



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:28 PM GMT

PDB ID : 1TT4  
Title : Structure of NP459575, a predicted glutathione synthase from *Salmonella typhimurium*  
Authors : Miller, D.J.; Shuvalova, L.; Anderson, W.F.; Midwest Center for Structural Genomics (MCSG)  
Deposited on : 2004-06-21  
Resolution : 2.80 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](mailto: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.

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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

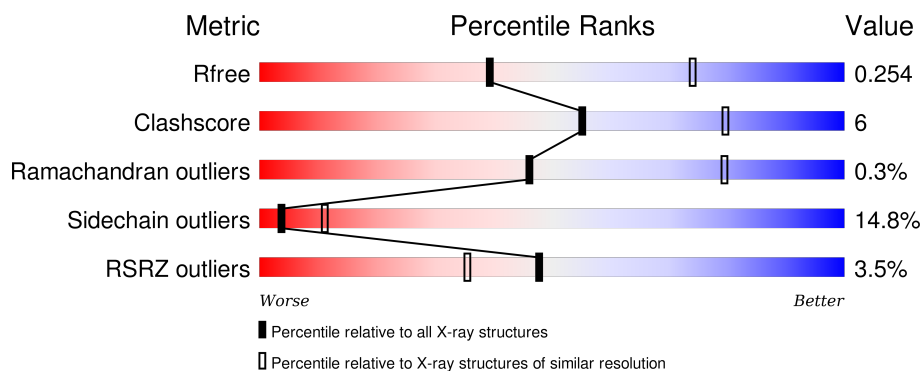
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

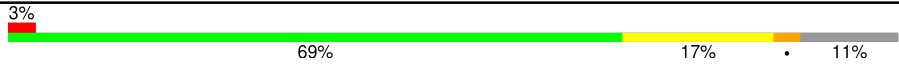

The reported resolution of this entry is 2.80 Å.

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	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	396	
1	B	396	

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	MG	B	401	-	-	-	X

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5624 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 cytoplasmic protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	354	Total	C	N	O	S	0	2	0
			2787	1765	489	517	16			
1	B	354	Total	C	N	O	S	0	1	0
			2786	1767	487	516	16			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-23	MET	-	INITIATING METHIONINE	UNP Q8ZR41
A	-22	HIS	-	EXPRESSION TAG	UNP Q8ZR41
A	-21	HIS	-	EXPRESSION TAG	UNP Q8ZR41
A	-20	HIS	-	EXPRESSION TAG	UNP Q8ZR41
A	-19	HIS	-	EXPRESSION TAG	UNP Q8ZR41
A	-18	HIS	-	EXPRESSION TAG	UNP Q8ZR41
A	-17	HIS	-	EXPRESSION TAG	UNP Q8ZR41
A	-16	SER	-	CLONING ARTIFACT	UNP Q8ZR41
A	-15	SER	-	CLONING ARTIFACT	UNP Q8ZR41
A	-14	GLY	-	CLONING ARTIFACT	UNP Q8ZR41
A	-13	VAL	-	CLONING ARTIFACT	UNP Q8ZR41
A	-12	ASP	-	CLONING ARTIFACT	UNP Q8ZR41
A	-11	LEU	-	CLONING ARTIFACT	UNP Q8ZR41
A	-10	GLY	-	CLONING ARTIFACT	UNP Q8ZR41
A	-9	THR	-	CLONING ARTIFACT	UNP Q8ZR41
A	-8	GLU	-	CLONING ARTIFACT	UNP Q8ZR41
A	-7	ASN	-	CLONING ARTIFACT	UNP Q8ZR41
A	-6	LEU	-	CLONING ARTIFACT	UNP Q8ZR41
A	-5	TYR	-	CLONING ARTIFACT	UNP Q8ZR41
A	-4	PHE	-	CLONING ARTIFACT	UNP Q8ZR41
A	-3	GLN	-	CLONING ARTIFACT	UNP Q8ZR41
A	-2	SER	-	CLONING ARTIFACT	UNP Q8ZR41
A	-1	ASN	-	CLONING ARTIFACT	UNP Q8ZR41
A	0	ALA	-	CLONING ARTIFACT	UNP Q8ZR41
B	-23	MET	-	INITIATING METHIONINE	UNP Q8ZR41

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-22	HIS	-	EXPRESSION TAG	UNP Q8ZR41
B	-21	HIS	-	EXPRESSION TAG	UNP Q8ZR41
B	-20	HIS	-	EXPRESSION TAG	UNP Q8ZR41
B	-19	HIS	-	EXPRESSION TAG	UNP Q8ZR41
B	-18	HIS	-	EXPRESSION TAG	UNP Q8ZR41
B	-17	HIS	-	EXPRESSION TAG	UNP Q8ZR41
B	-16	SER	-	CLONING ARTIFACT	UNP Q8ZR41
B	-15	SER	-	CLONING ARTIFACT	UNP Q8ZR41
B	-14	GLY	-	CLONING ARTIFACT	UNP Q8ZR41
B	-13	VAL	-	CLONING ARTIFACT	UNP Q8ZR41
B	-12	ASP	-	CLONING ARTIFACT	UNP Q8ZR41
B	-11	LEU	-	CLONING ARTIFACT	UNP Q8ZR41
B	-10	GLY	-	CLONING ARTIFACT	UNP Q8ZR41
B	-9	THR	-	CLONING ARTIFACT	UNP Q8ZR41
B	-8	GLU	-	CLONING ARTIFACT	UNP Q8ZR41
B	-7	ASN	-	CLONING ARTIFACT	UNP Q8ZR41
B	-6	LEU	-	CLONING ARTIFACT	UNP Q8ZR41
B	-5	TYR	-	CLONING ARTIFACT	UNP Q8ZR41
B	-4	PHE	-	CLONING ARTIFACT	UNP Q8ZR41
B	-3	GLN	-	CLONING ARTIFACT	UNP Q8ZR41
B	-2	SER	-	CLONING ARTIFACT	UNP Q8ZR41
B	-1	ASN	-	CLONING ARTIFACT	UNP Q8ZR41
B	0	ALA	-	CLONING ARTIFACT	UNP Q8ZR41

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

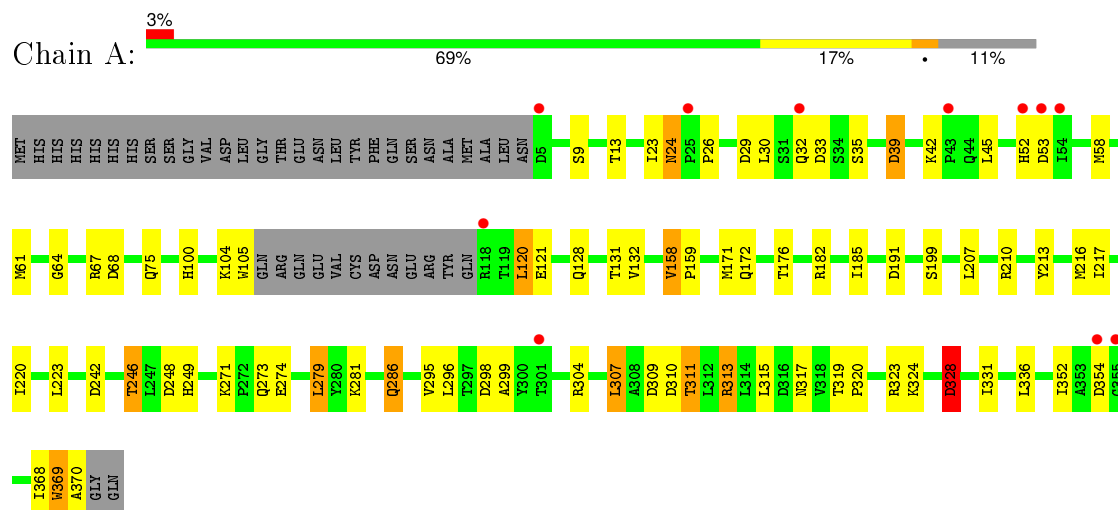
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	19	Total	O	0	0
			19	19		
4	B	20	Total	O	0	0
			20	20		

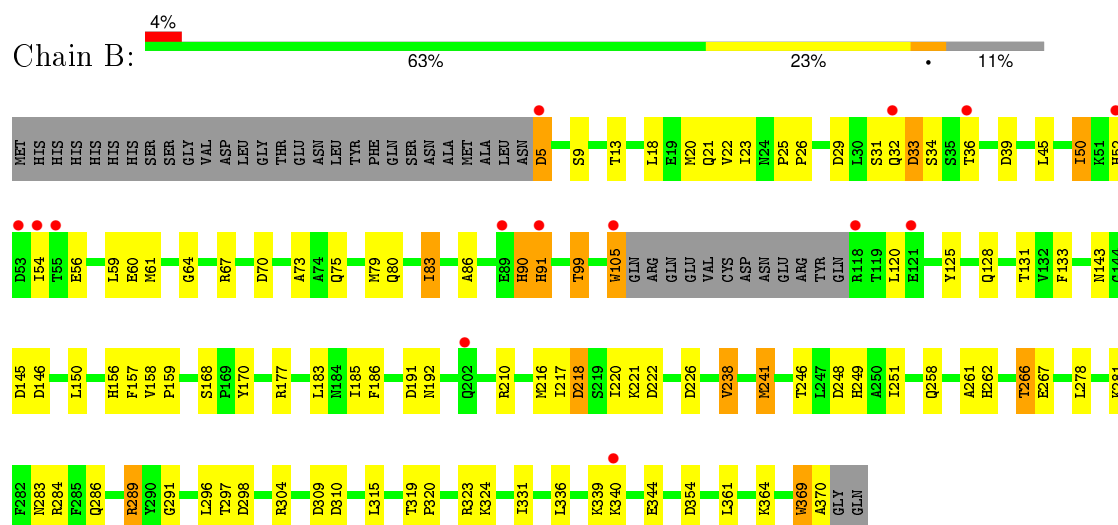
### 3 Residue-property plots [i](#)

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: putative cytoplasmic protein



- Molecule 1: putative cytoplasmic protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	188.93Å 188.93Å 188.93Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.92 – 2.80 19.91 – 2.80	Depositor EDS
% Data completeness (in resolution range)	95.0 (19.92-2.80) 100.0 (19.91-2.80)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	7.69 (at 2.79Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, $R_{free}$	0.209 , 0.258 0.203 , 0.254	Depositor DCC
$R_{free}$ test set	1391 reflections (5.31%)	DCC
Wilson B-factor (Å <sup>2</sup> )	55.4	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 38.7	EDS
Estimated twinning fraction	0.027 for -l,-k,-h	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 27578 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	5624	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.93% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: MG, 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	0.40	0/2855	0.76	11/3876 (0.3%)
1	B	0.42	0/2855	0.78	15/3877 (0.4%)
All	All	0.41	0/5710	0.77	26/7753 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	B	145	ASP	CB-CG-OD2	6.77	124.40	118.30
1	A	328	ASP	CB-CG-OD2	6.54	124.19	118.30
1	B	39	ASP	CB-CG-OD2	6.07	123.77	118.30
1	A	33	ASP	CB-CG-OD2	6.06	123.75	118.30
1	A	354	ASP	CB-CG-OD2	6.04	123.73	118.30
1	B	309	ASP	CB-CG-OD2	5.97	123.68	118.30
1	B	310	ASP	CB-CG-OD2	5.97	123.67	118.30
1	B	298	ASP	CB-CG-OD2	5.78	123.50	118.30
1	B	191	ASP	CB-CG-OD2	5.63	123.36	118.30
1	B	33	ASP	CB-CG-OD2	5.61	123.35	118.30
1	A	298	ASP	CB-CG-OD2	5.58	123.32	118.30
1	A	191	ASP	CB-CG-OD2	5.54	123.28	118.30
1	B	29	ASP	CB-CG-OD2	5.48	123.23	118.30
1	B	70	ASP	CB-CG-OD2	5.33	123.10	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	5	ASP	CB-CG-OD2	5.33	123.09	118.30
1	A	309	ASP	CB-CG-OD2	5.27	123.05	118.30
1	A	29	ASP	CB-CG-OD2	5.20	122.98	118.30
1	B	248	ASP	CB-CG-OD2	5.20	122.98	118.30
1	B	218	ASP	CB-CG-OD2	5.19	122.97	118.30
1	B	226	ASP	CB-CG-OD2	5.15	122.94	118.30
1	B	222	ASP	CB-CG-OD2	5.11	122.90	118.30
1	B	146	ASP	CB-CG-OD2	5.10	122.89	118.30
1	A	68	ASP	CB-CG-OD2	5.06	122.85	118.30
1	A	248	ASP	CB-CG-OD2	5.06	122.85	118.30
1	A	39	ASP	CB-CG-OD2	5.02	122.81	118.30
1	A	310	ASP	CB-CG-OD2	5.02	122.81	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	291	GLY	Peptide

## 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	2787	0	2699	32	0
1	B	2786	0	2701	44	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	19	0	0	0	0
4	B	20	0	0	1	0
All	All	5624	0	5400	69	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 6.

All (69) 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:52[B]:HIS:H	1:A:52[B]:HIS:CD2	1.95	0.85
1:B:262:HIS:O	1:B:266:THR:HB	1.80	0.82
1:B:79:MET:O	1:B:83:ILE:HG13	1.82	0.80
1:A:64:GLY:H	1:A:75:GLN:HE22	1.29	0.80
1:A:216:MET:CE	1:B:286:GLN:HE21	1.98	0.77
1:A:121:GLU:O	1:B:289:ARG:NH1	2.22	0.72
1:B:91:HIS:ND1	1:B:91:HIS:O	2.25	0.70
1:B:99:THR:HG21	1:B:170:TYR:O	1.94	0.68
1:B:238:VAL:HG22	1:B:241:MET:HE1	1.78	0.66
1:A:216:MET:HE2	1:B:286:GLN:HE21	1.62	0.65
1:A:216:MET:HE3	1:B:286:GLN:HE21	1.63	0.64
1:B:99:THR:CG2	1:B:170:TYR:O	2.49	0.60
1:B:99:THR:HG22	4:B:519:HOH:O	2.02	0.59
1:A:30:LEU:HG	1:A:100:HIS:CD2	2.37	0.59
1:B:91:HIS:CG	1:B:91:HIS:O	2.57	0.58
1:A:104:LYS:O	1:A:105:TRP:C	2.40	0.58
1:B:90:HIS:O	1:B:91:HIS:HB3	2.03	0.58
1:B:32:GLN:O	1:B:56:GLU:HB2	2.04	0.57
1:A:120:LEU:HD21	1:A:128:GLN:NE2	2.20	0.57
1:B:86:ALA:O	1:B:90:HIS:HB2	2.05	0.57
1:A:369:TRP:O	1:A:370:ALA:CB	2.53	0.57
1:A:286:GLN:HG2	1:A:295:VAL:H	1.72	0.55
1:B:64:GLY:H	1:B:75:GLN:HE22	1.55	0.54
1:A:279:LEU:HD22	1:A:299:ALA:HA	1.90	0.54
1:A:182:ARG:HD3	1:A:242:ASP:OD1	2.07	0.54
1:A:53:ASP:OD2	1:A:58:MET:SD	2.67	0.53
1:A:323:ARG:HA	1:A:328:ASP:OD1	2.10	0.52
1:B:23:ILE:O	1:B:25:PRO:HD3	2.10	0.51
1:A:52[B]:HIS:CD2	1:A:52[B]:HIS:N	2.72	0.51
1:B:22:VAL:HG21	1:B:34:SER:HB2	1.93	0.50
1:A:286:GLN:CG	1:A:295:VAL:H	2.23	0.50
1:B:105:TRP:HZ2	1:B:128:GLN:HG2	1.76	0.50
1:B:18:LEU:HB3	1:B:20:MET:HE2	1.94	0.50
1:A:158:VAL:N	1:A:159:PRO:CD	2.75	0.49
1:B:283:ASN:HD21	1:B:297:THR:H	1.58	0.49
1:A:207:LEU:CD1	1:B:278:LEU:HD11	2.41	0.49
1:B:73:ALA:HA	1:B:251:ILE:HD11	1.95	0.49
1:B:105:TRP:CZ2	1:B:128:GLN:HG2	2.48	0.49
1:B:61:MET:HE1	1:B:369:TRP:HZ2	1.78	0.49
1:B:156[B]:HIS:HD2	1:B:157:PHE:CE1	2.30	0.48
1:B:21:GLN:HG2	1:B:133:PHE:CZ	2.48	0.48
1:B:246:THR:HG22	1:B:249:HIS:ND1	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:249:HIS:ND1	1:B:344:GLU:OE1	2.40	0.47
1:A:286:GLN:HG2	1:A:295:VAL:N	2.29	0.47
1:B:59:LEU:HD21	1:B:83:ILE:HG21	1.97	0.46
1:B:319:THR:HB	1:B:320:PRO:HD3	1.98	0.45
1:A:217:ILE:HD13	1:A:223:LEU:HD21	1.98	0.45
1:A:307:LEU:O	1:A:311:THR:HG22	2.15	0.45
1:B:369:TRP:O	1:B:370:ALA:HB2	2.17	0.45
1:A:369:TRP:O	1:A:370:ALA:HB2	2.17	0.45
1:B:61:MET:HE1	1:B:369:TRP:CZ2	2.52	0.45
1:B:21:GLN:HG2	1:B:133:PHE:CE2	2.52	0.45
1:B:45:LEU:HD11	1:B:50:ILE:HG13	1.99	0.45
1:B:266:THR:HG22	1:B:267:GLU:HG2	1.99	0.44
1:A:246:THR:HG22	1:A:249:HIS:H	1.82	0.44
1:A:23:ILE:HG22	1:A:24:ASN:N	2.31	0.44
1:B:99:THR:HB	1:B:168:SER:HB2	1.99	0.44
1:A:319:THR:HB	1:A:320:PRO:HD3	1.98	0.44
1:B:158:VAL:N	1:B:159:PRO:CD	2.82	0.43
1:B:186:PHE:HB3	1:B:192:ASN:ND2	2.34	0.42
1:A:311:THR:O	1:A:315:LEU:HG	2.19	0.42
1:B:61:MET:CE	1:B:369:TRP:CZ2	3.02	0.42
1:A:246:THR:HG21	1:A:249:HIS:CD2	2.54	0.41
1:A:213:TYR:O	1:B:304:ARG:NH1	2.53	0.41
1:A:313:ARG:HH11	1:A:317:ASN:HD21	1.67	0.41
1:A:176:THR:HB	1:B:125:TYR:CD2	2.56	0.41
1:B:143:ASN:H	1:B:143:ASN:ND2	2.18	0.41
1:B:150:LEU:HD11	1:B:261:ALA:HA	2.02	0.41
1:A:307:LEU:O	1:A:311:THR:CG2	2.69	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	352/396 (89%)	335 (95%)	16 (4%)	1 (0%)	46	79
1	B	351/396 (89%)	335 (95%)	15 (4%)	1 (0%)	46	79
All	All	703/792 (89%)	670 (95%)	31 (4%)	2 (0%)	46	79

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	26	PRO
1	B	26	PRO

### 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	293/329 (89%)	254 (87%)	39 (13%)	5	14
1	B	293/329 (89%)	246 (84%)	47 (16%)	3	9
All	All	586/658 (89%)	500 (85%)	86 (15%)	4	11

All (86) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	SER
1	A	13	THR
1	A	24	ASN
1	A	32	GLN
1	A	35	SER
1	A	39	ASP
1	A	42	LYS
1	A	45	LEU
1	A	61	MET
1	A	67	ARG
1	A	120	LEU
1	A	131	THR
1	A	132	VAL
1	A	158	VAL

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Mol	Chain	Res	Type
1	A	171	MET
1	A	172	GLN
1	A	185	ILE
1	A	199	SER
1	A	210	ARG
1	A	220	ILE
1	A	246	THR
1	A	271	LYS
1	A	273	GLN
1	A	274	GLU
1	A	279	LEU
1	A	281	LYS
1	A	286	GLN
1	A	296	LEU
1	A	304	ARG
1	A	307	LEU
1	A	311	THR
1	A	313	ARG
1	A	324	LYS
1	A	328	ASP
1	A	331	ILE
1	A	336	LEU
1	A	352	ILE
1	A	368	ILE
1	A	369	TRP
1	B	5	ASP
1	B	9	SER
1	B	13	THR
1	B	31	SER
1	B	33	ASP
1	B	36	THR
1	B	50	ILE
1	B	52	HIS
1	B	54	ILE
1	B	60	GLU
1	B	67	ARG
1	B	80	GLN
1	B	83	ILE
1	B	90	HIS
1	B	91	HIS
1	B	99	THR
1	B	105	TRP

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Mol	Chain	Res	Type
1	B	120	LEU
1	B	131	THR
1	B	177	ARG
1	B	183	LEU
1	B	185	ILE
1	B	210	ARG
1	B	216	MET
1	B	217	ILE
1	B	218	ASP
1	B	220	ILE
1	B	221	LYS
1	B	238	VAL
1	B	241	MET
1	B	258	GLN
1	B	266	THR
1	B	281	LYS
1	B	284	ARG
1	B	289	ARG
1	B	296	LEU
1	B	315	LEU
1	B	323	ARG
1	B	324	LYS
1	B	331	ILE
1	B	336	LEU
1	B	339	LYS
1	B	340	LYS
1	B	354	ASP
1	B	361	LEU
1	B	364	LYS
1	B	369	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	ASN
1	A	75	GLN
1	A	103	GLN
1	A	249	HIS
1	A	317	ASN
1	B	21	GLN
1	B	24	ASN
1	B	75	GLN

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Mol	Chain	Res	Type
1	B	192	ASN
1	B	258	GLN
1	B	283	ASN
1	B	286	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 carbohydrates in this entry.

### 5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	500	-	4,4,4	0.23	0	6,6,6	0.35	0
3	SO4	B	501	2	4,4,4	0.17	0	6,6,6	0.31	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	500	-	-	0/0/0/0	0/0/0/0
3	SO4	B	501	2	-	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.

No monomer is involved in short contacts.

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	354/396 (89%)	-0.11	11 (3%)	52 40	29, 48, 75, 94	0
1	B	354/396 (89%)	-0.10	14 (3%)	42 30	23, 43, 78, 90	0
All	All	708/792 (89%)	-0.11	25 (3%)	48 35	23, 45, 77, 94	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	54	ILE	5.3
1	B	91	HIS	4.6
1	B	105	TRP	4.5
1	B	118	ARG	3.9
1	B	32	GLN	3.8
1	A	25	PRO	3.6
1	B	121	GLU	3.5
1	A	118	ARG	3.4
1	A	52[A]	HIS	3.4
1	B	5	ASP	3.3
1	B	52	HIS	3.2
1	B	53	ASP	3.2
1	A	355	GLY	2.9
1	B	89	GLU	2.9
1	B	340	LYS	2.9
1	A	53	ASP	2.8
1	A	301	THR	2.8
1	A	32	GLN	2.7
1	A	54	ILE	2.6
1	A	354	ASP	2.6
1	B	202	GLN	2.5
1	B	36	THR	2.5
1	B	55	THR	2.3
1	A	5	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	43	PRO	2.2

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	MG	B	401	1/1	0.77	0.20	4.75	43,43,43,43	0
3	SO4	A	500	5/5	0.83	0.28	1.17	71,85,86,88	0
3	SO4	B	501	5/5	0.89	0.21	0.43	54,64,77,80	0
2	MG	A	400	1/1	0.83	0.08	-2.46	38,38,38,38	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.