



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:15 PM GMT

PDB ID : 1O75  
Title : TP47, THE 47-KILODALTON LIPOPROTEIN OF TREPONEMA PAL-  
LIDUM  
Authors : Deka, R.K.; Machius, M.; Norgard, M.V.; Tomchick, D.R.  
Deposited on : 2002-10-23  
Resolution : 1.95 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

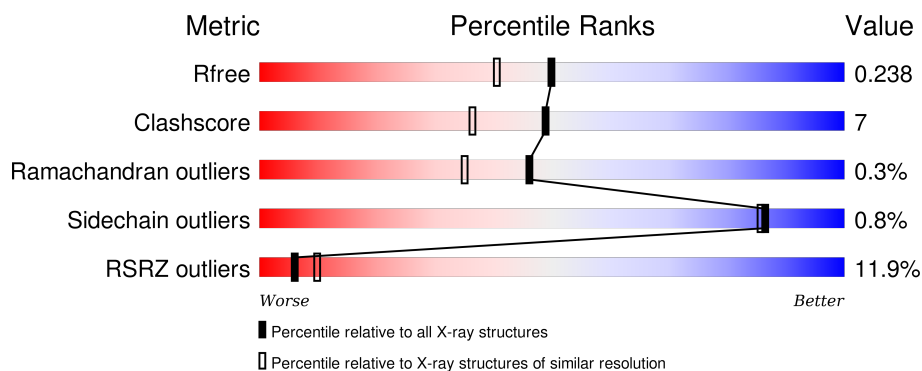
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.95 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	415	<div> <div>9%</div> <div>87%</div> <div>9%</div> <div>.</div> </div>
1	B	415	<div> <div>13%</div> <div>80%</div> <div>16%</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
2	XE	B	1415	-	-	X	-
2	XE	B	1417	-	-	X	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7017 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 47 KDA MEMBRANE ANTIGEN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	399	Total	C	N	O	S	0	9	0
			3175	1993	539	630	13			
1	B	402	Total	C	N	O	S	0	5	0
			3179	1998	542	626	13			

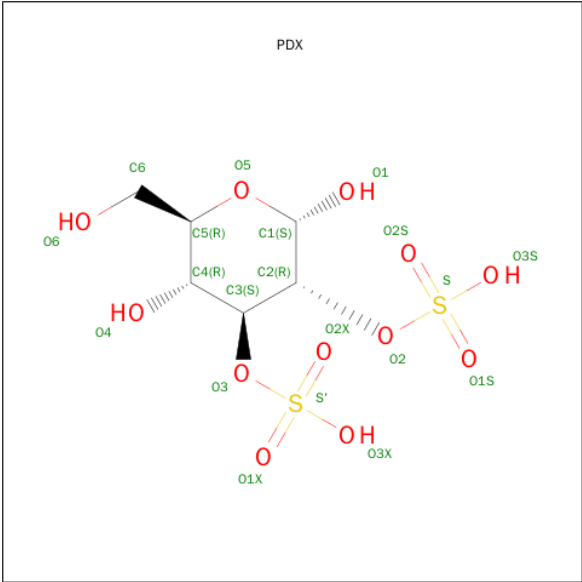
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	5	SER	HIS	ENGINEERED MUTATION	UNP P29723
A	9	SER	HIS	ENGINEERED MUTATION	UNP P29273
B	5	SER	HIS	ENGINEERED MUTATION	UNP P29723
B	9	SER	HIS	ENGINEERED MUTATION	UNP P29273

- Molecule 2 is XENON (three-letter code: XE) (formula: Xe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	3	Total	Xe	0	0
			3	3		
2	A	2	Total	Xe	0	0
			2	2		

- Molecule 3 is 2,3-DI-O-SULFO-ALPHA-D-GLUCOPYRANOSE (three-letter code: PDX) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>12</sub>S<sub>2</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	S	0	0
			19	6	11	2		
3	B	1	Total	C	O	S	0	0
			20	6	12	2		

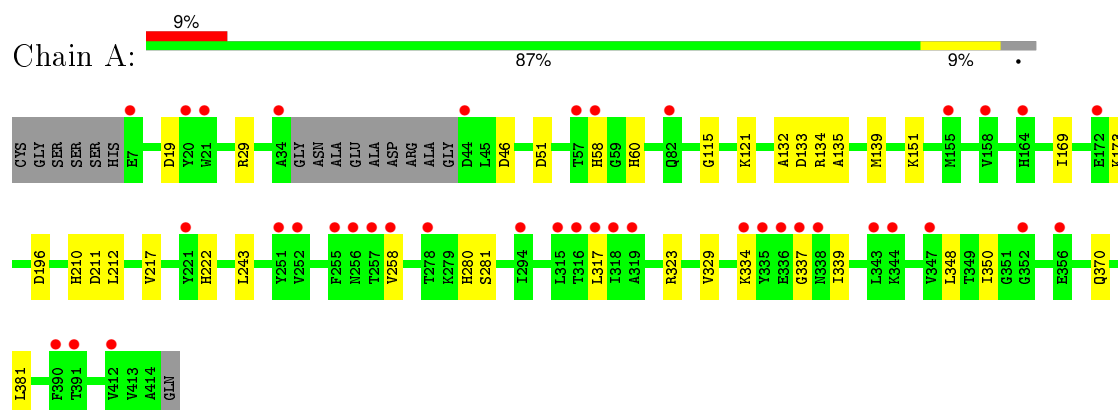
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	311	Total	O	0	0
			311	311		
4	B	308	Total	O	0	0
			308	308		

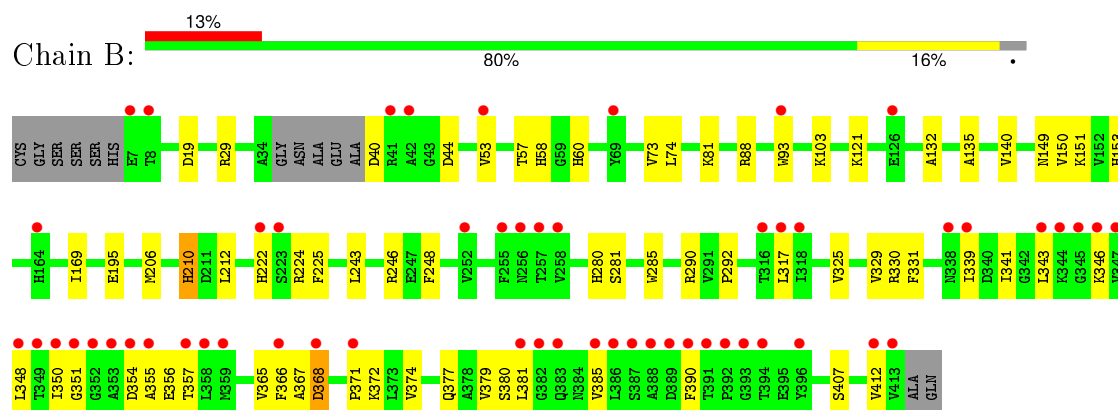
### 3 Residue-property plots

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: 47 KDA MEMBRANE ANTIGEN



#### • Molecule 1: 47 KDA MEMBRANE ANTIGEN



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	128.93Å 128.93Å 151.16Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	28.20 – 1.95 28.15 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.8 (28.20-1.95) 94.2 (28.15-1.95)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.17 (at 1.95Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.205 , 0.228 0.216 , 0.238	Depositor DCC
$R_{free}$ test set	5036 reflections (5.05%)	DCC
Wilson B-factor (Å <sup>2</sup> )	33.7	Xtriage
Anisotropy	0.034	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 50.9	EDS
Estimated twinning fraction	0.019 for -h,-k,l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 99701 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7017	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.83% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PDX, XE

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	0.51	0/3239	0.68	0/4372
1	B	0.51	0/3245	0.68	0/4380
All	All	0.51	0/6484	0.68	0/8752

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	3175	0	3036	26	0
1	B	3179	0	3042	67	0
2	A	2	0	0	1	0
2	B	3	0	0	7	0
3	A	19	0	7	0	0
3	B	20	0	11	1	0
4	A	311	0	0	2	0
4	B	308	0	0	1	0
All	All	7017	0	6096	90	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 7.



All (90) 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:339:ILE:HD11	1:B:350:ILE:HG23	1.44	0.98
1:B:151:LYS:HE2	1:B:153:HIS:CE1	2.06	0.91
1:B:169:ILE:HD13	1:B:195:GLU:HB2	1.54	0.88
1:B:350:ILE:HD12	1:B:379:VAL:HG11	1.57	0.85
1:B:19:ASP:OD1	1:B:29:ARG:HD3	1.79	0.83
1:A:211:ASP:OD2	1:A:323[A]:ARG:HD3	1.80	0.79
1:B:93[B]:TRP:CZ3	2:B:1417:XE:XE	3.20	0.72
1:B:350:ILE:HD12	1:B:379:VAL:CG1	2.20	0.70
1:B:350:ILE:CD1	1:B:379:VAL:HG11	2.27	0.64
1:B:169:ILE:CD1	1:B:195:GLU:HB2	2.25	0.64
1:A:222:HIS:O	1:B:371:PRO:HG3	1.97	0.63
1:B:339:ILE:HD12	1:B:351:GLY:O	1.98	0.63
1:B:53:VAL:O	1:B:151:LYS:HA	2.00	0.62
1:B:58:HIS:CD2	1:B:60:HIS:H	2.18	0.61
1:B:212:LEU:HD22	1:B:243:LEU:HD21	1.83	0.60
1:A:212:LEU:HD22	1:A:243:LEU:HD21	1.84	0.60
1:B:57:THR:CG2	1:B:150:VAL:HG23	2.32	0.60
1:A:370:GLN:HG2	1:B:225:PHE:CE1	2.37	0.59
1:A:217:VAL:HG21	1:A:329[A]:VAL:CG1	2.35	0.57
1:B:29:ARG:HG2	1:B:29:ARG:HH11	1.71	0.56
1:A:258:VAL:HG13	1:A:317:LEU:HD23	1.88	0.55
1:A:173:LYS:NZ	1:A:196:ASP:HB3	2.21	0.55
1:B:81:LYS:HE3	2:B:1417:XE:XE	2.85	0.54
1:A:169:ILE:HG13	4:A:2139:HOH:O	2.07	0.54
1:A:58:HIS:HD2	1:A:60:HIS:CG	2.25	0.54
1:B:331:PHE:CZ	2:B:1415:XE:XE	3.39	0.53
1:B:88:ARG:NH2	3:B:1414:PDX:H4	2.22	0.53
1:B:103:LYS:HG2	1:B:140:VAL:HG21	1.91	0.53
1:B:350:ILE:O	1:B:381:LEU:HD12	2.08	0.53
1:B:246:ARG:CZ	4:B:2209:HOH:O	2.57	0.53
1:A:173:LYS:HZ1	1:A:196:ASP:HB3	1.73	0.52
1:B:379:VAL:CG1	1:B:380:SER:N	2.74	0.51
1:A:19:ASP:OD1	1:A:29:ARG:HD3	2.09	0.51
1:B:341:ILE:HG23	1:B:348:LEU:HD11	1.92	0.51
1:B:222:HIS:HE1	1:B:407:SER:C	2.14	0.50
1:B:73:VAL:HG12	1:B:150:VAL:HG22	1.93	0.50
1:B:354:ASP:OD1	1:B:355:ALA:N	2.44	0.49
1:B:354:ASP:O	1:B:357:THR:HG22	2.12	0.49
1:B:224:ARG:NH1	1:B:285:TRP:CG	2.81	0.49
1:B:343:LEU:HD22	1:B:390:PHE:CE2	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:334:LYS:HE3	4:A:2097:HOH:O	2.12	0.48
1:B:354:ASP:OD1	1:B:356:GLU:N	2.42	0.48
1:A:350:ILE:O	1:A:381:LEU:HD12	2.13	0.48
1:B:93[A]:TRP:CD1	2:B:1417:XE:XE	3.44	0.48
1:B:40:ASP:HB3	1:B:44:ASP:HB2	1.96	0.48
1:B:151:LYS:HE2	1:B:153:HIS:HE1	1.72	0.48
1:B:93[B]:TRP:CE3	2:B:1417:XE:XE	3.45	0.47
1:A:46[B]:ASP:N	1:A:46[B]:ASP:OD2	2.46	0.47
1:B:331:PHE:HZ	2:B:1415:XE:XE	2.75	0.47
1:B:331:PHE:CE2	2:B:1415:XE:XE	3.45	0.47
1:B:58:HIS:CD2	1:B:60:HIS:HB2	2.50	0.47
1:B:248:PHE:CD1	1:B:292:PRO:HA	2.50	0.47
1:A:115:GLY:HA3	1:A:121:LYS:HE3	1.96	0.47
1:A:370:GLN:HG2	1:B:225:PHE:CZ	2.50	0.46
1:B:224:ARG:NH2	1:B:281:SER:O	2.48	0.46
1:A:339:ILE:HG23	1:A:339:ILE:O	2.16	0.46
1:B:374:VAL:HG13	1:B:377:GLN:CG	2.46	0.46
1:B:57:THR:HG23	1:B:150:VAL:HG23	1.98	0.46
1:A:280:HIS:O	1:A:281:SER:HB3	2.15	0.45
1:B:280:HIS:O	1:B:281:SER:HB3	2.17	0.45
1:B:248:PHE:CE1	1:B:292:PRO:HA	2.52	0.45
1:B:74:LEU:HD23	1:B:149:ASN:HD22	1.82	0.44
1:B:132:ALA:O	1:B:135:ALA:HB2	2.17	0.44
1:A:151:LYS:HB2	2:A:1416:XE:XE	2.96	0.44
1:A:339:ILE:HD11	1:A:350:ILE:CG2	2.48	0.44
1:B:354:ASP:C	1:B:357:THR:HG22	2.38	0.44
1:A:58:HIS:CD2	1:A:60:HIS:CG	3.05	0.44
1:B:74:LEU:HD23	1:B:149:ASN:ND2	2.33	0.43
1:B:341:ILE:HG23	1:B:348:LEU:CD1	2.48	0.43
1:A:132:ALA:O	1:A:135:ALA:HB2	2.18	0.43
1:B:212:LEU:CD2	1:B:243:LEU:HD21	2.48	0.43
1:B:366:PHE:HB3	1:B:372:LYS:HA	2.01	0.43
1:B:317:LEU:HB2	1:B:325:VAL:HB	2.00	0.42
1:A:133:ASP:O	1:A:134:ARG:HB2	2.19	0.42
1:B:365:VAL:HG11	1:B:390:PHE:CE1	2.55	0.42
1:A:217:VAL:HG21	1:A:329[B]:VAL:HG22	2.01	0.42
1:B:121:LYS:HD3	1:B:121:LYS:HA	1.88	0.42
1:B:222:HIS:CE1	1:B:407:SER:O	2.73	0.41
1:B:367:ALA:O	1:B:368:ASP:C	2.59	0.41
1:B:29:ARG:HG2	1:B:29:ARG:NH1	2.34	0.41
1:B:329:VAL:HG22	1:B:330:ARG:N	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:290:ARG:NH1	1:B:290:ARG:HG2	2.35	0.41
1:A:217:VAL:HG21	1:A:329[B]:VAL:CG2	2.50	0.41
1:B:88:ARG:HG2	1:B:88:ARG:O	2.20	0.41
1:B:346:LYS:O	1:B:385:VAL:HA	2.21	0.41
1:B:343:LEU:HG	1:B:412:VAL:HG11	2.03	0.40
1:B:210:HIS:CD2	1:B:210:HIS:C	2.94	0.40
1:A:348:LEU:C	1:A:348:LEU:HD23	2.41	0.40
1:B:290:ARG:HH11	1:B:290:ARG:HG2	1.86	0.40
1:B:222:HIS:CE1	1:B:407:SER:C	2.94	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	404/415 (97%)	391 (97%)	12 (3%)	1 (0%)	52	43
1	B	403/415 (97%)	389 (96%)	13 (3%)	1 (0%)	52	43
All	All	807/830 (97%)	780 (97%)	25 (3%)	2 (0%)	46	43

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	368	ASP
1	A	337	GLY

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	336/337 (100%)	333 (99%)	3 (1%)	84	83
1	B	334/337 (99%)	332 (99%)	2 (1%)	90	89
All	All	670/674 (99%)	665 (99%)	5 (1%)	86	88

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	ASP
1	A	139	MET
1	A	210	HIS
1	B	206	MET
1	B	210	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	HIS
1	A	82	GLN
1	A	149	ASN
1	B	149	ASN
1	B	153	HIS
1	B	185	ASN
1	B	222	HIS

### 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

Of 7 ligands modelled in this entry, 5 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	PDX	A	1417	3	18,19,20	1.78	4 (22%)	26,30,31	1.64	4 (15%)
3	PDX	B	1414	3	19,20,20	1.54	5 (26%)	27,31,31	0.90	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	PDX	A	1417	3	-	0/10/30/32	0/1/1/1
3	PDX	B	1414	3	-	0/12/32/32	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1414	PDX	O2-S	-2.93	1.47	1.57
3	B	1414	PDX	O1S-S	-2.85	1.34	1.45
3	A	1417	PDX	O1S-S	-2.76	1.34	1.45
3	B	1414	PDX	O2X-S'	-2.36	1.36	1.45
3	B	1414	PDX	O3-S'	-2.05	1.50	1.57
3	A	1417	PDX	O3-S'	-2.04	1.50	1.57
3	A	1417	PDX	O2-S	-2.01	1.50	1.57
3	B	1414	PDX	O3S-S	2.27	1.62	1.50
3	A	1417	PDX	O2S-S	4.87	1.63	1.45

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1417	PDX	O3-S'-O1X	-2.56	98.57	106.86
3	A	1417	PDX	O2-C2-C1	2.23	110.68	107.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1417	PDX	O4-C4-C5	2.80	116.41	109.84
3	A	1417	PDX	C3-O3-S'	5.81	129.83	118.77

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
3	B	1414	PDX	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	399/415 (96%)	0.53	39 (9%) 10 15	25, 39, 68, 76	0
1	B	402/415 (96%)	0.73	56 (13%) 4 6	24, 38, 76, 89	0
All	All	801/830 (96%)	0.63	95 (11%) 6 10	24, 38, 74, 89	0

All (95) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	381	LEU	7.6
1	B	390	PHE	6.6
1	A	221	TYR	5.7
1	B	391	THR	5.4
1	B	7	GLU	5.3
1	B	347	VAL	5.2
1	B	382	GLY	5.2
1	A	337	GLY	5.1
1	B	385	VAL	5.1
1	B	352	GLY	5.0
1	B	343	LEU	5.0
1	B	53	VAL	4.9
1	B	392	PRO	4.8
1	B	412	VAL	4.8
1	B	383	GLN	4.8
1	A	34	ALA	4.8
1	B	368	ASP	4.7
1	A	255	PHE	4.6
1	A	338	ASN	4.5
1	B	413	VAL	4.4
1	A	58	HIS	4.3
1	A	318	ILE	4.2
1	B	355	ALA	4.2
1	A	258	VAL	4.0

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Mol	Chain	Res	Type	RSRZ
1	B	359	MET	4.0
1	B	345	GLY	3.9
1	B	255	PHE	3.9
1	A	257	THR	3.9
1	B	354	ASP	3.8
1	B	344	LYS	3.7
1	B	357	THR	3.7
1	A	317	LEU	3.6
1	B	348	LEU	3.6
1	B	358	LEU	3.6
1	B	93[A]	TRP	3.6
1	B	393	GLY	3.5
1	B	386	LEU	3.5
1	B	258	VAL	3.4
1	B	394	THR	3.4
1	B	42	ALA	3.3
1	B	353	ALA	3.3
1	A	44	ASP	3.3
1	A	252	VAL	3.3
1	A	390	PHE	3.2
1	B	338	ASN	3.2
1	B	318	ILE	3.2
1	B	252	VAL	3.1
1	A	57	THR	3.1
1	B	389	ASP	3.1
1	A	344	LYS	3.0
1	B	351	GLY	3.0
1	B	164	HIS	3.0
1	B	69	TYR	2.9
1	B	223	SER	2.8
1	B	387	SER	2.8
1	B	366	PHE	2.8
1	A	172	GLU	2.8
1	A	251	TYR	2.8
1	A	412	VAL	2.7
1	A	82	GLN	2.7
1	B	346	LYS	2.7
1	A	347	VAL	2.7
1	B	350	ILE	2.7
1	A	256	ASN	2.7
1	B	396	TYR	2.7
1	B	257	THR	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	339	ILE	2.6
1	A	164	HIS	2.6
1	B	317	LEU	2.6
1	A	294	ILE	2.6
1	A	7	GLU	2.5
1	A	158	VAL	2.5
1	B	8	THR	2.5
1	A	352	GLY	2.5
1	A	319	ALA	2.5
1	A	336	GLU	2.5
1	B	388	ALA	2.5
1	B	41	ARG	2.5
1	A	316	THR	2.5
1	A	278	THR	2.4
1	A	335	TYR	2.4
1	B	126	GLU	2.4
1	A	21	TRP	2.3
1	A	356	GLU	2.2
1	A	334	LYS	2.2
1	A	155	MET	2.2
1	A	343	LEU	2.2
1	A	315	LEU	2.1
1	B	371	PRO	2.1
1	A	20	TYR	2.1
1	B	256	ASN	2.1
1	A	391	THR	2.1
1	B	316	THR	2.1
1	B	349	THR	2.1
1	B	222	HIS	2.0

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

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

## 6.3 Carbohydrates ⓘ

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	PDX	A	1417	19/20	0.94	0.16	0.73	52,55,57,58	0
3	PDX	B	1414	20/20	0.93	0.18	0.42	52,56,60,60	0
2	XE	B	1417	1/1	0.97	0.11	-0.97	58,58,58,58	1
2	XE	A	1415	1/1	1.00	0.08	-1.13	35,35,35,35	1
2	XE	B	1415	1/1	0.98	0.10	-1.37	34,34,34,34	1
2	XE	B	1416	1/1	0.98	0.06	-2.41	45,45,45,45	1
2	XE	A	1416	1/1	0.95	0.06	-2.61	56,56,56,56	1

## 6.5 Other polymers [i](#)

There are no such residues in this entry.