



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:07 AM GMT

PDB ID : 3KWF  
Title : human DPP-IV with carmegliptin (S)-1-((2S,3S,11bS)-2-Amino-9,10-dimethoxy-1,3,4,6,7,11b-hexahydro-2H-pyrido[2,1-a]isoquinolin-3-yl)-4-fluoromethyl-pyrrolidin-2-one  
Authors : Hennig, M.; Stihle, M.; Thoma, R.  
Deposited on : 2009-12-01  
Resolution : 2.40 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

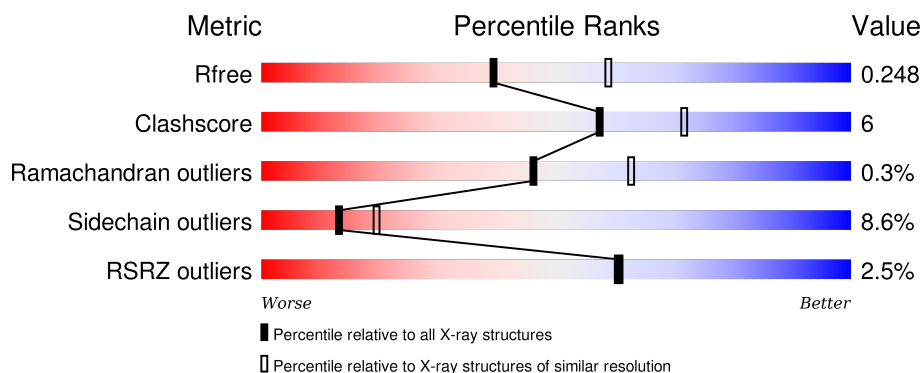
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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	728	<div> <div>2%</div> <div>82%</div> <div>15%</div> <div>.</div> </div>
1	B	728	<div> <div>3%</div> <div>82%</div> <div>15%</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	NAG	A	794	-	-	-	X

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 12464 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.

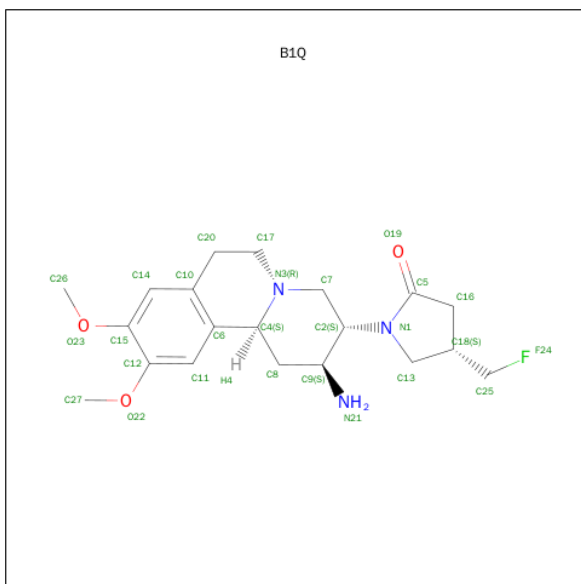
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	728	Total	C	N	O	S	0	0	0
			5963	3827	982	1128	26			
1	B	728	Total	C	N	O	S	0	0	0
			5963	3827	982	1128	26			

- Molecule 2 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).



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

- Molecule 3 is (4S)-1-[(2S,3S,11BS)-2-AMINO-9,10-DIMETHOXY-1,3,4,6,7,11B-HEXAHYDRO-2H-PYRIDO[2,1-A]ISOQUINOLIN-3-YL]-4-(FLUOROMETHYL)PYRROLIDIN-2-ONE (three-letter code: B1Q) (formula: C<sub>20</sub>H<sub>28</sub>FN<sub>3</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			27	20	1	3	3		
3	B	1	Total	C	F	N	O	0	0
			27	20	1	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	203	Total	O	0	0
			203	203		
4	B	225	Total	O	0	0
			225	225		





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	65.63Å 69.54Å 423.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.40 29.95 – 2.40	Depositor EDS
% Data completeness (in resolution range)	88.0 (15.00-2.40) 88.0 (29.95-2.40)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.52 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.200 , 0.246 0.202 , 0.248	Depositor DCC
$R_{free}$ test set	3403 reflections (5.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	38.9	Xtriage
Anisotropy	0.668	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 31.0	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 67948 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	12464	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.88% 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: NAG, B1Q

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	0.50	1/6135 (0.0%)	0.64	1/8344 (0.0%)
1	B	0.50	0/6135	0.64	1/8344 (0.0%)
All	All	0.50	1/12270 (0.0%)	0.64	2/16688 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	301	CYS	CB-SG	-6.59	1.71	1.82

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	60	LEU	CA-CB-CG	6.79	130.92	115.30
1	B	60	LEU	CA-CB-CG	6.54	130.35	115.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5963	0	5685	75	0
1	B	5963	0	5684	73	0
2	A	28	0	26	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	28	0	26	2	0
3	A	27	0	28	0	0
3	B	27	0	28	0	0
4	A	203	0	0	10	0
4	B	225	0	0	13	0
All	All	12464	0	11477	141	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 (141) 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:85:ASN:HD21	2:A:794:NAG:C1	1.24	1.49
1:B:229:ASN:HD21	2:B:796:NAG:C1	1.30	1.45
1:A:85:ASN:ND2	2:A:794:NAG:C1	1.93	1.29
1:A:229:ASN:HD21	2:A:796:NAG:C1	1.48	1.24
1:B:229:ASN:ND2	2:B:796:NAG:C1	2.15	1.10
1:B:600:THR:CG2	1:B:601:PHE:H	1.66	1.09
1:A:237:GLU:HG2	1:A:253:ARG:HG2	1.46	0.97
1:B:600:THR:HG23	1:B:601:PHE:H	1.29	0.96
1:A:229:ASN:ND2	2:A:796:NAG:C1	2.28	0.95
1:B:65:ASP:HB2	4:B:975:HOH:O	1.69	0.92
1:A:630:SER:HB3	1:A:740:HIS:NE2	1.87	0.90
1:B:600:THR:HG23	1:B:601:PHE:N	1.87	0.90
1:A:466:LYS:HG2	4:A:944:HOH:O	1.79	0.81
1:B:600:THR:HG22	1:B:601:PHE:H	1.50	0.77
1:B:56:LYS:HE3	4:B:767:HOH:O	1.83	0.77
1:B:403:GLU:H	1:B:420:ASN:HD21	1.32	0.77
1:A:630:SER:HB3	1:A:740:HIS:CE1	2.22	0.74
1:A:431:LEU:HD13	1:A:445:LEU:HD12	1.69	0.72
1:A:44:THR:HG22	1:A:46:THR:H	1.55	0.72
1:B:726:VAL:HG12	1:B:728:VAL:HG23	1.73	0.70
1:A:736:THR:HG21	1:B:717:ALA:O	1.93	0.68
1:A:39:SER:N	4:A:871:HOH:O	2.26	0.67
1:B:237:GLU:HG2	1:B:253:ARG:HG2	1.77	0.67
1:A:403:GLU:H	1:A:420:ASN:HD21	1.42	0.66
1:B:466:LYS:HB2	4:B:975:HOH:O	1.96	0.66
1:B:735:TYR:OH	1:B:750:HIS:HD2	1.79	0.64
1:A:693:GLU:O	1:A:696:LYS:HG3	1.97	0.64
1:A:450:ASN:OD1	4:A:25:HOH:O	2.14	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:640:LEU:HD11	1:B:650:GLY:HA3	1.80	0.63
1:B:510:PRO:HD3	1:B:569:SER:HB2	1.79	0.63
1:B:657:SER:H	1:B:715:GLN:NE2	1.97	0.63
1:B:54:ARG:HD2	4:B:767:HOH:O	1.98	0.63
1:A:726:VAL:HG12	1:A:728:VAL:HG23	1.79	0.62
1:B:600:THR:CG2	1:B:601:PHE:N	2.35	0.62
1:B:45:LEU:HG	1:B:49:LEU:HD22	1.83	0.61
1:B:126:HIS:HD2	4:B:24:HOH:O	1.83	0.61
1:A:44:THR:HG22	1:A:46:THR:N	2.16	0.60
1:B:327:ILE:HD13	1:B:389:ILE:HG12	1.84	0.60
1:B:598:LEU:O	1:B:682:HIS:HE1	1.85	0.59
1:A:85:ASN:ND2	2:A:794:NAG:O5	2.35	0.59
1:B:42:THR:HG23	1:B:570:THR:OG1	2.03	0.59
1:B:42:THR:HG22	4:B:780:HOH:O	2.03	0.57
1:B:134:ILE:HB	1:B:143:ILE:HG12	1.85	0.57
1:A:640:LEU:HD11	1:A:650:GLY:HA3	1.85	0.56
1:A:657:SER:H	1:A:715:GLN:NE2	2.03	0.56
1:B:145:GLU:HG3	4:B:858:HOH:O	2.05	0.56
1:B:224:ALA:HB1	1:B:268:PHE:CZ	2.41	0.56
1:A:718:GLN:OE1	1:A:721:LYS:NZ	2.39	0.56
1:A:253:ARG:HH12	1:B:253:ARG:NH1	2.04	0.55
1:B:302:ASP:HB3	1:B:314:GLN:HB2	1.89	0.55
1:B:403:GLU:H	1:B:420:ASN:ND2	2.03	0.54
1:A:253:ARG:NH1	1:B:253:ARG:HH12	2.06	0.53
1:A:481:THR:OG1	1:A:483:HIS:HE1	1.91	0.53
1:B:340:LEU:HB2	1:B:343:ARG:HD2	1.90	0.53
1:B:377:ASN:HD22	1:B:377:ASN:C	2.11	0.53
1:A:295:ILE:HG23	4:A:834:HOH:O	2.07	0.53
1:B:562:ASN:HD22	1:B:562:ASN:C	2.11	0.53
1:A:499:ALA:O	1:A:502:LYS:HG3	2.08	0.53
1:A:65:ASP:HB2	4:A:944:HOH:O	2.10	0.52
1:A:229:ASN:ND2	2:A:796:NAG:O5	2.37	0.52
1:B:486:VAL:HG23	4:B:911:HOH:O	2.09	0.52
1:A:750:HIS:HE1	1:B:728:VAL:O	1.93	0.51
1:B:105:TYR:HB2	1:B:114:ILE:HD11	1.91	0.51
1:A:305:TRP:CZ3	1:A:311:ILE:HG12	2.45	0.51
1:A:114:ILE:HG23	1:A:135:TYR:HB3	1.92	0.51
1:A:139:LYS:O	1:A:140:ARG:HG3	2.11	0.50
1:A:105:TYR:HB2	1:A:114:ILE:HD11	1.94	0.50
1:A:500:LEU:HD22	1:A:504:LEU:HG	1.94	0.50
1:B:563:TRP:CH2	1:B:759:ILE:HD12	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:196:ASN:OD1	1:A:227:GLN:HG3	2.11	0.50
1:A:327:ILE:HD13	1:A:389:ILE:HG12	1.93	0.50
1:A:401:THR:O	1:A:401:THR:HG22	2.12	0.50
1:A:289:ALA:HB3	1:A:294:LEU:HD13	1.93	0.50
1:B:401:THR:HG22	1:B:401:THR:O	2.12	0.49
1:A:44:THR:CG2	4:A:942:HOH:O	2.61	0.49
1:B:177:GLU:HB2	1:B:180:LEU:HD22	1.95	0.49
1:A:502:LYS:HD3	4:A:900:HOH:O	2.13	0.48
1:B:306:ALA:HB3	1:B:310:ARG:HG2	1.96	0.48
1:B:154:TRP:CE2	1:B:212:SER:HB2	2.49	0.48
1:A:377:ASN:C	1:A:377:ASN:HD22	2.16	0.48
1:B:614:SER:HA	1:B:619:VAL:CG2	2.43	0.47
1:B:598:LEU:O	1:B:682:HIS:CE1	2.67	0.47
1:A:470:LEU:HD12	1:A:483:HIS:CE1	2.50	0.47
1:B:414:TYR:CE1	1:B:433:LYS:HE2	2.50	0.47
1:A:82:GLU:HB2	1:A:467:TYR:OH	2.15	0.46
1:A:66:HIS:HB2	4:A:828:HOH:O	2.13	0.46
1:B:657:SER:H	1:B:715:GLN:HE21	1.63	0.46
1:B:133:ASP:OD1	1:B:147:ARG:NH1	2.48	0.46
1:A:626:ILE:O	1:A:650:GLY:HA2	2.16	0.46
1:A:377:ASN:ND2	1:A:381:TYR:H	2.14	0.46
1:B:504:LEU:HA	1:B:507:VAL:HG13	1.98	0.45
1:B:221:THR:O	1:B:273:THR:HB	2.17	0.45
1:A:45:LEU:HG	1:A:49:LEU:HD22	1.99	0.45
1:B:543:LEU:HD12	1:B:567:LEU:HD13	1.99	0.45
1:B:756:SER:O	1:B:760:LYS:HG3	2.16	0.45
1:A:519:LEU:HA	1:A:519:LEU:HD23	1.75	0.45
1:A:44:THR:HG23	4:A:942:HOH:O	2.17	0.44
1:A:420:ASN:HD22	1:A:420:ASN:C	2.21	0.44
1:A:728:VAL:O	1:B:750:HIS:HE1	2.01	0.44
1:B:595:ASN:O	1:B:597:ARG:NE	2.49	0.44
1:B:649:CYS:HB3	1:B:699:GLU:HB2	2.00	0.44
1:A:403:GLU:H	1:A:420:ASN:ND2	2.11	0.44
1:A:351:THR:OG1	1:A:592:HIS:HD2	2.01	0.44
1:A:55:LEU:HD22	1:A:478:PRO:HG2	2.00	0.44
1:A:360:SER:O	1:A:373:LYS:NZ	2.48	0.44
1:A:143:ILE:HD13	1:A:178:PRO:HB2	2.00	0.44
1:B:242:SER:HB3	1:B:246:LEU:HD12	2.00	0.43
1:A:60:LEU:HB2	1:A:68:TYR:CD1	2.53	0.43
1:A:657:SER:H	1:A:715:GLN:HE21	1.66	0.43
1:B:144:THR:HB	4:B:858:HOH:O	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:364:PHE:HE1	1:A:389:ILE:HD11	1.84	0.43
1:B:236:ILE:HG13	1:B:712:HIS:CE1	2.54	0.43
1:A:594:ILE:HD11	1:A:602:GLU:OE1	2.19	0.43
1:A:206:GLU:OE2	1:A:663:ASP:OD2	2.35	0.43
1:B:402:TRP:CD1	1:B:421:GLU:HG3	2.54	0.43
1:B:40:ARG:NH1	4:B:978:HOH:O	2.52	0.43
1:B:42:THR:CG2	1:B:570:THR:OG1	2.66	0.42
1:A:177:GLU:HB2	1:A:180:LEU:HD22	1.99	0.42
1:B:49:LEU:HG	1:B:749:GLN:HG2	2.01	0.42
1:B:219:ASN:ND2	4:B:812:HOH:O	2.51	0.42
1:A:156:THR:HG23	1:A:216:TRP:HE1	1.84	0.42
1:A:735:TYR:OH	1:A:750:HIS:HD2	2.01	0.42
1:A:463:LYS:C	1:A:465:ALA:H	2.23	0.42
1:A:253:ARG:HH12	1:B:253:ARG:HH12	1.60	0.42
1:A:554:LYS:HB3	1:A:577:SER:HB3	2.01	0.42
1:A:301:CYS:SG	1:A:316:LEU:HB2	2.60	0.41
1:A:152:THR:HG21	1:A:155:VAL:HG22	2.02	0.41
1:B:152:THR:HG21	1:B:155:VAL:HG22	2.02	0.41
1:A:592:HIS:HE1	4:A:865:HOH:O	2.03	0.41
1:A:397:ILE:HG22	1:A:439:TYR:CE2	2.55	0.41
1:B:597:ARG:HA	1:B:682:HIS:CE1	2.55	0.41
1:B:127:SER:HB3	1:B:211:TYR:CD1	2.55	0.41
1:B:626:ILE:O	1:B:650:GLY:HA2	2.20	0.41
1:B:48:TYR:CE1	1:B:562:ASN:HA	2.56	0.41
1:A:302:ASP:HB3	1:A:314:GLN:HB2	2.03	0.41
1:A:258:LYS:HD2	1:B:247:GLN:HG2	2.03	0.41
1:A:271:VAL:HG22	1:A:284:SER:HA	2.02	0.41
1:A:381:TYR:CZ	1:A:401:THR:HG23	2.56	0.40
1:B:693:GLU:HG2	4:B:905:HOH:O	2.21	0.40
1:B:433:LYS:HE3	4:B:28:HOH:O	2.21	0.40
1:B:114:ILE:HG23	1:B:135:TYR: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	726/728 (100%)	689 (95%)	34 (5%)	3 (0%)	39	56
1	B	726/728 (100%)	688 (95%)	36 (5%)	2 (0%)	46	63
All	All	1452/1456 (100%)	1377 (95%)	70 (5%)	5 (0%)	46	63

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	74	ASN
1	A	139	LYS
1	A	40	ARG
1	A	630	SER
1	B	109	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	653/653 (100%)	596 (91%)	57 (9%)	13	19
1	B	653/653 (100%)	598 (92%)	55 (8%)	14	20
All	All	1306/1306 (100%)	1194 (91%)	112 (9%)	13	19

All (112) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	SER
1	A	40	ARG
1	A	42	THR
1	A	49	LEU
1	A	51	ASN
1	A	60	LEU
1	A	74	ASN
1	A	87	SER

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Mol	Chain	Res	Type
1	A	88	VAL
1	A	90	LEU
1	A	145	GLU
1	A	180	LEU
1	A	202	VAL
1	A	223	LEU
1	A	230	ASP
1	A	246	LEU
1	A	280	THR
1	A	294	LEU
1	A	300	LEU
1	A	311	ILE
1	A	312	SER
1	A	316	LEU
1	A	326	ASP
1	A	343	ARG
1	A	358	ARG
1	A	361	GLU
1	A	366	LEU
1	A	377	ASN
1	A	379	GLU
1	A	385	CYS
1	A	389	ILE
1	A	410	LEU
1	A	420	ASN
1	A	431	LEU
1	A	452	GLU
1	A	463	LYS
1	A	466	LYS
1	A	472	CYS
1	A	479	LEU
1	A	482	LEU
1	A	500	LEU
1	A	502	LYS
1	A	505	GLN
1	A	514	LEU
1	A	543	LEU
1	A	544	LEU
1	A	566	TYR
1	A	594	ILE
1	A	621	ASN
1	A	627	TRP

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Mol	Chain	Res	Type
1	A	660	GLU
1	A	684	ARG
1	A	696	LYS
1	A	704	HIS
1	A	715	GLN
1	A	726	VAL
1	A	736	THR
1	B	40	ARG
1	B	42	THR
1	B	49	LEU
1	B	56	LYS
1	B	60	LEU
1	B	63	ILE
1	B	74	ASN
1	B	88	VAL
1	B	94	THR
1	B	110	ASP
1	B	129	THR
1	B	140	ARG
1	B	156	THR
1	B	202	VAL
1	B	212	SER
1	B	214	LEU
1	B	223	LEU
1	B	230	ASP
1	B	236	ILE
1	B	244	GLU
1	B	246	LEU
1	B	295	ILE
1	B	300	LEU
1	B	301	CYS
1	B	311	ILE
1	B	326	ASP
1	B	336	ARG
1	B	343	ARG
1	B	375	ILE
1	B	377	ASN
1	B	385	CYS
1	B	389	ILE
1	B	392	LYS
1	B	420	ASN
1	B	472	CYS

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Mol	Chain	Res	Type
1	B	473	SER
1	B	479	LEU
1	B	489	LYS
1	B	500	LEU
1	B	507	VAL
1	B	514	LEU
1	B	515	ASP
1	B	543	LEU
1	B	558	VAL
1	B	562	ASN
1	B	566	TYR
1	B	600	THR
1	B	621	ASN
1	B	627	TRP
1	B	680	LEU
1	B	684	ARG
1	B	696	LYS
1	B	704	HIS
1	B	715	GLN
1	B	726	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	85	ASN
1	A	103	ASN
1	A	169	ASN
1	A	229	ASN
1	A	263	ASN
1	A	298	HIS
1	A	377	ASN
1	A	420	ASN
1	A	430	ASN
1	A	435	GLN
1	A	483	HIS
1	A	508	GLN
1	A	592	HIS
1	A	621	ASN
1	A	679	ASN
1	A	715	GLN
1	A	750	HIS
1	B	123	GLN

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Mol	Chain	Res	Type
1	B	126	HIS
1	B	169	ASN
1	B	229	ASN
1	B	263	ASN
1	B	298	HIS
1	B	377	ASN
1	B	383	HIS
1	B	420	ASN
1	B	483	HIS
1	B	562	ASN
1	B	621	ASN
1	B	679	ASN
1	B	682	HIS
1	B	715	GLN
1	B	750	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 [i](#)

6 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	B1Q	A	1	-	29,30,30	0.70	0	30,44,44	1.68	7 (23%)
2	NAG	A	794	-	14,14,15	0.53	0	15,19,21	0.89	1 (6%)
2	NAG	A	796	-	14,14,15	0.50	0	15,19,21	0.99	1 (6%)
3	B1Q	B	1	-	29,30,30	0.76	0	30,44,44	1.63	8 (26%)
2	NAG	B	794	1	14,14,15	0.56	0	15,19,21	0.84	0
2	NAG	B	796	-	14,14,15	0.52	0	15,19,21	1.34	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
3	B1Q	A	1	-	-	0/10/47/47	0/4/4/4
2	NAG	A	794	-	-	0/6/23/26	0/1/1/1
2	NAG	A	796	-	-	0/6/23/26	0/1/1/1
3	B1Q	B	1	-	-	0/10/47/47	0/4/4/4
2	NAG	B	794	1	-	0/6/23/26	0/1/1/1
2	NAG	B	796	-	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1	B1Q	O19-C5-C16	-2.72	123.76	127.41
2	B	796	NAG	C1-O5-C5	-2.67	108.86	112.25
3	A	1	B1Q	C7-C2-N1	-2.52	109.15	112.40
3	A	1	B1Q	O19-C5-C16	-2.47	124.09	127.41
2	B	796	NAG	C2-N2-C7	-2.26	120.14	123.04
3	B	1	B1Q	O22-C12-C11	-2.26	120.36	124.21
3	A	1	B1Q	O23-C15-C14	-2.22	120.42	124.21
3	B	1	B1Q	O19-C5-N1	2.04	127.44	124.82
2	A	796	NAG	C3-C4-C5	2.04	113.76	110.20
3	B	1	B1Q	C8-C4-N3	2.06	112.84	109.67
2	A	794	NAG	C2-N2-C7	2.09	125.72	123.04
3	A	1	B1Q	O19-C5-N1	2.20	127.64	124.82
3	B	1	B1Q	O22-C12-C15	2.66	119.19	115.40
3	A	1	B1Q	C26-O23-C15	2.86	121.89	117.54
2	B	796	NAG	C3-C4-C5	2.90	115.25	110.20
3	B	1	B1Q	C27-O22-C12	3.04	122.15	117.54
3	B	1	B1Q	O23-C15-C12	3.07	119.78	115.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1	B1Q	O23-C15-C12	3.50	120.39	115.40
3	B	1	B1Q	C26-O23-C15	4.06	123.70	117.54
3	A	1	B1Q	C27-O22-C12	4.50	124.37	117.54

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	794	NAG	3	0
2	A	796	NAG	3	0
2	B	796	NAG	2	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	728/728 (100%)	-0.12	14 (1%) 70 69	25, 39, 59, 73	0
1	B	728/728 (100%)	-0.08	22 (3%) 54 53	25, 39, 57, 72	0
All	All	1456/1456 (100%)	-0.10	36 (2%) 61 60	25, 39, 59, 73	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	766	PRO	6.3
1	A	279	VAL	5.4
1	B	279	VAL	5.1
1	A	278	SER	5.0
1	B	280	THR	3.7
1	B	140	ARG	3.6
1	B	73	GLU	3.4
1	B	392	LYS	3.4
1	B	97	GLU	3.3
1	A	333	SER	3.2
1	B	96	ASP	3.2
1	A	277	SER	3.0
1	A	105	TYR	3.0
1	B	702	LEU	2.9
1	A	766	PRO	2.9
1	A	39	SER	2.8
1	A	73	GLU	2.7
1	B	651	ILE	2.6
1	B	332	GLU	2.5
1	B	716	SER	2.5
1	B	703	ILE	2.4
1	A	97	GLU	2.4
1	B	331	ASP	2.4
1	B	110	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	165	ALA	2.3
1	B	378	GLU	2.3
1	B	652	ALA	2.3
1	A	378	GLU	2.3
1	B	138	ASN	2.2
1	B	301	CYS	2.2
1	A	138	ASN	2.2
1	B	701	LEU	2.2
1	A	521	GLU	2.1
1	A	280	THR	2.1
1	B	278	SER	2.1
1	B	334	SER	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 ⓘ

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
2	NAG	A	794	14/15	0.92	0.16	2.06	46,48,50,50	0
2	NAG	A	796	14/15	0.81	0.19	2.00	44,50,53,53	0
2	NAG	B	796	14/15	0.90	0.18	0.43	49,54,55,55	0
3	B1Q	A	1	27/27	0.94	0.12	-0.36	24,28,30,31	0
3	B1Q	B	1	27/27	0.96	0.13	-0.56	24,28,29,30	0
2	NAG	B	794	14/15	0.97	0.11	-1.15	41,43,46,46	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.