



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:47 AM UTC

PDB ID : 4AHQ / pdb_00004ahq
Title : Crystal Structure of N-acetylneuraminic acid lyase mutant K165C from Staphylococcus aureus
Authors : Timms, N.; Polyakova, A.; Windle, C.L.; Trinh, C.H.; Nelson, A.; Trinh, A.R.; Berry, A.
Deposited on : 2012-02-06
Resolution : 1.95 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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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

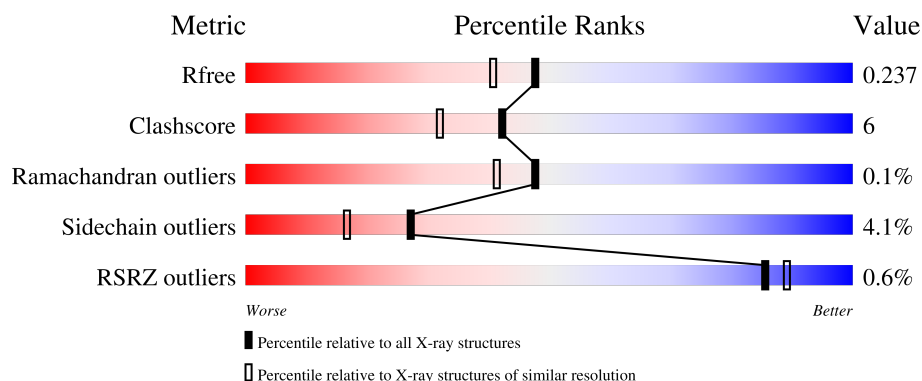
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	298	86% 11% ..
1	B	298	84% 13% ..
1	C	298	89% 8% ..
1	D	298	83% 11% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10090 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 N-ACETYLNEURAMINATE LYASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	293	Total	C	N	O	S	0	1	0
			2319	1490	385	440	4			
1	B	293	Total	C	N	O	S	0	0	0
			2302	1475	384	439	4			
1	C	292	Total	C	N	O	S	0	0	0
			2296	1472	382	438	4			
1	D	291	Total	C	N	O	S	0	2	0
			2305	1481	380	440	4			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	HIS	-	expression tag	UNP Q2G160
A	-3	HIS	-	expression tag	UNP Q2G160
A	-2	HIS	-	expression tag	UNP Q2G160
A	-1	HIS	-	expression tag	UNP Q2G160
A	0	HIS	-	expression tag	UNP Q2G160
A	1	HIS	-	expression tag	UNP Q2G160
A	165	CYS	LYS	engineered mutation	UNP Q2G160
B	-4	HIS	-	expression tag	UNP Q2G160
B	-3	HIS	-	expression tag	UNP Q2G160
B	-2	HIS	-	expression tag	UNP Q2G160
B	-1	HIS	-	expression tag	UNP Q2G160
B	0	HIS	-	expression tag	UNP Q2G160
B	1	HIS	-	expression tag	UNP Q2G160
B	165	CYS	LYS	engineered mutation	UNP Q2G160
C	-4	HIS	-	expression tag	UNP Q2G160
C	-3	HIS	-	expression tag	UNP Q2G160
C	-2	HIS	-	expression tag	UNP Q2G160
C	-1	HIS	-	expression tag	UNP Q2G160
C	0	HIS	-	expression tag	UNP Q2G160
C	1	HIS	-	expression tag	UNP Q2G160
C	165	CYS	LYS	engineered mutation	UNP Q2G160

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-4	HIS	-	expression tag	UNP Q2G160
D	-3	HIS	-	expression tag	UNP Q2G160
D	-2	HIS	-	expression tag	UNP Q2G160
D	-1	HIS	-	expression tag	UNP Q2G160
D	0	HIS	-	expression tag	UNP Q2G160
D	1	HIS	-	expression tag	UNP Q2G160
D	165	CYS	LYS	engineered mutation	UNP Q2G160

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	237	Total O 237 237	0	0
2	B	226	Total O 226 226	0	0
2	C	174	Total O 174 174	0	0
2	D	231	Total O 231 231	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	82.16Å 110.02Å 133.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.27 – 1.95 39.27 – 1.95	Depositor EDS
% Data completeness (in resolution range)	97.5 (39.27-1.95) 97.5 (39.27-1.95)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.59 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.191 , 0.238 0.189 , 0.237	Depositor DCC
R_{free} test set	3476 reflections (4.01%)	wwPDB-VP
Wilson B-factor (Å ²)	26.3	Xtriage
Anisotropy	0.634	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 39.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10090	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.48% 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	0.81	0/2367	0.97	1/3203 (0.0%)
1	B	0.79	0/2345	0.97	2/3174 (0.1%)
1	C	0.78	0/2339	0.98	2/3167 (0.1%)
1	D	0.81	1/2355 (0.0%)	1.01	2/3188 (0.1%)
All	All	0.80	1/9406 (0.0%)	0.98	7/12732 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	261	HIS	CG-CD2	5.60	1.42	1.35

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	251	ILE	N-CA-C	7.72	117.83	110.42
1	B	251	ILE	N-CA-C	6.82	116.91	110.30
1	C	251	ILE	N-CA-C	6.12	116.79	110.36
1	C	229	ILE	N-CA-C	5.49	116.27	110.72
1	D	11	ALA	N-CA-C	-5.35	99.02	108.23
1	B	278	GLU	N-CA-C	5.31	117.81	111.71
1	D	251	ILE	N-CA-C	5.07	115.68	110.36

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	2319	0	2294	29	0
1	B	2302	0	2271	36	0
1	C	2296	0	2264	21	0
1	D	2305	0	2275	35	0
2	A	237	0	0	1	0
2	B	226	0	0	5	0
2	C	174	0	0	1	0
2	D	231	0	0	3	0
All	All	10090	0	9104	115	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 6.

All (115) 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:139:ILE:HB	2:D:2123:HOH:O	1.30	1.29
1:C:215:ARG:HH11	1:C:215:ARG:HG2	1.27	0.99
1:D:215:ARG:HH11	1:D:215:ARG:HG2	1.33	0.92
1:D:33:ASN:HD21	1:D:262:ARG:HH11	1.28	0.82
1:B:274:LYS:HE3	1:C:117:GLU:HG3	1.62	0.81
1:B:138:ALA:O	1:B:167:THR:HB	1.82	0.79
1:B:33:ASN:HD21	1:B:262:ARG:HH11	1.30	0.75
1:B:256:LYS:HE2	2:B:2203:HOH:O	1.90	0.72
1:A:168:ALA:O	1:A:190[A]:PHE:HE2	1.74	0.70
1:B:167:THR:O	1:B:167:THR:CG2	2.38	0.70
1:C:215:ARG:HG2	1:C:215:ARG:NH1	2.05	0.70
1:D:215:ARG:HH11	1:D:215:ARG:CG	2.04	0.69
1:C:33:ASN:HD21	1:C:262:ARG:HH11	1.40	0.68
1:A:117:GLU:H	1:A:117:GLU:CD	2.01	0.67
1:B:280:HIS:HD2	2:B:2094:HOH:O	1.76	0.67
1:D:178:ARG:O	1:D:178:ARG:HD3	1.95	0.67
1:D:107:VAL:HG23	1:D:139:ILE:HD11	1.76	0.66
1:A:168:ALA:O	1:A:190[A]:PHE:CE2	2.48	0.65
1:D:257:GLU:OE2	1:D:260:ARG:NH1	2.30	0.65
1:D:167:THR:O	1:D:190[B]:PHE:CE2	2.50	0.64
1:A:225:ARG:HA	2:A:2182:HOH:O	1.97	0.64
1:C:215:ARG:HH11	1:C:215:ARG:CG	2.08	0.64
1:B:178:ARG:O	1:B:178:ARG:HD3	1.98	0.64
1:C:277:ASN:HD22	1:C:278:GLU:N	1.99	0.61
1:C:178:ARG:NH2	1:C:203:ASP:OD2	2.34	0.59
1:A:107:VAL:HG23	1:A:139:ILE:HD11	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:LYS:HD2	1:A:68:LYS:O	2.04	0.58
1:D:159:GLU:HB3	2:D:2057:HOH:O	2.04	0.56
1:A:25:GLN:H	1:A:25:GLN:CD	2.12	0.56
1:D:178:ARG:HD3	1:D:178:ARG:C	2.29	0.55
1:D:277:ASN:HD22	1:D:277:ASN:C	2.15	0.55
1:B:167:THR:O	1:B:167:THR:HG23	2.05	0.55
1:B:33:ASN:HD21	1:B:262:ARG:NH1	2.02	0.55
1:B:209:THR:HG21	1:B:243:ILE:HG12	1.87	0.55
1:D:68:LYS:HD2	1:D:77:LEU:CD1	2.36	0.54
1:B:178:ARG:NH2	1:B:203:ASP:OD2	2.39	0.54
1:D:68:LYS:HE3	1:D:99:LEU:O	2.08	0.54
1:A:179:LYS:HE2	2:C:2143:HOH:O	2.07	0.54
1:A:153:SER:HA	1:A:181:PHE:CZ	2.43	0.53
1:B:178:ARG:HD3	1:B:178:ARG:C	2.31	0.53
1:D:28:LYS:CB	2:D:2024:HOH:O	2.56	0.53
1:B:175:GLU:OE1	1:B:176:ARG:CD	2.57	0.53
1:A:178:ARG:O	1:A:178:ARG:HD3	2.09	0.53
1:D:221:PHE:O	1:D:225:ARG:HG3	2.09	0.53
1:B:140:PRO:HD2	1:B:167:THR:HG22	1.92	0.51
1:D:277:ASN:HD22	1:D:278:GLU:N	2.08	0.51
1:B:175:GLU:OE1	1:B:176:ARG:HD3	2.11	0.51
1:A:107:VAL:HG23	1:A:139:ILE:CD1	2.41	0.51
1:A:178:ARG:HD3	1:A:178:ARG:C	2.33	0.50
1:D:18:GLU:H	1:D:18:GLU:CD	2.20	0.50
1:A:178:ARG:NH2	1:A:203:ASP:OD2	2.43	0.50
1:A:25:GLN:CD	1:A:25:GLN:N	2.70	0.50
1:C:277:ASN:HD22	1:C:277:ASN:C	2.18	0.49
1:C:175:GLU:OE1	1:C:176:ARG:CD	2.60	0.49
1:B:25:GLN:HB3	2:B:2022:HOH:O	2.12	0.49
1:B:118:GLU:HG2	1:C:274:LYS:HG3	1.95	0.48
1:C:178:ARG:C	1:C:178:ARG:HD3	2.36	0.48
1:C:175:GLU:OE1	1:C:176:ARG:HD2	2.14	0.48
1:C:178:ARG:HD3	1:C:178:ARG:O	2.14	0.47
1:A:79:ALA:HB2	1:A:101:TYR:CD2	2.49	0.47
1:B:167:THR:O	1:B:167:THR:HG22	2.09	0.47
1:A:28:LYS:O	1:A:32:GLN:HG3	2.14	0.47
1:D:215:ARG:CG	1:D:215:ARG:NH1	2.73	0.47
1:B:131:GLN:O	1:B:160:LYS:HE2	2.15	0.47
1:B:58:GLU:H	1:B:58:GLU:CD	2.23	0.46
1:B:51:GLU:OE2	1:B:256:LYS:NZ	2.48	0.46
1:D:178:ARG:NH2	1:D:203:ASP:OD2	2.47	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:SER:HA	1:A:181:PHE:HZ	1.79	0.46
1:B:260:ARG:NH2	1:B:267:GLY:O	2.48	0.46
1:A:56:ASN:OD1	1:A:59:GLN:HG3	2.16	0.45
1:B:175:GLU:OE1	1:B:176:ARG:HD2	2.16	0.45
1:B:25:GLN:CD	1:B:25:GLN:H	2.25	0.45
1:A:175:GLU:OE1	1:A:176:ARG:HD2	2.16	0.45
1:D:209:THR:HG21	1:D:243:ILE:HG12	1.98	0.45
1:C:176:ARG:HD2	1:C:176:ARG:HA	1.76	0.45
1:C:215:ARG:NH1	1:C:215:ARG:CG	2.73	0.45
1:D:33:ASN:HD21	1:D:262:ARG:NH1	2.06	0.45
1:B:139:ILE:HA	1:B:167:THR:HG21	1.98	0.45
1:B:221:PHE:O	1:B:225:ARG:HG3	2.18	0.44
1:D:168:ALA:O	1:D:190[A]:PHE:HE2	2.01	0.44
1:A:175:GLU:OE1	1:A:176:ARG:CD	2.65	0.44
1:A:68:LYS:HD3	1:A:77:LEU:CD1	2.48	0.44
1:A:242:ILE:HA	1:A:287:LEU:HD11	2.00	0.43
1:C:33:ASN:HD21	1:C:262:ARG:NH1	2.10	0.43
1:A:199:ILE:HG21	1:A:229:ILE:HD11	2.00	0.43
1:B:199:ILE:HG12	1:D:199:ILE:HG12	1.99	0.43
1:D:253:PRO:HB3	1:D:269:PRO:HG3	1.99	0.43
1:B:226:GLN:HG3	2:B:2183:HOH:O	2.18	0.43
1:B:277:ASN:ND2	1:B:279:ALA:H	2.16	0.43
1:D:277:ASN:ND2	1:D:279:ALA:H	2.16	0.43
1:B:127:ILE:HG21	1:B:158:HIS:CE1	2.53	0.43
1:A:54:LEU:HD13	1:D:112:TYR:HE2	1.83	0.43
1:D:71:VAL:HB	1:D:75:VAL:HG21	2.00	0.42
1:A:54:LEU:HD13	1:D:112:TYR:CE2	2.55	0.42
1:A:178:ARG:HD3	1:A:178:ARG:HA	1.73	0.42
1:B:217:ALA:HA	1:B:220:ILE:HD12	2.01	0.42
1:C:79:ALA:HB2	1:C:101:TYR:CD2	2.54	0.42
1:B:112:TYR:CE2	1:C:54:LEU:HD13	2.55	0.42
1:C:175:GLU:OE1	1:C:176:ARG:HD3	2.20	0.41
1:D:246:VAL:HG13	1:D:251:ILE:HG13	2.01	0.41
1:D:246:VAL:HG22	1:D:254:THR:HG22	2.02	0.41
1:D:277:ASN:C	1:D:277:ASN:ND2	2.75	0.41
1:C:278:GLU:OE1	1:C:281:ARG:HB2	2.21	0.41
1:B:178:ARG:HD3	1:B:178:ARG:HA	1.83	0.41
1:A:35:ILE:O	1:A:39:GLU:HA	2.21	0.41
1:B:65:LYS:O	1:B:69:GLU:HG3	2.21	0.41
1:C:241:ASP:OD2	1:C:291:TYR:OH	2.35	0.41
1:D:156:PHE:HB3	1:D:184:LYS:HE3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:51:GLU:HB3	1:A:54:LEU:HD22	2.03	0.41
1:D:68:LYS:HD2	1:D:77:LEU:HD12	2.03	0.41
1:D:176:ARG:HA	1:D:176:ARG:HD2	1.58	0.41
1:B:131:GLN:HG2	2:B:2122:HOH:O	2.20	0.40
1:D:4:ASP:OD1	1:D:6:LYS:HG3	2.21	0.40
1:A:117:GLU:CD	1:A:117:GLU:N	2.75	0.40
1:B:156:PHE:HB3	1:B:184:LYS:HE3	2.03	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	292/298 (98%)	290 (99%)	2 (1%)	0	100	100
1	B	291/298 (98%)	286 (98%)	5 (2%)	0	100	100
1	C	290/298 (97%)	283 (98%)	6 (2%)	1 (0%)	36	28
1	D	291/298 (98%)	289 (99%)	2 (1%)	0	100	100
All	All	1164/1192 (98%)	1148 (99%)	15 (1%)	1 (0%)	48	41

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	111	TYR

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	244/254 (96%)	233 (96%)	11 (4%)	24	14
1	B	241/254 (95%)	234 (97%)	7 (3%)	37	29
1	C	241/254 (95%)	230 (95%)	11 (5%)	24	13
1	D	243/254 (96%)	231 (95%)	12 (5%)	22	11
All	All	969/1016 (95%)	928 (96%)	41 (4%)	27	16

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	GLN
1	A	54	LEU
1	A	117	GLU
1	A	139	ILE
1	A	141	ASP
1	A	145	VAL
1	A	162	VAL
1	A	178	ARG
1	A	231	GLU
1	A	246	VAL
1	A	262	ARG
1	B	25	GLN
1	B	139	ILE
1	B	141	ASP
1	B	167	THR
1	B	178	ARG
1	B	246	VAL
1	B	261	HIS
1	C	2	ASN
1	C	117	GLU
1	C	128	GLU
1	C	139	ILE
1	C	141	ASP
1	C	145	VAL
1	C	149	ILE
1	C	150	GLU
1	C	178	ARG
1	C	215	ARG
1	C	230	GLN
1	D	18	GLU

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Mol	Chain	Res	Type
1	D	65	LYS
1	D	139	ILE
1	D	159	GLU
1	D	162	VAL
1	D	167	THR
1	D	176	ARG
1	D	190[A]	PHE
1	D	190[B]	PHE
1	D	215	ARG
1	D	246	VAL
1	D	261	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	228	GLN
1	A	277	ASN
1	A	280	HIS
1	B	33	ASN
1	B	261	HIS
1	B	277	ASN
1	B	280	HIS
1	C	25	GLN
1	C	33	ASN
1	C	52	ASN
1	C	131	GLN
1	C	228	GLN
1	C	277	ASN
1	C	280	HIS
1	D	33	ASN
1	D	277	ASN

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 [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	293/298 (98%)	-0.27	2 (0%) 84 88	16, 30, 42, 74	1 (0%)
1	B	293/298 (98%)	-0.26	4 (1%) 73 79	21, 29, 40, 61	0
1	C	292/298 (97%)	-0.17	1 (0%) 90 92	25, 32, 47, 82	0
1	D	291/298 (97%)	-0.43	0 100 100	13, 27, 37, 69	2 (0%)
All	All	1169/1192 (98%)	-0.28	7 (0%) 85 89	13, 29, 43, 82	3 (0%)

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	HIS	3.9
1	B	1	HIS	3.3
1	A	2	ASN	3.3
1	B	190	PHE	3.3
1	C	2	ASN	2.3
1	B	183	ASP	2.1
1	B	2	ASN	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.