:242:VAL:O	2.16	0.45
1:B:133:CYS:HB2	2:B:340:HOH:O	2.16	0.45
1:C:250:VAL:HG12	1:C:252:VAL:HG13	1.99	0.45
1:A:263:VAL:HG23	1:A:299:VAL:HG11	1.98	0.45
1:A:78:GLU:HG3	1:B:213:LYS:HG2	1.98	0.44
1:C:320:LYS:HA	1:C:324:THR:HB	2.00	0.44
1:A:176:GLN:HG2	1:A:186:ALA:HB1	2.00	0.44
1:A:250:VAL:HG12	1:A:252:VAL:HG13	2.00	0.44
1:C:89[B]:GLU:HG2	2:C:468:HOH:O	2.17	0.44
1:B:19:GLU:O	1:B:23:GLU:HG3	2.17	0.44
1:A:263:VAL:CG2	1:A:299:VAL:HG11	2.48	0.43
1:A:262:GLN:O	1:A:266:MET:HG3	2.19	0.43
1:C:154:ILE:HG12	1:C:252:VAL:HG12	2.01	0.43
1:A:158:THR:HG22	1:A:248:GLU:HG2	2.01	0.43
1:C:22:LEU:C	1:C:22:LEU:HD23	2.44	0.43
1:A:114:VAL:HG13	1:A:143:LYS:HD3	2.01	0.43
1:C:106:ASP:OD2	1:C:125:ARG:HD3	2.20	0.42
1:C:84:ALA:HB3	1:C:85:PRO:HD3	2.01	0.42
1:A:288:ALA:HB2	1:A:314:VAL:HG12	2.02	0.42
1:A:19:GLU:OE1	1:A:48:TYR:OH	2.32	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:146:TYR:CE1	1:B:231:ILE:HD11	2.54	0.41
1:C:259:ASP:H	1:C:262:GLN:NE2	2.18	0.41
1:B:88:ALA:HB1	1:B:126:ASN:HB2	2.02	0.41
1:A:143:LYS:HB3	1:A:144:PRO:HD3	2.02	0.40
1:B:93:VAL:HG11	1:B:333:LEU:HD21	2.02	0.40
1:A:41:GLU:O	1:A:44:GLU:HB3	2.21	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	334/336 (99%)	318 (95%)	15 (4%)	1 (0%)	36	42
1	B	333/336 (99%)	314 (94%)	17 (5%)	2 (1%)	21	23
1	C	335/336 (100%)	321 (96%)	14 (4%)	0	100	100
All	All	1002/1008 (99%)	953 (95%)	46 (5%)	3 (0%)	36	42

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	41	GLU
1	A	322	ALA
1	B	44	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	279/280 (100%)	263 (94%)	16 (6%)	18	23
1	B	278/280 (99%)	269 (97%)	9 (3%)	34	47
1	C	280/280 (100%)	270 (96%)	10 (4%)	31	42
All	All	837/840 (100%)	802 (96%)	35 (4%)	26	36

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	LYS
1	A	53	LYS
1	A	68	VAL
1	A	78	GLU
1	A	82	LYS
1	A	132	ASN
1	A	179	LYS
1	A	187	GLU
1	A	203	ILE
1	A	213	LYS
1	A	231	ILE
1	A	263	VAL
1	A	286	ARG
1	A	296	VAL
1	A	307	SER
1	A	311	LEU
1	B	68	VAL
1	B	82	LYS
1	B	117	GLU
1	B	189	ASN
1	B	263	VAL
1	B	286	ARG
1	B	298	ARG
1	B	303	ILE
1	B	311	LEU
1	C	41	GLU
1	C	43	SER
1	C	68	VAL
1	C	82	LYS
1	C	117	GLU
1	C	231	ILE
1	C	263	VAL

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Mol	Chain	Res	Type
1	C	286	ARG
1	C	296	VAL
1	C	311	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	67	GLN
1	A	130	ASN
1	A	155	ASN
1	A	176	GLN
1	A	202	GLN
1	A	234	ASN
1	A	246	HIS
1	A	306	HIS
1	B	160	GLN
1	B	176	GLN
1	B	189	ASN
1	B	194	GLN
1	B	202	GLN
1	B	234	ASN
1	B	246	HIS
1	C	67	GLN
1	C	130	ASN
1	C	155	ASN
1	C	160	GLN
1	C	176	GLN
1	C	193	GLN
1	C	234	ASN
1	C	246	HIS
1	C	284	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

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	336/336 (100%)	0.10	7 (2%) 63 60	22, 37, 60, 75	2 (0%)
1	B	335/336 (99%)	0.45	26 (7%) 19 16	29, 44, 75, 90	2 (0%)
1	C	336/336 (100%)	-0.02	4 (1%) 76 74	23, 35, 56, 67	3 (0%)
All	All	1007/1008 (99%)	0.18	37 (3%) 45 42	22, 39, 64, 90	7 (0%)

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	43	SER	4.8
1	B	166	GLY	3.4
1	B	244	TYR	3.3
1	B	178	ALA	3.2
1	B	168	ALA	3.1
1	B	255[A]	ARG	3.1
1	B	173	LEU	3.0
1	B	175	GLY	2.9
1	B	208	ASP	2.9
1	A	180	LEU	2.9
1	A	255[A]	ARG	2.8
1	B	181	LEU	2.7
1	A	181	LEU	2.7
1	B	306	HIS	2.6
1	A	53	LYS	2.6
1	B	103	TYR	2.6
1	B	184	TYR	2.6
1	B	186	ALA	2.5
1	B	191	PHE	2.4
1	B	43	SER	2.4
1	A	184	TYR	2.4
1	C	184	TYR	2.4
1	B	196	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	189	ASN	2.3
1	C	255[A]	ARG	2.3
1	B	188	THR	2.3
1	B	45	GLY	2.3
1	B	171	ASP	2.2
1	A	271	ASP	2.2
1	B	39	ALA	2.2
1	B	42	ARG	2.2
1	B	170	ILE	2.2
1	B	167	LYS	2.2
1	B	180	LEU	2.1
1	B	177	THR	2.0
1	A	178	ALA	2.0
1	C	183	GLY	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.