



# Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 12:38 pm BST

PDB ID : 6IKO  
Title : Crystal structure of mouse GAS7cb  
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Deposited on : 2018-10-16  
Resolution : 3.76 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](mailto: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.

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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	:	2.11
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)	:	2.11

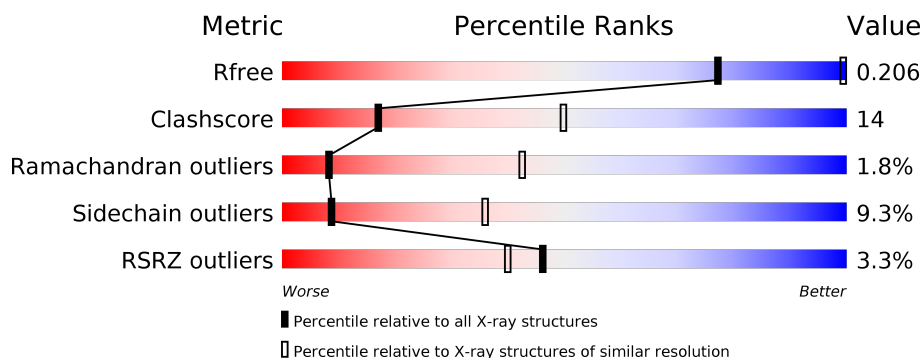
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.76 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	1039 (3.94-3.58)
Clashscore	141614	1051 (3.92-3.60)
Ramachandran outliers	138981	1015 (3.92-3.60)
Sidechain outliers	138945	1011 (3.92-3.60)
RSRZ outliers	127900	1050 (3.96-3.56)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	496	<div> <div> <div></div> <div>41%</div> <div>16%</div> <div>•</div> <div>39%</div> </div> </div>
1	B	496	<div> <div> <div>3%</div> <div>37%</div> <div>20%</div> <div>•</div> <div>39%</div> </div> </div>

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 4987 atoms, of which 0 are hydrogens and 0 are deuteriums.

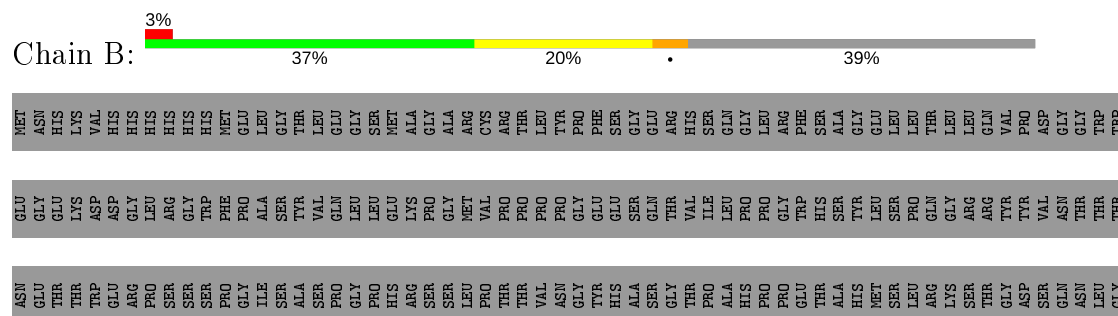
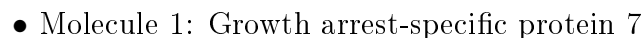
In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

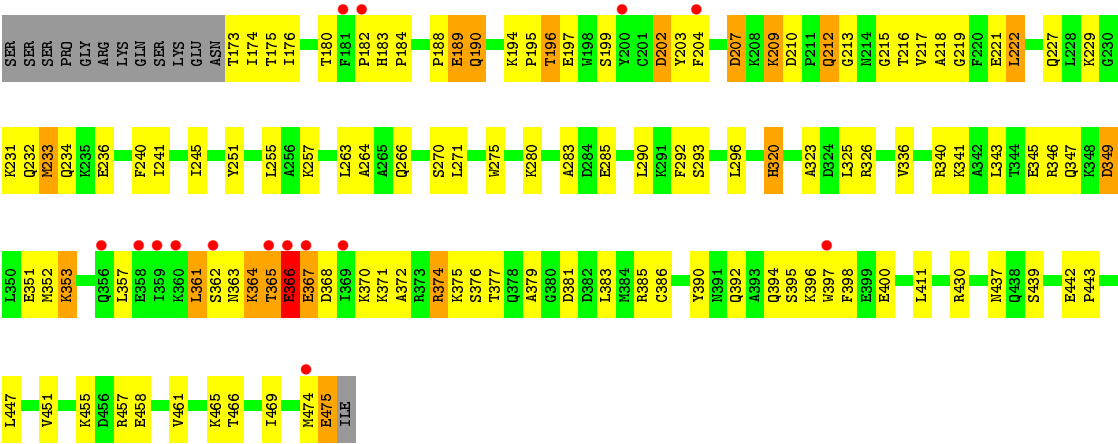
- Molecule 1 is a protein called Growth arrest-specific protein 7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	303	Total	C	N	O	S	0	0	0
			2493	1556	444	477	16			
1	B	303	Total	C	N	O	S	0	0	0
			2494	1557	443	478	16			



- Molecule 1: Growth arrest-specific protein 7





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	194.18Å 194.18Å 194.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.77 – 3.76 48.55 – 3.76	Depositor EDS
% Data completeness (in resolution range)	99.8 (45.77-3.76) 99.9 (48.55-3.76)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.76 (at 3.77Å)	Xtriage
Refinement program	PHENIX 1.8.2_1309	Depositor
R, $R_{free}$	0.173 , 0.205 0.175 , 0.206	Depositor DCC
$R_{free}$ test set	2547 reflections (10.13%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	185.9	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 117.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.019 for l,-k,h	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4987	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	155.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.85% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.60	0/2534	0.79	2/3400 (0.1%)
1	B	0.56	0/2535	0.74	0/3401
All	All	0.58	0/5069	0.76	2/6801 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	295	LYS	CD-CE-NZ	5.69	124.78	111.70
1	A	383	LEU	CA-CB-CG	-5.23	103.28	115.30

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	2493	0	2496	71	0
1	B	2494	0	2496	81	0
All	All	4987	0	4992	140	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 14.

All (140) 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:B:174:ILE:HG12	1:B:175:THR:H	1.49	0.77
1:A:222:LEU:H	1:A:222:LEU:HD12	1.55	0.72
1:A:196:THR:OG1	1:A:197:GLU:N	2.23	0.71
1:A:266:GLN:HB2	1:B:233:MET:HE2	1.74	0.69
1:B:196:THR:OG1	1:B:197:GLU:N	2.26	0.68
1:A:336:VAL:HG11	1:B:474:MET:HE1	1.74	0.68
1:A:346:ARG:HE	1:A:375:LYS:HB3	1.59	0.67
1:B:381:ASP:O	1:B:385:ARG:NH1	2.27	0.67
1:A:209:LYS:HA	1:A:216:THR:HA	1.78	0.65
1:B:222:LEU:HD12	1:B:222:LEU:H	1.61	0.65
1:A:285:GLU:HG3	1:A:437:ASN:OD1	1.98	0.64
1:B:180:THR:HG23	1:B:390:TYR:OH	1.99	0.63
1:A:357:LEU:HG	1:A:361:LEU:HD12	1.81	0.63
1:B:285:GLU:HG3	1:B:437:ASN:OD1	2.00	0.62
1:B:241:ILE:O	1:B:245:ILE:HG13	2.00	0.62
1:B:188:PRO:O	1:B:190:GLN:N	2.34	0.61
1:B:209:LYS:HA	1:B:216:THR:HA	1.82	0.61
1:B:264:ALA:HB3	1:B:275:TRP:NE1	2.16	0.60
1:B:218:ALA:HB3	1:B:221:GLU:HG2	1.83	0.60
1:A:466:THR:OG1	1:A:467:GLY:N	2.30	0.60
1:B:174:ILE:HG12	1:B:175:THR:N	2.16	0.60
1:A:274:ALA:HB1	1:A:448:LEU:HD23	1.83	0.60
1:B:365:THR:O	1:B:367:GLU:N	2.35	0.60
1:A:207:ASP:N	1:A:207:ASP:OD1	2.35	0.59
1:B:353:LYS:HG2	1:B:368:ASP:O	2.01	0.59
1:A:371:LYS:HA	1:A:374:ARG:NH1	2.18	0.58
1:A:188:PRO:O	1:A:190:GLN:N	2.37	0.57
1:B:371:LYS:O	1:B:374:ARG:HB2	2.05	0.56
1:B:370:LYS:O	1:B:374:ARG:NH1	2.39	0.56
1:A:369:ILE:O	1:A:373:ARG:HB3	2.06	0.56
1:A:428:GLN:NE2	1:A:432:GLU:OE1	2.37	0.55
1:A:176:ILE:HG13	1:A:176:ILE:O	2.05	0.55
1:B:346:ARG:HD3	1:B:375:LYS:HB3	1.89	0.55
1:B:381:ASP:O	1:B:385:ARG:HD3	2.07	0.55
1:A:174:ILE:HG12	1:A:175:THR:H	1.73	0.54
1:B:207:ASP:OD2	1:B:326:ARG:NH2	2.40	0.54
1:B:395:SER:HA	1:B:398:PHE:HB3	1.88	0.54
1:A:241:ILE:O	1:A:245:ILE:HG13	2.07	0.54
1:B:251:TYR:CE1	1:B:255:LEU:HD11	2.42	0.54
1:A:245:ILE:HG23	1:A:293:SER:HB2	1.90	0.54
1:A:442:GLU:N	1:A:443:PRO:HD2	2.22	0.53
1:B:216:THR:O	1:B:216:THR:OG1	2.22	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:279:LYS:HG2	1:B:240:PHE:CZ	2.45	0.52
1:A:292:PHE:CZ	1:A:296:LEU:HD11	2.45	0.52
1:A:207:ASP:OD2	1:A:326:ARG:NH2	2.43	0.52
1:A:346:ARG:HD3	1:A:375:LYS:O	2.10	0.52
1:A:251:TYR:CE1	1:A:255:LEU:HD11	2.44	0.52
1:A:216:THR:O	1:A:216:THR:OG1	2.18	0.51
1:A:354:THR:HG22	1:A:369:ILE:HD13	1.91	0.51
1:A:418:MET:HE3	1:A:419:ILE:HG13	1.91	0.51
1:B:183:HIS:ND1	1:B:184:PRO:HD2	2.26	0.51
1:A:472:VAL:HG13	1:B:182:PRO:HD3	1.92	0.51
1:B:207:ASP:N	1:B:207:ASP:OD1	2.44	0.51
1:B:207:ASP:HB3	1:B:217:VAL:O	2.10	0.51
1:B:270:SER:OG	1:B:451:VAL:HA	2.11	0.50
1:B:343:LEU:O	1:B:347:GLN:HG3	2.11	0.49
1:B:437:ASN:C	1:B:439:SER:H	2.15	0.49
1:A:264:ALA:HB3	1:A:275:TRP:NE1	2.27	0.49
1:A:354:THR:HA	1:A:369:ILE:HD11	1.93	0.49
1:B:442:GLU:N	1:B:443:PRO:HD2	2.28	0.49
1:A:461:VAL:O	1:A:465:LYS:HB3	2.12	0.49
1:A:371:LYS:HA	1:A:374:ARG:HH11	1.78	0.49
1:A:222:LEU:CD1	1:A:222:LEU:H	2.14	0.49
1:A:175:THR:OG1	1:A:180:THR:HG22	2.13	0.48
1:B:341:LYS:HE3	1:B:341:LYS:HB3	1.64	0.48
1:B:202:ASP:HB3	1:B:203:TYR:CD2	2.48	0.48
1:B:292:PHE:CZ	1:B:296:LEU:HD11	2.47	0.48
1:A:455:LYS:HB2	1:A:455:LYS:NZ	2.29	0.48
1:A:263:LEU:HD22	1:B:236:GLU:HB3	1.95	0.47
1:A:383:LEU:HG	1:A:383:LEU:O	2.14	0.47
1:A:463:GLU:HG2	1:A:463:GLU:O	2.14	0.47
1:B:347:GLN:HG2	1:B:376:SER:OG	2.15	0.47
1:B:204:PHE:HB3	1:B:219:GLY:O	2.15	0.47
1:A:257:LYS:HB2	1:A:257:LYS:HE3	1.70	0.47
1:A:416:VAL:HG13	1:B:271:LEU:HD13	1.97	0.47
1:A:202:ASP:HB3	1:A:203:TYR:CD2	2.49	0.47
1:A:183:HIS:ND1	1:A:184:PRO:HD2	2.31	0.46
1:A:251:TYR:CZ	1:A:255:LEU:HD11	2.50	0.46
1:B:349:ASP:HB3	1:B:372:ALA:O	2.15	0.46
1:A:351:GLU:HG3	1:A:352:MET:N	2.29	0.46
1:B:293:SER:O	1:B:296:LEU:HB2	2.16	0.46
1:A:233:MET:HE2	1:B:266:GLN:HB2	1.98	0.45
1:A:275:TRP:NE1	1:A:279:LYS:HE2	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:357:LEU:HB2	1:A:369:ILE:HD11	1.98	0.45
1:B:455:LYS:O	1:B:458:GLU:HB3	2.17	0.45
1:A:371:LYS:HE3	1:A:371:LYS:HB3	1.76	0.44
1:A:207:ASP:HB3	1:A:217:VAL:O	2.18	0.44
1:A:293:SER:O	1:A:296:LEU:HB2	2.17	0.44
1:B:212:GLN:HB3	1:B:213:GLY:H	1.49	0.44
1:B:229:LYS:O	1:B:232:GLN:HB3	2.16	0.44
1:A:285:GLU:OE2	1:B:430:ARG:NH1	2.51	0.44
1:A:370:LYS:O	1:A:374:ARG:HD3	2.17	0.44
1:B:357:LEU:O	1:B:361:LEU:HB2	2.18	0.44
1:A:222:LEU:N	1:A:222:LEU:HD12	2.29	0.44
1:A:426:TYR:OH	1:A:430:ARG:HD3	2.17	0.44
1:B:190:GLN:HA	1:B:190:GLN:HE21	1.83	0.44
1:B:394:GLN:C	1:B:396:LYS:H	2.21	0.44
1:B:280:LYS:O	1:B:283:ALA:N	2.51	0.43
1:B:340:ARG:HA	1:B:383:LEU:HD13	2.00	0.43
1:B:461:VAL:O	1:B:465:LYS:HB3	2.18	0.43
1:A:280:LYS:O	1:A:283:ALA:N	2.50	0.43
1:B:437:ASN:O	1:B:439:SER:N	2.51	0.43
1:B:257:LYS:HE3	1:B:257:LYS:HB2	1.82	0.43
1:A:234:GLN:OE1	1:A:307:PHE:HB3	2.19	0.43
1:A:457:ARG:HH21	1:B:197:GLU:HB2	1.84	0.43
1:B:196:THR:HG1	1:B:197:GLU:H	1.62	0.43
1:A:236:GLU:HB3	1:B:263:LEU:HD22	2.01	0.43
1:A:253:LYS:HB3	1:A:253:LYS:HE2	1.90	0.43
1:B:475:GLU:H	1:B:475:GLU:HG2	1.67	0.43
1:B:352:MET:O	1:B:353:LYS:NZ	2.36	0.42
1:A:234:GLN:HG2	1:A:304:LEU:O	2.19	0.42
1:A:346:ARG:NE	1:A:375:LYS:HB3	2.32	0.42
1:A:180:THR:HG23	1:A:390:TYR:OH	2.19	0.42
1:B:466:THR:O	1:B:466:THR:OG1	2.38	0.42
1:B:194:LYS:HA	1:B:195:PRO:HD3	1.73	0.41
1:B:364:LYS:HD3	1:B:366:GLU:OE1	2.19	0.41
1:B:343:LEU:HA	1:B:379:ALA:HB1	2.01	0.41
1:B:263:LEU:HD23	1:B:263:LEU:HA	1.85	0.41
1:A:197:GLU:HB2	1:B:457:ARG:HH21	1.86	0.41
1:B:210:ASP:HB2	1:B:215:GLY:O	2.20	0.41
1:B:325:LEU:HB3	1:B:397:TRP:HB2	2.03	0.41
1:B:336:VAL:HG22	1:B:386:CYS:HB3	2.02	0.41
1:A:268:GLU:OE1	1:B:229:LYS:NZ	2.46	0.41
1:A:337:GLU:C	1:A:339:ALA:H	2.23	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:394:GLN:C	1:A:396:LYS:H	2.25	0.41
1:B:361:LEU:HD23	1:B:362:SER:OG	2.21	0.41
1:A:212:GLN:HB3	1:A:213:GLY:H	1.60	0.41
1:B:357:LEU:HD13	1:B:368:ASP:OD2	2.21	0.41
1:B:343:LEU:HA	1:B:379:ALA:CB	2.51	0.40
1:B:363:ASN:C	1:B:365:THR:H	2.25	0.40
1:B:381:ASP:HB3	1:B:385:ARG:NH1	2.36	0.40
1:B:411:LEU:HD23	1:B:411:LEU:HA	1.84	0.40
1:A:187:MET:HA	1:A:188:PRO:HD2	1.92	0.40
1:B:326:ARG:HG2	1:B:397:TRP:NE1	2.36	0.40
1:B:437:ASN:C	1:B:439:SER:N	2.75	0.40
1:A:190:GLN:HA	1:A:190:GLN:HE21	1.85	0.40
1:A:302:LYS:N	1:A:303:PRO:HD2	2.36	0.40
1:B:231:LYS:O	1:B:234:GLN:HB3	2.21	0.40
1:B:290:LEU:HD12	1:B:293:SER:OG	2.21	0.40
1:B:320:HIS:HA	1:B:323:ALA:HB3	2.02	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	301/496 (61%)	261 (87%)	35 (12%)	5 (2%)	9	43
1	B	301/496 (61%)	250 (83%)	45 (15%)	6 (2%)	7	41
All	All	602/992 (61%)	511 (85%)	80 (13%)	11 (2%)	8	42

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	189	GLU
1	A	212	GLN

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Mol	Chain	Res	Type
1	B	189	GLU
1	B	212	GLN
1	B	366	GLU
1	A	361	LEU
1	B	364	LYS
1	B	196	THR
1	B	367	GLU
1	A	196	THR
1	A	211	PRO

### 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	275/438 (63%)	250 (91%)	25 (9%)	9	37
1	B	275/438 (63%)	249 (90%)	26 (10%)	8	35
All	All	550/876 (63%)	499 (91%)	51 (9%)	9	36

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	176	ILE
1	A	189	GLU
1	A	190	GLN
1	A	202	ASP
1	A	207	ASP
1	A	209	LYS
1	A	216	THR
1	A	222	LEU
1	A	227	GLN
1	A	228	LEU
1	A	233	MET
1	A	268	GLU
1	A	349	ASP
1	A	351	GLU

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Mol	Chain	Res	Type
1	A	361	LEU
1	A	365	THR
1	A	367	GLU
1	A	374	ARG
1	A	377	THR
1	A	385	ARG
1	A	392	GLN
1	A	410	ARG
1	A	442	GLU
1	A	447	LEU
1	A	455	LYS
1	B	173	THR
1	B	176	ILE
1	B	189	GLU
1	B	190	GLN
1	B	199	SER
1	B	202	ASP
1	B	207	ASP
1	B	209	LYS
1	B	222	LEU
1	B	227	GLN
1	B	233	MET
1	B	320	HIS
1	B	345	GLU
1	B	349	ASP
1	B	351	GLU
1	B	353	LYS
1	B	361	LEU
1	B	365	THR
1	B	366	GLU
1	B	374	ARG
1	B	377	THR
1	B	392	GLN
1	B	400	GLU
1	B	447	LEU
1	B	469	ILE
1	B	475	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	227	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 carbohydrates 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	303/496 (61%)	0.07	5 (1%) 70 65	94, 139, 209, 266	0
1	B	303/496 (61%)	0.35	15 (4%) 28 26	91, 150, 259, 300	0
All	All	606/992 (61%)	0.21	20 (3%) 46 40	91, 144, 249, 300	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	474	MET	5.1
1	B	359	ILE	3.4
1	B	358	GLU	3.4
1	A	200	TYR	3.3
1	B	369	ILE	3.3
1	B	362	SER	3.2
1	B	200	TYR	3.1
1	B	356	GLN	3.1
1	B	367	GLU	2.7
1	B	182	PRO	2.5
1	B	204	PHE	2.5
1	A	182	PRO	2.3
1	B	474	MET	2.3
1	B	366	GLU	2.2
1	B	365	THR	2.1
1	A	473	ASP	2.1
1	A	471	PRO	2.1
1	B	397	TRP	2.0
1	B	360	LYS	2.0
1	B	181	PHE	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 carbohydrates 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.