



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2018 – 12:15 pm GMT

PDB ID : 3S3P  
Title : Transglutaminase 2 in complex with a novel inhibitor  
Authors : Lindemann, I.; Heine, A.; Klebe, G.  
Deposited on : 2011-05-18  
Resolution : 2.50 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  
Mogul : 1.7.3 (157068), CSD as539be (2018)  
Xtriage (Phenix) : 1.13  
EDS : trunk31020  
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)  
Refmac : 5.8.0158  
CCP4 : 7.0 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk31020

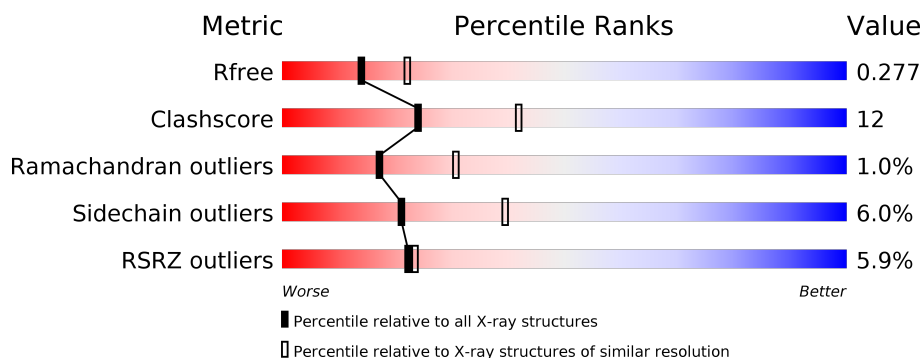
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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))
$R_{free}$	111664	4155 (2.50-2.50)
Clashscore	122126	4827 (2.50-2.50)
Ramachandran outliers	120053	4735 (2.50-2.50)
Sidechain outliers	120020	4737 (2.50-2.50)
RSRZ outliers	108989	4058 (2.50-2.50)

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. 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	694	<div> <div>5%</div> <div> <div></div> <div>65%</div> <div>21%</div> <div>•</div> <div>12%</div> </div> </div>
2	B	5	<div> <div>20%</div> <div>80%</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4810 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 Protein-glutamine gamma-glutamyltransferase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	611	Total	C	N	O	S	0	1	0
			4682	2983	788	886	25			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	MET	-	INITIATING METHIONINE	UNP P21980
A	-5	ALA	-	EXPRESSION TAG	UNP P21980
A	-4	HIS	-	EXPRESSION TAG	UNP P21980
A	-3	HIS	-	EXPRESSION TAG	UNP P21980
A	-2	HIS	-	EXPRESSION TAG	UNP P21980
A	-1	HIS	-	EXPRESSION TAG	UNP P21980
A	0	HIS	-	EXPRESSION TAG	UNP P21980
A	1	HIS	-	EXPRESSION TAG	UNP P21980

- Molecule 2 is a protein called Peptide inhibitor.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	5	Total	C	N	O	0	0	0
			48	33	5	10			

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		

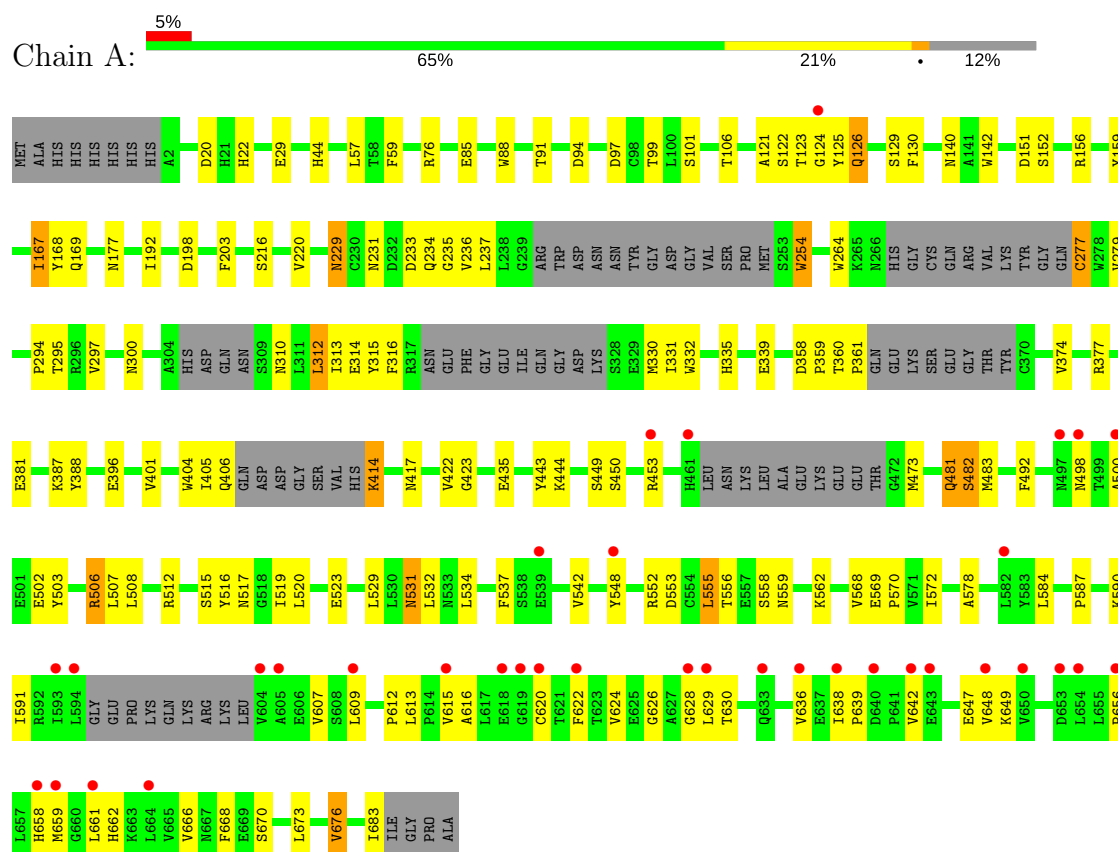
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	50	Total	O	0	0
			50	50		

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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: Protein-glutamine gamma-glutamyltransferase 2



- Molecule 2: Peptide inhibitor



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	71.00Å 71.00Å 310.36Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.73 – 2.50 46.73 – 2.50	Depositor EDS
% Data completeness (in resolution range)	87.8 (46.73-2.50) 87.6 (46.73-2.50)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.49 (at 2.48Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
R, $R_{free}$	0.226 , 0.286 0.219 , 0.277	Depositor DCC
$R_{free}$ test set	1242 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	43.1	Xtriage
Anisotropy	0.816	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 48.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	4810	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.91% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PHQ, SO4, XW1

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	0.38	0/4782	0.55	1/6515 (0.0%)
2	B	2.36	2/25 (8.0%)	1.36	0/32
All	All	0.42	2/4807 (0.0%)	0.55	1/6547 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	3	GLN	CD-NE2	6.80	1.49	1.32
2	B	3	GLN	C-N	6.31	1.46	1.34

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	277	CYS	CA-CB-SG	5.16	123.29	114.00

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	4682	0	4492	110	0
2	B	48	0	47	8	0
3	A	30	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	50	0	0	1	0
All	All	4810	0	4539	111	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 12.

All (111) 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:229[A]:ASN:H	1:A:229[A]:ASN:HD22	1.10	0.92
1:A:229[A]:ASN:ND2	1:A:229[A]:ASN:H	1.70	0.88
1:A:555:LEU:HD12	1:A:556:THR:H	1.37	0.87
1:A:330:MET:HG2	1:A:331:ILE:H	1.44	0.82
1:A:414:LYS:HG2	1:A:414:LYS:O	1.80	0.82
1:A:401:VAL:HG12	1:A:417:ASN:HB3	1.71	0.73
1:A:169:GLN:HB3	2:B:1:PHQ:H71	1.69	0.73
1:A:229[B]:ASN:ND2	1:A:361:PRO:HG3	2.05	0.72
1:A:624:VAL:HG22	1:A:666:VAL:HG22	1.73	0.71
1:A:552:ARG:HG3	1:A:553:ASP:H	1.55	0.70
1:A:94:ASP:HB3	1:A:101:SER:HB2	1.73	0.70
1:A:520:LEU:HD21	1:A:562:LYS:HD2	1.75	0.69
1:A:449:SER:O	1:A:453:ARG:HG3	1.94	0.68
1:A:659:MET:O	1:A:683:ILE:HG21	1.95	0.67
1:A:555:LEU:HD12	1:A:556:THR:N	2.10	0.67
1:A:552:ARG:HG3	1:A:553:ASP:N	2.12	0.65
1:A:126:GLN:HE21	1:A:126:GLN:HA	1.64	0.62
1:A:330:MET:SD	2:B:4:PRO:HB3	2.40	0.61
1:A:152:SER:O	1:A:156:ARG:HG3	2.00	0.61
1:A:167:ILE:HG23	1:A:297:VAL:HG11	1.83	0.61
1:A:404:TRP:O	1:A:578:ALA:HA	2.01	0.61
1:A:44:HIS:CD2	1:A:99:THR:HG22	2.36	0.60
1:A:481:GLN:HG2	1:A:482:SER:N	2.16	0.60
1:A:558:SER:O	1:A:559:ASN:HB3	2.03	0.59
1:A:167:ILE:HD13	1:A:297:VAL:HG11	1.85	0.59
1:A:358:ASP:OD1	1:A:360:THR:HB	2.04	0.58
1:A:515:SER:HB2	1:A:517:ASN:OD1	2.04	0.58
1:A:168:TYR:CE2	1:A:177:ASN:HB3	2.39	0.57
1:A:587:PRO:HG3	1:A:612:PRO:HG3	1.85	0.57
1:A:169:GLN:CB	2:B:1:PHQ:H71	2.34	0.56
1:A:404:TRP:CZ3	1:A:414:LYS:HB3	2.40	0.56
1:A:558:SER:O	1:A:559:ASN:CB	2.54	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:335:HIS:ND1	2:B:2:XW1:H93	2.23	0.54
1:A:628:GLY:HA3	1:A:662:HIS:CE1	2.43	0.53
1:A:332:TRP:NE1	2:B:2:XW1:H94A	2.23	0.53
1:A:473:MET:SD	1:A:534:LEU:HD13	2.48	0.53
1:A:330:MET:HG2	1:A:331:ILE:N	2.19	0.52
1:A:335:HIS:CE1	2:B:2:XW1:H93	2.45	0.52
1:A:569:GLU:HG3	1:A:572:ILE:HG12	1.90	0.52
1:A:515:SER:HA	1:A:558:SER:CB	2.40	0.52
1:A:517:ASN:OD1	1:A:519:ILE:HG12	2.10	0.52
1:A:661:LEU:HD12	1:A:662:HIS:N	2.25	0.51
1:A:313:ILE:H	1:A:313:ILE:HD12	1.76	0.50
1:A:642:VAL:HG21	1:A:648:VAL:HB	1.94	0.50
1:A:237:LEU:HD11	1:A:279:VAL:HG12	1.94	0.50
1:A:512:ARG:HG2	1:A:523:GLU:HA	1.93	0.50
1:A:236:VAL:HA	1:A:264:TRP:CD1	2.46	0.49
1:A:310:ASN:O	1:A:314:GLU:HG3	2.12	0.49
1:A:548:TYR:CE2	1:A:552:ARG:HD3	2.47	0.49
1:A:123:THR:O	1:A:125:TYR:N	2.45	0.49
1:A:374:VAL:HB	1:A:388:TYR:O	2.13	0.49
1:A:300:ASN:OD1	1:A:423:GLY:HA2	2.13	0.48
1:A:88:TRP:CZ3	1:A:106:THR:HG22	2.48	0.48
1:A:516:TYR:CD2	1:A:558:SER:HA	2.49	0.48
1:A:628:GLY:HA3	1:A:662:HIS:ND1	2.29	0.48
1:A:649:LYS:NZ	1:A:649:LYS:HB2	2.28	0.48
1:A:559:ASN:ND2	1:A:584:LEU:O	2.39	0.47
1:A:591:ILE:HD11	1:A:666:VAL:CG2	2.45	0.47
1:A:159:TYR:O	1:A:294:PRO:HA	2.15	0.47
1:A:591:ILE:HD11	1:A:666:VAL:HG21	1.97	0.47
1:A:377:ARG:O	1:A:381:GLU:HB2	2.15	0.46
1:A:569:GLU:HG2	1:A:572:ILE:HD11	1.96	0.46
1:A:670:SER:HB3	1:A:673:LEU:O	2.16	0.46
1:A:123:THR:O	1:A:123:THR:HG22	2.16	0.46
1:A:500:ALA:HA	1:A:537:PHE:CE1	2.51	0.46
1:A:360:THR:HG23	1:A:361:PRO:HD2	1.97	0.46
1:A:59:PHE:O	1:A:76:ARG:HA	2.16	0.46
1:A:234:GLN:HA	1:A:235:GLY:HA3	1.62	0.45
1:A:668:PHE:HB3	1:A:676:VAL:HG13	1.98	0.45
1:A:330:MET:CG	1:A:331:ILE:H	2.19	0.45
1:A:620:CYS:HB2	1:A:638:ILE:HB	1.98	0.45
1:A:254:TRP:N	1:A:254:TRP:CD1	2.85	0.45
2:B:1:PHQ:H81	2:B:1:PHQ:C1	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:312:LEU:O	1:A:315:TYR:HB3	2.17	0.45
1:A:177:ASN:ND2	1:A:177:ASN:O	2.49	0.44
1:A:229[B]:ASN:HD21	1:A:361:PRO:HG3	1.80	0.44
1:A:503:TYR:HD2	1:A:569:GLU:OE2	2.00	0.44
1:A:506:ARG:HG2	1:A:568:VAL:HB	2.00	0.44
1:A:553:ASP:OD1	1:A:553:ASP:N	2.51	0.44
1:A:295:THR:HA	1:A:339:GLU:O	2.18	0.44
1:A:443:TYR:C	1:A:444:LYS:HG3	2.39	0.44
1:A:216:SER:O	1:A:220:VAL:HG23	2.17	0.43
1:A:129:SER:O	1:A:130:PHE:HB3	2.19	0.43
1:A:335:HIS:CD2	1:A:396:GLU:OE1	2.71	0.43
1:A:405:ILE:O	1:A:406:GLN:C	2.57	0.43
1:A:569:GLU:HA	1:A:570:PRO:HD3	1.85	0.43
1:A:613:LEU:HB3	1:A:615:VAL:HG12	2.01	0.43
1:A:661:LEU:HD12	1:A:662:HIS:O	2.18	0.43
1:A:506:ARG:H	1:A:506:ARG:HG2	1.70	0.42
1:A:622:PHE:CD2	1:A:638:ILE:HD11	2.54	0.42
1:A:140:ASN:OD1	1:A:142:TRP:HB2	2.19	0.42
1:A:531:ASN:H	1:A:531:ASN:ND2	2.16	0.42
1:A:332:TRP:HE1	2:B:2:XW1:H94A	1.81	0.42
1:A:231:ASN:C	1:A:233:ASP:H	2.23	0.42
1:A:609:LEU:HD13	1:A:668:PHE:CG	2.55	0.42
1:A:507:LEU:HD22	1:A:532:LEU:HD22	2.01	0.42
1:A:569:GLU:O	1:A:569:GLU:HG3	2.20	0.42
1:A:590:LYS:O	1:A:607:VAL:HA	2.20	0.42
1:A:638:ILE:HA	1:A:639:PRO:HD2	1.86	0.42
1:A:198:ASP:O	1:A:203:PHE:CD2	2.73	0.41
1:A:628:GLY:O	1:A:629:LEU:HD23	2.21	0.41
1:A:387:LYS:HA	1:A:388:TYR:HA	1.78	0.41
1:A:569:GLU:HG3	1:A:572:ILE:CG1	2.50	0.41
1:A:615:VAL:HG22	1:A:616:ALA:H	1.84	0.41
1:A:192:ILE:HG23	1:A:264:TRP:CZ3	2.55	0.41
1:A:531:ASN:N	1:A:531:ASN:ND2	2.66	0.41
1:A:492:PHE:HA	1:A:542:VAL:O	2.20	0.41
1:A:22:HIS:HA	4:A:828:HOH:O	2.20	0.40
1:A:57:LEU:HD22	1:A:121:ALA:HA	2.04	0.40
1:A:358:ASP:HA	1:A:359:PRO:HD2	1.90	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	594/694 (86%)	555 (93%)	33 (6%)	6 (1%)	17	31
2	B	2/5 (40%)	2 (100%)	0	0	100	100
All	All	596/699 (85%)	557 (94%)	33 (6%)	6 (1%)	17	31

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	124	GLY
1	A	97	ASP
1	A	656	PRO
1	A	312	LEU
1	A	316	PHE
1	A	626	GLY

### 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	497/603 (82%)	466 (94%)	31 (6%)	20	38
2	B	3/3 (100%)	3 (100%)	0	100	100
All	All	500/606 (82%)	469 (94%)	31 (6%)	21	38

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	ASP
1	A	29	GLU
1	A	85	GLU
1	A	91	THR
1	A	122	SER
1	A	126	GLN
1	A	151	ASP
1	A	167	ILE
1	A	229[A]	ASN
1	A	229[B]	ASN
1	A	254	TRP
1	A	277	CYS
1	A	414	LYS
1	A	422	VAL
1	A	435	GLU
1	A	450	SER
1	A	481	GLN
1	A	482	SER
1	A	483	MET
1	A	498	ASN
1	A	502	GLU
1	A	506	ARG
1	A	508	LEU
1	A	529	LEU
1	A	531	ASN
1	A	555	LEU
1	A	630	THR
1	A	636	VAL
1	A	647	GLU
1	A	658	HIS
1	A	676	VAL

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

Mol	Chain	Res	Type
1	A	44	HIS
1	A	126	GLN
1	A	484	ASN
1	A	531	ASN
1	A	586	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	XW1	B	2	1,2	12,12,13	1.45	2 (16%)	10,13,15	1.63	3 (30%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	XW1	B	2	1,2	-	0/10/12/14	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2	XW1	O92-C1	2.17	1.39	1.33
2	B	2	XW1	CA-C	3.36	1.54	1.50

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	XW1	O-C-CA	-2.12	119.14	124.96
2	B	2	XW1	O92-C1-C17	2.34	118.68	111.92
2	B	2	XW1	O92-C93-C94	3.16	120.13	108.39

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2	XW1	4	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	SO4	A	701	-	4,4,4	0.28	0	6,6,6	0.51	0
3	SO4	A	702	-	4,4,4	0.18	0	6,6,6	0.32	0
3	SO4	A	703	-	4,4,4	0.21	0	6,6,6	0.17	0
3	SO4	A	704	-	4,4,4	0.13	0	6,6,6	0.40	0
3	SO4	A	705	-	4,4,4	0.21	0	6,6,6	0.20	0
3	SO4	A	706	-	4,4,4	0.22	0	6,6,6	0.15	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SO4	A	701	-	-	0/0/0/0	0/0/0/0
3	SO4	A	702	-	-	0/0/0/0	0/0/0/0
3	SO4	A	703	-	-	0/0/0/0	0/0/0/0
3	SO4	A	704	-	-	0/0/0/0	0/0/0/0
3	SO4	A	705	-	-	0/0/0/0	0/0/0/0
3	SO4	A	706	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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	611/694 (88%)	0.30	36 (5%) 22 23	39, 57, 87, 103	0
2	B	3/5 (60%)	0.36	0 100 100	69, 69, 70, 72	0
All	All	614/699 (87%)	0.30	36 (5%) 22 23	39, 57, 87, 103	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	629	LEU	6.2
1	A	604	VAL	4.9
1	A	654	LEU	4.7
1	A	594	LEU	4.4
1	A	658	HIS	4.2
1	A	605	ALA	4.0
1	A	124	GLY	3.8
1	A	618	GLU	3.3
1	A	620	CYS	3.2
1	A	622	PHE	3.1
1	A	656	PRO	3.1
1	A	628	GLY	3.1
1	A	453	ARG	2.8
1	A	650	VAL	2.8
1	A	593	ILE	2.8
1	A	461	HIS	2.6
1	A	642	VAL	2.6
1	A	640	ASP	2.6
1	A	582	LEU	2.5
1	A	615	VAL	2.4
1	A	661	LEU	2.4
1	A	664	LEU	2.4
1	A	636	VAL	2.4
1	A	619	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	643	GLU	2.3
1	A	659	MET	2.3
1	A	638	ILE	2.3
1	A	633	GLN	2.3
1	A	539	GLU	2.3
1	A	609	LEU	2.3
1	A	653	ASP	2.1
1	A	500	ALA	2.1
1	A	648	VAL	2.1
1	A	548	TYR	2.1
1	A	497	ASN	2.1
1	A	498	ASN	2.1

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	XW1	B	2	13/14	0.94	0.20	56,61,70,75	0

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	SO4	A	703	5/5	0.81	0.21	76,81,91,95	5
3	SO4	A	706	5/5	0.95	0.11	61,64,70,79	5
3	SO4	A	705	5/5	0.96	0.09	67,72,83,90	5
3	SO4	A	701	5/5	0.97	0.11	48,50,64,65	0
3	SO4	A	702	5/5	0.97	0.13	50,55,61,65	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	SO4	A	704	5/5	0.98	0.09	54,57,64,65	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.