Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 06:18 PM GMT

PDB ID	:	4L72
Title	:	Crystal structure of MERS-CoV complexed with human DPP4
Authors	:	Wang, X.Q.; Wang, N.S.
Deposited on	:	2013-06-13
Resolution	:	3.00 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
A user guide is available at
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

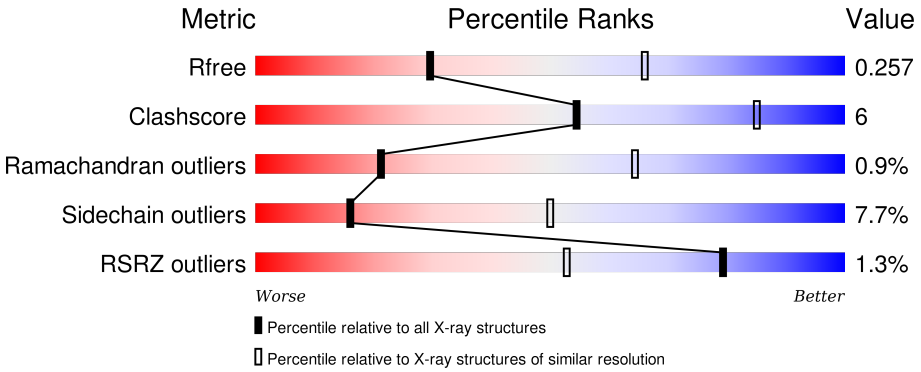
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

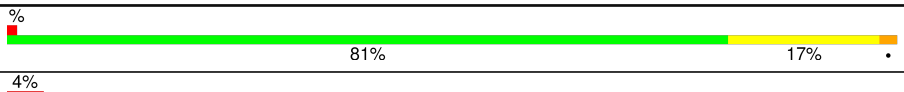
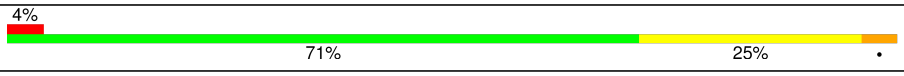
The reported resolution of this entry is 3.00 Å.

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	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

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 >=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 <=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	728	
2	B	204	

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	NAG	A	812	-	-	-	X
4	NAG	A	805	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7705 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.

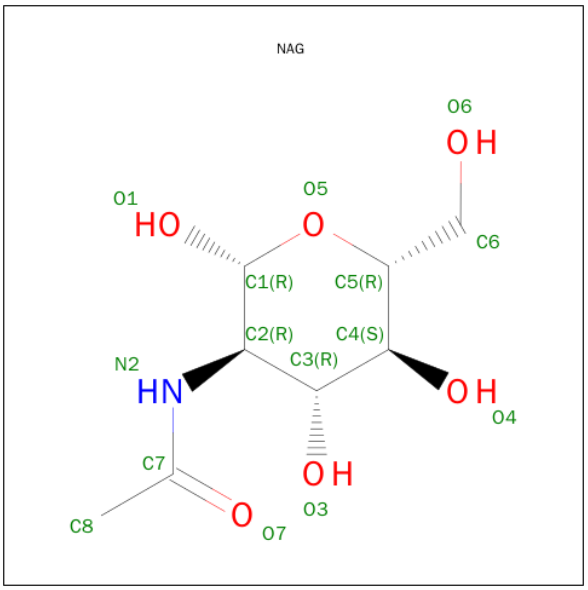
- Molecule 1 is a protein called Dipeptidyl peptidase 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	728	Total	C	N	O	S	0	0	0
			5963	3827	982	1128	26			

- Molecule 2 is a protein called MERS-CoV RBD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	204	Total	C	N	O	S	0	0	0
			1577	1004	251	311	11			

- Molecule 3 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	2	Total	C	N	O	0	0
			28	16	2	10		
4	A	2	Total	C	N	O	0	0
			28	16	2	10		

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

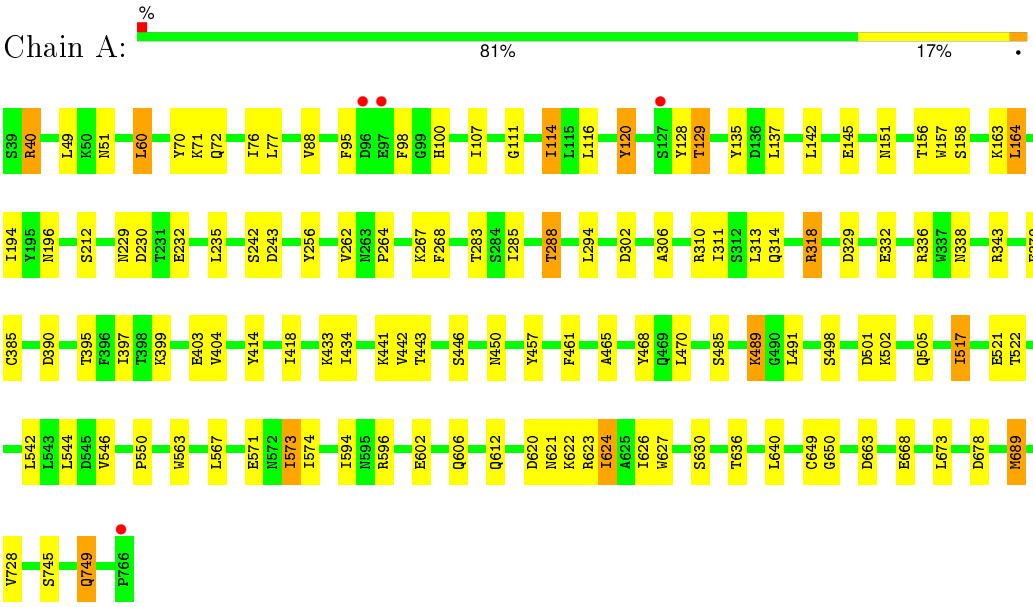
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	3	Total	C	N	O	0	0
			39	22	2	15		

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.

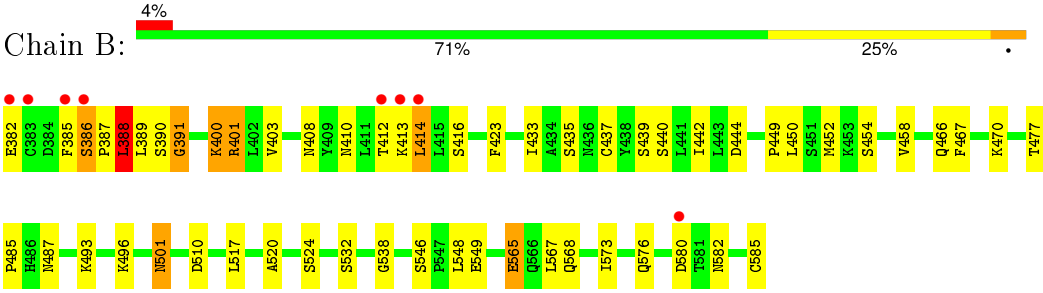
- Molecule 1: Dipeptidyl peptidase 4

Chain A:



- Molecule 2: MERS-CoV RBD

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α, β, γ	110.61Å 110.61Å 527.71Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.90 – 3.00 48.99 – 3.01	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.90-3.00) 99.9 (48.99-3.01)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	8.23 (at 3.01Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.205 , 0.253 0.214 , 0.257	Depositor DCC
R_{free} test set	1979 reflections (5.28%)	DCC
Wilson B-factor (Å^2)	46.3	Xtriage
Anisotropy	0.368	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 24.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Outliers	0 of 39452 reflections	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	7705	wwPDB-VP
Average B, all atoms (Å^2)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.26% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: BMA, NAG

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	0.49	0/6135	0.62	1/8344 (0.0%)
2	B	0.44	0/1615	0.64	1/2205 (0.0%)
All	All	0.48	0/7750	0.63	2/10549 (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
2	B	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	388	LEU	CA-CB-CG	6.45	130.12	115.30
1	A	60	LEU	CA-CB-CG	6.22	129.62	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	386	SER	Peptide

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen

atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5963	0	5676	68	0
2	B	1577	0	1534	31	0
3	A	70	0	65	0	0
4	A	56	0	50	0	0
5	A	39	0	34	4	0
All	All	7705	0	7359	97	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (97) 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:596:ARG:NH2	1:A:678:ASP:OD1	2.12	0.82
1:A:517:ILE:HD12	1:A:612:GLN:HG2	1.70	0.73
1:A:621:ASN:HA	1:A:624:ILE:HD11	1.69	0.72
1:A:329:ASP:OD1	1:A:343:ARG:NH1	2.25	0.69
1:A:640:LEU:HD11	1:A:650:GLY:HA3	1.75	0.68
1:A:302:ASP:HB3	1:A:314:GLN:HB2	1.78	0.64
2:B:493:LYS:NZ	2:B:565:GLU:O	2.31	0.64
1:A:397:ILE:HD12	1:A:434:ILE:HD13	1.81	0.63
1:A:267:LYS:HD3	2:B:538:GLY:HA3	1.82	0.62
2:B:400:LYS:HE2	2:B:401:ARG:H	1.66	0.60
1:A:242:SER:OG	1:A:243:ASP:N	2.27	0.60
1:A:384:ILE:HG13	1:A:404:VAL:HG21	1.84	0.59
2:B:410:ASN:HB3	2:B:413:LYS:HB2	1.85	0.57
1:A:158:SER:HB3	1:A:163:LYS:HB2	1.87	0.56
2:B:386:SER:OG	2:B:388:LEU:HB3	2.06	0.55
1:A:288:THR:HG21	2:B:501:ASN:ND2	2.21	0.55
1:A:49:LEU:HD22	1:A:749:GLN:HA	1.89	0.55
1:A:310:ARG:NH1	1:A:329:ASP:OD2	2.40	0.54
2:B:390:SER:OG	2:B:391:GLY:N	2.38	0.54
2:B:388:LEU:HD23	2:B:389:LEU:HD13	1.90	0.53
1:A:465:ALA:O	1:A:485:SER:OG	2.22	0.53
1:A:620:ASP:OD2	1:A:623:ARG:HD3	2.08	0.53
1:A:229:ASN:HD22	5:A:806:NAG:H83	1.74	0.53
1:A:40:ARG:H	1:A:40:ARG:HE	1.56	0.53
1:A:267:LYS:NZ	5:A:806:NAG:H81	2.25	0.52
1:A:535:ASP:HB3	1:A:538:LYS:HG3	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:386:SER:HG	2:B:388:LEU:HB3	1.74	0.51
1:A:567:LEU:HB3	1:A:573:ILE:HG23	1.91	0.51
1:A:626:ILE:HG23	1:A:636:THR:HG23	1.94	0.50
2:B:408:ASN:HB2	2:B:585:CYS:O	2.11	0.50
2:B:450:LEU:HD23	2:B:568:GLN:NE2	2.27	0.49
1:A:188:THR:HG23	1:A:194:ILE:HG21	1.92	0.49
1:A:98:PHE:CD2	1:A:100:HIS:HB2	2.46	0.49
1:A:726:VAL:HG23	1:A:728:VAL:HG23	1.95	0.49
1:A:318:ARG:HD3	1:A:668:GLU:OE1	2.13	0.49
2:B:483:THR:O	2:B:485:PRO:HD3	2.13	0.49
2:B:440:SER:HB3	2:B:576:GLN:HB3	1.94	0.49
1:A:95:PHE:CE1	1:A:116:LEU:HD11	2.48	0.48
2:B:401:ARG:HD2	2:B:442:ILE:HG23	1.93	0.48
2:B:386:SER:HB3	2:B:388:LEU:HD22	1.94	0.48
2:B:385:PHE:HB2	2:B:386:SER:HB2	1.95	0.48
1:A:76:ILE:HD13	1:A:107:ILE:HD11	1.96	0.48
1:A:77:LEU:HD23	1:A:88:VAL:HA	1.97	0.47
1:A:72:GLN:HB3	1:A:77:LEU:HD12	1.97	0.47
2:B:470:LYS:O	2:B:520:ALA:HA	2.15	0.47
2:B:466:GLN:HG2	2:B:517:LEU:HD22	1.97	0.47
1:A:135:TYR:HD1	1:A:142:LEU:HD12	1.79	0.46
1:A:649:CYS:HB3	1:A:699:GLU:HB2	1.98	0.46
1:A:571:GLU:HB2	1:A:573:ILE:HG22	1.96	0.46
1:A:192:ASP:O	1:A:193:ILE:HD13	2.16	0.46
1:A:550:PRO:HG3	1:A:594:ILE:HD11	1.98	0.46
1:A:188:THR:HG22	1:A:196:ASN:OD1	2.16	0.46
1:A:129:THR:HG23	1:A:151:ASN:HA	1.97	0.46
1:A:546:VAL:HG12	1:A:627:TRP:O	2.16	0.46
2:B:450:LEU:HD13	2:B:450:LEU:HA	1.84	0.46
1:A:382:ARG:H	1:A:403:GLU:HG2	1.81	0.46
2:B:501:ASN:C	2:B:501:ASN:HD22	2.19	0.45
1:A:745:SER:O	1:A:749:GLN:HG2	2.16	0.45
1:A:114:ILE:HG23	1:A:135:TYR:HB3	1.98	0.45
2:B:410:ASN:HB3	2:B:413:LYS:CB	2.46	0.45
2:B:478:CYS:HB2	2:B:573:ILE:HB	1.99	0.45
2:B:467:PHE:O	2:B:524:SER:HB2	2.16	0.45
1:A:414:TYR:CD1	1:A:433:LYS:HD2	2.51	0.44
1:A:157:TRP:CE3	1:A:164:LEU:HD13	2.52	0.44
2:B:401:ARG:HD3	2:B:444:ASP:OD1	2.17	0.44
2:B:546:SER:OG	2:B:549:GLU:HB2	2.17	0.44
1:A:544:LEU:HD12	1:A:544:LEU:HA	1.74	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:689:MET:HB3	1:A:689:MET:HE2	1.75	0.43
2:B:385:PHE:HA	2:B:386:SER:HA	1.55	0.43
1:A:158:SER:CB	1:A:163:LYS:HB2	2.49	0.43
2:B:403:VAL:HG13	2:B:442:ILE:HD11	2.00	0.42
1:A:542:LEU:HD12	1:A:574:ILE:O	2.18	0.42
2:B:386:SER:CB	2:B:388:LEU:H	2.33	0.42
1:A:283:THR:HG22	1:A:285:ILE:HD12	2.00	0.42
2:B:401:ARG:NH1	2:B:444:ASP:OD1	2.49	0.42
1:A:379:GLU:OE2	1:A:379:GLU:N	2.53	0.42
1:A:306:ALA:HB3	1:A:310:ARG:HB3	2.00	0.42
1:A:70:TYR:CG	1:A:71:LYS:N	2.87	0.42
1:A:446:SER:HB2	1:A:457:TYR:CE2	2.54	0.42
1:A:267:LYS:HZ2	5:A:806:NAG:H81	1.84	0.42
1:A:622:LYS:O	1:A:623:ARG:HG3	2.20	0.42
1:A:268:PHE:CE2	1:A:313:LEU:HD21	2.55	0.42
1:A:544:LEU:HD21	1:A:606:GLN:HG3	2.02	0.42
1:A:120:TYR:CE1	1:A:128:TYR:CD2	3.08	0.41
1:A:489:LYS:HD3	1:A:489:LYS:HA	1.61	0.41
1:A:256:TYR:CZ	1:A:663:ASP:HB3	2.56	0.41
1:A:229:ASN:ND2	5:A:806:NAG:H83	2.35	0.41
1:A:230:ASP:OD1	1:A:264:PRO:HB3	2.20	0.41
1:A:461:PHE:CD1	1:A:468:TYR:HB3	2.55	0.41
1:A:111:GLY:O	1:A:137:LEU:HD12	2.20	0.41
2:B:388:LEU:HD22	2:B:414:LEU:HD21	2.03	0.41
2:B:449:PRO:HG2	2:B:452:MET:HE2	2.02	0.41
1:A:294:LEU:HD23	1:A:294:LEU:HA	1.78	0.41
1:A:120:TYR:CD2	1:A:120:TYR:C	2.94	0.40
1:A:501:ASP:O	1:A:505:GLN:HG3	2.22	0.40
1:A:232:GLU:HB2	1:A:262:VAL:HG11	2.04	0.40
1:A:602:GLU:OE1	1:A:602:GLU:N	2.46	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	726/728 (100%)	676 (93%)	49 (7%)	1 (0%)	56 90
2	B	202/204 (99%)	179 (89%)	16 (8%)	7 (4%)	4 24
All	All	928/932 (100%)	855 (92%)	65 (7%)	8 (1%)	21 64

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	582	ASN
2	B	414	LEU
2	B	580	ASP
1	A	536	LYS
2	B	567	LEU
2	B	487	ASN
2	B	387	PRO
2	B	391	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	653/653 (100%)	610 (93%)	43 (7%)	21 57
2	B	186/186 (100%)	164 (88%)	22 (12%)	6 26
All	All	839/839 (100%)	774 (92%)	65 (8%)	16 50

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	ARG
1	A	51	ASN
1	A	60	LEU
1	A	114	ILE
1	A	120	TYR

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Mol	Chain	Res	Type
1	A	129	THR
1	A	145	GLU
1	A	156	THR
1	A	164	LEU
1	A	188	THR
1	A	212	SER
1	A	235	LEU
1	A	288	THR
1	A	311	ILE
1	A	318	ARG
1	A	332	GLU
1	A	336	ARG
1	A	338	ASN
1	A	385	CYS
1	A	390	ASP
1	A	395	THR
1	A	399	LYS
1	A	418	ILE
1	A	441	LYS
1	A	442	VAL
1	A	443	THR
1	A	450	ASN
1	A	470	LEU
1	A	489	LYS
1	A	491	LEU
1	A	498	SER
1	A	502	LYS
1	A	517	ILE
1	A	521	GLU
1	A	522	THR
1	A	537	SER
1	A	563	TRP
1	A	573	ILE
1	A	624	ILE
1	A	630	SER
1	A	673	LEU
1	A	689	MET
1	A	749	GLN
2	B	382	GLU
2	B	388	LEU
2	B	400	LYS
2	B	401	ARG

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Mol	Chain	Res	Type
2	B	412	THR
2	B	416	SER
2	B	423	PHE
2	B	433	ILE
2	B	435	SER
2	B	437	CYS
2	B	439	SER
2	B	454	SER
2	B	458	VAL
2	B	477	THR
2	B	483	THR
2	B	484	VAL
2	B	496	LYS
2	B	501	ASN
2	B	510	ASP
2	B	532	SER
2	B	548	LEU
2	B	565	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	501	ASN

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)

7 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	$\#\|Z\| > 2$
4	NAG	A	804	1,4	14,14,15	0.57	0	15,19,21	1.40	3 (20%)
4	NAG	A	805	4	14,14,15	0.60	0	15,19,21	1.47	2 (13%)
5	NAG	A	806	1,5	14,14,15	0.53	0	15,19,21	1.33	1 (6%)
5	NAG	A	807	5	14,14,15	0.62	0	15,19,21	1.18	0
5	BMA	A	808	5	11,11,12	1.77	2 (18%)	14,15,17	1.78	4 (28%)
4	NAG	A	809	4	14,14,15	0.51	0	15,19,21	1.33	1 (6%)
4	NAG	A	810	1,4	14,14,15	0.49	0	15,19,21	2.37	3 (20%)

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
4	NAG	A	804	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	805	4	-	0/6/23/26	0/1/1/1
5	NAG	A	806	1,5	-	0/6/23/26	0/1/1/1
5	NAG	A	807	5	-	0/6/23/26	0/1/1/1
5	BMA	A	808	5	-	0/2/19/22	0/1/1/1
4	NAG	A	809	4	-	0/6/23/26	0/1/1/1
4	NAG	A	810	1,4	-	0/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	808	BMA	C4-C5	2.93	1.59	1.53
5	A	808	BMA	C1-C2	3.19	1.59	1.52

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	809	NAG	C2-N2-C7	-3.73	118.24	123.04
5	A	808	BMA	C2-C3-C4	-3.57	104.97	111.04
4	A	804	NAG	C1-O5-C5	-2.91	108.55	112.25
4	A	804	NAG	C2-N2-C7	-2.32	120.06	123.04
5	A	808	BMA	O2-C2-C3	-2.18	105.73	110.12
5	A	808	BMA	O3-C3-C2	2.14	113.87	110.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	805	NAG	O4-C4-C3	2.30	115.51	110.34
4	A	804	NAG	C4-C3-C2	2.68	115.40	111.23
4	A	810	NAG	O6-C6-C5	2.75	120.41	111.33
4	A	810	NAG	O5-C5-C6	3.04	113.93	107.35
5	A	808	BMA	C1-O5-C5	3.06	116.13	112.25
4	A	805	NAG	C1-O5-C5	3.19	116.29	112.25
5	A	806	NAG	C1-O5-C5	3.20	116.31	112.25
4	A	810	NAG	C1-O5-C5	7.30	121.51	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	806	NAG	4	0

5.6 Ligand geometry (i)

5 ligands 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	NAG	A	801	1	14,14,15	0.43	0	15,19,21	1.34	1 (6%)
3	NAG	A	802	1	14,14,15	0.76	0	15,19,21	1.26	1 (6%)
3	NAG	A	803	1	14,14,15	0.69	0	15,19,21	1.49	2 (13%)
3	NAG	A	811	1	14,14,15	0.89	0	15,19,21	1.49	2 (13%)
3	NAG	A	812	1	14,14,15	0.58	0	15,19,21	2.00	2 (13%)

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	NAG	A	801	1	-	0/6/23/26	0/1/1/1
3	NAG	A	802	1	-	0/6/23/26	0/1/1/1
3	NAG	A	803	1	-	0/6/23/26	0/1/1/1
3	NAG	A	811	1	-	2/6/23/26	0/1/1/1
3	NAG	A	812	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	801	NAG	C2-N2-C7	-4.15	117.71	123.04
3	A	811	NAG	C4-C3-C2	-3.23	106.20	111.23
3	A	812	NAG	C2-N2-C7	-2.82	119.42	123.04
3	A	811	NAG	O3-C3-C2	2.13	113.33	109.11
3	A	803	NAG	C4-C3-C2	3.08	116.01	111.23
3	A	803	NAG	C1-O5-C5	3.32	116.46	112.25
3	A	802	NAG	C1-O5-C5	3.76	117.02	112.25
3	A	812	NAG	C1-O5-C5	6.33	120.28	112.25

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	811	NAG	O7-C7-N2-C2
3	A	811	NAG	C8-C7-N2-C2

There are no ring outliers.

No monomer is involved in short contacts.

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^{th} 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^2)	Q<0.9
1	A	728/728 (100%)	-0.48	4 (0%) 91 76	7, 18, 44, 76	0
2	B	204/204 (100%)	-0.09	8 (3%) 43 18	16, 37, 63, 79	0
All	All	932/932 (100%)	-0.40	12 (1%) 79 53	7, 20, 54, 79	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	127	SER	3.1
2	B	414	LEU	3.0
2	B	413	LYS	2.9
2	B	385	PHE	2.8
2	B	386	SER	2.7
1	A	96	ASP	2.5
1	A	97	GLU	2.5
2	B	580	ASP	2.4
1	A	766	PRO	2.4
2	B	383	CYS	2.2
2	B	412	THR	2.1
2	B	382	GLU	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

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^{th} 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^2)	Q<0.9
4	NAG	A	805	14/15	0.87	0.26	2.75	15,15,15,15	0
5	NAG	A	806	14/15	0.97	0.12	-1.46	15,15,15,15	0
4	NAG	A	810	14/15	0.89	0.13	-	15,15,15,15	0
5	NAG	A	807	14/15	0.93	0.22	-	15,15,15,15	0
5	BMA	A	808	11/12	0.75	0.29	-	15,15,15,15	0
4	NAG	A	809	14/15	0.92	0.26	-	15,15,15,15	0
4	NAG	A	804	14/15	0.93	0.20	-	15,15,15,15	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^{th} 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^2)	Q<0.9
3	NAG	A	812	14/15	0.81	0.44	5.11	15,15,15,15	0
3	NAG	A	811	14/15	0.88	0.19	1.17	15,15,15,15	0
3	NAG	A	801	14/15	0.86	0.26	-0.06	15,15,15,15	0
3	NAG	A	803	14/15	0.81	0.32	-	15,15,15,15	0
3	NAG	A	802	14/15	0.82	0.39	-	15,15,15,15	0

6.5 Other polymers (i)

There are no such residues in this entry.