



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:27 AM GMT

PDB ID : 2X51  
Title : M6 DELTA INSERT1  
Authors : Squires, G.; Houdusse, A.  
Deposited on : 2010-02-04  
Resolution : 2.20 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

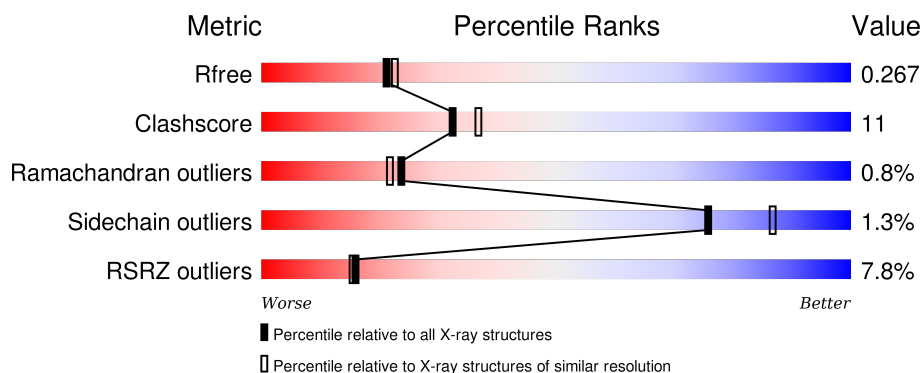
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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	789	<div> <div>8%</div> <div>72%</div> <div>20%</div> <div>7%</div> </div>
2	B	149	<div> <div>3%</div> <div>79%</div> <div>18%</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	GOL	A	1817	-	-	-	X
5	CA	B	1150	-	-	-	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7401 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 MYOSIN-VI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	737	Total	C	N	O	S	0	0	0
			5950	3790	1028	1103	29			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	.	-	LYS	DELETION	UNP Q29122
A	547	VAL	GLY	ENGINEERED MUTATION	UNP Q29122
A	572	ARG	ALA	ENGINEERED MUTATION	UNP Q29122
A	573	ASP	TYR	ENGINEERED MUTATION	UNP Q29122
A	714	LEU	VAL	ENGINEERED MUTATION	UNP Q29122
A	721	TYR	SER	ENGINEERED MUTATION	UNP Q29122
A	722	MET	LEU	ENGINEERED MUTATION	UNP Q29122

- Molecule 2 is a protein called CALMODULIN.

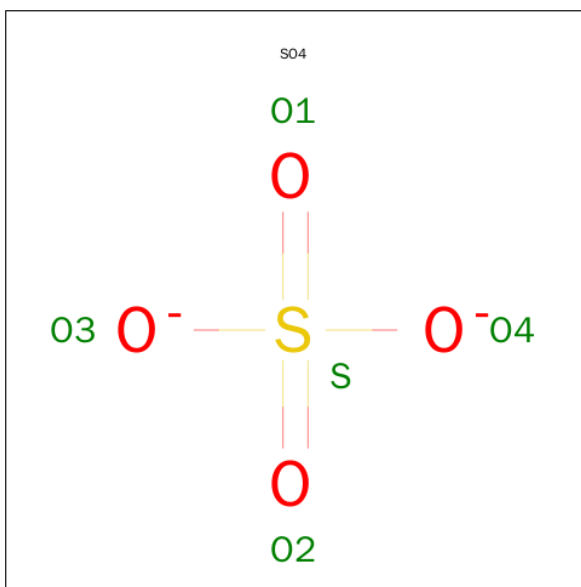
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	146	Total	C	N	O	S	0	0	0
			1150	706	185	250	9			

- Molecule 3 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
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		

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



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

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	3	Total	Ca	0	0
			3	3		

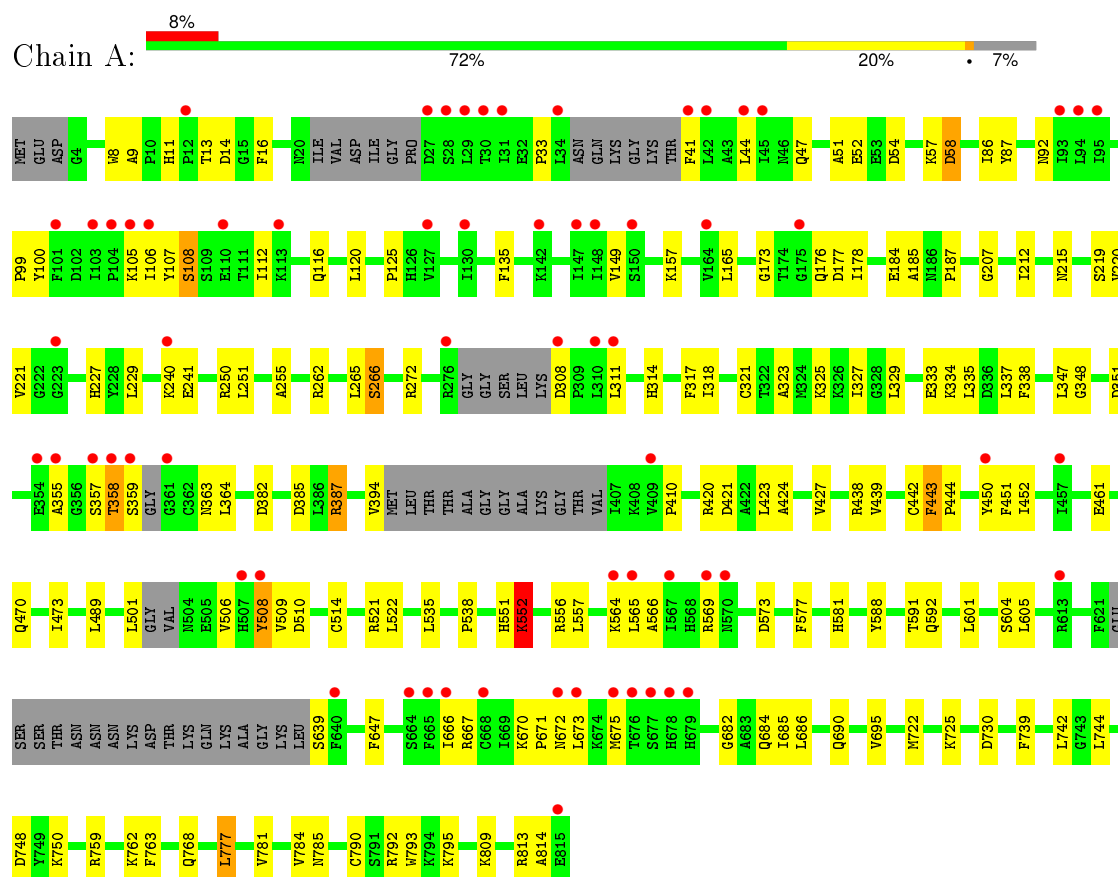
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	198	Total	O	0	0
			198	198		
6	B	53	Total	O	0	0
			53	53		

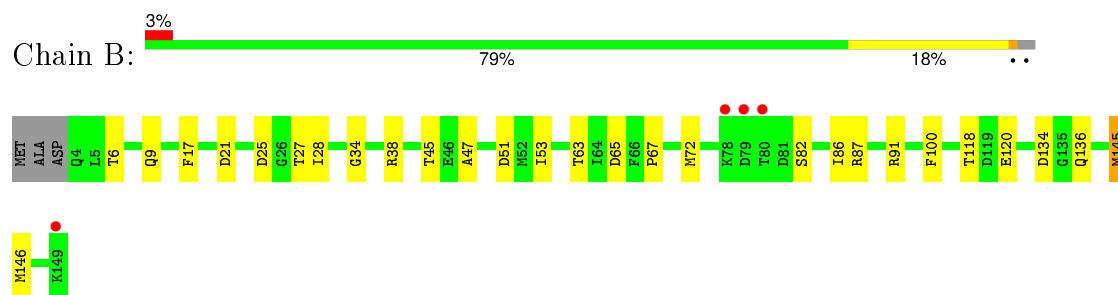
### 3 Residue-property plots [i](#)

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: MYOSIN-VI



#### • Molecule 2: CALMODULIN



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	163.89Å 61.02Å 133.27Å 90.00° 116.19° 90.00°	Depositor
Resolution (Å)	26.82 – 2.20 29.56 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.9 (26.82-2.20) 99.9 (29.56-2.20)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.83 (at 2.20Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.216 , 0.272 0.211 , 0.267	Depositor DCC
$R_{free}$ test set	3014 reflections (5.00%)	DCC
Wilson B-factor (Å <sup>2</sup> )	36.6	Xtriage
Anisotropy	0.786	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 43.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 60348 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	7401	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 3.95% 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: GOL, CA, 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	0.41	0/6071	0.53	0/8176
2	B	0.43	0/1162	0.53	0/1559
All	All	0.41	0/7233	0.53	0/9735

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	5950	0	5866	125	0
2	B	1150	0	1083	28	0
3	A	42	0	56	5	0
4	A	5	0	0	0	0
5	B	3	0	0	0	0
6	A	198	0	0	7	0
6	B	53	0	0	2	0
All	All	7401	0	7005	149	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 11.

All (149) 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:790:CYS:HA	2:B:145:MET:HE3	1.53	0.91
1:A:722:MET:SD	1:A:777:LEU:HD21	2.12	0.90
1:A:11:HIS:HD2	1:A:14:ASP:H	1.20	0.84
2:B:45:THR:HG21	6:B:2021:HOH:O	1.81	0.79
1:A:327:ILE:HG23	1:A:442:CYS:SG	2.26	0.75
2:B:25:ASP:OD1	2:B:27:THR:HG22	1.87	0.75
1:A:272:ARG:NH2	1:A:351:ASP:OD1	2.20	0.74
1:A:508:TYR:CZ	1:A:510:ASP:HB3	2.24	0.72
1:A:364:LEU:HD13	1:A:387:ARG:HG3	1.70	0.72
1:A:510:ASP:HB2	6:A:2098:HOH:O	1.89	0.71
1:A:323:ALA:O	1:A:327:ILE:HG12	1.91	0.70
1:A:461:GLU:OE2	3:A:1820:GOL:H2	1.91	0.70
1:A:739:PHE:CD2	1:A:744:LEU:HD12	2.27	0.69
1:A:11:HIS:HD2	1:A:14:ASP:N	1.90	0.69
1:A:581:HIS:HE1	1:A:588:TYR:OH	1.75	0.68
1:A:671:PRO:HA	1:A:685:ILE:HD11	1.76	0.68
1:A:809:LYS:HG3	2:B:72:MET:O	1.94	0.67
1:A:105:LYS:O	1:A:108:SER:HB3	1.96	0.65
2:B:45:THR:HG22	2:B:47:ALA:H	1.61	0.65
1:A:722:MET:SD	1:A:777:LEU:CD2	2.85	0.64
1:A:44:LEU:HB2	1:A:47:GLN:OE1	1.99	0.63
1:A:489:LEU:CD1	1:A:506:VAL:HB	2.28	0.63
1:A:690:GLN:HG2	1:A:695:VAL:HG21	1.79	0.63
1:A:329:LEU:HD12	1:A:438:ARG:HH21	1.62	0.63
1:A:149:VAL:HG11	1:A:157:LYS:HG3	1.80	0.62
1:A:165:LEU:HD22	1:A:212:ILE:HD11	1.81	0.62
1:A:686:LEU:O	1:A:690:GLN:HG3	1.98	0.62
1:A:443:PHE:N	1:A:444:PRO:HD3	2.15	0.61
1:A:185:ALA:HB1	1:A:439:VAL:HG13	1.81	0.61
1:A:551:HIS:HE1	6:A:2101:HOH:O	1.83	0.61
1:A:11:HIS:CD2	1:A:14:ASP:H	2.12	0.61
1:A:394:VAL:HG22	1:A:604:SER:HB3	1.84	0.60
1:A:52:GLU:HB3	1:A:57:LYS:HE2	1.82	0.60
1:A:581:HIS:CE1	1:A:588:TYR:OH	2.54	0.59
2:B:45:THR:HG22	2:B:47:ALA:N	2.17	0.59
1:A:262:ARG:HH21	1:A:335:LEU:HD21	1.67	0.59
1:A:240:LYS:O	1:A:241:GLU:HB2	2.03	0.59
1:A:666:ILE:O	1:A:667:ARG:HD2	2.03	0.59
1:A:364:LEU:CD1	1:A:387:ARG:HG3	2.33	0.58
1:A:86:ILE:HG21	1:A:99:PRO:HG3	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:215:ASN:HB3	1:A:221:VAL:HG11	1.86	0.58
1:A:501:LEU:HD22	1:A:750:LYS:HB3	1.86	0.58
1:A:591:THR:O	1:A:592:GLN:HB2	2.03	0.58
2:B:21:ASP:CA	2:B:28:ILE:HD11	2.34	0.57
1:A:54:ASP:HB3	1:A:57:LYS:HD3	1.86	0.57
1:A:666:ILE:C	1:A:667:ARG:HD2	2.24	0.57
1:A:672:ASN:HA	1:A:684:GLN:NE2	2.20	0.56
1:A:793:TRP:CD1	2:B:145:MET:HE2	2.39	0.56
1:A:184:GLU:O	1:A:187:PRO:HD2	2.06	0.56
1:A:321:CYS:O	1:A:325:LYS:HG3	2.05	0.55
1:A:514:CYS:HB2	1:A:556:ARG:HD2	1.87	0.55
1:A:251:LEU:HD12	1:A:255:ALA:HB2	1.87	0.55
1:A:759:ARG:HD2	6:A:2164:HOH:O	2.06	0.55
1:A:489:LEU:HD11	1:A:506:VAL:HB	1.88	0.55
3:A:1817:GOL:H11	6:A:2039:HOH:O	2.07	0.55
1:A:508:TYR:OH	1:A:510:ASP:HB3	2.08	0.54
1:A:792:ARG:HH11	1:A:792:ARG:HG2	1.71	0.54
1:A:566:ALA:HA	1:A:569:ARG:HD2	1.88	0.54
2:B:34:GLY:O	2:B:38:ARG:HG3	2.08	0.54
1:A:308:ASP:HB3	1:A:311:LEU:HD22	1.90	0.53
1:A:308:ASP:O	1:A:311:LEU:HB2	2.08	0.53
1:A:790:CYS:HA	2:B:145:MET:CE	2.32	0.53
1:A:58:ASP:C	1:A:58:ASP:OD1	2.46	0.53
1:A:813:ARG:HH12	2:B:51:ASP:HB3	1.75	0.52
2:B:118:THR:OG1	2:B:120:GLU:HG2	2.10	0.52
1:A:509:VAL:O	1:A:509:VAL:HG13	2.09	0.52
1:A:730:ASP:OD1	1:A:795:LYS:NZ	2.43	0.52
1:A:591:THR:HG23	6:A:2128:HOH:O	2.09	0.52
1:A:461:GLU:OE2	1:A:470:GLN:HG3	2.10	0.51
1:A:364:LEU:HD13	1:A:387:ARG:CG	2.39	0.51
1:A:573:ASP:HB2	3:A:1822:GOL:H2	1.92	0.51
1:A:639:SER:OG	1:A:639:SER:O	2.25	0.51
2:B:6:THR:H	2:B:9:GLN:NE2	2.09	0.51
1:A:112:ILE:O	1:A:116:GLN:HG3	2.11	0.51
1:A:725:LYS:HE2	1:A:785:ASN:OD1	2.10	0.50
2:B:53:ILE:HA	2:B:72:MET:CE	2.42	0.50
1:A:265:LEU:O	1:A:266:SER:HB3	2.11	0.50
2:B:82:SER:O	2:B:86:ILE:HG12	2.12	0.49
1:A:41:PHE:CD1	1:A:41:PHE:N	2.80	0.49
1:A:357:SER:O	1:A:358:THR:C	2.51	0.49
2:B:87:ARG:O	2:B:91:ARG:HG3	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:173:GLY:HA2	1:A:220:VAL:HB	1.94	0.49
1:A:327:ILE:CG2	1:A:442:CYS:SG	2.99	0.48
2:B:53:ILE:HA	2:B:72:MET:HE1	1.96	0.48
1:A:14:ASP:HA	1:A:682:GLY:HA3	1.96	0.47
1:A:423:LEU:O	1:A:427:VAL:HG23	2.13	0.47
1:A:178:ILE:HD12	1:A:220:VAL:CG1	2.44	0.47
1:A:329:LEU:CD1	1:A:438:ARG:HH21	2.25	0.47
2:B:27:THR:OG1	2:B:63:THR:HB	2.15	0.47
1:A:557:LEU:HD11	1:A:577:PHE:HB2	1.96	0.47
1:A:768:GLN:HA	1:A:768:GLN:OE1	2.15	0.47
2:B:86:ILE:HD12	2:B:146:MET:CE	2.45	0.47
1:A:8:TRP:CD1	1:A:51:ALA:HB2	2.49	0.47
1:A:359:SER:O	1:A:410:PRO:HG2	2.14	0.46
1:A:54:ASP:O	1:A:57:LYS:HG2	2.15	0.46
1:A:41:PHE:N	1:A:41:PHE:HD1	2.14	0.46
1:A:420:ARG:HD2	1:A:421:ASP:OD1	2.15	0.46
2:B:65:ASP:OD1	2:B:67:PRO:HD2	2.15	0.46
1:A:748:ASP:HB3	1:A:762:LYS:HG3	1.98	0.45
1:A:106:ILE:HG13	1:A:107:TYR:CD2	2.50	0.45
2:B:38:ARG:HD3	6:B:2014:HOH:O	2.15	0.45
1:A:473:ILE:HG12	3:A:1818:GOL:O3	2.16	0.45
1:A:9:ALA:O	1:A:16:PHE:HA	2.17	0.45
1:A:722:MET:HE1	1:A:781:VAL:CG2	2.46	0.45
1:A:100:TYR:CE2	1:A:675:MET:HA	2.52	0.45
1:A:157:LYS:HE2	1:A:157:LYS:HB3	1.71	0.45
1:A:358:THR:O	1:A:359:SER:O	2.34	0.44
1:A:793:TRP:CD1	2:B:145:MET:CE	3.00	0.44
1:A:348:GLY:HA2	1:A:420:ARG:HD3	1.99	0.44
1:A:333:GLU:O	1:A:337:LEU:HG	2.18	0.44
1:A:508:TYR:OH	1:A:510:ASP:CB	2.66	0.44
1:A:809:LYS:O	1:A:813:ARG:HG3	2.18	0.44
1:A:601:LEU:HD11	1:A:605:LEU:HG	2.00	0.44
1:A:135:PHE:HA	1:A:452:ILE:HD11	2.00	0.44
1:A:334:LYS:HG2	1:A:338:PHE:CE2	2.52	0.44
1:A:521:ARG:O	1:A:522:LEU:HB2	2.18	0.44
1:A:355:ALA:CB	1:A:363:ASN:HD21	2.30	0.43
1:A:87:TYR:O	1:A:125:PRO:HB2	2.18	0.43
1:A:92:ASN:OD1	1:A:120:LEU:HD23	2.18	0.43
1:A:552:LYS:HD3	1:A:552:LYS:HA	1.68	0.43
1:A:742:LEU:HB2	1:A:744:LEU:HG	2.01	0.43
1:A:149:VAL:CG1	1:A:157:LYS:HG3	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:394:VAL:CG2	1:A:604:SER:HB3	2.49	0.43
2:B:17:PHE:CZ	2:B:28:ILE:HG12	2.54	0.43
2:B:21:ASP:N	2:B:28:ILE:HD11	2.34	0.43
1:A:11:HIS:CD2	1:A:13:THR:H	2.37	0.43
1:A:240:LYS:O	1:A:241:GLU:CB	2.63	0.42
2:B:21:ASP:HA	2:B:28:ILE:HD11	2.01	0.42
1:A:250:ARG:HB3	1:A:317:PHE:HB2	2.01	0.42
1:A:314:HIS:CE1	1:A:318:ILE:HD11	2.53	0.42
1:A:489:LEU:HD12	1:A:506:VAL:HB	1.97	0.42
1:A:54:ASP:OD2	1:A:57:LYS:HB3	2.19	0.42
3:A:1817:GOL:H12	6:A:2079:HOH:O	2.19	0.42
1:A:347:LEU:HD23	1:A:424:ALA:HB2	2.01	0.42
1:A:777:LEU:HD22	6:A:2183:HOH:O	2.19	0.42
1:A:207:GLY:O	1:A:227:HIS:HA	2.19	0.42
1:A:535:LEU:O	1:A:538:PRO:HD3	2.20	0.42
1:A:792:ARG:NH1	1:A:792:ARG:HG2	2.33	0.42
1:A:219:SER:O	1:A:221:VAL:HG13	2.20	0.42
2:B:134:ASP:C	2:B:134:ASP:OD1	2.58	0.42
2:B:28:ILE:HD13	2:B:28:ILE:HA	1.83	0.41
1:A:229:LEU:HD22	1:A:647:PHE:CE2	2.55	0.41
1:A:784:VAL:O	1:A:784:VAL:CG1	2.69	0.41
1:A:382:ASP:HB3	1:A:385:ASP:OD2	2.21	0.41
1:A:670:LYS:HE2	1:A:673:LEU:HA	2.02	0.41
1:A:564:LYS:O	1:A:565:LEU:HD23	2.21	0.41
1:A:451:PHE:C	1:A:451:PHE:CD1	2.94	0.41
1:A:100:TYR:HD1	1:A:100:TYR:HA	1.76	0.40
2:B:100:PHE:HB3	2:B:136:GLN:HB3	2.03	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	721/789 (91%)	682 (95%)	32 (4%)	7 (1%)	19	16
2	B	144/149 (97%)	144 (100%)	0	0	100	100
All	All	865/938 (92%)	826 (96%)	32 (4%)	7 (1%)	24	22

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	266	SER
1	A	176	GLN
1	A	552	LYS
1	A	814	ALA
1	A	358	THR
1	A	443	PHE
1	A	33	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	654/698 (94%)	645 (99%)	9 (1%)	74	85
2	B	126/128 (98%)	125 (99%)	1 (1%)	86	93
All	All	780/826 (94%)	770 (99%)	10 (1%)	76	87

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	ASP
1	A	108	SER
1	A	177	ASP
1	A	387	ARG
1	A	450	TYR
1	A	508	TYR
1	A	552	LYS
1	A	763	PHE
1	A	777	LEU

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Mol	Chain	Res	Type
2	B	145	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	HIS
1	A	17	GLN
1	A	116	GLN
1	A	363	ASN
1	A	481	GLN
1	A	485	ASN
1	A	581	HIS
1	A	684	GLN
2	B	9	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 11 ligands modelled in this entry, 3 are monoatomic - leaving 8 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	GOL	A	1816	-	5,5,5	0.31	0	5,5,5	0.35	0
3	GOL	A	1817	-	5,5,5	0.77	0	5,5,5	0.66	0
3	GOL	A	1818	-	5,5,5	0.40	0	5,5,5	0.34	0
3	GOL	A	1819	-	5,5,5	0.30	0	5,5,5	0.36	0
3	GOL	A	1820	-	5,5,5	0.23	0	5,5,5	0.66	0
3	GOL	A	1821	-	5,5,5	0.35	0	5,5,5	0.25	0
3	GOL	A	1822	-	5,5,5	0.20	0	5,5,5	0.43	0
4	SO4	A	1823	-	4,4,4	0.13	0	6,6,6	0.16	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	GOL	A	1816	-	-	0/4/4/4	0/0/0/0
3	GOL	A	1817	-	-	0/4/4/4	0/0/0/0
3	GOL	A	1818	-	-	0/4/4/4	0/0/0/0
3	GOL	A	1819	-	-	0/4/4/4	0/0/0/0
3	GOL	A	1820	-	-	0/4/4/4	0/0/0/0
3	GOL	A	1821	-	-	0/4/4/4	0/0/0/0
3	GOL	A	1822	-	-	0/4/4/4	0/0/0/0
4	SO4	A	1823	-	-	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.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1817	GOL	2	0
3	A	1818	GOL	1	0
3	A	1820	GOL	1	0
3	A	1822	GOL	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, 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	737/789 (93%)	0.44	65 (8%) 12 11	28, 46, 77, 106	0
2	B	146/149 (97%)	0.09	4 (2%) 58 57	30, 45, 65, 88	0
All	All	883/938 (94%)	0.38	69 (7%) 16 15	28, 46, 76, 106	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	44	LEU	7.1
1	A	175	GLY	6.3
1	A	34	LEU	5.7
1	A	42	LEU	5.6
1	A	103	ILE	5.4
1	A	31	ILE	5.2
1	A	101	PHE	5.2
1	A	45	ILE	5.0
1	A	676	THR	4.6
1	A	673	LEU	4.5
1	A	308	ASP	4.4
1	A	666	ILE	4.2
1	A	276	ARG	4.1
1	A	110	GLU	4.0
1	A	29	LEU	3.9
2	B	80	THR	3.9
1	A	113	LYS	3.7
1	A	41	PHE	3.7
1	A	28	SER	3.7
1	A	357	SER	3.6
1	A	127	VAL	3.6
2	B	79	ASP	3.5
1	A	359	SER	3.5
1	A	567	ILE	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	569	ARG	3.2
1	A	105	LYS	3.2
1	A	565	LEU	3.2
1	A	409	VAL	3.1
1	A	815	GLU	3.1
1	A	12	PRO	3.0
2	B	78	LYS	3.0
1	A	94	LEU	3.0
1	A	675	MET	2.9
1	A	508	TYR	2.9
1	A	640	PHE	2.9
1	A	679	HIS	2.8
1	A	450	TYR	2.8
1	A	30	THR	2.7
1	A	106	ILE	2.7
1	A	147	ILE	2.7
1	A	240	LYS	2.7
1	A	148	ILE	2.6
1	A	223	GLY	2.6
1	A	677	SER	2.5
1	A	672	ASN	2.5
1	A	355	ALA	2.5
1	A	311	LEU	2.5
1	A	104	PRO	2.5
1	A	93	ILE	2.4
1	A	564	LYS	2.4
1	A	27	ASP	2.4
1	A	613	ARG	2.4
1	A	150	SER	2.4
1	A	570	ASN	2.3
1	A	678	HIS	2.3
1	A	130	ILE	2.3
1	A	507	HIS	2.3
1	A	457	ILE	2.2
1	A	95	ILE	2.2
1	A	668	CYS	2.2
1	A	354	GLU	2.2
1	A	361	GLY	2.1
1	A	142	LYS	2.1
1	A	164	VAL	2.1
1	A	664	SER	2.0
1	A	310	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	665	PHE	2.0
2	B	149	LYS	2.0
1	A	358	THR	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	GOL	A	1817	6/6	0.85	0.28	6.42	32,34,35,39	0
5	CA	B	1150	1/1	0.92	0.26	3.80	74,74,74,74	0
3	GOL	A	1819	6/6	0.89	0.16	0.43	43,50,55,66	0
3	GOL	A	1818	6/6	0.91	0.18	0.33	32,34,35,36	0
3	GOL	A	1821	6/6	0.90	0.16	0.13	47,51,54,55	0
3	GOL	A	1816	6/6	0.95	0.15	-0.93	37,42,43,43	0
5	CA	B	1151	1/1	0.98	0.09	-1.02	40,40,40,40	0
3	GOL	A	1820	6/6	0.94	0.10	-1.51	38,43,49,53	0
4	SO4	A	1823	5/5	0.95	0.08	-2.22	61,62,65,67	0
5	CA	B	1152	1/1	0.96	0.05	-2.83	46,46,46,46	0
3	GOL	A	1822	6/6	0.62	0.29	-	57,60,65,69	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.