



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 03:32 AM GMT

PDB ID : 3S3S  
Title : Transglutaminase 2 in complex with a novel inhibitor  
Authors : Lindemann, I.; Heine, A.; Klebe, G.  
Deposited on : 2011-05-18  
Resolution : 2.30 Å(reported)

This is a full wwPDB 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 <http://wwpdb.org/ValidationPDFNotes.html>

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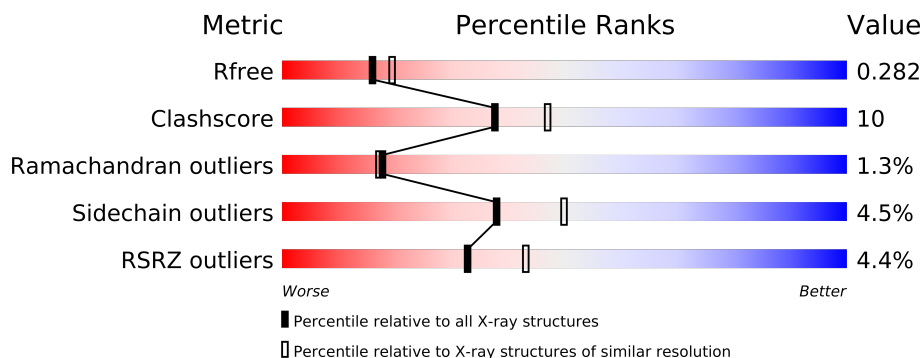
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.30 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))
$R_{free}$	66092	2929 (2.30-2.30)
Clashscore	79885	3679 (2.30-2.30)
Ramachandran outliers	78287	3642 (2.30-2.30)
Sidechain outliers	78261	3641 (2.30-2.30)
RSRZ outliers	66119	2930 (2.30-2.30)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	694	
2	B	5	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5013 atoms, of which 0 are hydrogen and 0 are deuterium.

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-glutamyltransferase2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	620	Total	C	N	O	S	0	0	0
			4784	3040	817	901	26			

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
			46	33	4	9			

- 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.

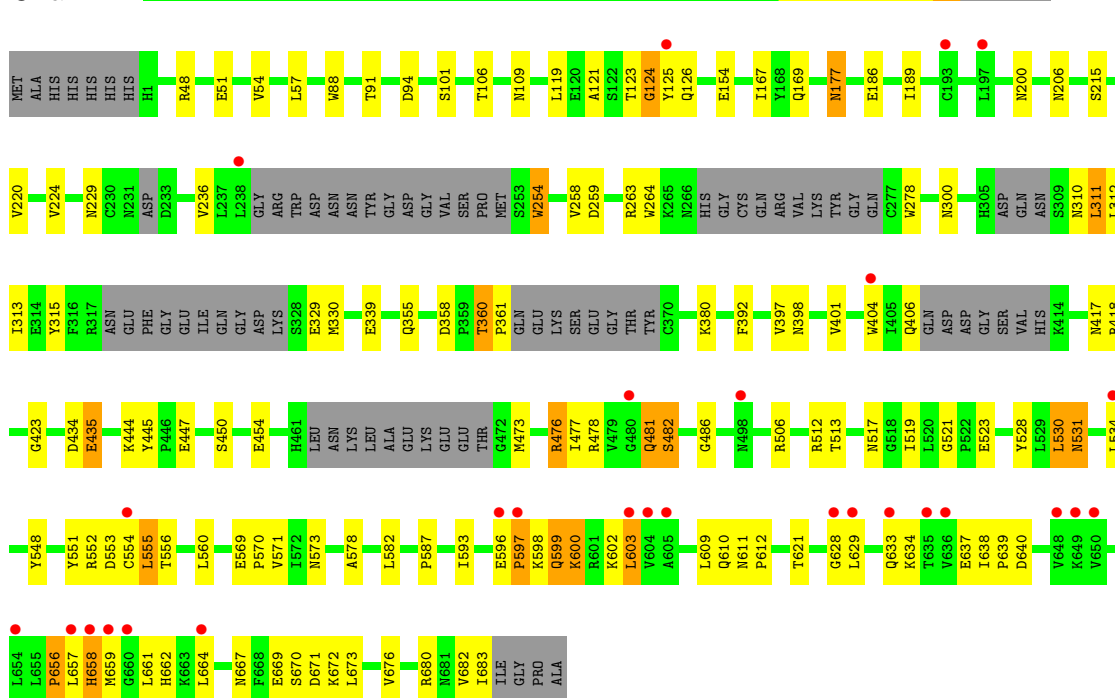
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	153	Total	O	0	0
			153	153		

### 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 errors 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-glutamyltransferase2

Chain A:



- Molecule 2: peptide inhibitor

Chain B:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	71.44Å 71.44Å 310.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.03 – 2.30 49.86 – 2.30	Depositor EDS
% Data completeness (in resolution range)	89.0 (48.03-2.30) 87.5 (49.86-2.30)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.51 (at 2.29Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
R, $R_{free}$	0.227 , 0.296 0.214 , 0.282	Depositor DCC
$R_{free}$ test set	1620 reflections (5.02%)	DCC
Wilson B-factor (Å <sup>2</sup> )	38.5	Xtriage
Anisotropy	0.629	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 45.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 34605 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	5013	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.77% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

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	0.40	0/4885	0.53	0/6651
2	B	2.74	3/23 (13.0%)	1.68	0/30
All	All	0.45	3/4908 (0.1%)	0.54	0/6681

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	3	VAL	C-N	7.95	1.49	1.34
2	B	5	LEU	C-O	6.62	1.35	1.23
2	B	4	PRO	C-N	6.15	1.48	1.34

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4784	0	4599	94	0
2	B	46	0	48	6	0
3	A	30	0	0	0	0
4	A	153	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5013	0	4647	95	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 10.

All (95) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:254:TRP:HZ3	1:A:263:ARG:HH21	1.20	0.90
1:A:621:THR:HG22	1:A:637:GLU:HA	1.56	0.87
1:A:587:PRO:HG3	1:A:612:PRO:HG3	1.54	0.87
1:A:124:GLY:O	1:A:125:TYR:HB2	1.85	0.77
1:A:57:LEU:HD22	1:A:119:LEU:HD11	1.67	0.77
1:A:401:VAL:HG12	1:A:417:ASN:HB3	1.67	0.75
1:A:569:GLU:OE1	1:A:571:VAL:HG22	1.87	0.75
1:A:593:ILE:HD11	1:A:680:ARG:HG2	1.70	0.74
1:A:476:ARG:HG3	1:A:478:ARG:HD3	1.69	0.74
1:A:94:ASP:HB3	1:A:101:SER:HB2	1.71	0.73
1:A:602:LYS:O	1:A:603:LEU:HD22	1.90	0.72
1:A:57:LEU:HD23	1:A:121:ALA:HA	1.74	0.70
1:A:51:GLU:HA	4:A:814:HOH:O	1.92	0.69
1:A:628:GLY:HA3	1:A:662:HIS:ND1	2.08	0.68
1:A:629:LEU:HD21	1:A:682:VAL:HG21	1.74	0.68
1:A:667:ASN:HD21	1:A:669:GLU:HG2	1.59	0.65
1:A:531:ASN:HD22	1:A:531:ASN:H	1.43	0.64
1:A:661:LEU:HD12	1:A:662:HIS:N	2.13	0.63
1:A:670:SER:HB3	1:A:673:LEU:O	1.99	0.62
1:A:531:ASN:HD22	1:A:531:ASN:N	1.97	0.61
1:A:517:ASN:OD1	1:A:519:ILE:HG12	2.01	0.61
1:A:597:PRO:HB2	1:A:683:ILE:HG12	1.84	0.59
1:A:154:GLU:HG2	4:A:841:HOH:O	2.02	0.59
1:A:481:GLN:HG2	1:A:482:SER:N	2.18	0.59
1:A:54:VAL:CG1	1:A:123:THR:HG23	2.33	0.59
1:A:671:ASP:OD1	1:A:672:LYS:HE3	2.03	0.59
1:A:486:GLY:HA2	1:A:548:TYR:CD1	2.38	0.58
1:A:312:LEU:O	1:A:315:TYR:HB3	2.03	0.58
1:A:552:ARG:HG3	1:A:553:ASP:N	2.17	0.58
1:A:569:GLU:CD	1:A:571:VAL:HG22	2.24	0.58
1:A:300:ASN:OD1	1:A:423:GLY:HA2	2.05	0.56
1:A:48:ARG:NH1	1:A:51:GLU:HG3	2.20	0.55
1:A:169:GLN:HB2	2:B:1:PHQ:H71	1.88	0.55
1:A:598:LYS:C	1:A:600:LYS:N	2.59	0.54
1:A:177:ASN:ND2	1:A:177:ASN:O	2.37	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:513:THR:HG21	1:A:554:CYS:O	2.06	0.54
1:A:548:TYR:HA	1:A:551:TYR:CE2	2.43	0.54
2:B:1:PHQ:H81	2:B:1:PHQ:C1	2.39	0.52
1:A:603:LEU:HD23	1:A:656:PRO:HG2	1.93	0.51
1:A:477:ILE:O	1:A:478:ARG:HD2	2.11	0.51
1:A:531:ASN:N	1:A:531:ASN:ND2	2.59	0.50
1:A:638:ILE:HG23	1:A:639:PRO:HD2	1.92	0.50
1:A:593:ILE:N	1:A:593:ILE:HD12	2.27	0.49
1:A:661:LEU:HD12	1:A:662:HIS:H	1.77	0.49
1:A:169:GLN:CG	2:B:1:PHQ:H71	2.42	0.49
1:A:229:ASN:OD1	1:A:361:PRO:HG3	2.12	0.49
1:A:552:ARG:NH1	1:A:553:ASP:HB3	2.26	0.49
1:A:473:MET:SD	1:A:534:LEU:HD13	2.53	0.49
1:A:124:GLY:O	1:A:125:TYR:CB	2.56	0.49
1:A:598:LYS:C	1:A:600:LYS:H	2.15	0.49
1:A:659:MET:O	1:A:683:ILE:HG21	2.13	0.49
1:A:506:ARG:NH2	1:A:570:PRO:HG3	2.27	0.48
1:A:512:ARG:HG2	1:A:523:GLU:HA	1.95	0.48
1:A:310:ASN:OD1	1:A:312:LEU:HB2	2.14	0.47
1:A:513:THR:OG1	1:A:521:GLY:HA3	2.14	0.47
1:A:167:ILE:HG21	1:A:278:TRP:HB2	1.97	0.47
1:A:418:ARG:HH12	1:A:447:GLU:CD	2.18	0.47
1:A:236:VAL:HA	1:A:264:TRP:CD1	2.51	0.46
1:A:358:ASP:OD1	1:A:360:THR:HB	2.16	0.46
1:A:313:ILE:HD12	1:A:313:ILE:N	2.32	0.45
1:A:629:LEU:HD13	1:A:664:LEU:HD13	1.97	0.45
1:A:311:LEU:HA	1:A:311:LEU:HD23	1.78	0.45
1:A:169:GLN:HE21	2:B:1:PHQ:C6	2.29	0.45
1:A:609:LEU:HD23	1:A:609:LEU:C	2.37	0.45
1:A:434:ASP:C	1:A:435:GLU:HG2	2.36	0.45
1:A:313:ILE:H	1:A:313:ILE:HD12	1.81	0.44
1:A:528:TYR:OH	1:A:530:LEU:HD22	2.18	0.44
1:A:658:HIS:HB3	1:A:659:MET:H	1.53	0.44
1:A:633:GLN:HG2	1:A:634:LYS:N	2.33	0.44
1:A:560:LEU:HA	1:A:582:LEU:O	2.17	0.44
1:A:598:LYS:O	1:A:600:LYS:N	2.51	0.43
1:A:599:GLN:HG2	1:A:659:MET:HA	1.99	0.43
1:A:404:TRP:O	1:A:578:ALA:HA	2.17	0.43
1:A:54:VAL:HG12	1:A:123:THR:HG23	1.99	0.43
1:A:639:PRO:HG2	1:A:640:ASP:H	1.84	0.43
1:A:169:GLN:HG3	2:B:1:PHQ:H71	1.99	0.43
1:A:126:GLN:HE21	1:A:126:GLN:HA	1.84	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:513:THR:HG22	1:A:555:LEU:HD13	2.00	0.43
1:A:330:MET:SD	2:B:4:PRO:HB3	2.59	0.42
1:A:548:TYR:HA	1:A:551:TYR:CZ	2.53	0.42
1:A:397:VAL:HG23	1:A:398:ASN:ND2	2.34	0.42
1:A:597:PRO:CB	1:A:683:ILE:HG12	2.49	0.42
1:A:186:GLU:OE1	1:A:258:VAL:HG21	2.19	0.42
1:A:392:PHE:C	1:A:392:PHE:CD2	2.93	0.41
1:A:596:GLU:O	1:A:603:LEU:HD21	2.20	0.41
1:A:88:TRP:CZ3	1:A:106:THR:HG22	2.55	0.41
1:A:380:LYS:HE2	1:A:445:TYR:CE2	2.56	0.41
1:A:339:GLU:HA	1:A:355:GLN:O	2.21	0.41
1:A:109:ASN:O	1:A:215:SER:HB2	2.21	0.41
1:A:329:GLU:OE2	1:A:506:ARG:NH2	2.54	0.40
1:A:220:VAL:O	1:A:224:VAL:HG23	2.22	0.40
1:A:682:VAL:O	1:A:683:ILE:HB	2.22	0.40
1:A:611:ASN:OD1	1:A:612:PRO:HD2	2.21	0.40
1:A:189:ILE:HD13	1:A:189:ILE:HA	1.88	0.40
1:A:555:LEU:HD12	1:A:556:THR:H	1.86	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

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	602/694 (87%)	565 (94%)	29 (5%)	8 (1%)	18	17
2	B	2/5 (40%)	2 (100%)	0	0	100	100
All	All	604/699 (86%)	567 (94%)	29 (5%)	8 (1%)	18	17

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	124	GLY
1	A	600	LYS

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Mol	Chain	Res	Type
1	A	573	ASN
1	A	599	GLN
1	A	656	PRO
1	A	657	LEU
1	A	206	ASN
1	A	597	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	510/603 (85%)	488 (96%)	22 (4%)	40	52
2	B	3/3 (100%)	2 (67%)	1 (33%)	0	0
All	All	513/606 (85%)	490 (96%)	23 (4%)	38	50

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	91	THR
1	A	177	ASN
1	A	200	ASN
1	A	254	TRP
1	A	259	ASP
1	A	311	LEU
1	A	360	THR
1	A	406	GLN
1	A	435	GLU
1	A	444	LYS
1	A	450	SER
1	A	454	GLU
1	A	476	ARG
1	A	481	GLN
1	A	482	SER
1	A	530	LEU
1	A	531	ASN
1	A	555	LEU
1	A	603	LEU

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Mol	Chain	Res	Type
1	A	610	GLN
1	A	658	HIS
1	A	676	VAL
2	B	3	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	69	GLN
1	A	126	GLN
1	A	169	GLN
1	A	266	ASN
1	A	398	ASN
1	A	441	HIS
1	A	531	ASN
1	A	533	ASN
1	A	559	ASN
1	A	599	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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	5.61	5 (41%)	11,13,15	1.91	3 (27%)

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 (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2	XW1	O-C	18.51	1.24	1.11
2	B	2	XW1	C13-CA	3.87	1.56	1.53
2	B	2	XW1	CA-C	2.57	1.53	1.48
2	B	2	XW1	O92-C93	-2.41	1.38	1.46
2	B	2	XW1	O92-C1	2.17	1.40	1.33

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	XW1	C-CA-N	-4.74	109.10	113.83
2	B	2	XW1	O92-C93-C94	2.50	118.14	108.43
2	B	2	XW1	O92-C1-C17	2.14	118.68	111.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

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	688	-	4,4,4	0.34	0	6,6,6	0.30	0
3	SO4	A	689	-	4,4,4	0.29	0	6,6,6	0.18	0
3	SO4	A	690	-	4,4,4	0.11	0	6,6,6	0.13	0
3	SO4	A	691	-	4,4,4	0.19	0	6,6,6	0.32	0
3	SO4	A	692	-	4,4,4	0.14	0	6,6,6	0.10	0
3	SO4	A	693	-	4,4,4	0.20	0	6,6,6	0.14	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	688	-	-	0/0/0/0	0/0/0/0
3	SO4	A	689	-	-	0/0/0/0	0/0/0/0
3	SO4	A	690	-	-	0/0/0/0	0/0/0/0
3	SO4	A	691	-	-	0/0/0/0	0/0/0/0
3	SO4	A	692	-	-	0/0/0/0	0/0/0/0
3	SO4	A	693	-	-	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.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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	620/694 (89%)	0.22	28 (4%) 32 42	27, 44, 77, 102	0
2	B	5/5 (100%)	0.04	0 100 100	46, 48, 52, 62	0
All	All	625/699 (89%)	0.22	28 (4%) 33 42	27, 45, 76, 102	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	597	PRO	7.3
1	A	629	LEU	5.3
1	A	480	GLY	3.7
1	A	658	HIS	3.5
1	A	654	LEU	3.5
1	A	648	VAL	3.2
1	A	633	GLN	3.1
1	A	604	VAL	3.0
1	A	603	LEU	3.0
1	A	635	THR	2.9
1	A	650	VAL	2.8
1	A	659	MET	2.7
1	A	605	ALA	2.7
1	A	649	LYS	2.7
1	A	660	GLY	2.6
1	A	664	LEU	2.6
1	A	125	TYR	2.5
1	A	554	CYS	2.4
1	A	596	GLU	2.4
1	A	404	TRP	2.3
1	A	534	LEU	2.3
1	A	636	VAL	2.3
1	A	628	GLY	2.3
1	A	193	CYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	657	LEU	2.1
1	A	238	LEU	2.1
1	A	498	ASN	2.0
1	A	197	LEU	2.0

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

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	XW1	B	2	13/14	0.18	1.62	40,46,54,54	0

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	SO4	A	693	5/5	0.13	-0.18	45,48,58,66	5
3	SO4	A	691	5/5	0.10	-0.56	49,49,50,51	0
3	SO4	A	692	5/5	0.16	-0.73	61,66,75,80	0
3	SO4	A	688	5/5	0.11	-0.82	44,46,48,50	0
3	SO4	A	689	5/5	0.08	-2.05	47,48,53,55	0
3	SO4	A	690	5/5	0.20	-	56,61,74,75	5



## 6.5 Other polymers ⓘ

There are no such residues in this entry.