



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:43 PM GMT

PDB ID : 1Q9W  
Title : S45-18 Fab pentasaccharide bisphosphate complex  
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Deposited on : 2003-08-26  
Resolution : 1.75 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

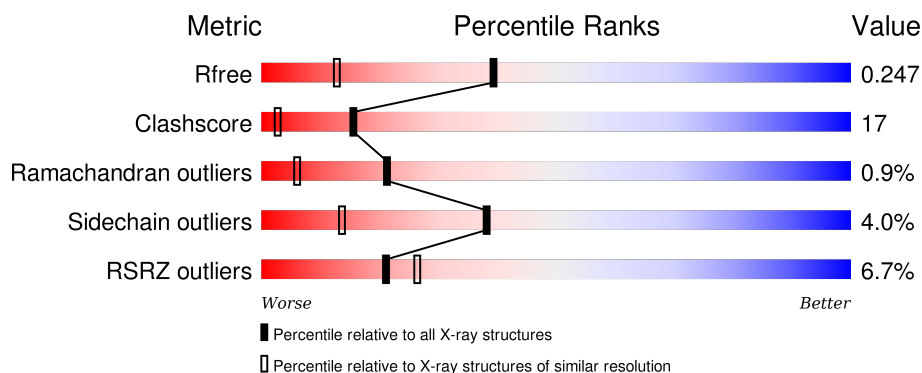
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.75 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	1609 (1.76-1.76)
Clashscore	102246	1730 (1.76-1.76)
Ramachandran outliers	100387	1711 (1.76-1.76)
Sidechain outliers	100360	1711 (1.76-1.76)
RSRZ outliers	91569	1610 (1.76-1.76)

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  $\geq 3$ , 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	219	<div> <div>6%</div> <div>74%</div> <div>25%</div> <div>.</div> </div>
1	C	219	<div> <div>5%</div> <div>72%</div> <div>25%</div> <div>.</div> </div>
2	B	226	<div> <div>8%</div> <div>75%</div> <div>22%</div> <div>.</div> </div>
2	D	226	<div> <div>8%</div> <div>72%</div> <div>27%</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	KDO	B	303	-	-	-	X
3	GP4	B	304	X	-	-	-
4	KDO	C	313	-	-	-	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7744 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 S45-18 Fab (IgG1k) light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	219	Total	C	N	O	S	0	0	0
			1707	1066	289	344	8			
1	C	219	Total	C	N	O	S	0	0	0
			1707	1066	289	344	8			

- Molecule 2 is a protein called S45-18 Fab (IgG1k) heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	226	Total	C	N	O	S	0	0	0
			1718	1085	284	340	9			
2	D	226	Total	C	N	O	S	0	0	0
			1718	1085	284	340	9			

- Molecule 3 is a polymer of unknown type called SUGAR (5-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	5	Total	C	N	O	P	0	0
			76	36	2	36	2		

- Molecule 4 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	3	Total	C	O	0	0
			46	24	22		

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	2	Total 2	Mg 2	0	0

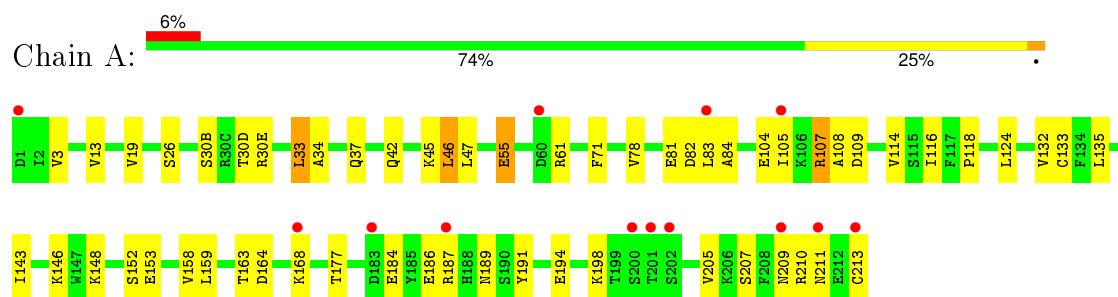
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	160	Total 160	O 160	0	0
6	B	220	Total 220	O 220	0	0
6	C	176	Total 176	O 176	0	0
6	D	211	Total 211	O 211	0	0

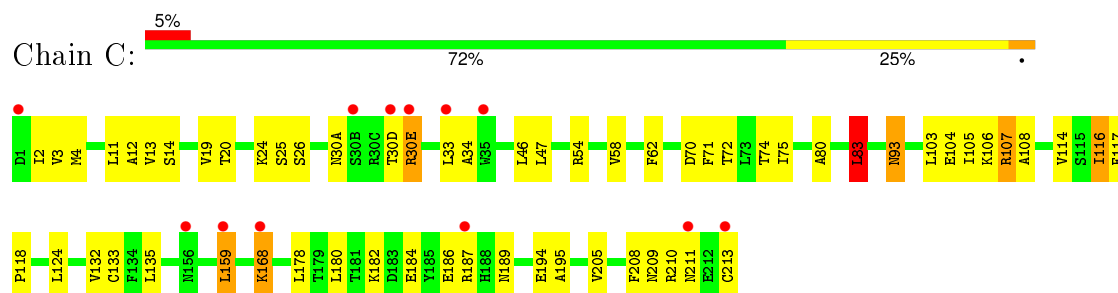
### 3 Residue-property plots [i](#)

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.

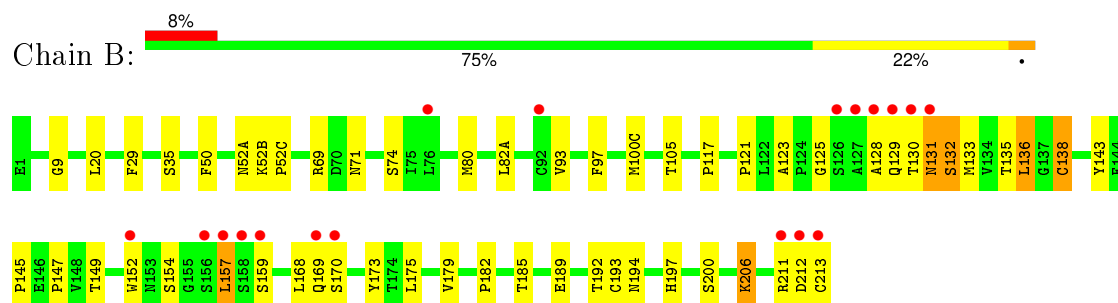
- Molecule 1: S45-18 Fab (IgG1k) light chain



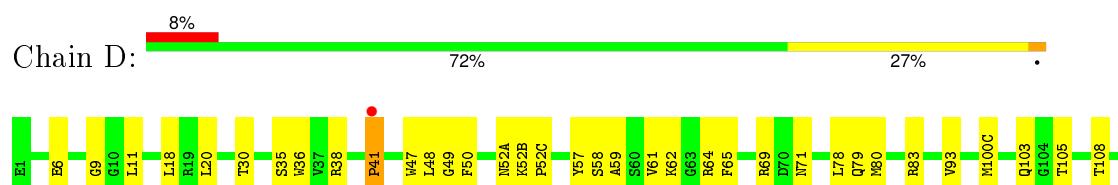
- Molecule 1: S45-18 Fab (IgG1k) light chain



- Molecule 2: S45-18 Fab (IgG1k) heavy chain



- Molecule 2: S45-18 Fab (IgG1k) heavy chain





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	71.20Å 113.90Å 133.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.90 – 1.75 19.90 – 1.75	Depositor EDS
% Data completeness (in resolution range)	90.5 (19.90-1.75) 90.4 (19.90-1.75)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.33 (at 1.74Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.212 , 0.247 0.212 , 0.247	Depositor DCC
$R_{free}$ test set	9994 reflections (10.02%)	DCC
Wilson B-factor (Å <sup>2</sup> )	23.4	Xtriage
Anisotropy	0.450	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 50.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 106072 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7744	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.83% 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: GP4, MG, KDO, GP1

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	0.42	0/1744	0.71	1/2361 (0.0%)
1	C	0.42	0/1744	0.72	1/2361 (0.0%)
2	B	0.46	0/1763	0.79	1/2406 (0.0%)
2	D	0.47	0/1763	0.78	0/2406
All	All	0.44	0/7014	0.75	3/9534 (0.0%)

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
3	B	1	0

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	138	CYS	CA-CB-SG	5.98	124.77	114.00
1	A	33	LEU	CA-CB-CG	-5.97	101.57	115.30
1	C	83	LEU	CA-CB-CG	5.20	127.27	115.30

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	B	304	GP4	C1A

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	1707	0	1653	63	0
1	C	1707	0	1653	71	0
2	B	1718	0	1671	42	0
2	D	1718	0	1671	66	0
3	B	76	0	54	3	0
4	C	46	0	35	1	0
5	A	2	0	0	0	0
5	B	1	0	0	0	0
5	C	2	0	0	0	0
6	A	160	0	0	3	0
6	B	220	0	0	1	0
6	C	176	0	0	2	0
6	D	211	0	0	4	0
All	All	7744	0	6737	236	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 17.

All (236) 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)
2:D:154:SER:H	2:D:194:ASN:HD21	1.04	0.99
1:C:114:VAL:HG22	1:C:135:LEU:HD13	1.45	0.98
2:D:105:THR:HG23	6:D:377:HOH:O	1.76	0.85
1:C:210:ARG:HA	1:C:213:CYS:SG	2.15	0.85
1:C:116:ILE:HD13	1:C:117:PHE:N	1.92	0.84
1:C:2:ILE:HG22	1:C:4:MET:CE	2.06	0.84
2:D:154:SER:H	2:D:194:ASN:ND2	1.76	0.82
1:A:194:GLU:HG2	1:A:205:VAL:HG12	1.61	0.81
2:B:69:ARG:HE	2:B:71:ASN:HD21	1.30	0.79
1:C:168:LYS:HZ3	1:C:168:LYS:HB2	1.47	0.79
1:A:210:ARG:HA	1:A:213:CYS:SG	2.22	0.78
1:C:30(E):ARG:HD2	1:C:30(E):ARG:H	1.48	0.78
1:C:117:PHE:HB2	1:C:132:VAL:HG13	1.65	0.77
2:D:48:LEU:HD22	2:D:61:VAL:HG11	1.65	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:163:THR:HG23	6:A:269:HOH:O	1.83	0.77
2:D:119:VAL:HG21	2:D:204:VAL:HG11	1.67	0.77
2:B:154:SER:H	2:B:194:ASN:HD21	1.34	0.76
1:C:2:ILE:HG22	1:C:4:MET:HE1	1.68	0.76
2:B:168:LEU:HB2	2:B:173:TYR:CE1	2.20	0.75
1:C:30(A):ASN:O	1:C:30(D):THR:O	2.06	0.74
2:D:9:GLY:H	2:D:105:THR:HG21	1.52	0.73
1:C:189:ASN:ND2	1:C:211:ASN:HB3	2.04	0.72
1:C:132:VAL:HG11	2:D:122:LEU:HD13	1.71	0.72
1:A:159:LEU:HD21	2:B:169:GLN:NE2	2.05	0.72
2:B:9:GLY:HA3	2:B:105:THR:HG22	1.72	0.71
2:D:30:THR:OG1	2:D:52(B):LYS:HB2	1.90	0.71
2:B:125:GLY:HA2	2:B:212:ASP:O	1.91	0.70
2:B:157:LEU:HD23	2:B:179:VAL:HG21	1.73	0.69
2:B:136:LEU:HD22	2:B:136:LEU:N	2.08	0.68
1:C:168:LYS:H	1:C:168:LYS:HZ2	1.41	0.68
1:A:118:PRO:HG3	2:B:213:CYS:HB2	1.74	0.68
1:C:80:ALA:HA	1:C:105:ILE:HD13	1.76	0.68
2:B:9:GLY:H	2:B:105:THR:HG21	1.59	0.68
1:A:189:ASN:ND2	1:A:211:ASN:HB3	2.09	0.68
2:D:69:ARG:HE	2:D:71:ASN:HD21	1.41	0.67
1:A:83:LEU:HD23	1:A:105:ILE:HD13	1.78	0.66
1:A:13:VAL:HG11	1:A:19:VAL:HG22	1.78	0.66
1:A:184:GLU:HA	1:A:187:ARG:NH1	2.11	0.65
2:D:65:PHE:HD1	2:D:78:LEU:HD11	1.60	0.65
1:C:13:VAL:HG11	1:C:19:VAL:HG22	1.78	0.65
1:C:62:PHE:CD2	1:C:75:ILE:HD12	2.31	0.65
1:C:117:PHE:HB2	1:C:132:VAL:CG1	2.27	0.65
2:D:103:GLN:OE1	6:D:259:HOH:O	2.14	0.64
1:A:46:LEU:HD13	1:A:55:GLU:HG2	1.77	0.64
1:C:107:ARG:HD3	1:C:108:ALA:O	1.97	0.64
1:C:213:CYS:SG	2:D:213:CYS:C	2.76	0.63
1:C:12:ALA:HB2	1:C:104:GLU:HG3	1.80	0.63
2:D:79:GLN:HG2	6:D:233:HOH:O	1.98	0.63
1:C:209:ASN:O	1:C:210:ARG:HB3	1.96	0.63
1:C:135:LEU:HD11	1:C:195:ALA:HB2	1.80	0.63
1:C:2:ILE:CG2	1:C:4:MET:HE1	2.28	0.63
2:D:93:VAL:CG2	2:D:100(C):MET:HB3	2.28	0.63
1:A:61:ARG:HD3	6:A:370:HOH:O	1.99	0.63
1:A:163:THR:HG22	1:A:164:ASP:O	1.98	0.62
1:A:163:THR:HG22	1:A:164:ASP:N	2.15	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:107:ARG:HD3	1:A:108:ALA:O	2.00	0.62
2:D:35:SER:HB2	2:D:93:VAL:CG1	2.30	0.62
1:C:116:ILE:HG12	1:C:133:CYS:SG	2.40	0.62
1:C:3:VAL:O	1:C:4:MET:HE2	1.98	0.62
2:D:93:VAL:HG21	2:D:100(C):MET:HB3	1.81	0.61
1:A:61:ARG:NH2	1:A:82:ASP:OD1	2.33	0.61
2:D:119:VAL:HG21	2:D:204:VAL:CG1	2.30	0.61
2:B:93:VAL:HG21	2:B:100(C):MET:HB3	1.83	0.61
1:A:114:VAL:HG22	1:A:135:LEU:CD2	2.31	0.61
1:A:146:LYS:HE2	1:A:153:GLU:OE2	2.01	0.61
2:B:152:TRP:O	2:B:192:THR:O	2.17	0.61
2:D:65:PHE:CD1	2:D:78:LEU:HD11	2.35	0.60
1:A:105:ILE:HD12	1:A:105:ILE:N	2.16	0.60
2:D:78:LEU:CD1	2:D:80:MET:HG2	2.32	0.60
2:B:69:ARG:NE	2:B:71:ASN:HD21	2.00	0.60
2:D:9:GLY:HA3	2:D:105:THR:HG22	1.84	0.60
1:A:114:VAL:HG22	1:A:135:LEU:HD22	1.83	0.60
2:D:57:TYR:HB2	2:D:62:LYS:HD2	1.85	0.59
2:D:64:ARG:HG2	2:D:64:ARG:HH11	1.68	0.59
1:C:11:LEU:CD1	1:C:103:LEU:HD13	2.33	0.59
2:B:145:PRO:O	2:B:197:HIS:HE1	1.86	0.58
1:A:118:PRO:HG3	2:B:213:CYS:CB	2.32	0.58
1:C:114:VAL:HG22	1:C:135:LEU:CD1	2.26	0.58
2:D:145:PRO:O	2:D:197:HIS:HE1	1.85	0.58
1:C:189:ASN:ND2	1:C:209:ASN:O	2.37	0.58
1:C:168:LYS:HZ2	1:C:168:LYS:N	1.99	0.58
1:A:30(B):SER:HA	1:A:30(E):ARG:HH21	1.69	0.58
1:A:37:GLN:CB	1:A:47:LEU:HD11	2.33	0.58
1:C:30(E):ARG:N	1:C:30(E):ARG:HD2	2.16	0.57
2:B:93:VAL:CG2	2:B:100(C):MET:HB3	2.33	0.57
1:C:3:VAL:HB	1:C:26:SER:OG	2.04	0.57
2:D:11:LEU:HD12	2:D:108:THR:O	2.05	0.56
2:B:123:ALA:HB3	2:B:212:ASP:HB2	1.86	0.56
1:C:213:CYS:SG	2:D:213:CYS:O	2.64	0.56
1:C:20:THR:HG23	1:C:72:THR:CG2	2.36	0.56
4:C:312:KDO:H5	4:C:313:KDO:O1A	2.05	0.56
1:C:116:ILE:HD13	1:C:117:PHE:H	1.71	0.55
2:B:128:ALA:O	2:B:129:GLN:HB2	2.06	0.55
1:A:213:CYS:HB3	2:B:213:CYS:O	2.06	0.55
1:C:2:ILE:HG22	1:C:4:MET:HE3	1.85	0.55
1:C:83:LEU:HD22	6:C:453:HOH:O	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:45:LYS:O	1:A:47:LEU:HD12	2.06	0.55
1:C:168:LYS:H	1:C:168:LYS:NZ	2.05	0.55
1:C:118:PRO:CG	2:D:213:CYS:HB3	2.37	0.54
1:C:24:LYS:HE2	1:C:70:ASP:OD1	2.06	0.54
1:A:33:LEU:HG	1:A:34:ALA:N	2.21	0.54
2:D:119:VAL:CG2	2:D:204:VAL:HG11	2.34	0.54
2:B:212:ASP:O	2:B:213:CYS:HB3	2.06	0.54
1:A:158:VAL:O	1:A:159:LEU:HD12	2.08	0.53
2:D:69:ARG:NE	2:D:71:ASN:HD21	2.06	0.53
2:B:52(B):LYS:HB3	2:B:52(C):PRO:HD3	1.89	0.53
1:A:189:ASN:O	1:A:209:ASN:O	2.25	0.53
1:A:189:ASN:HD21	1:A:211:ASN:HB3	1.72	0.53
2:D:35:SER:O	2:D:93:VAL:HG12	2.08	0.53
1:C:178:LEU:C	1:C:178:LEU:HD13	2.29	0.53
2:D:195:VAL:HB	2:D:204:VAL:HG13	1.91	0.53
2:D:58:SER:O	2:D:62:LYS:HG2	2.09	0.53
2:D:48:LEU:CD2	2:D:61:VAL:HG11	2.38	0.53
1:C:14:SER:N	1:C:106:LYS:HE2	2.23	0.52
1:A:37:GLN:N	1:A:47:LEU:HD11	2.24	0.52
2:D:132:SER:O	2:D:133:MET:HB2	2.08	0.52
1:C:194:GLU:HG2	1:C:205:VAL:HG12	1.90	0.52
1:A:135:LEU:HD13	1:A:143:ILE:HD13	1.91	0.52
2:D:182:PRO:HG2	2:D:185:THR:OG1	2.10	0.52
1:C:189:ASN:HD21	1:C:211:ASN:HB3	1.72	0.52
1:A:158:VAL:C	1:A:159:LEU:HD12	2.30	0.52
1:C:178:LEU:CD1	1:C:180:LEU:HG	2.40	0.52
1:A:189:ASN:ND2	1:A:209:ASN:O	2.43	0.51
1:C:13:VAL:HA	1:C:106:LYS:HE2	1.91	0.51
2:B:35:SER:HB2	2:B:93:VAL:CG1	2.40	0.51
2:B:69:ARG:HE	2:B:71:ASN:ND2	2.05	0.51
2:D:195:VAL:HB	2:D:204:VAL:CG1	2.39	0.51
1:A:33:LEU:HD22	1:A:71:PHE:CG	2.46	0.51
1:C:118:PRO:HB3	1:C:208:PHE:CZ	2.45	0.51
1:A:148:LYS:HA	1:A:152:SER:O	2.10	0.51
1:C:74:THR:C	1:C:75:ILE:HD13	2.31	0.51
2:D:78:LEU:C	2:D:78:LEU:HD13	2.31	0.51
2:D:197:HIS:HD2	2:D:200:SER:OG	1.93	0.50
2:D:210:PRO:O	2:D:211:ARG:HB3	2.12	0.50
2:D:136:LEU:HD22	2:D:136:LEU:N	2.27	0.50
1:C:159:LEU:HG	1:C:159:LEU:O	2.10	0.50
3:B:302:KDO:H5	3:B:303:KDO:O1A	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:131:ASN:O	2:B:132:SER:HB2	2.11	0.49
1:A:109:ASP:OD2	1:A:198:LYS:HE3	2.12	0.49
2:B:169:GLN:O	2:B:170:SER:HB2	2.11	0.49
2:B:29:PHE:CD2	2:B:74:SER:HA	2.47	0.49
1:C:12:ALA:CB	1:C:104:GLU:HG3	2.41	0.49
2:D:52(A):ASN:CG	2:D:52(C):PRO:HD2	2.33	0.49
1:A:3:VAL:H	1:A:26:SER:HB2	1.76	0.49
1:A:19:VAL:CG2	1:A:78:VAL:HG21	2.43	0.49
1:A:163:THR:CG2	1:A:164:ASP:N	2.75	0.49
1:C:189:ASN:HD21	1:C:211:ASN:CB	2.26	0.48
2:B:97:PHE:HB2	3:B:302:KDO:H32	1.95	0.48
2:D:64:ARG:CZ	2:D:64:ARG:HB2	2.41	0.48
1:A:104:GLU:C	1:A:105:ILE:HD12	2.34	0.48
2:B:80:MET:HB3	2:B:82(A):LEU:HD21	1.96	0.48
2:B:149:THR:HG22	6:B:436:HOH:O	2.13	0.48
2:B:117:PRO:HB3	2:B:143:TYR:HB3	1.94	0.48
2:D:175:LEU:C	2:D:175:LEU:HD12	2.34	0.48
2:D:59:ALA:HA	2:D:62:LYS:CG	2.43	0.47
2:B:154:SER:H	2:B:194:ASN:ND2	2.06	0.47
1:C:13:VAL:C	1:C:106:LYS:HE2	2.34	0.47
1:C:83:LEU:HD21	6:C:390:HOH:O	2.14	0.47
1:C:33:LEU:HD22	1:C:71:PHE:CG	2.49	0.47
1:C:13:VAL:CA	1:C:106:LYS:HE2	2.45	0.47
1:C:182:LYS:O	1:C:186:GLU:HG3	2.14	0.47
1:C:19:VAL:HG12	1:C:20:THR:N	2.29	0.47
2:D:61:VAL:HB	2:D:65:PHE:CD2	2.49	0.47
2:B:135:THR:C	2:B:136:LEU:HD22	2.34	0.47
2:D:130:THR:HG22	2:D:131:ASN:N	2.30	0.47
1:A:30(D):THR:O	1:A:30(E):ARG:HB2	2.14	0.47
2:D:103:GLN:CD	6:D:236:HOH:O	2.53	0.47
1:C:168:LYS:N	1:C:168:LYS:NZ	2.63	0.46
2:D:36:TRP:CE2	2:D:78:LEU:HB2	2.50	0.46
2:B:185:THR:O	2:B:189:GLU:HB2	2.15	0.46
2:D:69:ARG:HE	2:D:71:ASN:ND2	2.11	0.46
1:C:20:THR:HG23	1:C:72:THR:HG23	1.97	0.46
1:A:47:LEU:N	1:A:47:LEU:HD12	2.31	0.46
2:D:6:GLU:HB2	2:D:105:THR:OG1	2.16	0.46
2:B:197:HIS:HD2	2:B:200:SER:OG	1.99	0.46
2:D:121:PRO:HD3	2:D:206:LYS:HG2	1.98	0.46
2:D:185:THR:O	2:D:189:GLU:N	2.36	0.45
1:C:135:LEU:HD11	1:C:195:ALA:CB	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:136:LEU:N	2:B:136:LEU:CD2	2.78	0.45
1:A:209:ASN:O	1:A:210:ARG:CB	2.62	0.45
1:A:118:PRO:CG	2:B:213:CYS:HB2	2.44	0.45
1:C:54:ARG:HG2	1:C:58:VAL:HB	1.98	0.45
1:A:116:ILE:HG13	1:A:133:CYS:SG	2.57	0.45
1:A:19:VAL:HG23	1:A:78:VAL:HG21	1.99	0.45
2:D:64:ARG:NH2	2:D:83:ARG:HH21	2.15	0.44
1:C:4:MET:HE2	1:C:25:SER:HA	1.98	0.44
2:D:38:ARG:CD	2:D:48:LEU:HD21	2.48	0.44
1:A:3:VAL:HB	1:A:26:SER:OG	2.17	0.44
1:C:189:ASN:O	1:C:209:ASN:O	2.36	0.44
2:D:35:SER:HB2	2:D:93:VAL:HG13	1.99	0.44
1:C:118:PRO:HB3	1:C:208:PHE:CE2	2.53	0.44
1:A:83:LEU:C	1:A:83:LEU:HD13	2.38	0.44
1:A:83:LEU:O	1:A:84:ALA:HB2	2.18	0.44
1:C:168:LYS:HB2	1:C:168:LYS:NZ	2.25	0.44
2:D:47:TRP:CH2	2:D:49:GLY:HA2	2.53	0.44
2:B:133:MET:SD	2:B:182:PRO:HA	2.58	0.43
1:C:93:ASN:C	1:C:93:ASN:HD22	2.20	0.43
2:D:64:ARG:CG	2:D:64:ARG:HH11	2.29	0.43
1:A:19:VAL:HG23	1:A:78:VAL:CG2	2.49	0.43
2:B:182:PRO:HG2	2:B:185:THR:OG1	2.18	0.43
1:C:168:LYS:CB	1:C:168:LYS:HZ3	2.26	0.43
2:D:59:ALA:HA	2:D:62:LYS:HG3	2.01	0.43
1:C:186:GLU:HA	1:C:210:ARG:NH2	2.34	0.43
1:A:37:GLN:N	1:A:47:LEU:CD1	2.81	0.43
1:A:105:ILE:CD1	1:A:105:ILE:N	2.82	0.42
1:A:42:GLN:HG3	6:A:228:HOH:O	2.19	0.42
2:B:131:ASN:O	2:B:132:SER:CB	2.67	0.42
2:D:154:SER:N	2:D:194:ASN:ND2	2.57	0.42
2:D:126:SER:N	2:D:213:CYS:OXT	2.52	0.42
1:C:11:LEU:CD1	1:C:103:LEU:CD1	2.97	0.42
2:D:78:LEU:HD12	2:D:80:MET:HG2	2.00	0.42
2:D:189:GLU:OE2	2:D:189:GLU:HA	2.20	0.42
1:C:83:LEU:C	1:C:83:LEU:HD13	2.40	0.42
2:B:52(A):ASN:CG	2:B:52(C):PRO:HD2	2.40	0.42
2:D:47:TRP:CZ2	2:D:49:GLY:HA2	2.54	0.42
1:C:33:LEU:HG	1:C:34:ALA:N	2.34	0.41
1:A:132:VAL:HG22	1:A:177:THR:HG23	2.00	0.41
2:D:210:PRO:O	2:D:211:ARG:CB	2.68	0.41
2:B:175:LEU:C	2:B:175:LEU:HD12	2.40	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:LEU:HB2	1:A:105:ILE:HD13	2.01	0.41
1:A:81:GLU:CD	1:A:81:GLU:H	2.24	0.41
1:A:168:LYS:NZ	1:A:168:LYS:HB3	2.35	0.41
1:A:191:TYR:O	1:A:207:SER:HB2	2.19	0.41
1:C:184:GLU:OE2	1:C:187:ARG:CZ	2.69	0.41
3:B:301:KDO:H5	3:B:302:KDO:C1	2.50	0.41
1:C:178:LEU:HD11	1:C:180:LEU:HG	2.02	0.41
2:D:52(B):LYS:HB3	2:D:52(C):PRO:HD3	2.02	0.41
1:A:46:LEU:HD13	1:A:55:GLU:CG	2.46	0.40
1:A:146:LYS:NZ	1:A:153:GLU:HB2	2.35	0.40
1:A:189:ASN:HD21	1:A:211:ASN:CB	2.34	0.40
1:A:186:GLU:O	1:A:210:ARG:NH2	2.54	0.40
1:A:37:GLN:HB2	1:A:47:LEU:HD11	2.02	0.40
2:D:181:VAL:HB	2:D:182:PRO:HD2	2.03	0.40
1:A:135:LEU:HD13	1:A:143:ILE:CD1	2.51	0.40
2:D:64:ARG:NH2	2:D:83:ARG:NH2	2.70	0.40
2:B:121:PRO:HB3	2:B:206:LYS:HG2	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	217/219 (99%)	209 (96%)	8 (4%)	0	100	100
1	C	217/219 (99%)	205 (94%)	12 (6%)	0	100	100
2	B	224/226 (99%)	208 (93%)	12 (5%)	4 (2%)	11	1
2	D	224/226 (99%)	214 (96%)	6 (3%)	4 (2%)	11	1
All	All	882/890 (99%)	836 (95%)	38 (4%)	8 (1%)	21	6

All (8) Ramachandran outliers are listed below:



Mol	Chain	Res	Type
2	B	132	SER
2	B	211	ARG
2	D	41	PRO
2	D	128	ALA
2	D	211	ARG
2	B	130	THR
2	B	159	SER
2	D	133	MET

### 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	194/194 (100%)	190 (98%)	4 (2%)	61	37
1	C	194/194 (100%)	184 (95%)	10 (5%)	29	7
2	B	196/196 (100%)	187 (95%)	9 (5%)	33	10
2	D	196/196 (100%)	188 (96%)	8 (4%)	37	13
All	All	780/780 (100%)	749 (96%)	31 (4%)	38	14

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	LEU
1	A	55	GLU
1	A	107	ARG
1	A	124	LEU
2	B	20	LEU
2	B	50	PHE
2	B	131	ASN
2	B	136	LEU
2	B	138	CYS
2	B	147	PRO
2	B	157	LEU
2	B	193	CYS
2	B	206	LYS
1	C	30(E)	ARG

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Mol	Chain	Res	Type
1	C	46	LEU
1	C	47	LEU
1	C	83	LEU
1	C	93	ASN
1	C	107	ARG
1	C	116	ILE
1	C	124	LEU
1	C	159	LEU
1	C	168	LYS
2	D	18	LEU
2	D	20	LEU
2	D	41	PRO
2	D	50	PHE
2	D	129	GLN
2	D	136	LEU
2	D	147	PRO
2	D	148	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	GLN
1	A	136	ASN
1	A	137	ASN
1	A	155	GLN
1	A	189	ASN
1	A	209	ASN
2	B	71	ASN
2	B	162	HIS
2	B	169	GLN
2	B	194	ASN
2	B	197	HIS
1	C	27	GLN
1	C	93	ASN
1	C	136	ASN
1	C	137	ASN
1	C	189	ASN
1	C	209	ASN
2	D	71	ASN
2	D	162	HIS
2	D	194	ASN
2	D	197	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

8 carbohydrates are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z  > 2$	Counts	RMSZ	# $ Z  > 2$
3	KDO	B	301	3	12,15,16	2.07	3 (25%)	12,21,24	1.27	2 (16%)
3	KDO	B	302	3	12,15,16	2.04	3 (25%)	12,21,24	1.38	3 (25%)
3	KDO	B	303	3	12,15,16	1.74	2 (16%)	12,21,24	0.97	1 (8%)
3	GP4	B	304	3	15,15,16	1.27	1 (6%)	18,22,24	1.29	2 (11%)
3	GP1	B	305	3	15,16,16	1.46	2 (13%)	20,24,24	0.65	0
4	KDO	C	311	4	13,16,16	2.24	4 (30%)	15,24,24	1.31	1 (6%)
4	KDO	C	312	4	12,15,16	2.08	3 (25%)	12,21,24	1.45	2 (16%)
4	KDO	C	313	4	12,15,16	1.84	3 (25%)	12,21,24	0.94	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	KDO	B	301	3	-	0/6/26/30	0/1/1/1
3	KDO	B	302	3	-	0/6/26/30	0/1/1/1
3	KDO	B	303	3	-	0/6/26/30	0/1/1/1
3	GP4	B	304	3	1/1/5/6	0/7/24/27	0/1/1/1
3	GP1	B	305	3	-	0/6/27/27	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	KDO	C	311	4	-	0/6/30/30	0/1/1/1
4	KDO	C	312	4	-	0/6/26/30	0/1/1/1
4	KDO	C	313	4	-	0/6/26/30	0/1/1/1

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	305	GP1	P4B-O1B	-2.61	1.52	1.60
3	B	304	GP4	P4A-O4A	-2.36	1.52	1.60
3	B	303	KDO	O6-C6	2.03	1.47	1.43
3	B	302	KDO	O5-C5	2.07	1.47	1.43
3	B	302	KDO	O6-C6	2.08	1.47	1.43
4	C	313	KDO	C3-C4	2.12	1.56	1.52
4	C	312	KDO	O5-C5	2.16	1.48	1.43
3	B	301	KDO	O6-C2	2.25	1.50	1.43
4	C	311	KDO	O6-C6	2.34	1.48	1.44
4	C	313	KDO	O6-C6	2.49	1.48	1.43
4	C	312	KDO	O6-C6	2.50	1.48	1.43
3	B	305	GP1	C3B-C2B	3.18	1.57	1.53
4	C	311	KDO	C4-C5	3.42	1.57	1.52
3	B	301	KDO	O6-C6	3.63	1.49	1.43
4	C	311	KDO	O2-C2	4.14	1.44	1.40
3	B	303	KDO	C4-C5	4.49	1.58	1.52
4	C	311	KDO	O6-C2	4.60	1.47	1.42
4	C	313	KDO	C4-C5	4.81	1.59	1.52
3	B	302	KDO	C4-C5	5.35	1.59	1.52
3	B	301	KDO	C4-C5	5.43	1.60	1.52
4	C	312	KDO	C4-C5	5.69	1.60	1.52

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	311	KDO	C3-C4-C5	-3.81	107.34	110.84
4	C	312	KDO	C3-C4-C5	-2.61	107.00	110.56
3	B	302	KDO	C3-C4-C5	-2.27	107.47	110.56
3	B	303	KDO	C3-C4-C5	-2.20	107.56	110.56
3	B	302	KDO	O4-C4-C3	-2.00	105.16	110.06
3	B	301	KDO	O6-C2-C3	2.04	113.78	109.86
3	B	304	GP4	C1A-C2A-C3A	2.12	112.70	109.27
3	B	301	KDO	O5-C5-C4	2.13	113.85	110.00
3	B	302	KDO	O6-C2-C3	2.62	114.90	109.86
4	C	312	KDO	O6-C2-C3	2.83	115.29	109.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	304	GP4	C1A-O6A-C5A	3.63	116.86	112.25

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	B	304	GP4	C1A

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	301	KDO	1	0
3	B	302	KDO	3	0
3	B	303	KDO	1	0
4	C	312	KDO	1	0
4	C	313	KDO	1	0

## 5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 5 are monoatomic - leaving 0 for Mogul analysis.

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	219/219 (100%)	0.45	13 (5%) 26 31	16, 29, 48, 70	0
1	C	219/219 (100%)	0.37	12 (5%) 29 34	18, 29, 42, 55	0
2	B	226/226 (100%)	0.47	18 (7%) 15 19	17, 27, 55, 77	0
2	D	226/226 (100%)	0.42	17 (7%) 17 22	17, 25, 50, 70	0
All	All	890/890 (100%)	0.43	60 (6%) 21 26	16, 28, 50, 77	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	213	CYS	10.6
2	B	130	THR	10.2
2	B	213	CYS	8.9
2	B	128	ALA	7.2
2	B	129	GLN	6.3
1	A	213	CYS	6.1
2	D	132	SER	5.6
2	B	131	ASN	5.5
2	D	130	THR	5.5
2	B	127	ALA	5.0
2	D	128	ALA	4.9
2	D	129	GLN	4.8
2	D	131	ASN	4.7
1	A	211	ASN	4.7
2	D	211	ARG	4.2
1	C	1	ASP	4.2
2	D	138	CYS	4.1
2	B	170	SER	4.0
2	B	212	ASP	3.8
2	B	158	SER	3.8
1	C	33	LEU	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	201	THR	3.6
1	C	159	LEU	3.5
1	C	213	CYS	3.5
1	A	202	SER	3.3
2	D	189	GLU	3.3
2	B	126	SER	3.3
1	C	30(E)	ARG	3.3
1	A	1	ASP	3.2
1	C	156	ASN	3.1
1	C	211	ASN	3.1
2	B	159	SER	3.1
2	B	211	ARG	3.0
2	D	184	SER	2.9
2	D	127	ALA	2.9
1	A	187	ARG	2.7
1	C	168	LYS	2.6
1	C	30(B)	SER	2.6
1	A	209	ASN	2.6
1	C	187	ARG	2.5
2	B	152	TRP	2.5
2	B	169	GLN	2.5
2	D	212	ASP	2.5
2	B	92	CYS	2.3
1	A	183	ASP	2.3
2	B	157	LEU	2.2
1	A	200	SER	2.2
2	D	136	LEU	2.2
2	B	156	SER	2.2
2	D	158	SER	2.2
2	B	76	LEU	2.2
1	C	30(D)	THR	2.1
2	D	188	SER	2.1
1	C	35	TRP	2.1
2	D	122	LEU	2.1
1	A	168	LYS	2.1
2	D	41	PRO	2.1
1	A	105	ILE	2.0
1	A	60	ASP	2.0
1	A	83	LEU	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](#)

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	KDO	B	303	15/16	0.91	0.14	2.74	17,20,29,37	0
4	KDO	C	313	15/16	0.87	0.13	2.48	25,27,35,39	0
4	KDO	C	312	15/16	0.84	0.13	0.60	28,35,42,45	0
3	KDO	B	301	15/16	0.89	0.11	0.57	22,25,32,35	0
3	GP4	B	304	15/16	0.90	0.16	-	27,32,47,47	0
3	GP1	B	305	16/16	0.94	0.10	-	27,29,36,38	0
3	KDO	B	302	15/16	0.91	0.12	-	20,24,29,32	0
4	KDO	C	311	16/16	0.80	0.21	-	40,49,52,53	0

## 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
5	MG	A	214	1/1	1.00	0.10	1.40	16,16,16,16	0
5	MG	B	306	1/1	0.98	0.06	-0.79	23,23,23,23	0
5	MG	C	314	1/1	0.98	0.08	-0.84	24,24,24,24	0
5	MG	A	215	1/1	0.99	0.07	-1.27	26,26,26,26	0
5	MG	C	315	1/1	0.99	0.05	-2.12	32,32,32,32	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.