



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:36 PM GMT

PDB ID : 4PEF  
Title : Dbr1 in complex with sulfate  
Authors : Montemayor, E.J.; Katolik, A.; Clark, N.E.; Taylor, A.B.; Schuermann, J.P.; Combs, D.J.; Johnsson, R.; Holloway, S.P.; Stevens, S.W.; Damha, M.J.; Hart, P.J.  
Deposited on : 2014-04-23  
Resolution : 1.96 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

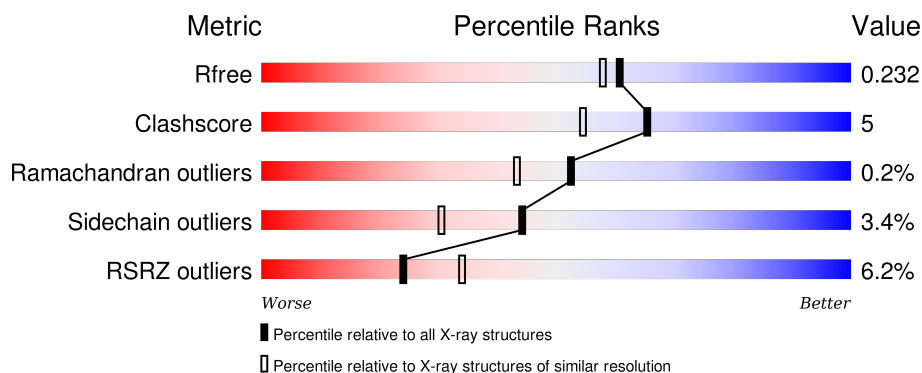
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.96 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	1833 (1.96-1.96)
Clashscore	102246	1953 (1.96-1.96)
Ramachandran outliers	100387	1936 (1.96-1.96)
Sidechain outliers	100360	1936 (1.96-1.96)
RSRZ outliers	91569	1835 (1.96-1.96)

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	360	<div> <div>5%</div> <div>86%</div> <div>11%</div> <div>• •</div> </div>
1	B	360	<div> <div>3%</div> <div>84%</div> <div>13%</div> <div>• •</div> </div>
1	C	360	<div> <div>6%</div> <div>85%</div> <div>11%</div> <div>• •</div> </div>
1	D	360	<div> <div>8%</div> <div>85%</div> <div>11%</div> <div>•</div> </div>
1	E	360	<div> <div>8%</div> <div>83%</div> <div>13%</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	E	402	-	-	-	X
4	GOL	A	404	-	-	-	X
4	GOL	D	404	-	-	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 15545 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 lariat debranching enzyme, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	349	Total	C	N	O	S	0	3	0
			2885	1873	470	527	15			
1	B	349	Total	C	N	O	S	0	3	0
			2876	1868	469	523	16			
1	C	349	Total	C	N	O	S	0	1	0
			2867	1863	468	521	15			
1	D	349	Total	C	N	O	S	0	2	0
			2870	1865	468	521	16			
1	E	349	Total	C	N	O	S	0	2	0
			2870	1865	468	521	16			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	355	HIS	-	expression tag	UNP C4M1P9
A	356	HIS	-	expression tag	UNP C4M1P9
A	357	HIS	-	expression tag	UNP C4M1P9
A	358	HIS	-	expression tag	UNP C4M1P9
A	359	HIS	-	expression tag	UNP C4M1P9
A	360	HIS	-	expression tag	UNP C4M1P9
B	355	HIS	-	expression tag	UNP C4M1P9
B	356	HIS	-	expression tag	UNP C4M1P9
B	357	HIS	-	expression tag	UNP C4M1P9
B	358	HIS	-	expression tag	UNP C4M1P9
B	359	HIS	-	expression tag	UNP C4M1P9
B	360	HIS	-	expression tag	UNP C4M1P9
C	355	HIS	-	expression tag	UNP C4M1P9
C	356	HIS	-	expression tag	UNP C4M1P9
C	357	HIS	-	expression tag	UNP C4M1P9
C	358	HIS	-	expression tag	UNP C4M1P9
C	359	HIS	-	expression tag	UNP C4M1P9
C	360	HIS	-	expression tag	UNP C4M1P9
D	355	HIS	-	expression tag	UNP C4M1P9

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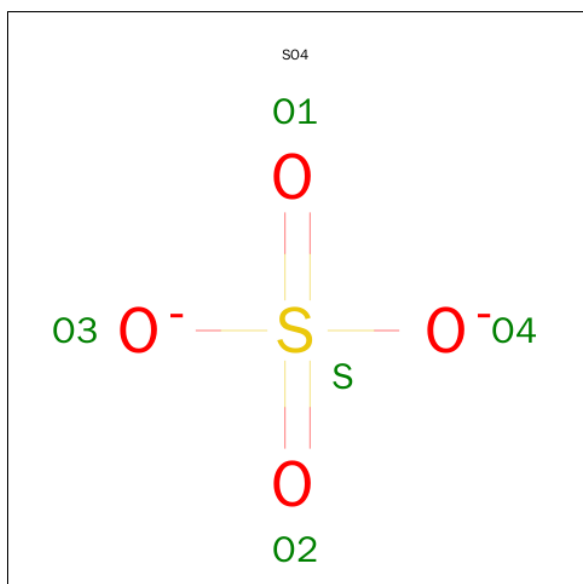
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Chain	Residue	Modelled	Actual	Comment	Reference
D	356	HIS	-	expression tag	UNP C4M1P9
D	357	HIS	-	expression tag	UNP C4M1P9
D	358	HIS	-	expression tag	UNP C4M1P9
D	359	HIS	-	expression tag	UNP C4M1P9
D	360	HIS	-	expression tag	UNP C4M1P9
E	355	HIS	-	expression tag	UNP C4M1P9
E	356	HIS	-	expression tag	UNP C4M1P9
E	357	HIS	-	expression tag	UNP C4M1P9
E	358	HIS	-	expression tag	UNP C4M1P9
E	359	HIS	-	expression tag	UNP C4M1P9
E	360	HIS	-	expression tag	UNP C4M1P9

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

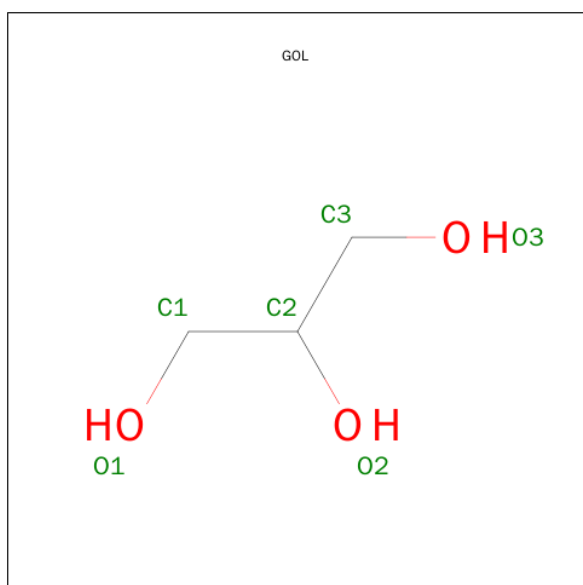
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Mn 1 1	0	0
2	A	1	Total Mn 1 1	0	0
2	D	1	Total Mn 1 1	0	0
2	C	1	Total Mn 1 1	0	0
2	E	1	Total Mn 1 1	0	0

- 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 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
3	C	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0
3	E	1	Total O S 5 4 1	0	0
3	E	1	Total O S 5 4 1	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 6 3 3	0	0
4	B	1	Total C O 6 3 3	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	D	1	Total	C	O	0	0
			6	3	3		

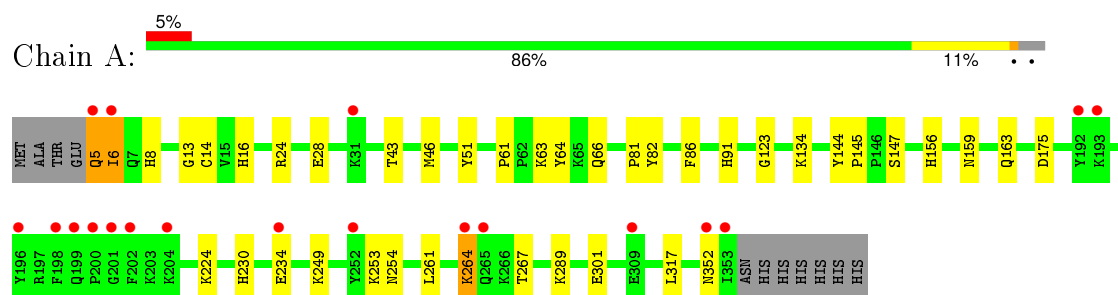
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	227	Total	O	0	0
			227	227		
5	B	236	Total	O	0	0
			236	236		
5	C	212	Total	O	0	0
			212	212		
5	D	220	Total	O	0	0
			220	220		
5	E	209	Total	O	0	0
			209	209		

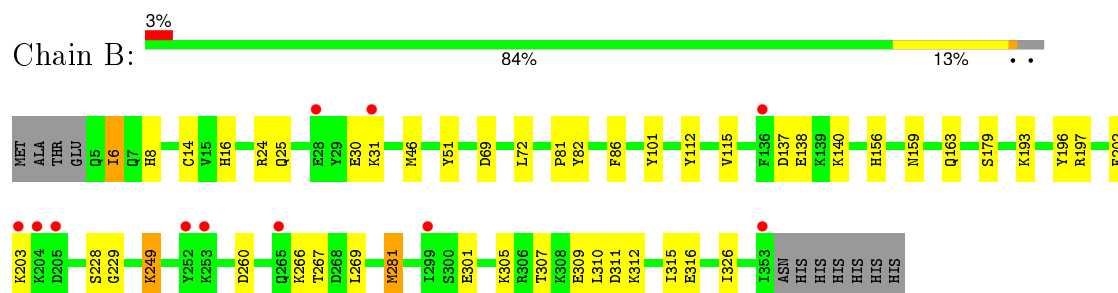
### 3 Residue-property plots [i](#)

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.

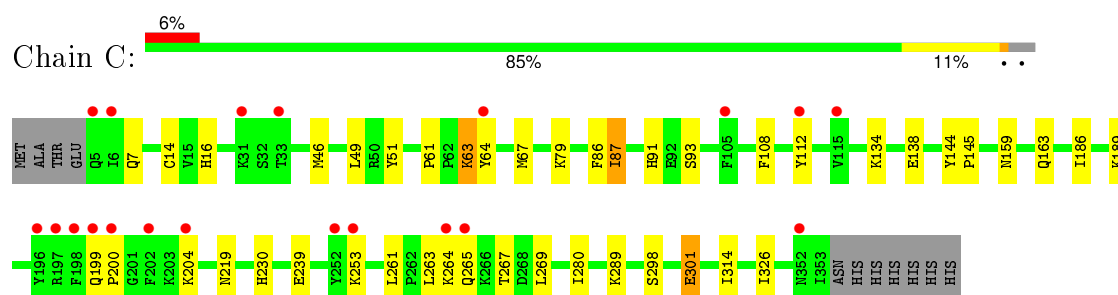
- Molecule 1: RNA lariat debranching enzyme, putative



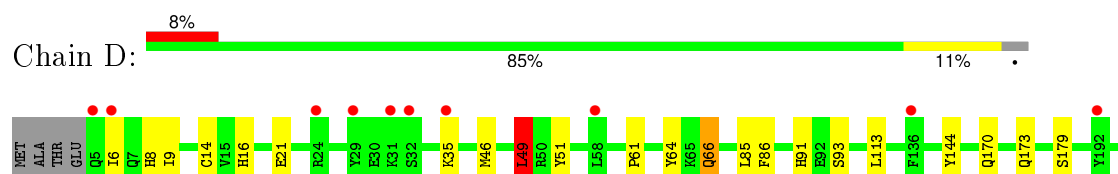
- Molecule 1: RNA lariat debranching enzyme, putative



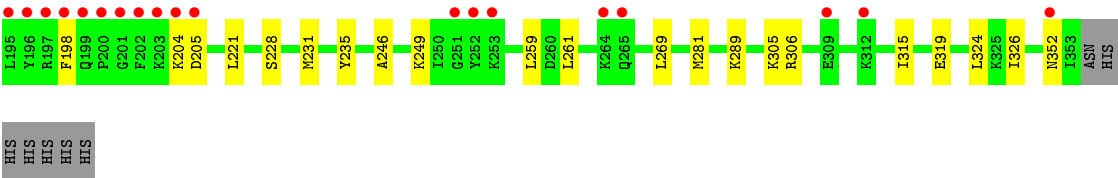
- Molecule 1: RNA lariat debranching enzyme, putative



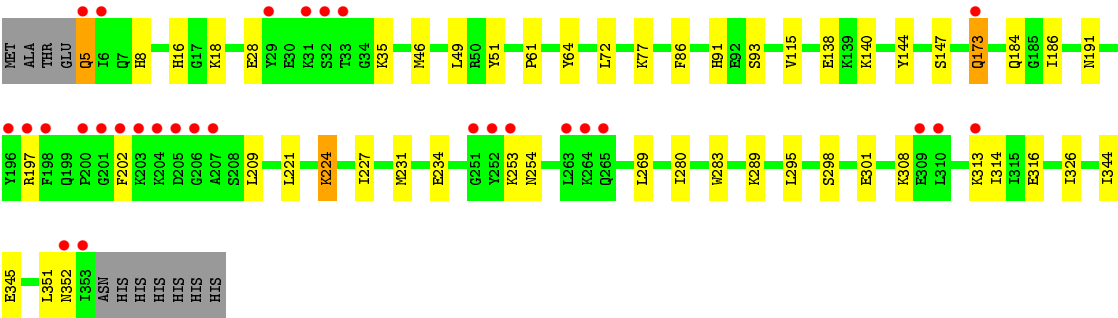
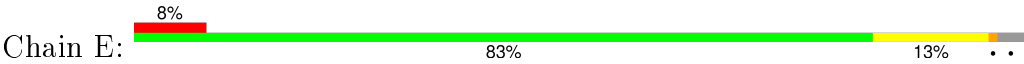
- Molecule 1: RNA lariat debranching enzyme, putative







- Molecule 1: RNA lariat debranching enzyme, putative



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.22Å 141.69Å 213.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.18 – 1.96 43.18 – 1.96	Depositor EDS
% Data completeness (in resolution range)	96.9 (43.18-1.96) 96.9 (43.18-1.96)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.78 (at 1.97Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, $R_{free}$	0.191 , 0.232 0.191 , 0.232	Depositor DCC
$R_{free}$ test set	7760 reflections (5.02%)	DCC
Wilson B-factor (Å <sup>2</sup> )	27.5	Xtriage
Anisotropy	0.120	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 46.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 154587 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	15545	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.70% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, MN, 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	0.39	0/2964	0.55	0/4006
1	B	0.42	0/2958	0.54	0/3998
1	C	0.40	0/2946	0.54	0/3982
1	D	0.38	0/2952	0.56	1/3990 (0.0%)
1	E	0.42	0/2952	0.57	0/3990
All	All	0.40	0/14772	0.55	1/19966 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	49	LEU	CA-CB-CG	5.19	127.24	115.30

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	2885	0	2848	22	0
1	B	2876	0	2847	35	0
1	C	2867	0	2838	29	0
1	D	2870	0	2843	25	0
1	E	2870	0	2843	33	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
3	A	10	0	0	0	0
3	B	10	0	0	0	0
3	C	10	0	0	0	0
3	D	10	0	0	0	0
3	E	10	0	0	0	0
4	A	6	0	8	0	0
4	B	6	0	8	1	0
4	D	6	0	8	0	0
5	A	227	0	0	2	0
5	B	236	0	0	6	0
5	C	212	0	0	3	0
5	D	220	0	0	4	0
5	E	209	0	0	4	0
All	All	15545	0	14243	144	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 5.

All (144) 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:281:MET:SD	5:B:622:HOH:O	2.31	0.89
1:D:14[A]:CYS:SG	5:D:542:HOH:O	2.29	0.74
1:A:24:ARG:NH1	1:A:28[A]:GLU:OE2	2.22	0.72
1:A:264:LYS:H	1:A:264:LYS:HE2	1.53	0.72
1:B:6:ILE:HD11	1:B:8:HIS:CE1	2.26	0.71
1:B:14[B]:CYS:HB3	1:B:16[B]:HIS:CD2	2.26	0.71
1:A:66:GLN:NE2	5:A:501:HOH:O	2.24	0.70
1:E:173:GLN:H	1:E:173:GLN:HE21	1.40	0.70
1:D:16[A]:HIS:HE1	1:D:91:HIS:HD2	1.45	0.64
1:A:123:GLY:HA3	1:A:264:LYS:HE3	1.81	0.63
1:D:8:HIS:CE1	1:D:35:LYS:HD2	2.33	0.63
1:B:138:GLU:HG2	1:B:159:ASN:HB2	1.83	0.61
1:B:14[B]:CYS:SG	5:B:516:HOH:O	2.55	0.61
1:B:137:ASP:HA	1:B:140:LYS:HD3	1.84	0.60
1:B:14[B]:CYS:HB3	1:B:16[B]:HIS:HD2	1.65	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:61:PRO:HG2	1:D:64:TYR:HD2	1.67	0.59
1:D:46:MET:HG3	1:D:86:PHE:CD2	2.36	0.59
1:B:307:THR:HG22	1:B:309:GLU:H	1.68	0.59
1:E:197:ARG:NH1	5:E:503:HOH:O	2.35	0.58
1:B:16[A]:HIS:CE1	1:B:249:LYS:HG3	2.37	0.58
1:A:16[B]:HIS:CE1	1:A:249:LYS:HG3	2.39	0.58
1:C:49:LEU:HD23	1:C:93:SER:HB2	1.86	0.58
1:E:16[B]:HIS:CE1	1:E:18:LYS:HD2	2.38	0.58
1:A:14:CYS:HB3	1:A:16[A]:HIS:CD2	2.39	0.57
1:B:312:LYS:NZ	1:B:316:GLU:OE1	2.37	0.57
1:E:173:GLN:H	1:E:173:GLN:NE2	2.02	0.56
1:E:46:MET:HG3	1:E:86:PHE:CD2	2.39	0.56
1:B:193:LYS:NZ	5:B:501:HOH:O	2.33	0.56
1:C:138:GLU:HG2	1:C:159:ASN:HB2	1.87	0.55
1:E:61:PRO:HG2	1:E:64:TYR:HD2	1.71	0.55
1:B:14[B]:CYS:SG	1:B:229:GLY:O	2.63	0.55
1:C:46:MET:HG3	1:C:86:PHE:CD1	2.41	0.55
1:A:6:ILE:HD12	1:A:8:HIS:CE1	2.41	0.55
1:E:77:LYS:NZ	5:E:705:HOH:O	2.39	0.55
1:E:49:LEU:HD23	1:E:93:SER:HB2	1.89	0.54
1:D:14[A]:CYS:HB3	1:D:16[A]:HIS:CD2	2.42	0.54
1:B:281:MET:HA	1:B:281:MET:CE	2.37	0.54
1:C:7:GLN:HB3	1:C:263:LEU:HD13	1.90	0.54
1:D:305:LYS:NZ	5:D:677:HOH:O	2.40	0.54
1:B:24:ARG:NH1	5:B:702:HOH:O	2.38	0.54
1:B:301:GLU:OE2	1:B:305:LYS:HE2	2.08	0.53
1:E:16[A]:HIS:HE1	1:E:91:HIS:CD2	2.26	0.53
1:B:307:THR:HG22	1:B:309:GLU:N	2.24	0.52
1:C:265:GLN:HG3	1:C:267:THR:OG1	2.09	0.52
1:C:163:GLN:NE2	5:C:710:HOH:O	2.41	0.52
1:A:16[A]:HIS:HE1	1:A:91:HIS:HD2	1.58	0.52
1:B:196:TYR:CD2	1:B:203:LYS:HG2	2.45	0.52
1:C:230:HIS:HA	5:C:709:HOH:O	2.08	0.52
1:D:14[B]:CYS:HB3	1:D:249:LYS:HA	1.92	0.51
1:C:61:PRO:HG2	1:C:64:TYR:CD2	2.46	0.51
1:C:49:LEU:HD12	1:C:67:MET:HA	1.93	0.50
1:C:16[A]:HIS:HE1	1:C:91:HIS:CD2	2.30	0.50
1:D:16[A]:HIS:HE1	1:D:91:HIS:CD2	2.28	0.50
1:D:259:LEU:HB3	1:D:261:LEU:HD13	1.92	0.50
1:E:313:LYS:HD3	1:E:351:LEU:HD23	1.94	0.50
1:E:280:ILE:HD13	1:E:344:ILE:HG23	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:173:GLN:NE2	5:D:675:HOH:O	2.42	0.50
1:E:173:GLN:HG3	1:E:224:LYS:HE3	1.94	0.49
1:B:193:LYS:O	1:B:197:ARG:HG3	2.11	0.49
1:A:234:GLU:HB2	1:A:254:ASN:HB2	1.94	0.49
1:C:63:LYS:CD	1:C:63:LYS:H	2.25	0.49
1:D:319:GLU:HG2	1:D:324:LEU:HG	1.94	0.49
1:C:298:SER:HB3	1:C:301:GLU:HB3	1.95	0.49
1:A:5:GLN:HG3	1:A:6:ILE:N	2.27	0.48
1:B:46:MET:HG3	1:B:86:PHE:CD1	2.48	0.48
1:B:137:ASP:OD2	1:B:156:HIS:ND1	2.40	0.48
1:B:14[B]:CYS:HB2	1:B:249:LYS:HB2	1.95	0.48
1:B:69:ASP:O	1:B:72:LEU:HB2	2.13	0.47
1:D:269:LEU:HG	1:D:326:ILE:HD12	1.97	0.47
1:C:144:TYR:HB3	1:C:289:LYS:O	2.15	0.47
1:E:283:TRP:CZ2	1:E:345:GLU:HB2	2.50	0.47
1:B:269:LEU:HG	1:B:326:ILE:HD12	1.96	0.47
1:E:269:LEU:HG	1:E:326:ILE:HD12	1.96	0.47
1:B:6:ILE:HD13	1:B:260:ASP:HB3	1.96	0.47
1:A:16[B]:HIS:NE2	1:A:249:LYS:HE3	2.30	0.47
1:B:81:PRO:HG2	1:B:82:TYR:CD2	2.49	0.47
1:A:46:MET:HG3	1:A:86:PHE:CD1	2.50	0.47
1:E:5:GLN:N	1:E:5:GLN:HE21	2.13	0.47
1:B:101:TYR:HB2	1:B:115:VAL:HG23	1.98	0.46
1:C:189:LYS:HG3	1:C:239:GLU:OE2	2.15	0.46
1:C:61:PRO:HB2	1:C:63:LYS:HD2	1.97	0.46
1:C:16[A]:HIS:HE1	1:C:91:HIS:HD2	1.64	0.46
1:B:30:GLU:OE1	1:B:82:TYR:OH	2.32	0.46
1:B:31:LYS:HD3	1:D:198:PHE:CZ	2.52	0.45
1:B:25:GLN:OE1	5:B:643:HOH:O	2.21	0.45
1:D:315:ILE:O	1:D:319:GLU:HG3	2.17	0.45
1:E:298:SER:HB3	1:E:301:GLU:HB2	1.98	0.45
1:B:86:PHE:CZ	1:B:112:TYR:HB2	2.51	0.45
1:A:61:PRO:HG2	1:A:64:TYR:HD2	1.81	0.45
1:E:16[B]:HIS:NE2	1:E:18:LYS:HD2	2.32	0.45
1:E:280:ILE:HG13	1:E:314:ILE:HG23	1.98	0.45
1:A:144:TYR:HB3	1:A:289:LYS:O	2.17	0.45
1:B:311:ASP:O	1:B:315:ILE:HG12	2.17	0.45
4:B:404:GOL:O1	5:B:728:HOH:O	2.20	0.44
1:A:16[B]:HIS:CE1	1:A:249:LYS:HE3	2.52	0.44
1:E:16[A]:HIS:HE1	1:E:91:HIS:HD2	1.64	0.44
1:E:295:LEU:HD23	1:E:295:LEU:HA	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:63:LYS:H	1:C:63:LYS:HD2	1.82	0.44
1:E:49:LEU:HD23	1:E:93:SER:CB	2.48	0.44
1:B:281:MET:HA	1:B:281:MET:HE3	2.00	0.44
1:E:191:ASN:OD1	5:E:501:HOH:O	2.21	0.44
1:D:16[A]:HIS:CE1	1:D:91:HIS:HD2	2.30	0.44
1:D:85:LEU:HB3	1:D:113:LEU:HD11	1.99	0.44
1:C:219:ASN:ND2	5:C:679:HOH:O	2.44	0.44
1:B:179:SER:O	1:B:228:SER:HA	2.18	0.44
1:D:61:PRO:HG2	1:D:64:TYR:CD2	2.50	0.43
1:B:307:THR:HG22	1:B:310:LEU:H	1.83	0.43
1:E:61:PRO:HG2	1:E:64:TYR:CD2	2.51	0.43
1:C:280:ILE:HG13	1:C:314:ILE:HG23	2.01	0.43
1:E:16[A]:HIS:CE1	1:E:91:HIS:HD2	2.36	0.43
1:A:264:LYS:H	1:A:264:LYS:CE	2.26	0.43
1:E:173:GLN:NE2	5:E:676:HOH:O	2.50	0.43
1:D:49:LEU:HD22	1:D:93:SER:HB2	2.00	0.43
1:C:86:PHE:CZ	1:C:112:TYR:HB2	2.54	0.43
1:C:14:CYS:SG	1:C:16[A]:HIS:CD2	3.12	0.43
1:E:202:PHE:N	1:E:202:PHE:CD1	2.86	0.43
1:D:9:ILE:HD12	1:D:261:LEU:HD22	2.00	0.42
1:C:79:LYS:HE3	1:C:108:PHE:CE2	2.54	0.42
1:E:184:GLN:HA	1:E:209:LEU:O	2.19	0.42
1:D:66:GLN:NE2	5:D:503:HOH:O	2.52	0.42
1:C:134:LYS:HA	1:C:134:LYS:HD3	1.87	0.42
1:A:134:LYS:HG3	1:A:156:HIS:NE2	2.35	0.42
1:C:199:GLN:HA	1:C:200:PRO:HD2	1.78	0.42
1:E:202:PHE:HD1	1:E:202:PHE:N	2.17	0.42
1:C:264:LYS:H	1:C:264:LYS:HD2	1.84	0.42
1:D:235:TYR:HB3	1:D:246:ALA:HB3	2.01	0.42
1:A:230:HIS:HA	5:A:614:HOH:O	2.20	0.41
1:E:234:GLU:HB2	1:E:254:ASN:HB2	2.02	0.41
1:D:179:SER:O	1:D:228:SER:HA	2.20	0.41
1:C:144:TYR:CG	1:C:145:PRO:HA	2.55	0.41
1:D:281:MET:HG3	1:D:306:ARG:HG2	2.03	0.41
1:B:202:PHE:N	1:B:202:PHE:CD1	2.89	0.41
1:E:8:HIS:CE1	1:E:35:LYS:HD3	2.56	0.41
1:C:87:ILE:H	1:C:87:ILE:HG13	1.71	0.41
1:E:16[A]:HIS:CE1	1:E:91:HIS:CD2	3.09	0.41
1:C:269:LEU:HG	1:C:326:ILE:HD12	2.02	0.41
1:A:81:PRO:HG2	1:A:82:TYR:CD2	2.56	0.41
1:E:308:LYS:HB3	1:E:308:LYS:HE2	1.74	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:186:ILE:HD11	1:C:239:GLU:HB2	2.04	0.41
1:A:144:TYR:CG	1:A:145:PRO:HA	2.56	0.40
1:D:144:TYR:HB3	1:D:289:LYS:O	2.21	0.40
1:A:13:GLY:HA2	1:A:43:THR:OG1	2.21	0.40
1:E:144:TYR:HB3	1:E:289:LYS:O	2.20	0.40
1:A:175:ASP:OD1	1:A:224:LYS:HE3	2.20	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	350/360 (97%)	339 (97%)	10 (3%)	1 (0%)	46	35
1	B	350/360 (97%)	342 (98%)	8 (2%)	0	100	100
1	C	348/360 (97%)	336 (97%)	12 (3%)	0	100	100
1	D	349/360 (97%)	334 (96%)	14 (4%)	1 (0%)	46	35
1	E	349/360 (97%)	336 (96%)	12 (3%)	1 (0%)	46	35
All	All	1746/1800 (97%)	1687 (97%)	56 (3%)	3 (0%)	52	43

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	352	ASN
1	E	352	ASN
1	A	352	ASN



### 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	319/326 (98%)	306 (96%)	13 (4%)	37	22
1	B	319/326 (98%)	312 (98%)	7 (2%)	60	51
1	C	317/326 (97%)	310 (98%)	7 (2%)	60	51
1	D	318/326 (98%)	308 (97%)	10 (3%)	47	34
1	E	318/326 (98%)	302 (95%)	16 (5%)	30	14
All	All	1591/1630 (98%)	1538 (97%)	53 (3%)	44	32

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	GLN
1	A	6	ILE
1	A	51	TYR
1	A	63	LYS
1	A	147	SER
1	A	159	ASN
1	A	163	GLN
1	A	253	LYS
1	A	261	LEU
1	A	264	LYS
1	A	267	THR
1	A	301	GLU
1	A	317	LEU
1	B	6	ILE
1	B	51	TYR
1	B	163	GLN
1	B	249	LYS
1	B	266	LYS
1	B	267	THR
1	B	281	MET
1	C	51	TYR
1	C	63	LYS
1	C	87	ILE

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Mol	Chain	Res	Type
1	C	204	LYS
1	C	253	LYS
1	C	261	LEU
1	C	301	GLU
1	D	6	ILE
1	D	21	GLU
1	D	49	LEU
1	D	51	TYR
1	D	66	GLN
1	D	170	GLN
1	D	204	LYS
1	D	205	ASP
1	D	221	LEU
1	D	231	MET
1	E	5	GLN
1	E	28	GLU
1	E	51	TYR
1	E	72	LEU
1	E	115	VAL
1	E	138	GLU
1	E	140	LYS
1	E	147	SER
1	E	173	GLN
1	E	186	ILE
1	E	221	LEU
1	E	224	LYS
1	E	227	ILE
1	E	231	MET
1	E	253	LYS
1	E	316	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	GLN
1	B	102	ASN
1	D	232	HIS

### 5.3.3 RNA ⓘ

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 18 ligands modelled in this entry, 5 are monoatomic - leaving 13 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	402	-	4,4,4	0.10	0	6,6,6	0.17	0
3	SO4	A	403	2	4,4,4	0.12	0	6,6,6	0.12	0
4	GOL	A	404	-	5,5,5	0.30	0	5,5,5	0.29	0
3	SO4	B	402	-	4,4,4	0.15	0	6,6,6	0.13	0
3	SO4	B	403	2	4,4,4	0.06	0	6,6,6	0.25	0
4	GOL	B	404	-	5,5,5	0.33	0	5,5,5	0.44	0
3	SO4	C	402	-	4,4,4	0.19	0	6,6,6	0.19	0
3	SO4	C	403	2	4,4,4	0.24	0	6,6,6	0.25	0
3	SO4	D	402	-	4,4,4	0.24	0	6,6,6	0.36	0
3	SO4	D	403	2	4,4,4	0.08	0	6,6,6	0.21	0
4	GOL	D	404	-	5,5,5	0.34	0	5,5,5	0.33	0
3	SO4	E	402	-	4,4,4	0.20	0	6,6,6	0.26	0
3	SO4	E	403	2	4,4,4	0.19	0	6,6,6	0.21	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	402	-	-	0/0/0/0	0/0/0/0
3	SO4	A	403	2	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	A	404	-	-	0/4/4/4	0/0/0/0
3	SO4	B	402	-	-	0/0/0/0	0/0/0/0
3	SO4	B	403	2	-	0/0/0/0	0/0/0/0
4	GOL	B	404	-	-	0/4/4/4	0/0/0/0
3	SO4	C	402	-	-	0/0/0/0	0/0/0/0
3	SO4	C	403	2	-	0/0/0/0	0/0/0/0
3	SO4	D	402	-	-	0/0/0/0	0/0/0/0
3	SO4	D	403	2	-	0/0/0/0	0/0/0/0
4	GOL	D	404	-	-	0/4/4/4	0/0/0/0
3	SO4	E	402	-	-	0/0/0/0	0/0/0/0
3	SO4	E	403	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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	404	GOL	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, 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	349/360 (96%)	0.39	19 (5%) 29 41	14, 25, 44, 59	0
1	B	349/360 (96%)	0.30	11 (3%) 51 61	13, 24, 41, 57	0
1	C	349/360 (96%)	0.60	20 (5%) 27 37	16, 27, 44, 56	0
1	D	349/360 (96%)	0.66	29 (8%) 14 22	14, 27, 55, 78	0
1	E	349/360 (96%)	0.62	29 (8%) 14 22	15, 29, 53, 69	0
All	All	1745/1800 (96%)	0.52	108 (6%) 24 34	13, 26, 49, 78	0

All (108) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	198	PHE	7.6
1	E	204	LYS	6.5
1	E	353	ILE	5.9
1	E	198	PHE	5.3
1	D	202	PHE	5.3
1	D	204	LYS	5.0
1	D	203	LYS	4.8
1	E	265	GLN	4.7
1	D	198	PHE	4.7
1	D	136	PHE	4.3
1	D	31	LYS	4.3
1	A	265	GLN	4.3
1	D	252	TYR	4.2
1	D	201	GLY	4.1
1	E	202	PHE	4.1
1	D	196	TYR	4.0
1	D	352	ASN	4.0
1	E	352	ASN	3.9
1	D	205	ASP	3.9
1	B	136	PHE	3.8

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Mol	Chain	Res	Type	RSRZ
1	C	264	LYS	3.7
1	A	198	PHE	3.7
1	D	265	GLN	3.7
1	E	252	TYR	3.6
1	C	6	ILE	3.6
1	A	204	LYS	3.6
1	E	203	LYS	3.6
1	A	252	TYR	3.6
1	A	6	ILE	3.6
1	D	6	ILE	3.5
1	E	6	ILE	3.5
1	D	264	LYS	3.4
1	A	353	ILE	3.4
1	E	310	LEU	3.4
1	D	200	PRO	3.4
1	E	205	ASP	3.4
1	E	29	TYR	3.4
1	B	205	ASP	3.3
1	D	309	GLU	3.2
1	D	195	LEU	3.2
1	E	5	GLN	3.2
1	E	309	GLU	3.2
1	A	192	TYR	3.1
1	D	199	GLN	3.1
1	E	33	THR	3.1
1	D	197	ARG	3.1
1	E	207	ALA	3.0
1	C	352	ASN	3.0
1	D	192	TYR	2.9
1	E	264	LYS	2.9
1	D	29	TYR	2.9
1	B	253	LYS	2.9
1	C	204	LYS	2.9
1	E	31	LYS	2.9
1	C	265	GLN	2.9
1	E	32	SER	2.8
1	C	202	PHE	2.8
1	C	199	GLN	2.8
1	D	5	GLN	2.8
1	C	252	TYR	2.8
1	D	35	LYS	2.8
1	C	197	ARG	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	353	ILE	2.8
1	B	252	TYR	2.8
1	A	193	LYS	2.8
1	E	196	TYR	2.7
1	B	203	LYS	2.7
1	C	253	LYS	2.7
1	B	265	GLN	2.7
1	D	24	ARG	2.7
1	A	201	GLY	2.6
1	A	352	ASN	2.6
1	E	253	LYS	2.6
1	E	206	GLY	2.5
1	A	31	LYS	2.5
1	D	32	SER	2.5
1	D	58	LEU	2.5
1	E	201	GLY	2.5
1	E	251	GLY	2.5
1	E	263	LEU	2.5
1	B	31	LYS	2.4
1	E	313	LYS	2.4
1	E	200	PRO	2.4
1	B	204	LYS	2.4
1	C	31	LYS	2.4
1	E	173	GLN	2.3
1	E	197	ARG	2.3
1	C	64	TYR	2.3
1	C	33	THR	2.3
1	C	5	GLN	2.3
1	A	202	PHE	2.2
1	D	312	LYS	2.2
1	A	309	GLU	2.2
1	D	251	GLY	2.2
1	A	199	GLN	2.2
1	A	196	TYR	2.2
1	A	234	GLU	2.1
1	A	264	LYS	2.1
1	C	196	TYR	2.1
1	D	253	LYS	2.1
1	B	28	GLU	2.1
1	A	200	PRO	2.1
1	C	115	VAL	2.0
1	A	5	GLN	2.0

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Mol	Chain	Res	Type	RSRZ
1	C	200	PRO	2.0
1	C	105	PHE	2.0
1	C	112	TYR	2.0
1	B	299	ILE	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, 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
3	SO4	E	402	5/5	0.95	0.17	6.48	41,42,47,50	0
4	GOL	A	404	6/6	0.81	0.29	5.82	43,44,50,52	0
4	GOL	D	404	6/6	0.85	0.21	3.60	34,43,46,49	0
4	GOL	B	404	6/6	0.91	0.14	1.51	39,42,45,46	0
3	SO4	D	402	5/5	0.97	0.12	0.61	30,38,42,47	0
3	SO4	B	402	5/5	0.98	0.09	-1.34	36,40,44,49	0
2	MN	A	401	1/1	0.99	0.08	-1.39	23,23,23,23	0
3	SO4	A	402	5/5	0.96	0.10	-1.56	39,41,44,45	0
3	SO4	B	403	5/5	0.97	0.09	-1.66	26,31,37,37	0
3	SO4	A	403	5/5	0.98	0.07	-1.68	29,38,40,40	0
3	SO4	E	403	5/5	0.97	0.08	-1.96	25,30,34,38	0
3	SO4	C	402	5/5	0.98	0.09	-2.12	38,39,42,44	0
2	MN	E	401	1/1	1.00	0.08	-2.34	21,21,21,21	0
2	MN	D	401	1/1	0.99	0.09	-2.71	27,27,27,27	0
3	SO4	C	403	5/5	0.97	0.09	-2.99	32,34,42,44	0
3	SO4	D	403	5/5	0.96	0.08	-3.02	31,36,40,50	0
2	MN	B	401	1/1	1.00	0.05	-3.83	20,20,20,20	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	MN	C	401	1/1	0.99	0.06	-7.24	22,22,22,22	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.