



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:17 PM UTC

PDB ID : 5J77 / pdb_00005j77
Title : Mutant glyceraldehyde dehydrogenase (F34M+S405N) from *Thermoplasma acidophilum*
Authors : Iermak, I.; Mesters, J.R.; Kuta Smatanova, I.
Deposited on : 2016-04-06
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

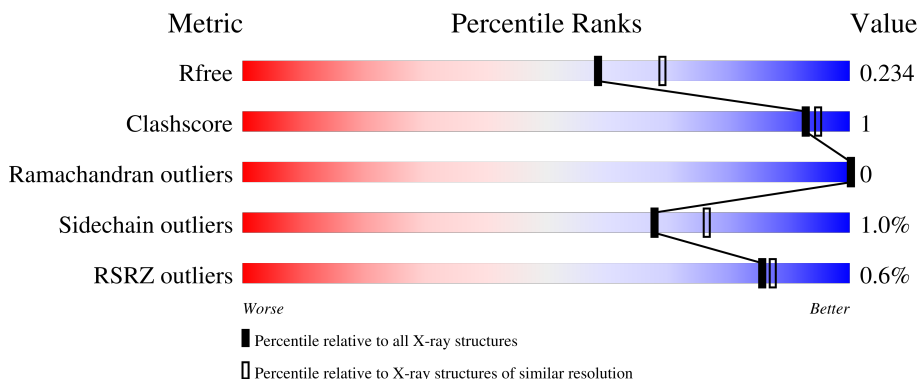
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

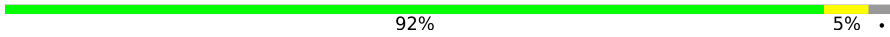
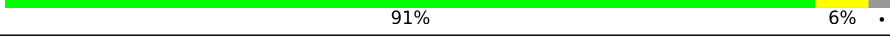
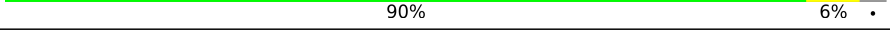
The reported resolution of this entry is 2.10 Å.

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	508	
1	B	508	
1	C	508	
1	D	508	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 15691 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 D-glyceraldehyde dehydrogenase (NADP(+)).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	492	Total	C	N	O	S	0	3	0
			3806	2430	630	730	16			
1	B	492	Total	C	N	O	S	0	5	0
			3818	2436	634	732	16			
1	C	492	Total	C	N	O	S	0	2	0
			3745	2397	610	724	14			
1	D	492	Total	C	N	O	S	0	2	0
			3744	2395	616	718	15			

There are 68 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	34	MET	PHE	engineered mutation	UNP Q9HK01
A	405	ASN	SER	engineered mutation	UNP Q9HK01
A	494	SER	-	expression tag	UNP Q9HK01
A	495	GLY	-	expression tag	UNP Q9HK01
A	496	ARG	-	expression tag	UNP Q9HK01
A	497	PRO	-	expression tag	UNP Q9HK01
A	498	VAL	-	expression tag	UNP Q9HK01
A	499	LEU	-	expression tag	UNP Q9HK01
A	500	GLY	-	expression tag	UNP Q9HK01
A	501	SER	-	expression tag	UNP Q9HK01
A	502	SER	-	expression tag	UNP Q9HK01
A	503	HIS	-	expression tag	UNP Q9HK01
A	504	HIS	-	expression tag	UNP Q9HK01
A	505	HIS	-	expression tag	UNP Q9HK01
A	506	HIS	-	expression tag	UNP Q9HK01
A	507	HIS	-	expression tag	UNP Q9HK01
A	508	HIS	-	expression tag	UNP Q9HK01
B	34	MET	PHE	engineered mutation	UNP Q9HK01
B	405	ASN	SER	engineered mutation	UNP Q9HK01
B	494	SER	-	expression tag	UNP Q9HK01
B	495	GLY	-	expression tag	UNP Q9HK01

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	496	ARG	-	expression tag	UNP Q9HK01
B	497	PRO	-	expression tag	UNP Q9HK01
B	498	VAL	-	expression tag	UNP Q9HK01
B	499	LEU	-	expression tag	UNP Q9HK01
B	500	GLY	-	expression tag	UNP Q9HK01
B	501	SER	-	expression tag	UNP Q9HK01
B	502	SER	-	expression tag	UNP Q9HK01
B	503	HIS	-	expression tag	UNP Q9HK01
B	504	HIS	-	expression tag	UNP Q9HK01
B	505	HIS	-	expression tag	UNP Q9HK01
B	506	HIS	-	expression tag	UNP Q9HK01
B	507	HIS	-	expression tag	UNP Q9HK01
B	508	HIS	-	expression tag	UNP Q9HK01
C	34	MET	PHE	engineered mutation	UNP Q9HK01
C	405	ASN	SER	engineered mutation	UNP Q9HK01
C	494	SER	-	expression tag	UNP Q9HK01
C	495	GLY	-	expression tag	UNP Q9HK01
C	496	ARG	-	expression tag	UNP Q9HK01
C	497	PRO	-	expression tag	UNP Q9HK01
C	498	VAL	-	expression tag	UNP Q9HK01
C	499	LEU	-	expression tag	UNP Q9HK01
C	500	GLY	-	expression tag	UNP Q9HK01
C	501	SER	-	expression tag	UNP Q9HK01
C	502	SER	-	expression tag	UNP Q9HK01
C	503	HIS	-	expression tag	UNP Q9HK01
C	504	HIS	-	expression tag	UNP Q9HK01
C	505	HIS	-	expression tag	UNP Q9HK01
C	506	HIS	-	expression tag	UNP Q9HK01
C	507	HIS	-	expression tag	UNP Q9HK01
C	508	HIS	-	expression tag	UNP Q9HK01
D	34	MET	PHE	engineered mutation	UNP Q9HK01
D	405	ASN	SER	engineered mutation	UNP Q9HK01
D	494	SER	-	expression tag	UNP Q9HK01
D	495	GLY	-	expression tag	UNP Q9HK01
D	496	ARG	-	expression tag	UNP Q9HK01
D	497	PRO	-	expression tag	UNP Q9HK01
D	498	VAL	-	expression tag	UNP Q9HK01
D	499	LEU	-	expression tag	UNP Q9HK01
D	500	GLY	-	expression tag	UNP Q9HK01
D	501	SER	-	expression tag	UNP Q9HK01
D	502	SER	-	expression tag	UNP Q9HK01
D	503	HIS	-	expression tag	UNP Q9HK01

Continued on next page...

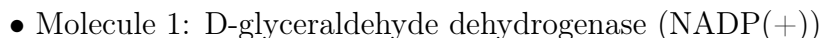
Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	504	HIS	-	expression tag	UNP Q9HK01
D	505	HIS	-	expression tag	UNP Q9HK01
D	506	HIS	-	expression tag	UNP Q9HK01
D	507	HIS	-	expression tag	UNP Q9HK01
D	508	HIS	-	expression tag	UNP Q9HK01

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	225	Total 225	O 225	0	0
2	B	190	Total 190	O 190	0	0
2	C	75	Total 75	O 75	0	0
2	D	88	Total 88	O 88	0	0

- Molecule 1: D-glyceraldehyde dehydrogenase (NADP(+))



[illegible]

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	108.07Å 158.42Å 130.05Å 90.00° 91.57° 90.00°	Depositor
Resolution (Å)	44.63 – 2.10 44.63 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.6 (44.63-2.10) 99.6 (44.63-2.10)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.80 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.176 , 0.226 0.188 , 0.234	Depositor DCC
R_{free} test set	6331 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	31.0	Xtriage
Anisotropy	0.488	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.019 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	15691	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.63% 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

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	1.24	5/3897 (0.1%)	1.10	10/5281 (0.2%)
1	B	1.17	3/3915 (0.1%)	1.02	4/5307 (0.1%)
1	C	1.04	0/3833	1.11	11/5208 (0.2%)
1	D	1.08	4/3832 (0.1%)	1.13	12/5204 (0.2%)
All	All	1.13	12/15477 (0.1%)	1.09	37/21000 (0.2%)

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	247[A]	GLU	CA-C	7.07	1.61	1.53
1	D	247[B]	GLU	CA-C	7.07	1.61	1.53
1	A	247[A]	GLU	CA-C	6.32	1.60	1.52
1	A	247[B]	GLU	CA-C	6.32	1.60	1.52
1	A	7	ILE	C-O	-5.93	1.17	1.24
1	D	39	ARG	CA-C	5.77	1.60	1.52
1	A	122	VAL	CA-C	5.74	1.59	1.52
1	B	30	VAL	C-O	-5.36	1.18	1.24
1	B	115	ALA	C-O	5.23	1.30	1.23
1	A	451	ARG	C-O	-5.15	1.19	1.24
1	B	70	ARG	C-O	-5.09	1.18	1.24
1	D	113	GLU	C-O	5.07	1.30	1.24

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	247[A]	GLU	CA-C-O	14.83	137.60	120.70
1	C	247[B]	GLU	CA-C-O	14.83	137.60	120.70
1	D	247[A]	GLU	CA-C-O	12.11	134.95	121.39
1	D	247[B]	GLU	CA-C-O	12.11	134.95	121.39
1	D	246	LEU	CA-C-N	10.81	138.78	122.77
1	D	246	LEU	C-N-CA	10.81	138.78	122.77
1	A	247[A]	GLU	CA-C-O	10.56	131.75	120.46
1	A	247[B]	GLU	CA-C-O	10.56	131.75	120.46

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	247[A]	GLU	N-CA-C	7.86	123.10	107.62
1	C	247[B]	GLU	N-CA-C	7.86	123.10	107.62
1	D	102	ASP	N-CA-C	-6.99	103.74	111.36
1	D	249	GLY	N-CA-C	6.95	121.43	112.54
1	A	149	ASN	N-CA-C	6.79	118.76	111.36
1	A	262	MET	N-CA-C	6.27	118.63	111.11
1	C	149	ASN	N-CA-C	6.20	118.12	111.36
1	A	405	ASN	N-CA-CB	-6.20	100.10	110.57
1	D	247[A]	GLU	N-CA-C	6.05	119.32	107.57
1	D	247[B]	GLU	N-CA-C	6.05	119.32	107.57
1	A	249	GLY	N-CA-C	6.03	118.96	112.33
1	D	291	GLU	CB-CA-C	-5.92	100.61	110.68
1	A	247[A]	GLU	N-CA-C	5.87	118.38	107.99
1	A	247[B]	GLU	N-CA-C	5.87	118.38	107.99
1	C	249	GLY	N-CA-C	5.78	121.36	112.51
1	B	322	ILE	N-CA-C	5.77	116.50	110.62
1	C	246	LEU	CA-C-N	5.72	131.60	123.14
1	C	246	LEU	C-N-CA	5.72	131.60	123.14
1	B	480	LYS	CA-C-N	-5.54	114.06	119.76
1	B	480	LYS	C-N-CA	-5.54	114.06	119.76
1	A	437	MET	N-CA-C	5.48	113.15	108.22
1	A	244	LEU	N-CA-C	5.39	117.75	109.07
1	C	137	VAL	CA-C-N	-5.29	114.51	119.85
1	C	137	VAL	C-N-CA	-5.29	114.51	119.85
1	D	94	VAL	N-CA-C	5.27	115.99	110.62
1	D	137	VAL	CA-C-N	-5.20	114.60	119.85
1	D	137	VAL	C-N-CA	-5.20	114.60	119.85
1	C	102	ASP	N-CA-C	-5.19	105.71	111.36
1	B	53	PHE	N-CA-C	5.04	117.16	111.11

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	3806	0	3713	10	0
1	B	3818	0	3716	9	0
1	C	3745	0	3583	14	0
1	D	3744	0	3599	14	0
2	A	225	0	0	0	0
2	B	190	0	0	0	1
2	C	75	0	0	0	0
2	D	88	0	0	0	0
All	All	15691	0	14611	44	1

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 1.

All (44) 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:188:PHE:O	1:C:193:VAL:HG23	1.97	0.64
1:D:247[B]:GLU:OE1	1:D:457:GLY:HA2	1.99	0.63
1:C:384:ALA:HB1	1:C:385:PRO:HD2	1.86	0.56
1:D:273:LYS:HE3	1:D:273:LYS:HA	1.89	0.55
1:D:142:VAL:HG22	1:D:220:LEU:HB3	1.88	0.55
1:C:291:GLU:CD	1:C:393:SER:HB3	2.34	0.52
1:A:444:GLN:HB2	1:B:123:VAL:HG11	1.92	0.52
1:D:70:ARG:NH1	1:D:191:ALA:O	2.43	0.51
1:C:34:MET:HE1	1:C:175:SER:O	2.10	0.51
1:B:172:LYS:HA	1:B:201:ILE:O	2.12	0.50
1:D:34:MET:HE1	1:D:175:SER:O	2.12	0.49
1:A:384:ALA:HB1	1:A:385:PRO:HD2	1.95	0.49
1:A:34:MET:HE1	1:A:175:SER:O	2.13	0.48
1:B:247[A]:GLU:HG3	1:B:458:GLU:OE1	2.14	0.48
1:D:172:LYS:HA	1:D:201:ILE:O	2.15	0.47
1:A:258:LYS:HB3	1:A:290:HIS:CG	2.50	0.47
1:A:274:TYR:CE1	1:A:365:PRO:HB3	2.50	0.47
1:A:257:TRP:HD1	1:A:259:ASP:OD1	1.98	0.46
1:C:273:LYS:HA	1:C:273:LYS:HE3	1.98	0.46
1:C:390:ARG:NH1	1:C:391:LYS:O	2.49	0.45
1:A:347:GLY:HA3	1:A:365:PRO:O	2.16	0.45
1:C:444:GLN:HB2	1:D:123:VAL:HG11	1.98	0.45
1:D:414:PHE:CE2	1:D:436:ASN:HA	2.52	0.45
1:B:303:GLU:OE2	1:B:306:ARG:NH1	2.51	0.44
1:A:483:HIS:ND1	1:A:489:ASP:OD2	2.43	0.44
1:B:390[A]:ARG:NH1	1:B:391:LYS:O	2.50	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:414:PHE:CE2	1:B:436:ASN:HA	2.52	0.44
1:C:473:ILE:N	1:C:473:ILE:HD12	2.33	0.44
1:C:123:VAL:HG11	1:D:444:GLN:HB2	1.99	0.43
1:A:273:LYS:HA	1:A:273:LYS:HE3	2.00	0.43
1:C:49:ALA:HA	1:C:196:GLY:O	2.19	0.43
1:C:417:ASP:OD1	1:C:417:ASP:C	2.63	0.42
1:D:384:ALA:HB1	1:D:385:PRO:HD2	2.00	0.42
1:D:15:SER:HB2	1:D:41:ASP:OD1	2.20	0.42
1:C:462:TYR:O	1:C:466:GLU:HG2	2.20	0.42
1:D:414:PHE:HA	1:D:436:ASN:OD1	2.20	0.42
1:B:56:TRP:CE2	1:B:64:ARG:HD2	2.54	0.41
1:C:302:VAL:HG13	1:C:346:PHE:HB2	2.02	0.41
1:B:142:VAL:HG22	1:B:220:LEU:HB3	2.03	0.41
1:B:310:LEU:HD12	1:B:310:LEU:N	2.35	0.41
1:D:449:GLY:HA3	1:D:453:THR:HG21	2.03	0.41
1:D:134:GLN:HA	1:D:471:LYS:O	2.21	0.41
1:C:151:PRO:HB2	1:C:180:SER:OG	2.21	0.41
1:A:247[B]:GLU:HG3	1:A:458:GLU:OE1	2.22	0.40

All (1) 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)
2:B:646:HOH:O	2:B:646:HOH:O[2_555]	2.16	0.04

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	493/508 (97%)	480 (97%)	13 (3%)	0	100	100
1	B	495/508 (97%)	480 (97%)	15 (3%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	492/508 (97%)	479 (97%)	13 (3%)	0	100	100
1	D	492/508 (97%)	476 (97%)	16 (3%)	0	100	100
All	All	1972/2032 (97%)	1915 (97%)	57 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	392/419 (94%)	386 (98%)	6 (2%)	57	65
1	B	392/419 (94%)	389 (99%)	3 (1%)	73	81
1	C	375/419 (90%)	372 (99%)	3 (1%)	73	81
1	D	376/419 (90%)	371 (99%)	5 (1%)	61	69
All	All	1535/1676 (92%)	1518 (99%)	17 (1%)	68	74

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	123	VAL
1	A	247[A]	GLU
1	A	247[B]	GLU
1	A	273	LYS
1	A	306	ARG
1	A	409	LEU
1	B	1	MET
1	B	273	LYS
1	B	409	LEU
1	C	273	LYS
1	C	383	PHE
1	C	461	LYS
1	D	57	ASN
1	D	247[A]	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	247[B]	GLU
1	D	273	LYS
1	D	409	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	13	ASN
1	C	57	ASN
1	C	109	GLN
1	D	57	ASN
1	D	373	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	492/508 (96%)	-0.23	1 (0%) 91 92	18, 37, 53, 69	7 (1%)
1	B	492/508 (96%)	-0.17	2 (0%) 88 90	19, 38, 58, 72	7 (1%)
1	C	492/508 (96%)	0.30	5 (1%) 79 81	27, 53, 69, 83	6 (1%)
1	D	492/508 (96%)	0.21	4 (0%) 82 84	27, 52, 65, 76	6 (1%)
All	All	1968/2032 (96%)	0.03	12 (0%) 85 87	18, 47, 64, 83	26 (1%)

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	392	ILE	3.8
1	C	483	HIS	3.2
1	D	207	GLU	3.1
1	C	352	SER	2.8
1	B	16	SER	2.7
1	A	338	LYS	2.6
1	C	491	LEU	2.5
1	C	1	MET	2.4
1	B	17	GLY	2.3
1	D	249	GLY	2.3
1	D	341	GLY	2.3
1	C	15	SER	2.1

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.