



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:16 AM GMT

PDB ID : 2Q3Z
Title : Transglutaminase 2 undergoes large conformational change upon activation
Authors : Strop, P.; Pinkas, D.M.; Brunger, A.T.; Khosla, C.
Deposited on : 2007-05-30
Resolution : 2.00 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

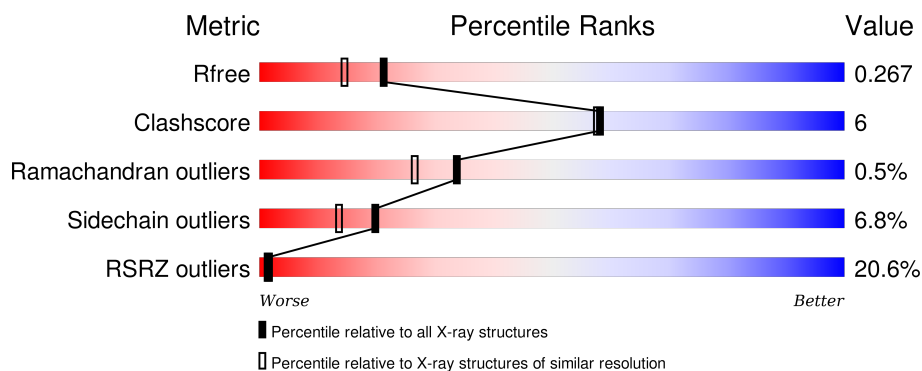
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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 ≥ 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	687	<div> <div>20%</div> <div>81%</div> <div>11%</div> <div>• 5%</div> </div>
2	X	7	<div> <div>14%</div> <div>86%</div> <div>14%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5519 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 Transglutaminase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	655	5186	3282	894	979	31	66	0	0

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	51	GLN	GLU	CONFLICT	UNP P21980
A	186	GLN	GLU	CONFLICT	UNP P21980
A	224	GLY	VAL	CONFLICT	UNP P21980
A	533	THR	ASN	CONFLICT	UNP P21980
A	655	VAL	LEU	CONFLICT	UNP P21980

- Molecule 2 is a protein called Polypeptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	X	7	46	33	6	7	0	0	1

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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		

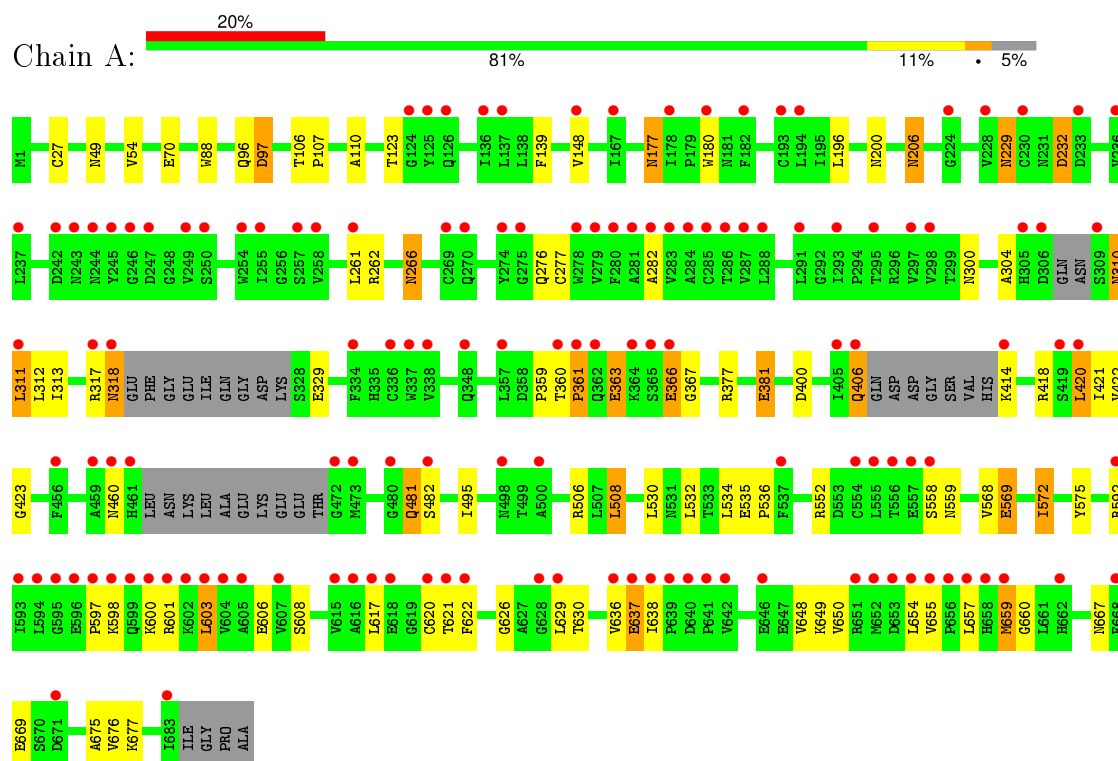
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	262	Total	O	0	0
			262	262		

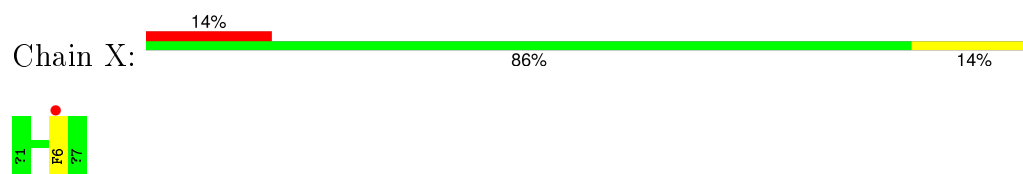
3 Residue-property plots [i](#)

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: Transglutaminase 2



• Molecule 2: Polypeptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	71.67Å 71.67Å 309.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.00 36.12 – 1.85	Depositor EDS
% Data completeness (in resolution range)	100.0 (25.00-2.00) 95.8 (36.12-1.85)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.88 (at 1.85Å)	Xtriage
Refinement program	REFMAC 5.3.0037	Depositor
R, R_{free}	0.228 , 0.266 0.228 , 0.267	Depositor DCC
R_{free} test set	3484 reflections (6.77%)	DCC
Wilson B-factor (Å ²)	30.3	Xtriage
Anisotropy	0.641	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 56.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	4 of 67283 reflections (0.006%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5519	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.09% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NH2, ONL, ACE, SO4

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	0.64	2/5297 (0.0%)	0.72	2/7186 (0.0%)
2	X	0.61	0/35	1.01	0/47
All	All	0.64	2/5332 (0.0%)	0.72	2/7233 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	329	GLU	CG-CD	-8.21	1.39	1.51
1	A	27	CYS	CB-SG	-5.75	1.72	1.81

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	97	ASP	CB-CG-OD1	-12.76	106.82	118.30
1	A	97	ASP	CB-CG-OD2	-11.78	107.70	118.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	363	GLU	Peptide

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Mol	Chain	Res	Type	Group
1	A	600	LYS	Peptide
1	A	659	MET	Peptide
1	A	97	ASP	Sidechain

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	5186	0	5097	60	0
2	X	46	0	43	4	0
3	A	25	0	0	0	0
4	A	262	0	0	0	0
All	All	5519	0	5140	60	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 (60) 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:636:VAL:HG21	1:A:650:VAL:HG11	1.49	0.95
1:A:360:THR:HG23	1:A:361:PRO:HD2	1.52	0.92
1:A:360:THR:CG2	1:A:361:PRO:HD2	2.10	0.81
1:A:603:LEU:HD12	1:A:654:LEU:HD12	1.64	0.77
1:A:638:ILE:HD13	1:A:648:VAL:HG21	1.66	0.76
1:A:622:PHE:HB2	1:A:636:VAL:HG22	1.74	0.68
1:A:318:ASN:HD22	1:A:318:ASN:H	1.43	0.64
1:A:54:VAL:HG12	1:A:123:THR:HG23	1.82	0.62
1:A:318:ASN:N	1:A:318:ASN:HD22	1.99	0.60
1:A:569:GLU:CG	1:A:572:ILE:HD12	2.32	0.60
1:A:420:LEU:HD23	2:X:6:PHE:CD2	2.38	0.58
1:A:232:ASP:N	1:A:232:ASP:OD2	2.37	0.57
1:A:569:GLU:HG3	1:A:572:ILE:HD12	1.87	0.57
1:A:304:ALA:HB2	2:X:6:PHE:CD2	2.40	0.57
1:A:568:VAL:HG13	1:A:575:TYR:CE2	2.41	0.55
1:A:139:PHE:HB3	1:A:148:VAL:HG21	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:659:MET:HG3	1:A:660:GLY:N	2.22	0.54
1:A:406:GLN:HA	1:A:406:GLN:HE21	1.71	0.54
1:A:266:ASN:HD22	1:A:266:ASN:N	2.06	0.53
1:A:377:ARG:O	1:A:381:GLU:HB2	2.08	0.53
1:A:420:LEU:HD11	1:A:422:VAL:HG13	1.89	0.53
1:A:311:LEU:HD22	1:A:508:LEU:HD13	1.92	0.52
1:A:667:ASN:HD21	1:A:675:ALA:HA	1.75	0.51
1:A:366:GLU:HG3	1:A:367:GLY:N	2.25	0.51
1:A:318:ASN:N	1:A:318:ASN:ND2	2.59	0.51
1:A:421:ILE:HD12	1:A:421:ILE:N	2.27	0.50
1:A:310:ASN:HB3	1:A:313:ILE:HD12	1.94	0.50
1:A:420:LEU:HD23	2:X:6:PHE:CE2	2.47	0.49
1:A:621:THR:HG22	1:A:637:GLU:HA	1.93	0.49
1:A:506:ARG:HD3	1:A:530:LEU:HD22	1.95	0.49
1:A:229:ASN:OD1	1:A:361:PRO:HD3	2.12	0.49
1:A:420:LEU:CD1	1:A:422:VAL:HG13	2.43	0.48
1:A:300:ASN:OD1	1:A:423:GLY:HA2	2.13	0.47
1:A:495:ILE:HG21	1:A:534:LEU:HD11	1.95	0.47
1:A:88:TRP:CZ3	1:A:106:THR:HG22	2.50	0.47
1:A:481:GLN:HG2	1:A:482:SER:N	2.30	0.47
1:A:535:GLU:HG3	1:A:536:PRO:HD2	1.96	0.47
1:A:304:ALA:HB2	2:X:6:PHE:HD2	1.78	0.47
1:A:262:ARG:O	1:A:266:ASN:ND2	2.48	0.46
1:A:620:CYS:HB2	1:A:638:ILE:HB	1.97	0.45
1:A:107:PRO:HD2	1:A:110:ALA:HB2	1.99	0.45
1:A:229:ASN:HD22	1:A:229:ASN:C	2.21	0.44
1:A:406:GLN:CA	1:A:406:GLN:HE21	2.30	0.44
1:A:276:GLN:O	1:A:277:CYS:C	2.56	0.43
1:A:420:LEU:HD11	1:A:422:VAL:CG1	2.48	0.43
1:A:196:LEU:HD23	1:A:196:LEU:C	2.39	0.43
1:A:608:SER:HB3	1:A:649:LYS:HG2	2.01	0.43
1:A:206:ASN:HD22	1:A:206:ASN:C	2.22	0.43
1:A:506:ARG:NH1	1:A:530:LEU:HD13	2.33	0.42
1:A:49:ASN:ND2	1:A:96:GLN:O	2.51	0.42
1:A:311:LEU:CD2	1:A:508:LEU:HD13	2.49	0.42
1:A:569:GLU:HG2	1:A:572:ILE:HD12	2.02	0.42
1:A:420:LEU:HD22	1:A:421:ILE:N	2.35	0.42
1:A:558:SER:O	1:A:559:ASN:HB2	2.19	0.42
1:A:54:VAL:CG1	1:A:123:THR:HG23	2.49	0.41
1:A:667:ASN:ND2	1:A:677:LYS:HD3	2.35	0.41
1:A:229:ASN:HB3	1:A:359:PRO:O	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:TRP:CE2	1:A:282:ALA:HB1	2.56	0.40
1:A:592:ARG:HB2	1:A:606:GLU:HB3	2.03	0.40
1:A:177:ASN:N	1:A:177:ASN:ND2	2.69	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	645/687 (94%)	610 (95%)	32 (5%)	3 (0%)	34	26
2	X	4/7 (57%)	4 (100%)	0	0	100	100
All	All	649/694 (94%)	614 (95%)	32 (5%)	3 (0%)	34	26

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	361	PRO
1	A	597	PRO
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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	570/596 (96%)	531 (93%)	39 (7%)	20	13
2	X	4/4 (100%)	4 (100%)	0	100	100
All	All	574/600 (96%)	535 (93%)	39 (7%)	20	13

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	70	GLU
1	A	177	ASN
1	A	200	ASN
1	A	206	ASN
1	A	229	ASN
1	A	232	ASP
1	A	261	LEU
1	A	266	ASN
1	A	310	ASN
1	A	311	LEU
1	A	312	LEU
1	A	317	ARG
1	A	318	ASN
1	A	363	GLU
1	A	366	GLU
1	A	381	GLU
1	A	400	ASP
1	A	406	GLN
1	A	414	LYS
1	A	418	ARG
1	A	420	LEU
1	A	460	ASN
1	A	481	GLN
1	A	508	LEU
1	A	532	LEU
1	A	552	ARG
1	A	569	GLU
1	A	572	ILE
1	A	598	LYS
1	A	601	ARG
1	A	603	LEU
1	A	617	LEU

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Mol	Chain	Res	Type
1	A	629	LEU
1	A	630	THR
1	A	637	GLU
1	A	655	VAL
1	A	657	LEU
1	A	669	GLU
1	A	676	VAL

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

Mol	Chain	Res	Type
1	A	51	GLN
1	A	126	GLN
1	A	134	HIS
1	A	169	GLN
1	A	177	ASN
1	A	206	ASN
1	A	266	ASN
1	A	267	HIS
1	A	318	ASN
1	A	406	GLN
1	A	531	ASN
1	A	573	ASN
1	A	610	GLN
1	A	633	GLN
1	A	662	HIS
1	A	667	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	ONL	X	3	1,2	7,8,9	0.62	0	6,9,11	1.65	1 (16%)

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	ONL	X	3	1,2	-	0/5/7/9	0/0/0/0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	X	3	ONL	CB-CG-CD	-3.82	110.25	114.52

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

5 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	0.36	0	6,6,6	0.37	0
3	SO4	A	689	-	4,4,4	0.19	0	6,6,6	0.20	0
3	SO4	A	690	-	4,4,4	0.15	0	6,6,6	0.14	0
3	SO4	A	691	-	4,4,4	0.53	0	6,6,6	0.42	0
3	SO4	A	692	-	4,4,4	0.24	0	6,6,6	0.10	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

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, 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	655/687 (95%)	1.10	135 (20%) ⓘ ⓘ	39, 47, 60, 70	21 (3%)
2	X	4/7 (57%)	2.55	1 (25%) ⓘ ⓘ	62, 63, 66, 71	0
All	All	659/694 (94%)	1.11	136 (20%) ⓘ ⓘ	39, 47, 61, 71	21 (3%)

All (136) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	654	LEU	7.8
1	A	593	ILE	7.1
1	A	601	ARG	6.8
1	A	554	CYS	6.5
2	X	6	PHE	6.4
1	A	603	LEU	6.2
1	A	597	PRO	6.0
1	A	659	MET	5.9
1	A	629	LEU	5.9
1	A	640	ASP	5.8
1	A	604	VAL	5.8
1	A	652	MET	5.7
1	A	602	LYS	5.6
1	A	125	TYR	5.4
1	A	365	SER	5.4
1	A	637	GLU	5.2
1	A	657	LEU	5.1
1	A	594	LEU	5.1
1	A	638	ILE	4.9
1	A	658	HIS	4.9
1	A	228	VAL	4.8
1	A	306	ASP	4.8
1	A	414	LYS	4.6
1	A	651	ARG	4.5

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Mol	Chain	Res	Type	RSRZ
1	A	283	VAL	4.4
1	A	309	SER	4.4
1	A	599	GLN	4.4
1	A	124	GLY	4.3
1	A	683	ILE	4.3
1	A	616	ALA	4.3
1	A	243	ASN	4.2
1	A	244	ASN	4.2
1	A	281	ALA	4.2
1	A	498	ASN	4.1
1	A	364	LYS	4.0
1	A	285	CYS	4.0
1	A	279	VAL	3.9
1	A	642	VAL	3.9
1	A	618	GLU	3.9
1	A	656	PRO	3.9
1	A	405	ILE	3.9
1	A	596	GLU	3.9
1	A	193	CYS	3.9
1	A	236	VAL	3.8
1	A	480	GLY	3.8
1	A	406	GLN	3.7
1	A	336	CYS	3.7
1	A	284	ALA	3.7
1	A	598	LYS	3.6
1	A	600	LYS	3.6
1	A	639	PRO	3.6
1	A	628	GLY	3.5
1	A	297	VAL	3.5
1	A	641	PRO	3.5
1	A	558	SER	3.5
1	A	287	VAL	3.5
1	A	615	VAL	3.5
1	A	456	PHE	3.5
1	A	537	PHE	3.5
1	A	655	VAL	3.5
1	A	278	TRP	3.5
1	A	461	HIS	3.5
1	A	556	THR	3.4
1	A	605	ALA	3.4
1	A	288	LEU	3.4
1	A	622	PHE	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	653	ASP	3.2
1	A	636	VAL	3.2
1	A	671	ASP	3.2
1	A	167	ILE	3.2
1	A	620	CYS	3.2
1	A	247	ASP	3.1
1	A	338	VAL	3.1
1	A	557	GLU	3.0
1	A	178	ILE	3.0
1	A	261	LEU	3.0
1	A	182	PHE	3.0
1	A	459	ALA	3.0
1	A	555	LEU	3.0
1	A	274	TYR	3.0
1	A	366	GLU	2.9
1	A	237	LEU	2.9
1	A	291	LEU	2.9
1	A	233	ASP	2.9
1	A	246	GLY	2.8
1	A	242	ASP	2.8
1	A	592	ARG	2.8
1	A	293	ILE	2.8
1	A	317	ARG	2.7
1	A	230	CYS	2.7
1	A	646	GLU	2.6
1	A	286	THR	2.6
1	A	357	LEU	2.6
1	A	595	GLY	2.6
1	A	348	GLN	2.6
1	A	419	SER	2.6
1	A	249	VAL	2.5
1	A	318	ASN	2.5
1	A	270	GLN	2.5
1	A	295	THR	2.5
1	A	668	PHE	2.5
1	A	482	SER	2.5
1	A	311	LEU	2.4
1	A	250	SER	2.4
1	A	337	TRP	2.4
1	A	137	LEU	2.4
1	A	420	LEU	2.4
1	A	269	CYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	180	TRP	2.4
1	A	148	VAL	2.3
1	A	282	ALA	2.3
1	A	126	GLN	2.3
1	A	245	TYR	2.3
1	A	194	LEU	2.3
1	A	224	GLY	2.3
1	A	362	GLN	2.3
1	A	617	LEU	2.2
1	A	361	PRO	2.2
1	A	334	PHE	2.2
1	A	258	VAL	2.2
1	A	472	GLY	2.2
1	A	255	ILE	2.2
1	A	254	TRP	2.2
1	A	257	SER	2.2
1	A	136	ILE	2.2
1	A	621	THR	2.1
1	A	607	VAL	2.1
1	A	280	PHE	2.1
1	A	298	VAL	2.1
1	A	662	HIS	2.1
1	A	460	ASN	2.1
1	A	473	MET	2.1
1	A	500	ALA	2.0
1	A	360	THR	2.0
1	A	305	HIS	2.0
1	A	275	GLY	2.0

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. 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, 95th 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	LLDF	B-factors(Å ²)	Q<0.9
2	ONL	X	3	9/10	0.91	0.12	-	55,57,62,63	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. 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, 95th 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	LLDF	B-factors(\AA^2)	Q<0.9
3	SO4	A	688	5/5	0.98	0.15	0.06	45,49,50,50	0
3	SO4	A	692	5/5	0.95	0.24	-	78,78,79,80	0
3	SO4	A	691	5/5	0.98	0.13	-	46,49,49,50	0
3	SO4	A	690	5/5	0.81	0.33	-	107,108,108,108	0
3	SO4	A	689	5/5	0.87	0.25	-	56,58,60,60	0

6.5 Other polymers [i](#)

There are no such residues in this entry.