



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:10 AM UTC

PDB ID : 3L09 / pdb_00003l09
Title : Crystal structure of Putative transcriptional regulator (JANN_22DEC04_C
ONTIG27_REVISIED_GENE3569) from Jannaschia sp. CCS1 at 2.81 Å
resolution
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : 2009-12-09
Resolution : 2.81 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

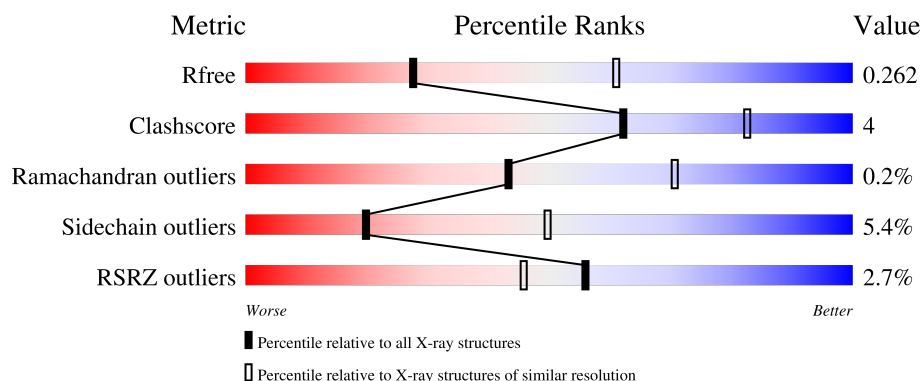
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.81 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}	180053	4591 (2.84-2.80)
Clashscore	190562	5010 (2.84-2.80)
Ramachandran outliers	187476	4916 (2.84-2.80)
Sidechain outliers	187428	4918 (2.84-2.80)
RSRZ outliers	180081	4594 (2.84-2.80)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	266	4% 77% 15% 6%
1	B	266	3% 80% 12% 6%
1	C	266	% 82% 12% 5%
1	D	266	2% 79% 14% 6%

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	A	266	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7745 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 Putative transcriptional regulator.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	250	Total	C	N	O	S	Se	0	0	0
			1883	1190	344	342	4	3			
1	B	251	Total	C	N	O	S	Se	0	2	0
			1922	1218	352	345	4	3			
1	C	253	Total	C	N	O	S	Se	0	2	0
			1942	1227	360	348	4	3			
1	D	251	Total	C	N	O	S	Se	0	1	0
			1914	1210	353	344	4	3			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP Q28UN6
A	1	VAL	-	expression tag	UNP Q28UN6
B	0	GLY	-	expression tag	UNP Q28UN6
B	1	VAL	-	expression tag	UNP Q28UN6
C	0	GLY	-	expression tag	UNP Q28UN6
C	1	VAL	-	expression tag	UNP Q28UN6
D	0	GLY	-	expression tag	UNP Q28UN6
D	1	VAL	-	expression tag	UNP Q28UN6

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

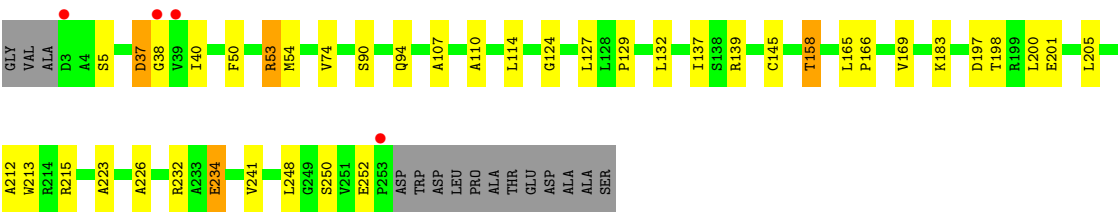
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	9	Total	O	0	0
			9	9		
3	B	5	Total	O	0	0
			5	5		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	7	Total	O	0	0
			7	7		
3	D	13	Total	O	0	0
			13	13		



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	69.95Å 81.72Å 199.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.93 – 2.81 46.93 – 2.81	Depositor EDS
% Data completeness (in resolution range)	(Not available) (46.93-2.81) 99.1 (46.93-2.81)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.41 (at 2.81Å)	Xtriage
Refinement program	BUSTER 2.8.0, PHENIX	Depositor
R, R_{free}	0.199 , 0.245 0.215 , 0.262	Depositor DCC
R_{free} test set	1430 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	51.2	Xtriage
Anisotropy	0.702	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 61.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7745	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% 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, 0.333 respectively for untwinned datasets, and 0.375, 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

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	0.88	1/1921 (0.1%)	1.42	7/2617 (0.3%)
1	B	0.83	0/1969	1.41	14/2681 (0.5%)
1	C	0.87	0/1988	1.34	8/2704 (0.3%)
1	D	0.87	1/1956 (0.1%)	1.39	10/2661 (0.4%)
All	All	0.86	2/7834 (0.0%)	1.39	39/10663 (0.4%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	40	ILE	CG1-CD1	5.50	1.73	1.51
1	A	40	ILE	CG1-CD1	5.06	1.71	1.51

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	37	ASP	CA-CB-CG	7.48	120.08	112.60
1	C	37	ASP	CA-CB-CG	7.30	119.90	112.60
1	C	38	GLY	N-CA-C	6.68	121.60	113.79
1	B	37	ASP	CA-CB-CG	6.55	119.15	112.60
1	B	38	GLY	N-CA-C	6.29	121.15	113.79
1	D	158	THR	CB-CA-C	6.00	121.23	109.35
1	D	38	GLY	N-CA-C	5.97	120.78	113.79
1	B	158	THR	CB-CA-C	5.92	121.08	109.35
1	C	158	THR	CB-CA-C	5.85	120.94	109.35
1	A	158	THR	CB-CA-C	5.74	120.70	109.35
1	A	94	GLN	CA-C-N	5.67	128.34	120.29
1	A	94	GLN	C-N-CA	5.67	128.34	120.29
1	D	94	GLN	CA-C-N	5.63	128.28	120.29
1	D	94	GLN	C-N-CA	5.63	128.28	120.29
1	D	226	ALA	CA-C-N	5.58	128.78	120.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	226	ALA	C-N-CA	5.58	128.78	120.31
1	B	94	GLN	CA-C-N	5.47	128.06	120.29
1	B	94	GLN	C-N-CA	5.47	128.06	120.29
1	A	6	ASP	CA-CB-CG	5.45	118.05	112.60
1	B	37	ASP	N-CA-C	5.43	118.91	112.72
1	A	253	PRO	N-CA-C	5.41	119.67	111.19
1	C	94	GLN	CA-C-N	5.34	127.88	120.29
1	C	94	GLN	C-N-CA	5.34	127.88	120.29
1	C	37	ASP	N-CA-C	5.20	118.65	112.72
1	B	201	GLU	CA-C-N	5.17	127.47	120.38
1	B	201	GLU	C-N-CA	5.17	127.47	120.38
1	D	37	ASP	N-CA-C	5.17	118.62	112.72
1	C	192	ILE	CA-C-N	5.15	129.23	120.88
1	C	192	ILE	C-N-CA	5.15	129.23	120.88
1	B	177	GLY	CA-C-N	5.14	131.36	121.54
1	B	177	GLY	C-N-CA	5.14	131.36	121.54
1	A	35	SER	CA-C-N	5.10	127.11	120.28
1	A	35	SER	C-N-CA	5.10	127.11	120.28
1	D	212	ALA	CA-C-N	5.05	127.30	120.44
1	D	212	ALA	C-N-CA	5.05	127.30	120.44
1	B	253	PRO	CA-C-N	5.02	127.27	120.44
1	B	253	PRO	C-N-CA	5.02	127.27	120.44
1	B	101	ARG	CA-C-N	5.01	126.99	120.88
1	B	101	ARG	C-N-CA	5.01	126.99	120.88

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	1883	0	1893	18	0
1	B	1922	0	1945	14	0
1	C	1942	0	1972	16	0
1	D	1914	0	1948	16	0
2	A	5	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	5	0	0	0	0
2	C	25	0	0	1	0
2	D	15	0	0	1	0
3	A	9	0	0	0	0
3	B	5	0	0	0	0
3	C	7	0	0	0	0
3	D	13	0	0	2	0
All	All	7745	0	7758	63	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 4.

All (63) 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:C:49:SER:HG	1:C:255:TRP:CD1	2.08	0.71
1:D:53:ARG:NH1	1:D:201:GLU:HG2	2.07	0.69
1:B:49:SER:HG	1:B:255:TRP:CD1	2.10	0.68
1:C:163:ARG:HD3	3:D:271:HOH:O	1.96	0.65
1:A:53:ARG:NH1	1:A:201:GLU:HG2	2.13	0.64
1:A:215:ARG:HD3	2:A:266:SO4:O2	2.02	0.60
1:B:178:CYS:HB3	1:B:181:GLU:HB2	1.84	0.59
1:B:185:LEU:HD11	1:B:212:ALA:HB1	1.86	0.57
1:D:232:ARG:HE	1:D:234:GLU:HG2	1.70	0.56
1:C:202[B]:ARG:HH21	1:C:251:VAL:HG22	1.70	0.55
1:D:114:LEU:HB3	1:D:158:THR:HG23	1.88	0.55
1:A:114:LEU:HB3	1:A:158:THR:HG23	1.87	0.55
1:C:77:ARG:HG2	1:C:84:PHE:HB2	1.90	0.54
1:B:129:PRO:HD2	1:B:132:LEU:HD12	1.88	0.54
1:C:129:PRO:HD2	1:C:132:LEU:HD12	1.89	0.54
1:C:232:ARG:HE	1:C:234:GLU:HG2	1.73	0.53
1:B:179:GLU:O	1:B:183:LYS:HG3	2.09	0.52
1:B:114:LEU:HB3	1:B:158:THR:HG23	1.90	0.52
1:D:215:ARG:HD3	2:D:266:SO4:O2	2.10	0.52
1:A:178:CYS:HB3	1:A:181:GLU:HB2	1.91	0.52
1:C:114:LEU:HB3	1:C:158:THR:HG23	1.91	0.51
1:B:173:VAL:HG11	1:B:227:ALA:HB3	1.93	0.51
1:C:163:ARG:CD	3:D:271:HOH:O	2.56	0.50
1:A:129:PRO:HD2	1:A:132:LEU:HD12	1.93	0.50
1:B:173:VAL:HG12	1:B:224:ALA:HA	1.95	0.49
1:A:189:LEU:HD22	1:A:244:LEU:HD13	1.92	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:211:HIS:NE2	2:A:266:SO4:O2	2.43	0.49
1:A:145:CYS:HB2	1:A:223:ALA:HA	1.95	0.49
1:A:107:ALA:HB3	1:A:232:ARG:HH22	1.78	0.48
1:D:129:PRO:HD2	1:D:132:LEU:HD12	1.94	0.48
1:A:232:ARG:HE	1:A:234:GLU:HG2	1.78	0.48
1:B:50:PHE:CZ	1:B:54:MSE:HE1	2.49	0.48
1:C:215:ARG:HD3	2:C:266:SO4:O4	2.14	0.47
1:C:202[B]:ARG:NH2	1:C:251:VAL:HG22	2.30	0.47
1:C:232:ARG:HE	1:C:234:GLU:CG	2.28	0.47
1:D:53:ARG:HH11	1:D:201:GLU:HG2	1.80	0.46
1:A:205:LEU:HD23	1:A:248:LEU:HD21	1.98	0.46
1:C:50:PHE:CZ	1:C:54:MSE:HE1	2.49	0.46
1:B:40:ILE:HD12	1:B:41:GLU:HG2	1.98	0.46
1:D:50:PHE:CZ	1:D:54:MSE:HE1	2.51	0.45
1:A:50:PHE:CZ	1:A:54:MSE:HE1	2.51	0.45
1:D:145:CYS:HB2	1:D:223:ALA:HA	1.98	0.45
1:A:165:LEU:HD22	1:A:169:VAL:HG11	2.00	0.44
1:D:205:LEU:HD23	1:D:248:LEU:HD21	1.99	0.44
1:B:205:LEU:HD23	1:B:248:LEU:HD21	2.00	0.43
1:C:165:LEU:HD22	1:C:169:VAL:HG11	2.00	0.43
1:B:165:LEU:HD22	1:B:169:VAL:HG11	2.01	0.42
1:A:189:LEU:HD22	1:A:244:LEU:CD1	2.49	0.42
1:A:229:GLY:O	1:A:232:ARG:HD3	2.19	0.42
1:C:107:ALA:HB1	1:C:110:ALA:HB2	2.01	0.42
1:C:205:LEU:HD23	1:C:248:LEU:HD21	2.01	0.42
1:A:124:GLY:HA2	1:A:127:LEU:HD12	2.01	0.42
1:D:107:ALA:HB1	1:D:110:ALA:HB2	2.02	0.42
1:D:198:THR:HG22	1:D:200:LEU:H	1.86	0.41
1:D:166:PRO:HG2	1:D:169:VAL:HG23	2.03	0.41
1:A:210:LEU:HD13	1:A:245:LEU:HD11	2.03	0.41
1:B:232:ARG:HE	1:B:234:GLU:CG	2.33	0.41
1:C:139:ARG:HH21	1:D:139:ARG:NH1	2.19	0.41
1:D:213:TRP:CG	1:D:241:VAL:HG11	2.56	0.41
1:B:49:SER:OG	1:B:255:TRP:CD1	2.71	0.41
1:D:124:GLY:HA2	1:D:127:LEU:HD12	2.03	0.40
1:D:165:LEU:HD22	1:D:169:VAL:HG11	2.03	0.40
1:A:53:ARG:HH11	1:A:201:GLU:HG2	1.84	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	248/266 (93%)	237 (96%)	10 (4%)	1 (0%)	30	58
1	B	251/266 (94%)	240 (96%)	10 (4%)	1 (0%)	30	58
1	C	253/266 (95%)	244 (96%)	9 (4%)	0	100	100
1	D	250/266 (94%)	241 (96%)	9 (4%)	0	100	100
All	All	1002/1064 (94%)	962 (96%)	38 (4%)	2 (0%)	43	71

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	6	ASP
1	B	178	CYS

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	195/209 (93%)	181 (93%)	14 (7%)	13	37
1	B	201/209 (96%)	192 (96%)	9 (4%)	24	57
1	C	203/209 (97%)	194 (96%)	9 (4%)	25	58
1	D	201/209 (96%)	190 (94%)	11 (6%)	19	49
All	All	800/836 (96%)	757 (95%)	43 (5%)	20	50

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	ARG
1	A	53	ARG
1	A	74	VAL
1	A	77	ARG
1	A	90	SER
1	A	137	ILE
1	A	145	CYS
1	A	149	GLU
1	A	188	THR
1	A	193	ASP
1	A	197	ASP
1	A	234	GLU
1	A	250	SER
1	A	252	GLU
1	B	37	ASP
1	B	40	ILE
1	B	74	VAL
1	B	137	ILE
1	B	153	GLU
1	B	197	ASP
1	B	234	GLU
1	B	250	SER
1	B	252	GLU
1	C	5	SER
1	C	37	ASP
1	C	53	ARG
1	C	74	VAL
1	C	90	SER
1	C	197	ASP
1	C	234	GLU
1	C	250	SER
1	C	252	GLU
1	D	5	SER
1	D	37	ASP
1	D	53	ARG
1	D	74	VAL
1	D	90	SER
1	D	137	ILE
1	D	183	LYS
1	D	197	ASP
1	D	234	GLU
1	D	250	SER
1	D	252	GLU

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

Mol	Chain	Res	Type
1	A	113	HIS
1	B	57	GLN
1	B	113	HIS
1	C	57	GLN
1	D	57	GLN
1	D	113	HIS
1	D	170	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

10 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	D	268	-	4,4,4	0.42	0	6,6,6	0.21	0
2	SO4	D	266	-	4,4,4	0.58	0	6,6,6	0.38	0
2	SO4	C	270	-	4,4,4	0.35	0	6,6,6	0.20	0
2	SO4	A	266	-	4,4,4	0.65	0	6,6,6	0.46	0
2	SO4	D	267	-	4,4,4	0.57	0	6,6,6	0.25	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	B	266	-	4,4,4	0.52	0	6,6,6	0.52	0
2	SO4	C	267	-	4,4,4	0.51	0	6,6,6	0.30	0
2	SO4	C	266	-	4,4,4	0.59	0	6,6,6	0.24	0
2	SO4	C	269	-	4,4,4	0.26	0	6,6,6	0.36	0
2	SO4	C	268	-	4,4,4	0.25	0	6,6,6	0.27	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.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	266	SO4	1	0
2	A	266	SO4	2	0
2	C	266	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	247/266 (92%)	0.07	11 (4%) 38 30	45, 70, 111, 138	0
1	B	248/266 (93%)	0.32	9 (3%) 46 37	32, 85, 138, 164	2 (0%)
1	C	250/266 (93%)	-0.22	3 (1%) 76 68	27, 52, 94, 111	2 (0%)
1	D	248/266 (93%)	-0.20	4 (1%) 70 61	24, 50, 85, 99	1 (0%)
All	All	993/1064 (93%)	-0.01	27 (2%) 56 46	24, 62, 121, 164	5 (0%)

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	5	SER	4.0
1	B	73	TRP	3.3
1	D	39	VAL	3.0
1	A	5	SER	2.9
1	B	123	ASP	2.9
1	B	122	PRO	2.7
1	A	121	ALA	2.6
1	A	254	ASP	2.6
1	A	37	ASP	2.6
1	A	123	ASP	2.6
1	A	38	GLY	2.5
1	C	3	ASP	2.4
1	C	40	ILE	2.4
1	B	79	LEU	2.4
1	A	122	PRO	2.4
1	D	3	ASP	2.4
1	D	38	GLY	2.3
1	A	126	SER	2.3
1	B	255	TRP	2.3
1	B	121	ALA	2.3
1	B	17	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	191	GLN	2.1
1	A	40	ILE	2.1
1	A	127	LEU	2.0
1	C	255	TRP	2.0
1	D	253	PRO	2.0
1	B	55	GLY	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 oligosaccharides 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, 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	B-factors(Å ²)	Q<0.9
2	SO4	C	268	5/5	0.73	0.19	56,59,60,61	5
2	SO4	C	270	5/5	0.76	0.17	71,75,76,76	5
2	SO4	C	267	5/5	0.86	0.15	46,50,51,52	5
2	SO4	B	266	5/5	0.88	0.11	80,84,85,86	0
2	SO4	D	268	5/5	0.88	0.13	45,49,50,52	5
2	SO4	D	267	5/5	0.89	0.08	93,97,98,99	0
2	SO4	D	266	5/5	0.92	0.30	93,98,99,100	0
2	SO4	C	266	5/5	0.93	0.11	41,46,46,47	5
2	SO4	A	266	5/5	0.95	0.10	42,46,47,49	5
2	SO4	C	269	5/5	0.97	0.07	48,52,53,56	0

6.5 Other polymers [i](#)

There are no such residues in this entry.