



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 10, 2016 – 07:52 AM EDT

PDB ID : 5CKV
Title : DAHP synthase from Mycobacterium tuberculosis, fully inhibited by tyrosine, phenylalanine, and tryptophan
Authors : Munack, S.; Kregel, U.
Deposited on : 2015-07-15
Resolution : 2.79 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027107
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) : rb-20027107

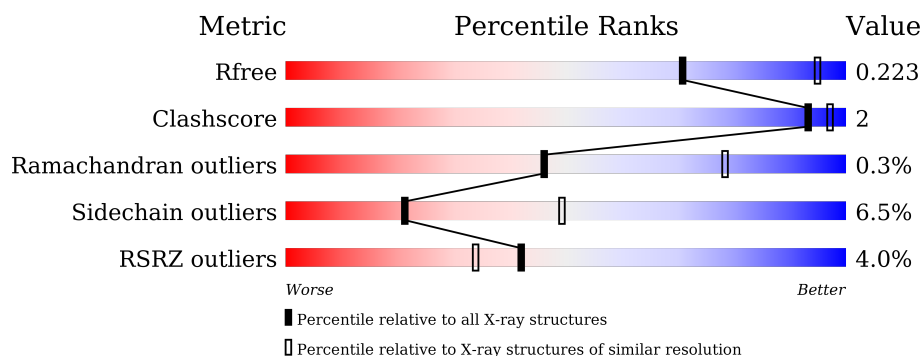
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.79 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	3004 (2.80-2.76)
Clashscore	102246	3480 (2.80-2.76)
Ramachandran outliers	100387	3423 (2.80-2.76)
Sidechain outliers	100360	3425 (2.80-2.76)
RSRZ outliers	91569	3016 (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 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	472	<div> <div>4%</div> <div>85%</div> <div>10%</div> <div>.</div> <div>.</div> </div>
1	B	472	<div> <div>3%</div> <div>88%</div> <div>8%</div> <div>.</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
3	SO4	A	502	-	-	-	X
4	GOL	A	505	-	-	-	X
4	GOL	A	506	-	-	-	X
4	GOL	B	506	-	-	-	X
4	GOL	B	507	-	-	-	X

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 13997 atoms, of which 6914 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 Phospho-2-dehydro-3-deoxyheptonate aldolase AroG.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	451	Total	C	H	N	O	S	0	0	0
			6868	2160	3420	632	639	17			
1	B	454	Total	C	H	N	O	S	0	0	0
			6874	2171	3406	631	649	17			

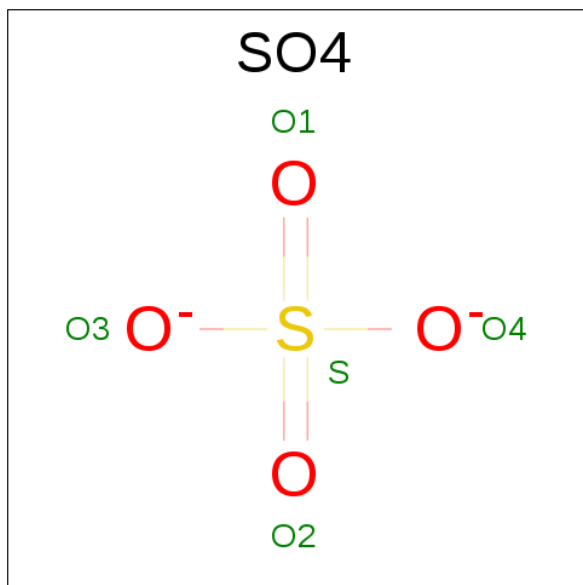
There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	MET	-	initiating methionine	UNP O53512
A	-8	HIS	-	expression tag	UNP O53512
A	-7	HIS	-	expression tag	UNP O53512
A	-6	HIS	-	expression tag	UNP O53512
A	-5	HIS	-	expression tag	UNP O53512
A	-4	HIS	-	expression tag	UNP O53512
A	-3	HIS	-	expression tag	UNP O53512
A	-2	SER	-	expression tag	UNP O53512
A	-1	SER	-	expression tag	UNP O53512
A	0	GLY	-	expression tag	UNP O53512
B	-9	MET	-	initiating methionine	UNP O53512
B	-8	HIS	-	expression tag	UNP O53512
B	-7	HIS	-	expression tag	UNP O53512
B	-6	HIS	-	expression tag	UNP O53512
B	-5	HIS	-	expression tag	UNP O53512
B	-4	HIS	-	expression tag	UNP O53512
B	-3	HIS	-	expression tag	UNP O53512
B	-2	SER	-	expression tag	UNP O53512
B	-1	SER	-	expression tag	UNP O53512
B	0	GLY	-	expression tag	UNP O53512

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

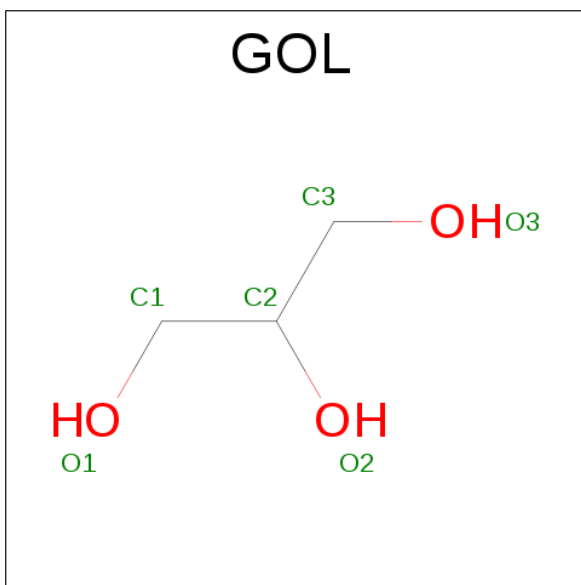
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Cl 1 1	0	0
2	A	1	Total Cl 1 1	0	0

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



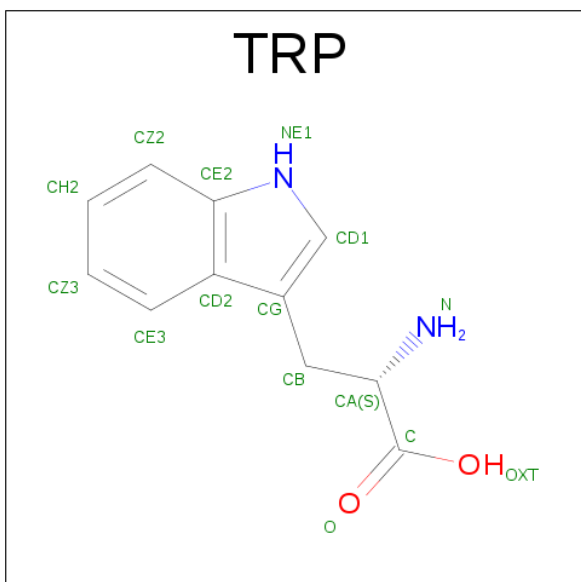
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			14	3	8	3		
4	A	1	Total	C	H	O	0	0
			14	3	8	3		
4	B	1	Total	C	H	O	0	0
			14	3	8	3		
4	B	1	Total	C	H	O	0	0
			14	3	8	3		
4	B	1	Total	C	H	O	0	0
			14	3	8	3		

- Molecule 5 is TRYPTOPHAN (three-letter code: TRP) (formula: $C_{11}H_{12}N_2O_2$).

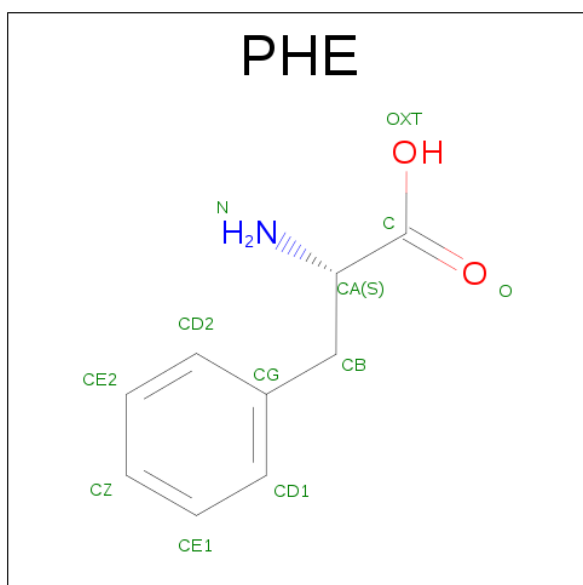


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	H	N	O	0	0
			27	11	12	2	2		
5	B	1	Total	C	H	N	O	0	0
			24	11	9	2	2		

- Molecule 6 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

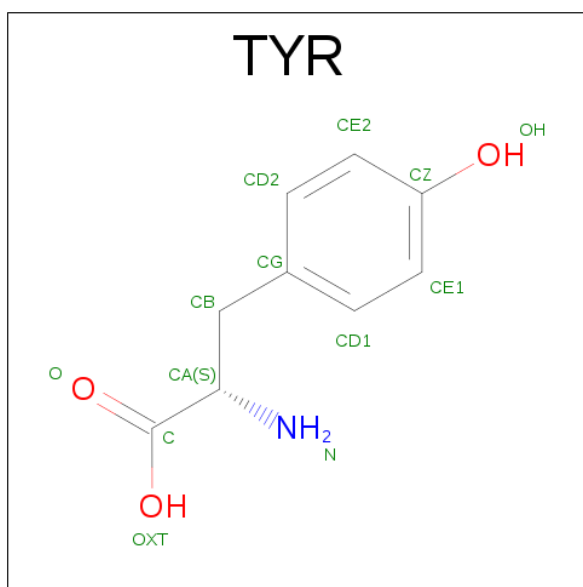
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Mn	0	0
			1	1		
6	A	1	Total	Mn	0	0
			1	1		

- Molecule 7 is PHENYLALANINE (three-letter code: PHE) (formula: C₉H₁₁NO₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total	C	H	N	O	0	0
			20	9	8	1	2		
7	B	1	Total	C	H	N	O	0	0
			20	9	8	1	2		

- Molecule 8 is TYROSINE (three-letter code: TYR) (formula: C₉H₁₁NO₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	B	1	Total	C	H	N	O	0	0
			24	9	11	1	3		

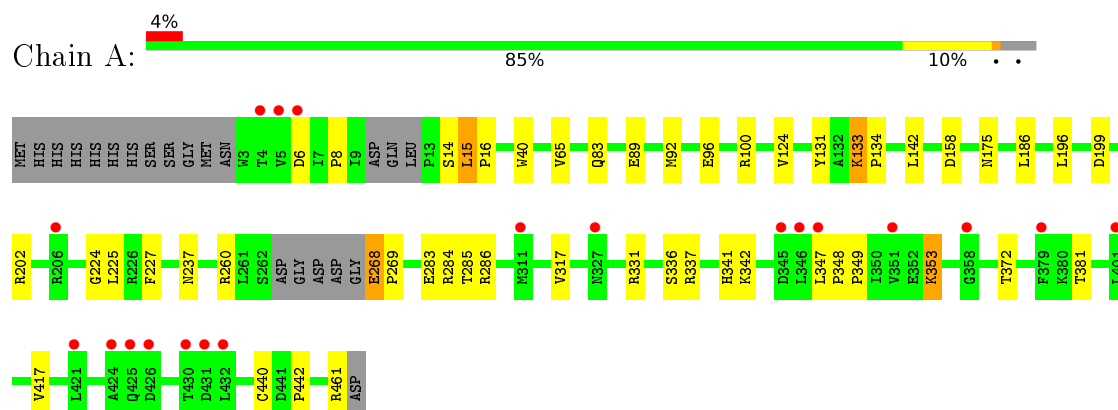
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	18	Total	O	0	0
			18	18		
9	B	23	Total	O	0	0
			23	23		

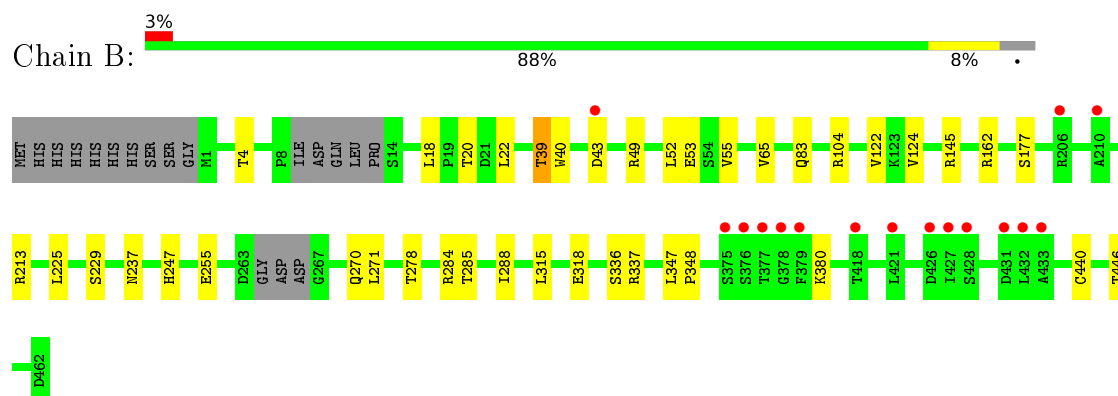
3 Residue-property plots [i](#)

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: Phospho-2-dehydro-3-deoxyheptonate aldolase AroG



- Molecule 1: Phospho-2-dehydro-3-deoxyheptonate aldolase AroG



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	204.76 Å 204.76 Å 66.77 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.30 – 2.79 47.30 – 2.79	Depositor EDS
% Data completeness (in resolution range)	99.4 (47.30-2.79) 99.4 (47.30-2.79)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.94 (at 2.77 Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.175 , 0.226 0.173 , 0.223	Depositor DCC
R_{free} test set	2006 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	55.7	Xtriage
Anisotropy	0.075	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 51.5	EDS
Estimated twinning fraction	0.033 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 40049 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13997	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.97% 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: GOL, MN, SO4, CL

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	0.33	0/3518	0.52	0/4786
1	B	0.33	0/3539	0.52	0/4818
All	All	0.33	0/7057	0.52	0/9604

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	3448	3420	3418	17	0
1	B	3468	3406	3407	9	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	15	0	0	0	0
3	B	10	0	0	0	0
4	A	12	16	16	0	0
4	B	18	24	24	0	0
5	A	15	12	9	0	0
5	B	15	9	9	0	0
6	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	1	0	0	0	0
7	A	12	8	8	1	0
7	B	12	8	8	0	0
8	B	13	11	8	0	0
9	A	18	0	0	0	0
9	B	23	0	0	0	0
All	All	7083	6914	6907	26	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 2.

All (26) 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:133:LYS:NZ	1:A:440:CYS:SG	2.46	0.88
1:A:96:GLU:OE2	1:A:100:ARG:NH1	2.35	0.59
1:A:15:LEU:HB3	1:A:16:PRO:HD2	1.95	0.49
1:B:315:LEU:O	1:B:318:GLU:N	2.46	0.48
1:A:15:LEU:HB3	1:A:16:PRO:CD	2.44	0.48
1:B:83:GLN:HA	1:B:124:VAL:O	2.15	0.46
1:A:196:LEU:O	1:A:199:ASP:HB2	2.16	0.46
1:A:134:PRO:HB2	1:A:284:ARG:HD2	1.97	0.46
1:A:317:VAL:HG21	1:A:353:LYS:HD3	1.98	0.46
1:B:18:LEU:HD21	1:B:271:LEU:HD21	1.97	0.46
1:B:39:THR:OG1	1:B:39:THR:O	2.31	0.45
1:A:283:GLU:OE2	1:A:286:ARG:NH2	2.48	0.45
1:B:52:LEU:O	1:B:55:VAL:HG22	2.17	0.45
1:A:175:ASN:HD21	7:A:509:PHE:N	2.15	0.45
1:A:83:GLN:HA	1:A:124:VAL:O	2.16	0.45
1:B:347:LEU:N	1:B:348:PRO:CD	2.80	0.44
1:A:199:ASP:OD1	1:A:202:ARG:NH2	2.51	0.44
1:A:89:GLU:HB3	1:A:131:TYR:CE1	2.54	0.43
1:A:268:GLU:N	1:A:268:GLU:OE1	2.52	0.43
1:A:224:GLY:O	1:A:227:PHE:HB3	2.19	0.43
1:A:348:PRO:N	1:A:349:PRO:HD2	2.34	0.43
1:B:145:ARG:O	1:B:162:ARG:NH1	2.44	0.43
1:A:347:LEU:N	1:A:348:PRO:CD	2.82	0.41
1:B:49:ARG:O	1:B:53:GLU:HG3	2.21	0.41
1:B:247:HIS:HE1	1:B:255:GLU:OE1	2.03	0.40
1:A:381:THR:HA	1:A:442:PRO:HG3	2.02	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	445/472 (94%)	418 (94%)	25 (6%)	2 (0%)	39	73
1	B	448/472 (95%)	425 (95%)	22 (5%)	1 (0%)	52	84
All	All	893/944 (95%)	843 (94%)	47 (5%)	3 (0%)	46	78

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	440	CYS
1	A	269	PRO
1	A	8	PRO

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	360/385 (94%)	336 (93%)	24 (7%)	20	47
1	B	360/385 (94%)	337 (94%)	23 (6%)	22	50
All	All	720/770 (94%)	673 (94%)	47 (6%)	21	49

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ASP
1	A	14	SER
1	A	15	LEU
1	A	40	TRP
1	A	65	VAL
1	A	92	MET
1	A	133	LYS
1	A	142	LEU
1	A	158	ASP
1	A	186	LEU
1	A	225	LEU
1	A	237	ASN
1	A	260	ARG
1	A	268	GLU
1	A	285	THR
1	A	331	ARG
1	A	336	SER
1	A	337	ARG
1	A	341	HIS
1	A	342	LYS
1	A	353	LYS
1	A	372	THR
1	A	417	VAL
1	A	461	ARG
1	B	4	THR
1	B	20	THR
1	B	22	LEU
1	B	39	THR
1	B	40	TRP
1	B	43	ASP
1	B	65	VAL
1	B	104	ARG
1	B	122	VAL
1	B	177	SER
1	B	213	ARG
1	B	225	LEU
1	B	229	SER
1	B	237	ASN
1	B	270	GLN
1	B	278	THR
1	B	284	ARG
1	B	285	THR
1	B	288	ILE

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Mol	Chain	Res	Type
1	B	336	SER
1	B	337	ARG
1	B	380	LYS
1	B	446	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 19 ligands modelled in this entry, 4 are monoatomic - leaving 15 for Mogul analysis.

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	502	-	4,4,4	0.16	0	6,6,6	0.15	0
3	SO4	A	503	-	4,4,4	0.11	0	6,6,6	0.09	0
3	SO4	A	504	-	4,4,4	0.21	0	6,6,6	0.14	0
4	GOL	A	505	-	5,5,5	0.36	0	5,5,5	0.22	0
4	GOL	A	506	-	5,5,5	0.34	0	5,5,5	0.22	0
5	TRP	A	507	-	11,16,16	0.55	0	9,22,22	0.87	0
7	PHE	A	509	-	9,12,12	0.38	0	10,15,15	0.24	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	B	503	-	4,4,4	0.15	0	6,6,6	0.23	0
3	SO4	B	504	-	4,4,4	0.18	0	6,6,6	0.14	0
4	GOL	B	505	-	5,5,5	0.34	0	5,5,5	0.20	0
4	GOL	B	506	-	5,5,5	0.32	0	5,5,5	0.17	0
4	GOL	B	507	-	5,5,5	0.38	0	5,5,5	0.36	0
8	TYR	B	508	-	10,13,13	0.38	0	12,17,17	0.25	0
5	TRP	B	509	-	11,16,16	0.63	0	9,22,22	0.87	0
7	PHE	B	510	-	9,12,12	0.32	0	10,15,15	0.18	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	502	-	-	0/0/0/0	0/0/0/0
3	SO4	A	503	-	-	0/0/0/0	0/0/0/0
3	SO4	A	504	-	-	0/0/0/0	0/0/0/0
4	GOL	A	505	-	-	0/4/4/4	0/0/0/0
4	GOL	A	506	-	-	0/4/4/4	0/0/0/0
5	TRP	A	507	-	-	0/3/8/8	0/2/2/2
7	PHE	A	509	-	-	0/4/8/8	0/1/1/1
3	SO4	B	503	-	-	0/0/0/0	0/0/0/0
3	SO4	B	504	-	-	0/0/0/0	0/0/0/0
4	GOL	B	505	-	-	0/4/4/4	0/0/0/0
4	GOL	B	506	-	-	0/4/4/4	0/0/0/0
4	GOL	B	507	-	-	0/4/4/4	0/0/0/0
8	TYR	B	508	-	-	0/4/8/8	0/1/1/1
5	TRP	B	509	-	-	0/3/8/8	0/2/2/2
7	PHE	B	510	-	-	0/4/8/8	0/1/1/1

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
7	A	509	PHE	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	451/472 (95%)	0.31	20 (4%) 38 30	27, 51, 80, 116	0
1	B	454/472 (96%)	0.20	16 (3%) 48 40	24, 44, 74, 105	0
All	All	905/944 (95%)	0.25	36 (3%) 42 34	24, 48, 77, 116	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	376	SER	4.8
1	A	351	VAL	3.7
1	B	375	SER	3.6
1	B	432	LEU	3.4
1	A	346	LEU	3.2
1	A	206	ARG	3.2
1	A	430	THR	3.1
1	A	311	MET	3.1
1	B	426	ASP	3.0
1	A	347	LEU	2.9
1	B	377	THR	2.9
1	B	379	PHE	2.9
1	A	6	ASP	2.9
1	A	432	LEU	2.7
1	B	210	ALA	2.7
1	A	5	VAL	2.6
1	A	327	ASN	2.6
1	A	358	GLY	2.6
1	A	426	ASP	2.6
1	A	401	LEU	2.5
1	A	379	PHE	2.5
1	B	206	ARG	2.4
1	B	421	LEU	2.4
1	A	425	GLN	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	421	LEU	2.3
1	B	43	ASP	2.3
1	B	418	THR	2.3
1	A	431	ASP	2.3
1	A	4	THR	2.2
1	B	427	ILE	2.2
1	B	428	SER	2.1
1	B	378	GLY	2.1
1	B	433	ALA	2.1
1	A	424	ALA	2.1
1	A	345	ASP	2.1
1	B	431	ASP	2.1

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(\AA^2)	Q<0.9
4	GOL	B	507	6/6	0.92	0.27	6.27	45,66,79,82	0
4	GOL	B	506	6/6	0.84	0.41	5.49	65,80,96,99	0
4	GOL	A	505	6/6	0.88	0.31	4.52	53,63,73,74	0
3	SO4	A	502	5/5	0.75	0.29	3.15	88,107,123,153	0
4	GOL	A	506	6/6	0.90	0.26	2.37	51,68,79,84	0
4	GOL	B	505	6/6	0.75	0.23	1.89	55,81,107,107	0
7	PHE	B	510	12/12	0.94	0.29	1.82	31,54,74,76	20
3	SO4	B	503	5/5	0.97	0.19	1.31	47,49,56,64	0
7	PHE	A	509	12/12	0.93	0.24	0.92	37,50,61,61	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	TRP	A	507	15/15	0.96	0.19	0.43	35,47,60,62	0
3	SO4	A	504	5/5	0.97	0.17	-0.16	51,56,63,77	0
3	SO4	A	503	5/5	0.91	0.18	-0.16	81,84,89,109	0
8	TYR	B	508	13/13	0.96	0.19	-0.20	33,45,61,72	0
5	TRP	B	509	15/15	0.97	0.15	-0.57	30,40,46,54	0
6	MN	A	508	1/1	0.98	0.11	-1.24	65,65,65,65	0
3	SO4	B	504	5/5	0.94	0.12	-1.50	84,86,98,131	0
6	MN	B	502	1/1	0.94	0.11	-1.54	59,59,59,59	1
2	CL	B	501	1/1	0.96	0.10	-2.42	54,54,54,54	0
2	CL	A	501	1/1	0.94	0.23	-	67,67,67,67	0

6.5 Other polymers [i](#)

There are no such residues in this entry.