



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:17 PM GMT

PDB ID : 3WPD
Title : Crystal structure of horse TLR9 in complex with inhibitory DNA4084
Authors : Ohto, U.; Tanji, H.; Shimizu, T.
Deposited on : 2014-01-11
Resolution : 2.75 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

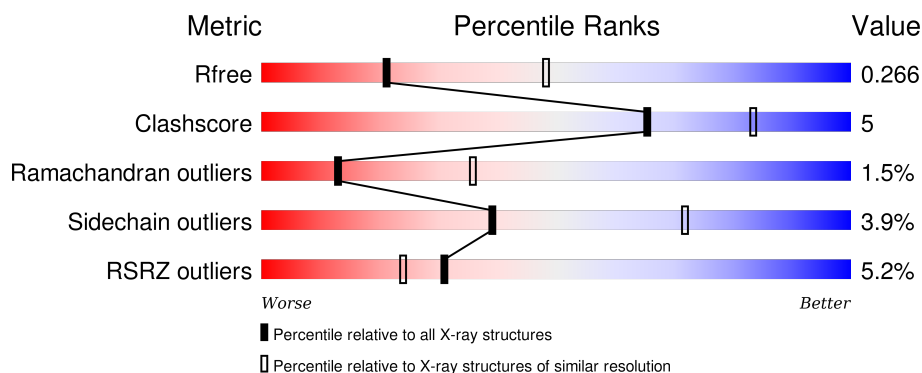
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.75 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	3340 (2.80-2.72)
Clashscore	102246	3829 (2.80-2.72)
Ramachandran outliers	100387	3767 (2.80-2.72)
Sidechain outliers	100360	3770 (2.80-2.72)
RSRZ outliers	91569	3352 (2.80-2.72)

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	802	
2	C	10	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	NAG	A	904	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6196 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 Toll-like receptor 9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	752	5909	3777	1040	1063	29	17	0	0

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	22	ARG	-	EXPRESSION TAG	UNP Q2EEY0
A	23	SER	-	EXPRESSION TAG	UNP Q2EEY0
A	24	PRO	-	EXPRESSION TAG	UNP Q2EEY0
A	25	TRP	-	EXPRESSION TAG	UNP Q2EEY0
A	818	GLU	-	EXPRESSION TAG	UNP Q2EEY0
A	819	PHE	-	EXPRESSION TAG	UNP Q2EEY0
A	820	LEU	-	EXPRESSION TAG	UNP Q2EEY0
A	821	VAL	-	EXPRESSION TAG	UNP Q2EEY0
A	822	PRO	-	EXPRESSION TAG	UNP Q2EEY0
A	823	ARG	-	EXPRESSION TAG	UNP Q2EEY0

- Molecule 2 is a DNA chain called DNA (5'-D(*CP*CP*TP*GP*GP*AP*TP*GP*GP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	C	10	206	98	40	59	9	0	0	0

- 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		

- Molecule 4 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 5 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	2	Total	C	N	O	0	0
			28	16	2	10		

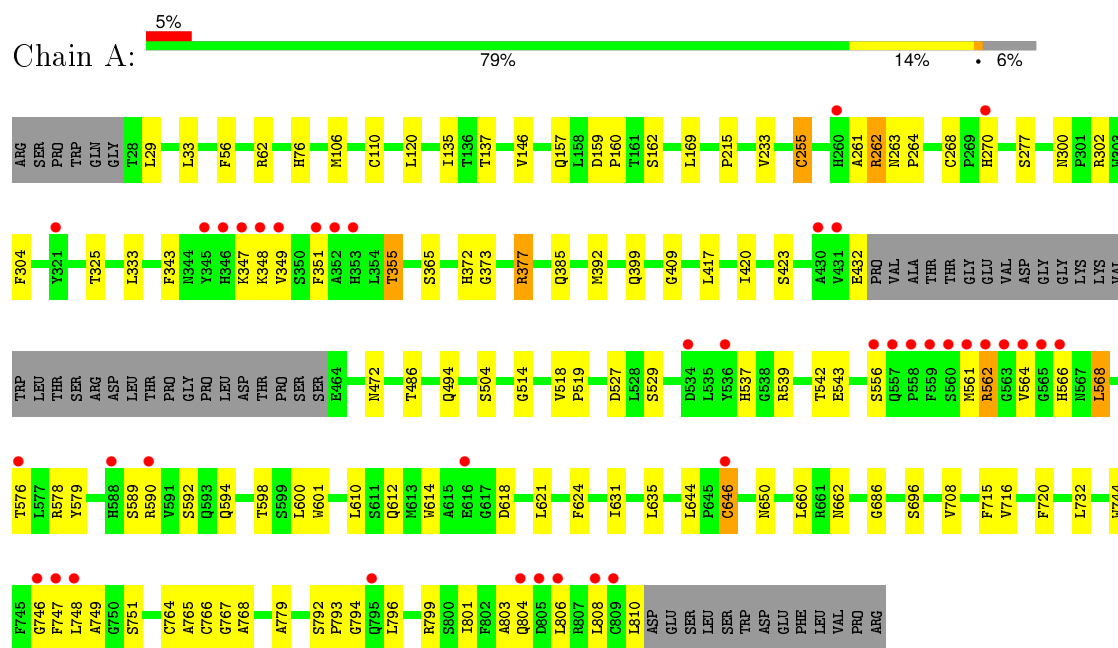
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	10	Total	O	0	0
			10	10		

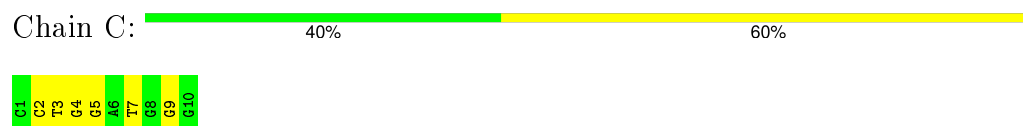
3 Residue-property plots

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: Toll-like receptor 9



• Molecule 2: DNA (5'-D(*CP*CP*TP*GP*GP*AP*TP*GP*GP*G)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	117.55Å 117.08Å 69.57Å 90.00° 105.46° 90.00°	Depositor
Resolution (Å)	58.61 – 2.75 58.54 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.9 (58.61-2.75) 99.9 (58.54-2.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.20 (at 2.77Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.201 , 0.258 0.205 , 0.266	Depositor DCC
R_{free} test set	1212 reflections (5.42%)	DCC
Wilson B-factor (Å ²)	37.1	Xtriage
Anisotropy	0.316	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 40.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	1 of 23565 reflections (0.004%)	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6196	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.03% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, 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	0.49	1/6048 (0.0%)	0.71	2/8228 (0.0%)
2	C	0.41	0/231	0.68	0/356
All	All	0.49	1/6279 (0.0%)	0.71	2/8584 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	432	GLU	CA-CB	-10.20	1.31	1.53

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	432	GLU	CB-CA-C	8.81	128.02	110.40
1	A	432	GLU	N-CA-CB	-7.37	97.34	110.60

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	5909	0	5941	57	1
2	C	206	0	114	5	0
3	A	15	0	0	1	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	28	0	26	0	0
5	A	28	0	25	0	0
6	A	10	0	0	0	0
All	All	6196	0	6106	61	1

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 5.

All (61) 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:261:ALA:HA	1:A:262:ARG:HB2	1.59	0.84
1:A:261:ALA:CA	1:A:262:ARG:HB2	2.22	0.70
1:A:56:PHE:O	1:A:62:ARG:NH1	2.26	0.69
1:A:696:SER:HB2	1:A:720:PHE:HA	1.77	0.65
1:A:747:PHE:O	1:A:749:ALA:N	2.35	0.60
1:A:76:HIS:CD2	1:A:110:CYS:SG	2.96	0.58
1:A:304:PHE:HB3	1:A:333:LEU:HD21	1.85	0.58
1:A:372:HIS:HB2	1:A:399:GLN:O	2.04	0.58
1:A:618:ASP:HA	1:A:621:LEU:HG	1.86	0.57
1:A:765:ALA:O	1:A:768:ALA:N	2.41	0.53
1:A:29:LEU:HD21	1:A:810:LEU:HD13	1.91	0.52
1:A:539:ARG:O	1:A:542:THR:OG1	2.24	0.52
1:A:803:ALA:O	1:A:806:LEU:N	2.32	0.51
1:A:537:HIS:HB3	3:A:901:SO4:O2	2.11	0.50
1:A:716:VAL:HB	1:A:744:TRP:CE2	2.46	0.50
1:A:578:ARG:HB2	1:A:579:TYR:CD1	2.45	0.50
1:A:261:ALA:HA	1:A:262:ARG:CB	2.35	0.50
1:A:263:ASN:ND2	2:C:7:DT:OP1	2.45	0.49
1:A:159:ASP:HB2	1:A:160:PRO:CD	2.43	0.49
2:C:4:DG:H5"	2:C:4:DG:N3	2.27	0.49
1:A:566:HIS:HB3	1:A:568:LEU:CD1	2.43	0.48
1:A:589:SER:OG	1:A:590:ARG:N	2.44	0.48
1:A:277:SER:CB	1:A:300:ASN:HD21	2.27	0.47
2:C:4:DG:H2"	2:C:5:DG:C8	2.50	0.47
1:A:646:CYS:O	1:A:650:ASN:ND2	2.48	0.47
1:A:417:LEU:HD13	1:A:420:ILE:HD11	1.95	0.47
1:A:796:LEU:HD12	1:A:796:LEU:N	2.31	0.46
1:A:601:TRP:O	1:A:631:ILE:HB	2.15	0.46
1:A:792:SER:HB2	1:A:793:PRO:HA	1.98	0.46
1:A:644:LEU:HB3	1:A:646:CYS:SG	2.57	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:409:GLY:HA3	1:A:494:GLN:HG3	1.98	0.45
1:A:233:VAL:HG12	1:A:255:CYS:HB3	1.97	0.45
1:A:365:SER:HA	1:A:392:MET:CE	2.46	0.45
1:A:796:LEU:CD2	1:A:806:LEU:HD21	2.46	0.45
1:A:796:LEU:HD23	1:A:806:LEU:HD11	1.99	0.45
1:A:277:SER:HB3	1:A:300:ASN:HD21	1.81	0.45
1:A:255:CYS:C	1:A:268:CYS:HB2	2.37	0.45
1:A:277:SER:HB2	1:A:302:ARG:HB2	1.99	0.45
1:A:610:LEU:HD22	1:A:614:TRP:CE2	2.52	0.44
1:A:399:GLN:HB2	1:A:423:SER:O	2.17	0.44
1:A:662:ASN:HA	1:A:686:GLY:O	2.17	0.44
1:A:708:VAL:CG2	1:A:732:LEU:HD23	2.47	0.44
1:A:159:ASP:C	1:A:159:ASP:OD1	2.56	0.43
1:A:343:PHE:HA	1:A:373:GLY:O	2.19	0.43
1:A:764:CYS:SG	1:A:801:ILE:HG22	2.59	0.43
1:A:766:CYS:HA	1:A:767:GLY:HA2	1.88	0.42
1:A:263:ASN:HB3	1:A:264:PRO:CD	2.50	0.42
2:C:2:DC:C2	2:C:9:DG:N2	2.87	0.42
1:A:527:ASP:OD1	1:A:527:ASP:C	2.58	0.41
1:A:504:SER:HA	1:A:529:SER:O	2.20	0.41
1:A:635:LEU:HB2	1:A:660:LEU:HD23	2.03	0.41
1:A:146:VAL:HA	1:A:169:LEU:HA	2.01	0.41
1:A:578:ARG:O	1:A:600:LEU:HD12	2.21	0.41
2:C:3:DT:H2'	2:C:4:DG:H5''	2.03	0.41
1:A:417:LEU:CD1	1:A:420:ILE:HD11	2.51	0.41
1:A:351:PHE:CE2	1:A:377:ARG:HG3	2.56	0.41
1:A:518:VAL:HB	1:A:519:PRO:HD3	2.02	0.40
1:A:610:LEU:HD11	1:A:624:PHE:CE2	2.56	0.40
1:A:561:MET:O	1:A:562:ARG:HB2	2.21	0.40
1:A:325:THR:HA	1:A:355:THR:O	2.21	0.40
1:A:514:GLY:HA2	1:A:543:GLU:OE1	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:GLN:OE1	1:A:539:ARG:NH1[3_455]	1.92	0.28

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	748/802 (93%)	666 (89%)	71 (10%)	11 (2%)	13	36

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	262	ARG
1	A	270	HIS
1	A	592	SER
1	A	748	LEU
1	A	779	ALA
1	A	794	GLY
1	A	215	PRO
1	A	804	GLN
1	A	349	VAL
1	A	746	GLY
1	A	564	VAL

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	668/712 (94%)	642 (96%)	26 (4%)	39	72

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	LEU
1	A	106	MET
1	A	120	LEU
1	A	135	ILE
1	A	137	THR
1	A	162	SER
1	A	255	CYS
1	A	347	LYS
1	A	348	LYS
1	A	355	THR
1	A	377	ARG
1	A	385	GLN
1	A	472	ASN
1	A	486	THR
1	A	556	SER
1	A	562	ARG
1	A	568	LEU
1	A	576	THR
1	A	594	GLN
1	A	598	THR
1	A	612	GLN
1	A	646	CYS
1	A	715	PHE
1	A	751	SER
1	A	799	ARG
1	A	808	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	76	HIS
1	A	157	GLN
1	A	260	HIS
1	A	530	HIS
1	A	566	HIS
1	A	650	ASN
1	A	797	GLN

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 ⓘ

2 carbohydrates 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	$\# Z > 2$
5	NAG	A	905	1,5	14,14,15	0.79	0	15,19,21	1.03	1 (6%)
5	NAG	A	906	5	14,14,15	0.60	0	15,19,21	1.44	2 (13%)

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
5	NAG	A	905	1,5	-	0/6/23/26	0/1/1/1
5	NAG	A	906	5	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	905	NAG	C2-N2-C7	-2.13	120.30	123.04
5	A	906	NAG	C4-C3-C2	3.04	115.95	111.23
5	A	906	NAG	C3-C4-C5	3.33	116.00	110.20

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.6 Ligand geometry

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	901	-	4,4,4	0.41	0	6,6,6	0.21	0
3	SO4	A	902	-	4,4,4	0.52	0	6,6,6	0.41	0
3	SO4	A	903	-	4,4,4	0.39	0	6,6,6	0.07	0
4	NAG	A	904	1	14,14,15	0.54	0	15,19,21	2.22	1 (6%)
4	NAG	A	907	1	14,14,15	0.51	0	15,19,21	1.18	1 (6%)

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	901	-	-	0/0/0/0	0/0/0/0
3	SO4	A	902	-	-	0/0/0/0	0/0/0/0
3	SO4	A	903	-	-	0/0/0/0	0/0/0/0
4	NAG	A	904	1	-	0/6/23/26	0/1/1/1
4	NAG	A	907	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	907	NAG	O7-C7-C8	-2.30	117.85	122.06
4	A	904	NAG	C1-O5-C5	7.66	121.97	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	901	SO4	1	0

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	752/802 (93%)	0.13	40 (5%) 30 23	14, 35, 96, 204	6 (0%)
2	C	10/10 (100%)	-0.31	0 100 100	30, 38, 46, 48	0
All	All	762/812 (93%)	0.12	40 (5%) 31 24	14, 35, 95, 204	6 (0%)

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	561	MET	11.9
1	A	563	GLY	11.8
1	A	349	VAL	8.2
1	A	534	ASP	7.4
1	A	564	VAL	7.3
1	A	559	PHE	7.2
1	A	560	SER	6.8
1	A	348	LYS	6.6
1	A	558	PRO	6.0
1	A	347	LYS	6.0
1	A	747	PHE	5.8
1	A	562	ARG	5.3
1	A	805	ASP	5.2
1	A	808	LEU	5.1
1	A	536	TYR	5.0
1	A	566	HIS	4.8
1	A	746	GLY	4.3
1	A	557	GLN	3.8
1	A	270	HIS	3.6
1	A	795	GLN	3.5
1	A	565	GLY	3.4
1	A	430	ALA	3.4
1	A	260	HIS	3.4
1	A	616	GLU	3.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	351	PHE	3.2
1	A	576	THR	3.1
1	A	748	LEU	2.9
1	A	590	ARG	2.9
1	A	352	ALA	2.8
1	A	646	CYS	2.6
1	A	346	HIS	2.6
1	A	809	CYS	2.4
1	A	588	HIS	2.4
1	A	804	GLN	2.3
1	A	431	VAL	2.2
1	A	556	SER	2.2
1	A	345	TYR	2.1
1	A	321	TYR	2.1
1	A	806	LEU	2.1
1	A	353	HIS	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](#)

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
5	NAG	A	905	14/15	0.94	0.14	-0.48	28,35,48,56	0
5	NAG	A	906	14/15	0.90	0.16	-	56,66,76,81	0

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
4	NAG	A	904	14/15	0.86	0.39	4.39	73,101,119,124	0
4	NAG	A	907	14/15	0.95	0.15	-0.05	38,42,47,49	0
3	SO4	A	903	5/5	0.91	0.16	-	80,84,97,97	5
3	SO4	A	902	5/5	0.92	0.21	-	51,59,75,76	0
3	SO4	A	901	5/5	0.95	0.16	-	59,68,75,77	0

6.5 Other polymers ⓘ

There are no such residues in this entry.