



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:00 PM GMT

PDB ID : 4K6J
Title : Human cohesin inhibitor WapL
Authors : Tomchick, D.R.; Yu, H.; Ouyang, Z.
Deposited on : 2013-04-16
Resolution : 2.62 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

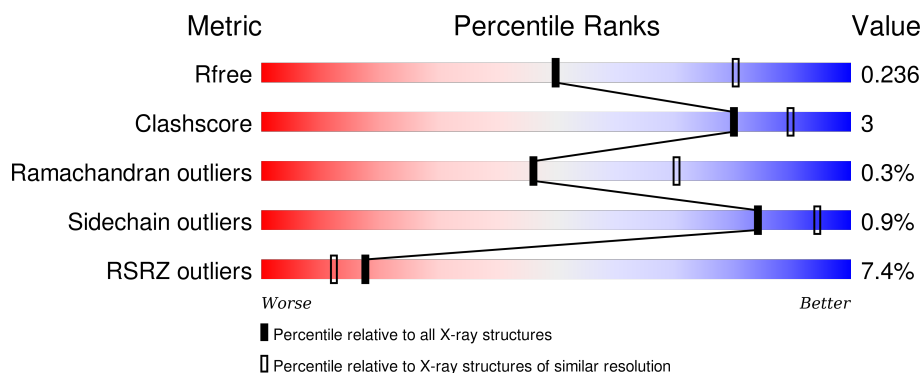
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.62 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	2700 (2.64-2.60)
Clashscore	102246	3065 (2.64-2.60)
Ramachandran outliers	100387	3015 (2.64-2.60)
Sidechain outliers	100360	3015 (2.64-2.60)
RSRZ outliers	91569	2706 (2.64-2.60)

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	568	<div> <div>5%</div> <div>79%</div> <div>8%</div> <div>13%</div> </div>
1	B	568	<div> <div>7%</div> <div>75%</div> <div>7%</div> <div>17%</div> </div>

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
2	SO4	B	1205	-	-	-	X
3	ACT	A	1208	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 15557 atoms, of which 7730 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 Wings apart-like protein homolog.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	497	Total	C	H	N	O	S	0	0	0
			7887	2451	3963	689	752	32			
1	B	471	Total	C	H	N	O	S	0	0	0
			7485	2330	3761	654	709	31			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	GLY	-	EXPRESSION TAG	UNP Q7Z5K2
A	-6	PRO	-	EXPRESSION TAG	UNP Q7Z5K2
A	-5	LEU	-	EXPRESSION TAG	UNP Q7Z5K2
A	-4	GLY	-	EXPRESSION TAG	UNP Q7Z5K2
A	-3	SER	-	EXPRESSION TAG	UNP Q7Z5K2
A	-2	GLY	-	EXPRESSION TAG	UNP Q7Z5K2
A	-1	ARG	-	EXPRESSION TAG	UNP Q7Z5K2
A	0	PRO	-	EXPRESSION TAG	UNP Q7Z5K2
B	-7	GLY	-	EXPRESSION TAG	UNP Q7Z5K2
B	-6	PRO	-	EXPRESSION TAG	UNP Q7Z5K2
B	-5	LEU	-	EXPRESSION TAG	UNP Q7Z5K2
B	-4	GLY	-	EXPRESSION TAG	UNP Q7Z5K2
B	-3	SER	-	EXPRESSION TAG	UNP Q7Z5K2
B	-2	GLY	-	EXPRESSION TAG	UNP Q7Z5K2
B	-1	ARG	-	EXPRESSION TAG	UNP Q7Z5K2
B	0	PRO	-	EXPRESSION TAG	UNP Q7Z5K2

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			7	2	3	2		
3	A	1	Total	C	H	O	0	0
			7	2	3	2		

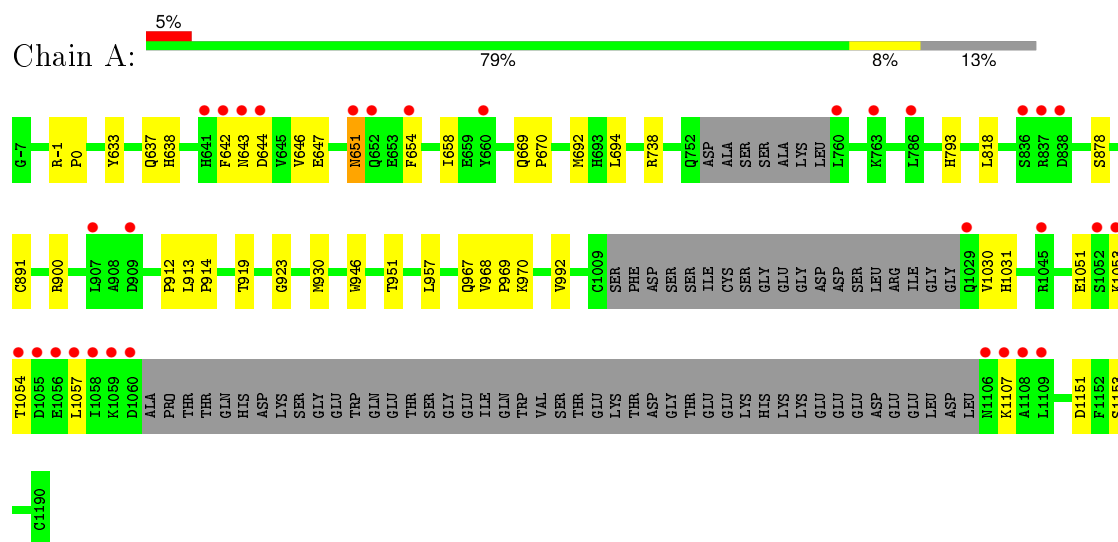
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	58	Total	O	0	0
			58	58		
4	B	48	Total	O	0	0
			48	48		

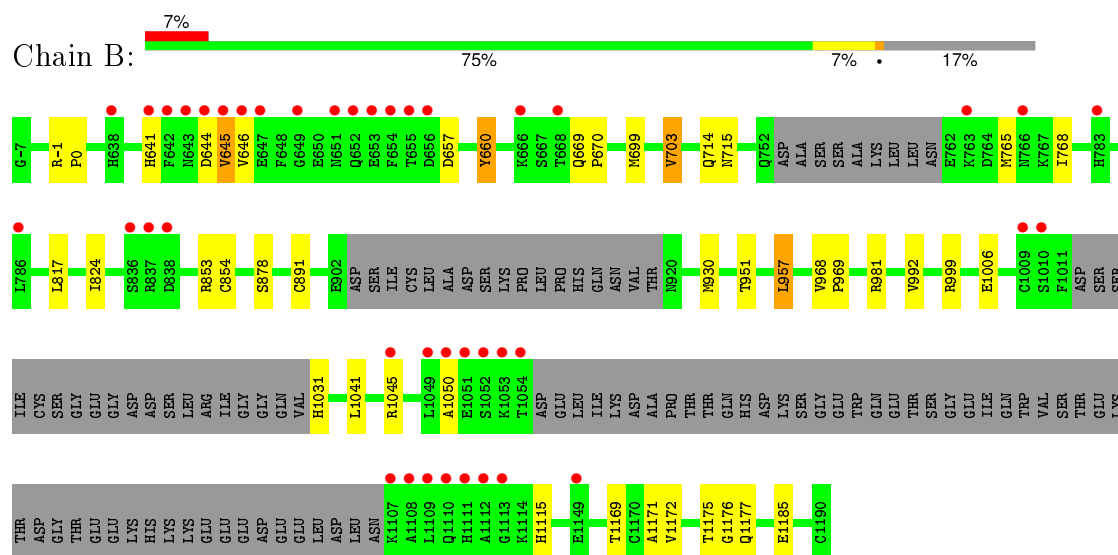
3 Residue-property plots

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: Wings apart-like protein homolog



- Molecule 1: Wings apart-like protein homolog



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	107.54Å 107.54Å 300.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.68 – 2.62 41.84 – 2.62	Depositor EDS
% Data completeness (in resolution range)	94.8 (29.68-2.62) 94.9 (41.84-2.62)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.42 (at 2.61Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.184 , 0.237 0.182 , 0.236	Depositor DCC
R_{free} test set	2584 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	34.7	Xtriage
Anisotropy	0.261	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 58.7	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	0 of 51187 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	15557	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.20% 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: SO4, ACT

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	0.22	0/3979	0.39	0/5369
1	B	0.22	0/3776	0.40	0/5089
All	All	0.22	0/7755	0.39	0/10458

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	3924	3963	3943	27	0
1	B	3724	3761	3742	27	0
2	A	35	0	0	0	0
2	B	30	0	0	1	0
3	A	8	6	6	0	0
4	A	58	0	0	0	0
4	B	48	0	0	1	0
All	All	7827	7730	7691	52	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 3.

All (52) 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:715:ASN:ND2	1:B:765:MET:SD	2.65	0.70
1:A:738:ARG:HG2	1:A:818:LEU:HD13	1.81	0.63
1:B:853:ARG:NH2	4:B:1309:HOH:O	2.31	0.63
1:B:-1:ARG:N	1:B:0:PRO:HD2	2.15	0.61
1:B:992:VAL:O	1:B:999:ARG:NH1	2.34	0.61
1:B:644:ASP:O	1:B:646:VAL:N	2.34	0.60
1:A:946:TRP:CD1	1:B:817:LEU:HD13	2.39	0.58
1:A:1051:GLU:HA	1:A:1054:THR:HG22	1.87	0.57
1:A:638:HIS:NE2	1:A:644:ASP:OD2	2.40	0.55
1:A:738:ARG:HG2	1:A:818:LEU:CD1	2.37	0.55
1:A:-1:ARG:N	1:A:0:PRO:CD	2.72	0.52
1:B:1006:GLU:HA	1:B:1031:HIS:HB3	1.92	0.52
1:A:646:VAL:CG1	1:A:647:GLU:N	2.73	0.51
1:A:1030:VAL:HG22	1:A:1031:HIS:N	2.27	0.50
1:B:951:THR:HG22	1:B:957:LEU:HD22	1.93	0.50
1:B:1050:ALA:HB3	1:B:1115:HIS:HB2	1.92	0.50
1:A:738:ARG:HD3	1:A:818:LEU:HD13	1.95	0.49
1:B:657:ASP:HA	1:B:660:TYR:CE2	2.49	0.48
1:B:968:VAL:N	1:B:969:PRO:CD	2.77	0.47
1:B:765:MET:O	1:B:768:ILE:N	2.48	0.47
1:B:981:ARG:NH2	2:B:1201:SO4:O3	2.48	0.47
1:A:968:VAL:N	1:A:969:PRO:CD	2.78	0.47
1:A:951:THR:HG22	1:A:957:LEU:HD22	1.96	0.46
1:A:891:CYS:HB3	1:A:930:MET:HG3	1.97	0.46
1:B:669:GLN:HG3	1:B:670:PRO:HD2	1.97	0.46
1:B:645:VAL:HG13	1:B:645:VAL:O	2.15	0.46
1:A:669:GLN:HG3	1:A:670:PRO:HD2	1.97	0.45
1:A:913:LEU:HB3	1:A:914:PRO:HD2	1.98	0.45
1:A:967:GLN:O	1:A:970:LYS:HG2	2.17	0.45
1:B:1175:THR:HG23	1:B:1176:GLY:N	2.31	0.44
1:A:900:ARG:HD2	1:A:923:GLY:HA3	1.99	0.44
1:A:642:PHE:O	1:A:643:ASN:HB2	2.17	0.44
1:B:657:ASP:HA	1:B:660:TYR:CD2	2.54	0.43
1:B:1172:VAL:HG13	1:B:1177:GLN:HG3	2.00	0.43
1:B:1041:LEU:O	1:B:1045:ARG:HB2	2.19	0.43
1:A:1151:ASP:CG	1:A:1153:SER:HG	2.22	0.43
1:A:633:TYR:CE2	1:A:637:GLN:HG3	2.54	0.43
1:A:646:VAL:HG12	1:A:647:GLU:N	2.33	0.42
1:A:651:ASN:O	1:A:654:PHE:HB3	2.19	0.42
1:B:824:ILE:CG2	1:B:854:CYS:HB3	2.49	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:891:CYS:HB3	1:B:930:MET:HG3	2.01	0.42
1:B:714:GLN:HG2	1:B:715:ASN:N	2.35	0.42
1:B:699:MET:O	1:B:703:VAL:HG13	2.20	0.42
1:B:644:ASP:HB3	1:B:646:VAL:HG23	2.02	0.42
1:B:1169:THR:HG22	1:B:1171:ALA:HB2	2.01	0.42
1:B:1172:VAL:HG13	1:B:1177:GLN:CG	2.50	0.42
1:A:646:VAL:HG11	1:A:692:MET:CE	2.50	0.41
1:A:658:ILE:HD11	1:A:694:LEU:HD13	2.01	0.41
1:A:1053:LYS:O	1:A:1057:LEU:HD13	2.20	0.41
1:A:912:PRO:HB2	1:A:919:THR:OG1	2.21	0.41
1:A:1057:LEU:HD21	1:A:1107:LYS:HB3	2.02	0.41
1:A:946:TRP:CD1	1:B:817:LEU:CD1	3.03	0.41

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	489/568 (86%)	471 (96%)	17 (4%)	1 (0%)	52	76
1	B	461/568 (81%)	444 (96%)	15 (3%)	2 (0%)	39	63
All	All	950/1136 (84%)	915 (96%)	32 (3%)	3 (0%)	46	70

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	645	VAL
1	A	651	ASN
1	B	641	HIS

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	443/504 (88%)	440 (99%)	3 (1%)	88	96
1	B	418/504 (83%)	413 (99%)	5 (1%)	78	92
All	All	861/1008 (85%)	853 (99%)	8 (1%)	84	94

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	793	HIS
1	A	878	SER
1	A	992	VAL
1	B	660	TYR
1	B	703	VAL
1	B	878	SER
1	B	957	LEU
1	B	1185	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	712	HIS

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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

15 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	$\# Z > 2$
2	SO4	A	1201	-	4,4,4	0.28	0	6,6,6	0.11	0
2	SO4	A	1202	-	4,4,4	0.24	0	6,6,6	0.13	0
2	SO4	A	1203	-	4,4,4	0.20	0	6,6,6	0.14	0
2	SO4	A	1204	-	4,4,4	0.11	0	6,6,6	0.06	0
2	SO4	A	1205	-	4,4,4	0.13	0	6,6,6	0.09	0
2	SO4	A	1206	-	4,4,4	0.18	0	6,6,6	0.08	0
2	SO4	A	1207	-	4,4,4	0.17	0	6,6,6	0.07	0
3	ACT	A	1208	-	1,3,3	1.44	0	0,3,3	0.00	-
3	ACT	A	1209	-	1,3,3	1.34	0	0,3,3	0.00	-
2	SO4	B	1201	-	4,4,4	0.24	0	6,6,6	0.09	0
2	SO4	B	1202	-	4,4,4	0.29	0	6,6,6	0.16	0
2	SO4	B	1203	-	4,4,4	0.19	0	6,6,6	0.12	0
2	SO4	B	1204	-	4,4,4	0.19	0	6,6,6	0.25	0
2	SO4	B	1205	-	4,4,4	0.20	0	6,6,6	0.09	0
2	SO4	B	1206	-	4,4,4	0.13	0	6,6,6	0.08	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
2	SO4	A	1201	-	-	0/0/0/0	0/0/0/0
2	SO4	A	1202	-	-	0/0/0/0	0/0/0/0
2	SO4	A	1203	-	-	0/0/0/0	0/0/0/0
2	SO4	A	1204	-	-	0/0/0/0	0/0/0/0
2	SO4	A	1205	-	-	0/0/0/0	0/0/0/0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SO4	A	1206	-	-	0/0/0/0	0/0/0/0
2	SO4	A	1207	-	-	0/0/0/0	0/0/0/0
3	ACT	A	1208	-	-	0/0/0/0	0/0/0/0
3	ACT	A	1209	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1201	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1202	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1203	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1204	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1205	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1206	-	-	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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1201	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	497/568 (87%)	0.07	31 (6%)	24 18	17, 38, 98, 147	0
1	B	471/568 (82%)	0.16	41 (8%)	13 9	19, 43, 104, 148	0
All	All	968/1136 (85%)	0.11	72 (7%)	17 12	17, 40, 101, 148	0

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	643	ASN	8.0
1	A	1060	ASP	7.6
1	A	1058	ILE	7.4
1	B	652	GLN	6.6
1	B	651	ASN	6.2
1	B	1109	LEU	6.1
1	B	641	HIS	5.8
1	B	653	GLU	5.6
1	A	1059	LYS	5.5
1	A	1057	LEU	5.4
1	B	1108	ALA	5.2
1	B	1052	SER	5.2
1	A	837	ARG	4.9
1	B	1111	HIS	4.7
1	A	644	ASP	4.5
1	B	642	PHE	4.5
1	A	1055	ASP	4.4
1	B	1110	GLN	4.3
1	B	1053	LYS	4.3
1	A	1056	GLU	4.2
1	B	1054	THR	4.2
1	B	1112	ALA	3.9
1	B	654	PHE	3.8
1	B	645	VAL	3.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	641	HIS	3.7
1	B	1049	LEU	3.5
1	B	783	HIS	3.4
1	B	644	ASP	3.2
1	B	647	GLU	3.2
1	B	837	ARG	3.1
1	A	760	LEU	3.1
1	A	1107	LYS	3.1
1	A	1054	THR	3.1
1	B	838	ASP	3.0
1	A	1106	ASN	3.0
1	A	1029	GLN	3.0
1	B	655	THR	3.0
1	A	763	LYS	3.0
1	A	838	ASP	2.9
1	B	656	ASP	2.9
1	A	786	LEU	2.8
1	B	1050	ALA	2.8
1	A	654	PHE	2.8
1	B	1113	GLY	2.8
1	A	907	LEU	2.7
1	A	836	SER	2.7
1	A	1052	SER	2.6
1	A	1053	LYS	2.6
1	A	1109	LEU	2.5
1	B	1051	GLU	2.5
1	B	1107	LYS	2.5
1	A	642	PHE	2.5
1	A	909	ASP	2.4
1	B	646	VAL	2.4
1	A	652	GLN	2.4
1	B	763	LYS	2.4
1	A	660	TYR	2.3
1	B	836	SER	2.3
1	B	1009	CYS	2.3
1	B	1149	GLU	2.3
1	B	786	LEU	2.3
1	B	666	LYS	2.3
1	B	643	ASN	2.3
1	B	1045	ARG	2.2
1	A	1108	ALA	2.2
1	B	668	THR	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	1045	ARG	2.1
1	B	649	GLY	2.1
1	B	1010	SER	2.1
1	A	651	ASN	2.1
1	B	766	ASN	2.0
1	B	638	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](#)

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(Å ²)	Q<0.9
3	ACT	A	1208	4/4	0.83	0.21	3.57	34,40,53,63	0
2	SO4	B	1205	5/5	0.97	0.17	2.11	47,53,57,73	0
3	ACT	A	1209	4/4	0.86	0.17	1.53	25,30,50,55	0
2	SO4	A	1204	5/5	0.90	0.18	0.82	66,79,92,94	0
2	SO4	A	1202	5/5	0.98	0.15	0.79	39,47,54,66	0
2	SO4	A	1207	5/5	0.97	0.16	0.07	49,53,68,69	0
2	SO4	A	1201	5/5	0.99	0.14	-0.07	34,38,44,44	0
2	SO4	B	1204	5/5	0.98	0.15	-0.52	34,45,64,75	0
2	SO4	A	1206	5/5	0.99	0.15	-0.55	50,57,59,59	0
2	SO4	B	1201	5/5	0.96	0.13	-0.59	34,40,63,69	0
2	SO4	A	1205	5/5	0.94	0.16	-0.72	63,66,78,86	0
2	SO4	B	1202	5/5	0.99	0.14	-0.73	27,35,39,49	0
2	SO4	B	1203	5/5	0.98	0.12	-0.75	42,53,54,56	0
2	SO4	B	1206	5/5	0.96	0.14	-0.79	45,59,64,71	0
2	SO4	A	1203	5/5	0.98	0.11	-1.04	46,50,59,66	0

6.5 Other polymers [i](#)

There are no such residues in this entry.