



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 3LES
Title : 2F5 Epitope scaffold ES2
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Deposited on : 2010-01-15
Resolution : 2.77 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.23.2
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.23.2

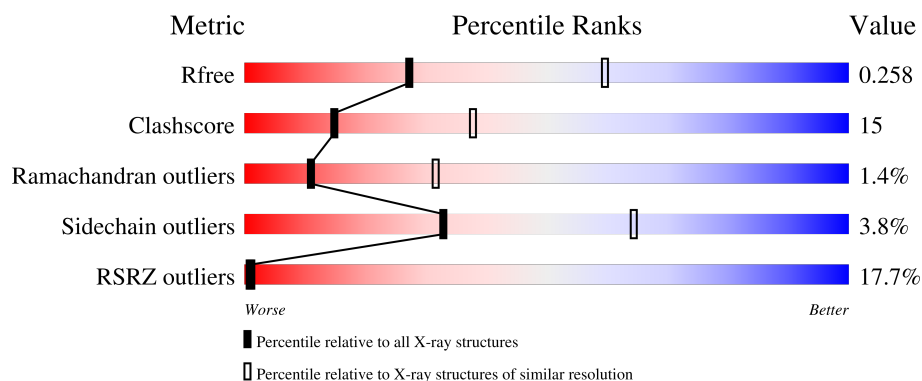
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.77 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}	130704	4107 (2.80-2.76)
Clashscore	141614	4575 (2.80-2.76)
Ramachandran outliers	138981	4487 (2.80-2.76)
Sidechain outliers	138945	4489 (2.80-2.76)
RSRZ outliers	127900	4027 (2.80-2.76)

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	179	
1	B	179	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2880 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 RNA polymerase sigma factor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	179	Total	C	N	O	S	0	0	0
			1424	907	255	261	1			
1	B	177	Total	C	N	O	S	0	0	0
			1408	898	250	259	1			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	105	GLU	GLN	engineered mutation	UNP Q9EZJ8
A	107	LEU	PRO	engineered mutation	UNP Q9EZJ8
A	108	GLU	LEU	engineered mutation	UNP Q9EZJ8
A	110	ASP	THR	engineered mutation	UNP Q9EZJ8
A	111	LYS	LEU	engineered mutation	UNP Q9EZJ8
A	112	TRP	GLU	engineered mutation	UNP Q9EZJ8
A	113	ALA	GLU	engineered mutation	UNP Q9EZJ8
A	115	LEU	ILE	engineered mutation	UNP Q9EZJ8
A	116	GLY	ASP	engineered mutation	UNP Q9EZJ8
A	117	ALA	LEU	engineered mutation	UNP Q9EZJ8
A	119	ALA	ARG	engineered mutation	UNP Q9EZJ8
A	155	ALA	ARG	engineered mutation	UNP Q9EZJ8
A	247	GLY	ARG	engineered mutation	UNP Q9EZJ8
A	249	ALA	LYS	engineered mutation	UNP Q9EZJ8
B	105	GLU	GLN	engineered mutation	UNP Q9EZJ8
B	107	LEU	PRO	engineered mutation	UNP Q9EZJ8
B	108	GLU	LEU	engineered mutation	UNP Q9EZJ8
B	110	ASP	THR	engineered mutation	UNP Q9EZJ8
B	111	LYS	LEU	engineered mutation	UNP Q9EZJ8
B	112	TRP	GLU	engineered mutation	UNP Q9EZJ8
B	113	ALA	GLU	engineered mutation	UNP Q9EZJ8
B	115	LEU	ILE	engineered mutation	UNP Q9EZJ8
B	116	GLY	ASP	engineered mutation	UNP Q9EZJ8
B	117	ALA	LEU	engineered mutation	UNP Q9EZJ8
B	119	ALA	ARG	engineered mutation	UNP Q9EZJ8

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Chain	Residue	Modelled	Actual	Comment	Reference
B	155	ALA	ARG	engineered mutation	UNP Q9EZJ8
B	247	GLY	ARG	engineered mutation	UNP Q9EZJ8
B	249	ALA	LYS	engineered mutation	UNP Q9EZJ8

- 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	B	1	Total	O	S	0	0
			5	4	1		

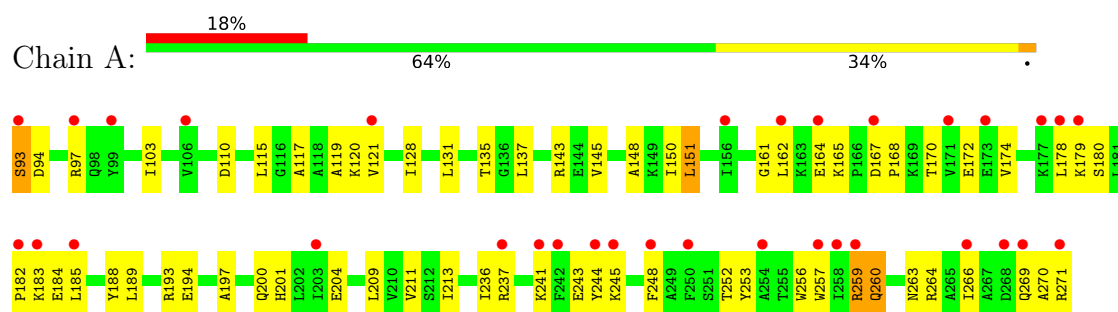
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	15	Total	O	0	0
			15	15		
3	B	13	Total	O	0	0
			13	13		

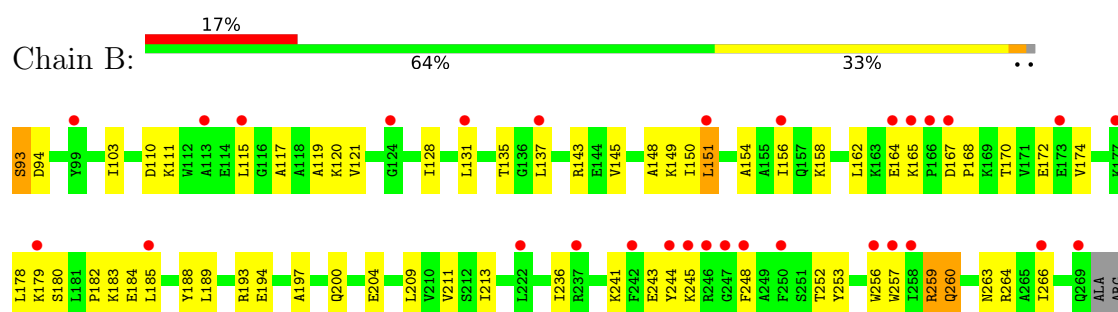
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: RNA polymerase sigma factor



• Molecule 1: RNA polymerase sigma factor



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	37.26Å 49.99Å 57.57Å 78.03° 89.83° 87.57°	Depositor
Resolution (Å)	48.85 – 2.77 48.86 – 2.77	Depositor EDS
% Data completeness (in resolution range)	85.6 (48.85-2.77) 85.7 (48.86-2.77)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.94 (at 2.77Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.227 , 0.260 0.226 , 0.258	Depositor DCC
R_{free} test set	1011 reflections (10.07%)	wwPDB-VP
Wilson B-factor (Å ²)	62.5	Xtriage
Anisotropy	0.720	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 89.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.034 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2880	wwPDB-VP
Average B, all atoms (Å ²)	100.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.39% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.26	0/1446	0.43	0/1944
1	B	0.26	0/1430	0.43	0/1923
All	All	0.26	0/2876	0.43	0/3867

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 [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	1424	0	1475	43	0
1	B	1408	0	1457	45	0
2	A	15	0	0	3	0
2	B	5	0	0	1	0
3	A	15	0	0	1	0
3	B	13	0	0	3	0
All	All	2880	0	2932	88	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 15.

All (88) 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:248:PHE:HE2	1:A:253:TYR:HB2	1.51	0.76
1:B:248:PHE:HE2	1:B:253:TYR:HB2	1.52	0.74
1:B:170:THR:O	1:B:174:VAL:HG23	1.91	0.71
1:A:263:ASN:HA	1:A:266:ILE:HG12	1.74	0.70
1:B:263:ASN:HA	1:B:266:ILE:HG12	1.74	0.69
1:A:201:HIS:NE2	2:A:2:SO4:O2	2.25	0.69
1:A:170:THR:O	1:A:174:VAL:HG23	1.93	0.69
1:A:135:THR:HG23	1:A:137:LEU:H	1.59	0.68
1:B:135:THR:HG23	1:B:137:LEU:H	1.58	0.68
1:A:182:PRO:HD2	1:A:185:LEU:HD22	1.79	0.64
1:B:150:ILE:HG21	1:B:197:ALA:HA	1.80	0.63
1:B:182:PRO:HD2	1:B:185:LEU:HD22	1.80	0.63
1:A:150:ILE:HG21	1:A:197:ALA:HA	1.79	0.63
1:A:209:LEU:O	1:A:213:ILE:HG13	1.97	0.62
1:B:209:LEU:O	1:B:213:ILE:HG13	1.99	0.62
1:A:168:PRO:O	1:A:172:GLU:HG3	2.00	0.61
1:A:97:ARG:HD3	3:A:16:HOH:O	2.00	0.61
1:B:260:GLN:HA	1:B:263:ASN:OD1	2.01	0.61
1:A:248:PHE:CE2	1:A:253:TYR:HB2	2.34	0.61
1:A:260:GLN:HA	1:A:263:ASN:OD1	2.00	0.61
1:B:168:PRO:O	1:B:172:GLU:HG3	2.01	0.60
1:B:248:PHE:CE2	1:B:253:TYR:HB2	2.36	0.60
1:A:260:GLN:O	1:A:264:ARG:HB2	2.02	0.59
1:B:260:GLN:O	1:B:264:ARG:HB2	2.03	0.59
1:B:162:LEU:HD21	1:B:236:ILE:HG13	1.86	0.58
1:A:167:ASP:HB2	1:A:170:THR:HB	1.88	0.56
1:A:103:ILE:HD11	1:A:211:VAL:HG21	1.87	0.55
1:B:167:ASP:HB2	1:B:170:THR:HB	1.87	0.55
1:B:182:PRO:HB2	1:B:185:LEU:HD13	1.88	0.55
1:A:182:PRO:HB2	1:A:185:LEU:HD13	1.87	0.55
1:A:248:PHE:CZ	1:A:252:THR:HB	2.43	0.54
1:B:158:LYS:NZ	3:B:23:HOH:O	2.40	0.54
1:B:248:PHE:CZ	1:B:252:THR:HB	2.43	0.54
1:B:103:ILE:HD11	1:B:211:VAL:HG21	1.90	0.53
1:A:162:LEU:HD21	1:A:236:ILE:HG13	1.90	0.53
1:A:145:VAL:O	1:A:148:ALA:HB3	2.09	0.53
1:B:120:LYS:HE2	1:B:194:GLU:HB3	1.91	0.53
1:A:244:TYR:OH	1:A:245:LYS:HE3	2.10	0.52
1:B:145:VAL:O	1:B:148:ALA:HB3	2.09	0.52
1:A:120:LYS:HE2	1:A:194:GLU:HB3	1.92	0.51
1:A:117:ALA:O	1:A:121:VAL:HG23	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:93:SER:O	1:B:94:ASP:C	2.49	0.50
1:B:156:ILE:HA	3:B:5:HOH:O	2.12	0.50
1:A:93:SER:O	1:A:94:ASP:C	2.50	0.50
1:B:117:ALA:O	1:B:121:VAL:HG23	2.12	0.50
1:B:189:LEU:O	1:B:193:ARG:HG3	2.12	0.50
1:A:189:LEU:O	1:A:193:ARG:HG3	2.12	0.49
1:B:178:LEU:C	1:B:180:SER:H	2.15	0.49
1:B:244:TYR:OH	1:B:245:LYS:HE3	2.11	0.49
1:A:131:LEU:O	1:A:135:THR:HG22	2.12	0.49
1:A:178:LEU:C	1:A:180:SER:H	2.15	0.49
1:B:110:ASP:OD1	1:B:110:ASP:C	2.51	0.49
1:A:165:LYS:HB3	1:A:165:LYS:NZ	2.28	0.48
1:B:165:LYS:NZ	1:B:165:LYS:HB3	2.29	0.48
1:B:241:LYS:HD2	1:B:257:TRP:CZ2	2.49	0.47
1:A:270:ALA:O	1:A:271:ARG:C	2.52	0.47
1:A:110:ASP:C	1:A:110:ASP:OD1	2.52	0.47
1:A:241:LYS:HD2	1:A:257:TRP:CZ2	2.50	0.47
1:B:154:ALA:HA	2:B:4:SO4:O4	2.15	0.47
1:A:200:GLN:O	1:A:204:GLU:HG3	2.15	0.46
1:B:149:LYS:NZ	3:B:27:HOH:O	2.44	0.46
1:A:237:ARG:HD3	2:A:1:SO4:O2	2.15	0.45
1:B:151:LEU:HA	1:B:151:LEU:HD23	1.70	0.45
1:A:161:GLY:N	2:A:3:SO4:O1	2.43	0.45
1:A:128:ILE:CD1	1:A:143:ARG:HG3	2.47	0.45
1:A:151:LEU:HD23	1:A:151:LEU:HA	1.68	0.45
1:B:162:LEU:CD2	1:B:236:ILE:HG13	2.47	0.44
1:B:200:GLN:O	1:B:204:GLU:HG3	2.18	0.44
1:B:135:THR:CG2	1:B:137:LEU:H	2.28	0.43
1:A:115:LEU:HD21	1:A:248:PHE:O	2.18	0.43
1:A:183:LYS:HG3	1:A:184:GLU:H	1.84	0.43
1:B:128:ILE:CD1	1:B:143:ARG:HG3	2.49	0.43
1:B:256:TRP:O	1:B:259:ARG:HB3	2.19	0.42
1:B:115:LEU:HD21	1:B:248:PHE:O	2.19	0.42
1:A:162:LEU:CD2	1:A:236:ILE:HG13	2.49	0.42
1:A:256:TRP:O	1:A:259:ARG:HB3	2.19	0.42
1:B:188:TYR:CD1	1:B:188:TYR:N	2.87	0.42
1:B:183:LYS:HG3	1:B:184:GLU:H	1.84	0.42
1:B:131:LEU:O	1:B:135:THR:HG22	2.19	0.41
1:B:178:LEU:O	1:B:180:SER:N	2.53	0.41
1:A:188:TYR:CD1	1:A:188:TYR:N	2.89	0.41
1:B:167:ASP:HB2	1:B:170:THR:CB	2.49	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:LEU:O	1:A:180:SER:N	2.54	0.41
1:B:188:TYR:HD1	1:B:188:TYR:H	1.69	0.41
1:A:167:ASP:HB2	1:A:170:THR:CB	2.49	0.41
1:B:162:LEU:HD11	1:B:236:ILE:HD11	2.03	0.41
1:A:119:ALA:HB2	1:A:244:TYR:CD1	2.56	0.40
1:B:119:ALA:HB2	1:B:244:TYR:CD1	2.56	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	177/179 (99%)	160 (90%)	15 (8%)	2 (1%)	14	38
1	B	175/179 (98%)	161 (92%)	11 (6%)	3 (2%)	9	27
All	All	352/358 (98%)	321 (91%)	26 (7%)	5 (1%)	11	31

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	179	LYS
1	A	179	LYS
1	A	259	ARG
1	B	111	LYS
1	B	259	ARG

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	145/145 (100%)	139 (96%)	6 (4%)	30	61
1	B	144/145 (99%)	139 (96%)	5 (4%)	36	67
All	All	289/290 (100%)	278 (96%)	11 (4%)	33	64

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	93	SER
1	A	151	LEU
1	A	164	GLU
1	A	243	GLU
1	A	260	GLN
1	A	269	GLN
1	B	93	SER
1	B	151	LEU
1	B	164	GLU
1	B	243	GLU
1	B	260	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	206	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry

4 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	B	4	-	4,4,4	0.16	0	6,6,6	0.24	0
2	SO4	A	1	-	4,4,4	0.13	0	6,6,6	0.08	0
2	SO4	A	3	-	4,4,4	0.17	0	6,6,6	0.50	0
2	SO4	A	2	-	4,4,4	0.10	0	6,6,6	0.39	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.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	4	SO4	1	0
2	A	1	SO4	1	0
2	A	3	SO4	1	0
2	A	2	SO4	1	0

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, 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	179/179 (100%)	1.15	33 (18%)  	59, 90, 153, 167	0
1	B	177/179 (98%)	1.10	30 (16%)  	59, 89, 151, 167	0
All	All	356/358 (99%)	1.12	63 (17%)  	59, 90, 153, 167	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	185	LEU	7.2
1	A	257	TRP	6.0
1	A	271	ARG	5.1
1	B	179	LYS	5.1
1	B	244	TYR	4.4
1	B	167	ASP	4.1
1	B	266	ILE	3.9
1	A	93	SER	3.8
1	B	245	LYS	3.7
1	B	164	GLU	3.5
1	B	222	LEU	3.5
1	B	248	PHE	3.4
1	A	177	LYS	3.4
1	B	165	LYS	3.4
1	A	266	ILE	3.4
1	A	269	GLN	3.3
1	A	242	PHE	3.1
1	A	248	PHE	3.0
1	A	244	TYR	3.0
1	A	164	GLU	3.0
1	A	237	ARG	2.9
1	A	173	GLU	2.8
1	A	258	ILE	2.8
1	A	97	ARG	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	256	TRP	2.7
1	A	245	LYS	2.7
1	B	166	PRO	2.7
1	A	179	LYS	2.7
1	B	269	GLN	2.6
1	B	137	LEU	2.6
1	A	268	ASP	2.6
1	A	254	ALA	2.6
1	A	182	PRO	2.6
1	A	183	LYS	2.6
1	A	185	LEU	2.5
1	A	178	LEU	2.5
1	A	156	ILE	2.5
1	B	151	LEU	2.5
1	B	246	ARG	2.5
1	A	167	ASP	2.5
1	B	113	ALA	2.4
1	B	250	PHE	2.3
1	B	258	ILE	2.3
1	B	131	LEU	2.3
1	B	99	TYR	2.3
1	A	250	PHE	2.3
1	A	259	ARG	2.2
1	B	115	LEU	2.2
1	B	177	LYS	2.1
1	B	237	ARG	2.1
1	A	162	LEU	2.1
1	A	106	VAL	2.1
1	A	99	TYR	2.1
1	B	124	GLY	2.1
1	A	121	VAL	2.1
1	B	156	ILE	2.1
1	B	173	GLU	2.1
1	A	241	LYS	2.0
1	B	242	PHE	2.0
1	B	257	TRP	2.0
1	A	171	VAL	2.0
1	A	203	ILE	2.0
1	B	247	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 monosaccharides 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(\AA^2)	Q<0.9
2	SO4	A	1	5/5	0.59	0.33	170,174,187,202	0
2	SO4	A	3	5/5	0.84	0.34	96,139,161,167	0
2	SO4	B	4	5/5	0.91	0.17	63,80,138,169	0
2	SO4	A	2	5/5	0.94	0.15	80,103,134,136	0

6.5 Other polymers [i](#)

There are no such residues in this entry.