



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 18, 2016 – 06:14 AM EDT

PDB ID : 5JW1  
Title : Crystal structure of Celecoxib bound to S121P murine COX-2 mutant  
Authors : Malkowski, M.G.; Orlando, B.J.  
Deposited on : 2016-05-11  
Resolution : 2.82 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 : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20027939  
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) : rb-20027939

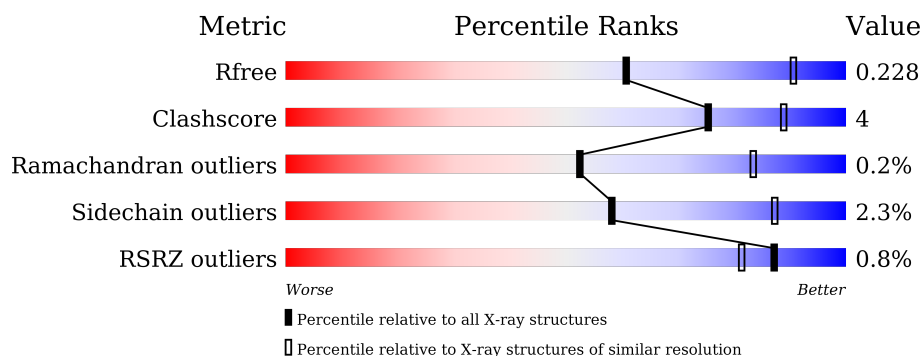
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.82 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	2676 (2.84-2.80)
Clashscore	102246	3124 (2.84-2.80)
Ramachandran outliers	100387	3072 (2.84-2.80)
Sidechain outliers	100360	3074 (2.84-2.80)
RSRZ outliers	91569	2690 (2.84-2.80)

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	552	<div> <div></div> <div>90%</div> <div>9%</div> <div>.</div> </div>
1	B	552	<div> <div></div> <div>89%</div> <div>10%</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	COH	A	601	-	-	-	X
2	COH	B	601	-	-	-	X
5	NAG	A	604	-	-	-	X
5	NAG	A	607	-	-	-	X
5	NAG	B	606	-	-	-	X



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.

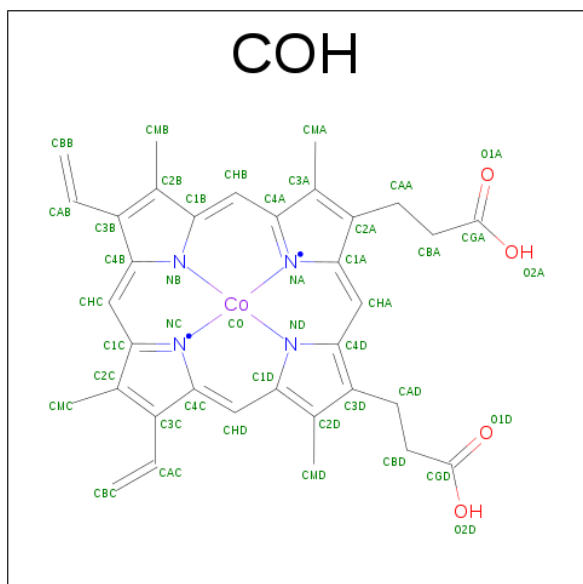
- Molecule 1 is a protein called Prostaglandin G/H synthase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	552	Total 4448	C 2876	N 738	O 809	S 25	0	1	0
1	B	551	Total 4435	C 2869	N 738	O 803	S 25	0	1	0

There are 6 discrepancies between the modelled and reference sequences:

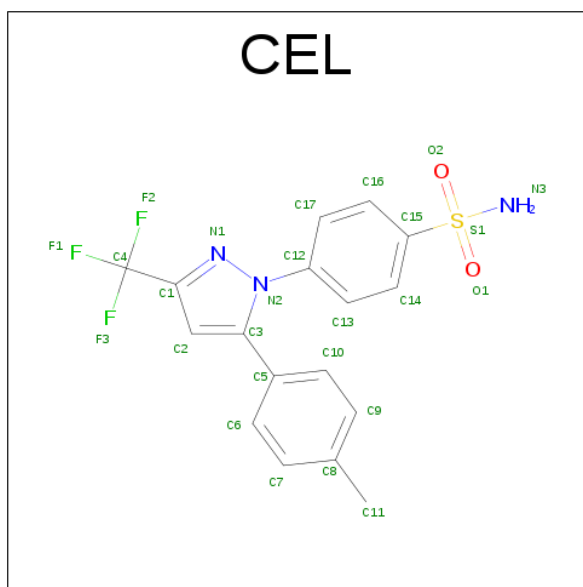
Chain	Residue	Modelled	Actual	Comment	Reference
A	33	HIS	-	expression tag	UNP Q05769
A	34	HIS	-	expression tag	UNP Q05769
A	122	PRO	SER	engineered mutation	UNP Q05769
B	33	HIS	-	expression tag	UNP Q05769
B	34	HIS	-	expression tag	UNP Q05769
B	122	PRO	SER	engineered mutation	UNP Q05769

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING CO (three-letter code: COH) (formula:  $C_{34}H_{32}CoN_4O_4$ ).



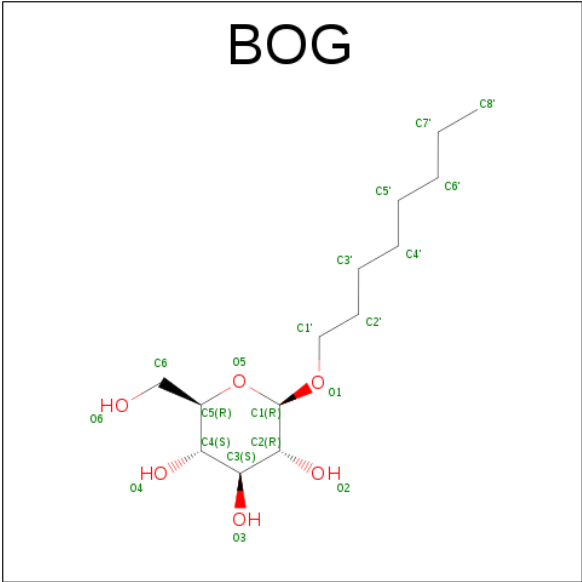
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Co	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Co	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is 4-[5-(4-METHYLPHENYL)-3-(TRIFLUOROMETHYL)-1H-PYRAZOL-1-YL]BENZENESULFONAMIDE (three-letter code: CEL) (formula: C<sub>17</sub>H<sub>14</sub>F<sub>3</sub>N<sub>3</sub>O<sub>2</sub>S).



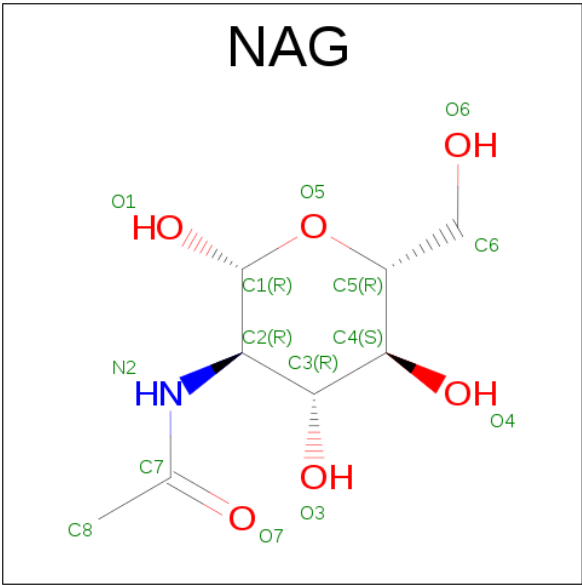
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total 26	C 17	F 3	N 3	O 2	S 1	0	0
3	B	1	Total 26	C 17	F 3	N 3	O 2	S 1	0	0

- Molecule 4 is B-OCTYLGLUCOSIDE (three-letter code: BOG) (formula: C<sub>14</sub>H<sub>28</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			20	14	6		

- Molecule 5 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



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

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

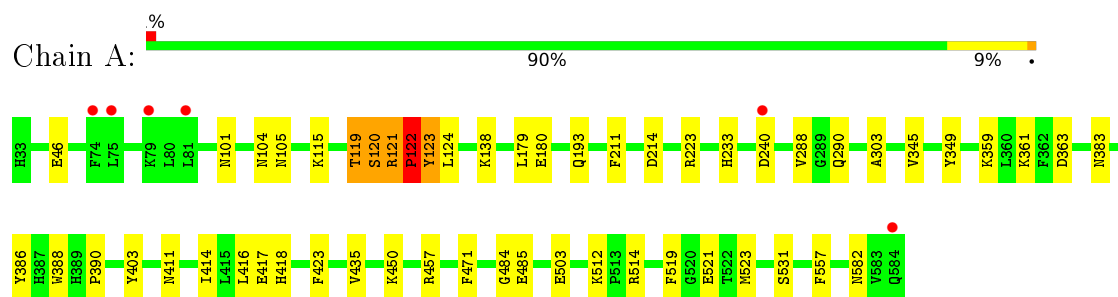
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	5	Total	O	0	0
			5	5		
6	B	10	Total	O	0	0
			10	10		

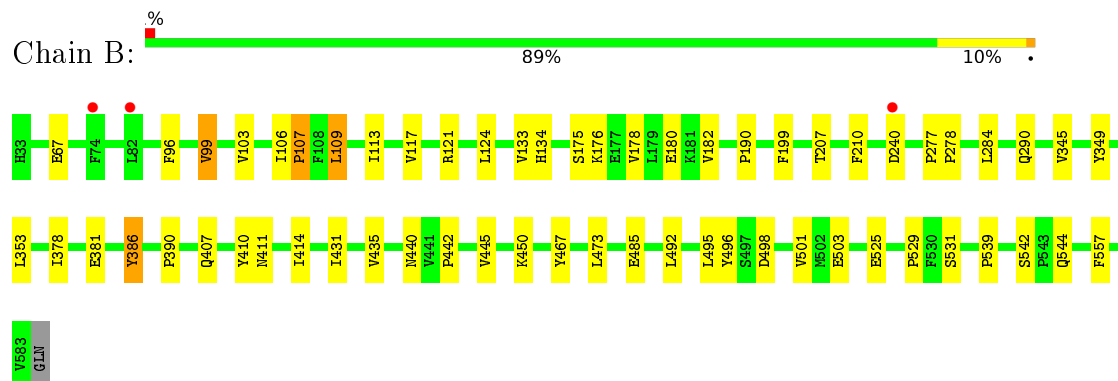
### 3 Residue-property plots

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 ( $\text{RSRZ} > 2$ ). 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: Prostaglandin G/H synthase 2



- Molecule 1: Prostaglandin G/H synthase 2





## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	120.36Å 132.20Å 180.39Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.82 30.09 – 2.82	Depositor EDS
% Data completeness (in resolution range)	93.1 (30.00-2.82) 89.1 (30.09-2.82)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.81 (at 2.80Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, $R_{free}$	0.189 , 0.226 0.197 , 0.228	Depositor DCC
$R_{free}$ test set	1726 reflections (5.87%)	DCC
Wilson B-factor (Å <sup>2</sup> )	49.1	Xtriage
Anisotropy	0.057	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 14.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	9168	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.03% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: COH, CEL, NAG, BOG

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	0.29	0/4581	0.42	2/6222 (0.0%)
1	B	0.30	1/4568 (0.0%)	0.42	1/6203 (0.0%)
All	All	0.30	1/9149 (0.0%)	0.42	3/12425 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	107	PRO	N-CD	5.04	1.54	1.47

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	121	ARG	C-N-CD	6.67	142.42	128.40
1	B	106	ILE	C-N-CD	5.42	139.79	128.40
1	A	122	PRO	CA-N-CD	-5.30	104.09	111.50

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	4448	0	4291	32	0
1	B	4435	0	4274	33	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	43	0	30	4	0
2	B	43	0	30	3	0
3	A	26	0	14	1	0
3	B	26	0	14	0	0
4	A	20	0	28	1	0
5	A	56	0	51	2	0
5	B	56	0	51	2	0
6	A	5	0	0	0	0
6	B	10	0	0	0	0
All	All	9168	0	8783	69	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 4.

All (69) 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:115:LYS:O	1:A:119:THR:OG1	1.95	0.84
2:A:601:COH:HMB1	2:A:601:COH:HBB1	1.68	0.75
1:B:96:PHE:HB3	1:B:99:VAL:CG1	2.18	0.74
2:A:601:COH:HHD	2:A:601:COH:HBC1	1.79	0.65
1:B:414:ILE:HG12	5:B:606:NAG:H82	1.78	0.65
1:B:121:ARG:HD3	1:B:529:PRO:HG3	1.79	0.64
2:B:601:COH:HHC	2:B:601:COH:HBB1	1.80	0.63
1:A:120:SER:OG	1:A:121:ARG:N	2.32	0.63
1:A:122:PRO:HD2	1:A:123:TYR:H	1.65	0.61
2:B:601:COH:HBC1	2:B:601:COH:HHD	1.82	0.61
1:A:104:ASN:HB3	1:A:359:LYS:HD3	1.83	0.61
1:A:414:ILE:HA	5:A:607:NAG:H82	1.84	0.60
1:A:101:ASN:O	1:A:105:ASN:ND2	2.34	0.59
1:A:124:LEU:HD13	1:A:471:PHE:HD1	1.69	0.58
1:A:124:LEU:HD13	1:A:471:PHE:CD1	2.40	0.57
1:A:211:PHE:HB3	2:A:601:COH:HBD1	1.87	0.56
1:B:176:LYS:NZ	1:B:180:GLU:OE2	2.38	0.55
1:B:96:PHE:O	1:B:99:VAL:HG13	2.07	0.55
1:B:473:LEU:HD21	1:B:525:GLU:HG3	1.88	0.54
1:A:514:ARG:HH21	1:A:521:GLU:HG3	1.73	0.54
1:B:390:PRO:HB3	1:B:440:ASN:HB3	1.90	0.53
1:A:122:PRO:HG2	1:A:123:TYR:CE2	2.43	0.53
1:B:381:GLU:HG2	1:B:467:TYR:CE2	2.46	0.51
1:B:175:SER:OG	1:B:450:LYS:NZ	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:411:ASN:HB3	1:B:414:ILE:HG13	1.92	0.50
1:A:484:GLY:HA3	1:A:512:LYS:HD3	1.94	0.50
1:A:179:LEU:HD22	1:A:450:LYS:HB2	1.93	0.49
1:B:96:PHE:HB3	1:B:99:VAL:HG11	1.95	0.49
1:B:539:PRO:O	1:B:542:SER:OG	2.25	0.48
1:A:214:ASP:HB2	1:A:223:ARG:HG3	1.96	0.48
1:B:442:PRO:HG2	1:B:445:VAL:HG22	1.95	0.47
1:A:361:LYS:HE2	1:A:363:ASP:HB2	1.96	0.47
1:A:138:LYS:HE2	1:B:544:GLN:O	2.14	0.47
1:A:383:ASN:OD1	2:A:601:COH:HAD2	2.14	0.47
1:B:407:GLN:O	1:B:411:ASN:ND2	2.49	0.46
1:A:193:GLN:O	1:A:582[B]:ASN:ND2	2.49	0.46
1:A:288:VAL:HG11	1:A:303:ALA:HB1	1.96	0.46
1:A:388:TRP:HZ2	3:A:602:CEL:H112	1.80	0.46
1:B:345:VAL:HA	1:B:349:TYR:HB3	1.98	0.45
1:A:46:GLU:OE1	1:A:138:LYS:NZ	2.47	0.45
1:B:190:PRO:HB2	1:B:431:ILE:HD13	1.99	0.44
1:B:178:VAL:HG22	1:B:495:LEU:HD13	1.98	0.44
1:B:67:GLU:HB3	5:B:603:NAG:H82	1.99	0.44
1:A:345:VAL:HA	1:A:349:TYR:HB3	1.99	0.44
1:A:180:GLU:HB3	4:A:603:BOG:H4'2	2.00	0.44
1:A:122:PRO:HG2	1:A:123:TYR:CD2	2.53	0.44
1:B:496:TYR:HE1	1:B:503:GLU:HG3	1.83	0.43
1:A:411:ASN:HB3	1:A:414:ILE:HG13	2.00	0.43
1:B:498:ASP:HB3	1:B:501:VAL:HG22	2.00	0.43
1:B:277:PRO:HA	1:B:278:PRO:HD3	1.91	0.43
1:B:210:PHE:HB2	1:B:378:ILE:HG13	2.01	0.43
1:B:390:PRO:HB2	1:B:435:VAL:HA	2.00	0.42
1:B:96:PHE:CB	1:B:99:VAL:CG1	2.94	0.42
1:B:284:LEU:HA	1:B:284:LEU:HD23	1.93	0.42
1:B:199:PHE:CZ	1:B:353:LEU:HD13	2.55	0.42
1:A:122:PRO:CD	1:A:123:TYR:H	2.31	0.42
1:B:121:ARG:HB3	1:B:124:LEU:HD12	2.02	0.42
1:A:414:ILE:HG12	5:A:607:NAG:H82	2.01	0.42
1:B:103:VAL:HG13	1:B:109:LEU:HD13	2.02	0.42
1:A:403:TYR:OH	1:A:418:HIS:NE2	2.45	0.42
1:B:207:THR:HG21	1:B:386:TYR:CE2	2.56	0.41
1:A:124:LEU:HD23	1:A:124:LEU:HA	1.91	0.41
1:A:519:PHE:HB3	1:A:523:MET:HB3	2.01	0.41
1:B:133:VAL:HG13	1:B:134:HIS:CD2	2.56	0.41
1:A:390:PRO:HB2	1:A:435:VAL:HA	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:457:ARG:NH2	1:A:503:GLU:OE1	2.42	0.41
1:B:182:VAL:HG21	1:B:492:LEU:HD21	2.01	0.41
2:B:601:COH:HBC1	2:B:601:COH:CHD	2.49	0.40
1:B:113:ILE:O	1:B:117:VAL:HG23	2.22	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	551/552 (100%)	528 (96%)	21 (4%)	2 (0%)	39	73
1	B	550/552 (100%)	526 (96%)	24 (4%)	0	100	100
All	All	1101/1104 (100%)	1054 (96%)	45 (4%)	2 (0%)	52	83

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	120	SER
1	A	122	PRO

### 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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	482/494 (98%)	470 (98%)	12 (2%)	55	85
1	B	477/494 (97%)	467 (98%)	10 (2%)	61	89
All	All	959/988 (97%)	937 (98%)	22 (2%)	58	87

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	119	THR
1	A	123	TYR
1	A	233	HIS
1	A	240	ASP
1	A	290	GLN
1	A	386	TYR
1	A	416	LEU
1	A	417	GLU
1	A	423	PHE
1	A	485	GLU
1	A	531	SER
1	A	557	PHE
1	B	99	VAL
1	B	107	PRO
1	B	109	LEU
1	B	240	ASP
1	B	290	GLN
1	B	386	TYR
1	B	410	TYR
1	B	485	GLU
1	B	531	SER
1	B	557	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

13 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$
2	COH	A	601	-	27,50,50	4.45	19 (70%)	23,82,82	4.45	10 (43%)
3	CEL	A	602	-	26,28,28	1.46	4 (15%)	39,43,43	2.86	12 (30%)
4	BOG	A	603	-	20,20,20	0.90	1 (5%)	25,25,25	0.89	1 (4%)
5	NAG	A	604	1	14,14,15	0.29	0	15,19,21	0.45	0
5	NAG	A	605	1,5	14,14,15	0.38	0	15,19,21	0.39	0
5	NAG	A	606	5	14,14,15	0.30	0	15,19,21	0.46	0
5	NAG	A	607	1	14,14,15	0.26	0	15,19,21	0.21	0
2	COH	B	601	-	27,50,50	4.46	19 (70%)	23,82,82	4.22	6 (26%)
3	CEL	B	602	-	26,28,28	1.46	3 (11%)	39,43,43	2.77	11 (28%)
5	NAG	B	603	1	14,14,15	0.19	0	15,19,21	0.38	0
5	NAG	B	604	1,5	14,14,15	0.33	0	15,19,21	0.39	0
5	NAG	B	605	5	14,14,15	0.31	0	15,19,21	0.33	0
5	NAG	B	606	1	14,14,15	0.28	0	15,19,21	0.25	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
2	COH	A	601	-	-	0/8/94/94	0/0/8/8

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	CEL	A	602	-	-	0/20/20/20	0/3/3/3
4	BOG	A	603	-	-	0/11/31/31	0/1/1/1
5	NAG	A	604	1	-	0/6/23/26	0/1/1/1
5	NAG	A	605	1,5	-	0/6/23/26	0/1/1/1
5	NAG	A	606	5	-	0/6/23/26	0/1/1/1
5	NAG	A	607	1	-	0/6/23/26	0/1/1/1
2	COH	B	601	-	-	0/8/94/94	0/0/8/8
3	CEL	B	602	-	-	0/20/20/20	0/3/3/3
5	NAG	B	603	1	-	0/6/23/26	0/1/1/1
5	NAG	B	604	1,5	-	0/6/23/26	0/1/1/1
5	NAG	B	605	5	-	0/6/23/26	0/1/1/1
5	NAG	B	606	1	-	0/6/23/26	0/1/1/1

All (46) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	COH	CHB-C4A	-7.86	1.38	1.53
2	B	601	COH	CHB-C4A	-7.75	1.38	1.53
2	B	601	COH	CHA-C1A	-7.72	1.38	1.53
2	A	601	COH	CHA-C1A	-7.65	1.38	1.53
2	A	601	COH	CHC-C1C	-7.60	1.38	1.53
2	B	601	COH	CHD-C4C	-7.59	1.38	1.53
2	A	601	COH	CHD-C4C	-7.47	1.39	1.53
2	B	601	COH	CHC-C1C	-7.32	1.39	1.53
2	B	601	COH	C3B-C2B	-5.15	1.33	1.40
3	A	602	CEL	C5-C3	-4.91	1.39	1.47
3	B	602	CEL	C5-C3	-4.89	1.39	1.47
2	A	601	COH	CHB-C1B	-4.84	1.38	1.51
2	B	601	COH	CHB-C1B	-4.81	1.38	1.51
2	B	601	COH	CHA-C4D	-4.72	1.38	1.51
2	A	601	COH	CHC-C4B	-4.70	1.38	1.51
2	B	601	COH	CHD-C1D	-4.68	1.38	1.51
2	A	601	COH	CHA-C4D	-4.62	1.39	1.51
2	A	601	COH	CHD-C1D	-4.61	1.39	1.51
2	B	601	COH	CHC-C4B	-4.50	1.39	1.51
3	B	602	CEL	C12-N2	-4.18	1.35	1.44
3	A	602	CEL	C12-N2	-4.18	1.35	1.44
2	A	601	COH	C3B-C2B	-4.13	1.35	1.40
2	B	601	COH	C4C-C3C	-3.19	1.46	1.50
2	A	601	COH	C4C-C3C	-2.98	1.46	1.50
3	A	602	CEL	C2-C1	-2.07	1.36	1.39
3	A	602	CEL	S1-N3	2.22	1.64	1.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	602	CEL	S1-N3	2.28	1.64	1.60
2	B	601	COH	CMC-C2C	2.36	1.54	1.50
2	A	601	COH	CMC-C2C	2.49	1.54	1.50
2	A	601	COH	CMA-C3A	2.63	1.55	1.50
4	A	603	BOG	O5-C1	2.68	1.48	1.41
2	B	601	COH	CMA-C3A	2.69	1.55	1.50
2	A	601	COH	CAA-C2A	3.01	1.55	1.51
2	B	601	COH	CAA-C2A	3.22	1.55	1.51
2	B	601	COH	C3B-CAB	3.56	1.55	1.47
2	A	601	COH	C3B-CAB	3.83	1.55	1.47
2	B	601	COH	C1B-C2B	5.00	1.45	1.38
2	A	601	COH	C4D-C3D	5.01	1.45	1.38
2	B	601	COH	CAC-C3C	5.07	1.55	1.45
2	A	601	COH	CAC-C3C	5.11	1.55	1.45
2	B	601	COH	C4D-C3D	5.19	1.45	1.38
2	B	601	COH	C1D-C2D	5.23	1.45	1.38
2	A	601	COH	C1B-C2B	5.29	1.45	1.38
2	B	601	COH	C3D-C2D	5.44	1.53	1.37
2	A	601	COH	C1D-C2D	5.46	1.45	1.38
2	A	601	COH	C3D-C2D	5.48	1.53	1.37

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602	CEL	O1-S1-O2	-12.97	101.44	118.87
3	B	602	CEL	O1-S1-O2	-12.96	101.45	118.87
2	A	601	COH	CMB-C2B-C1B	-3.88	123.17	127.13
3	B	602	CEL	C2-C1-N1	-2.95	106.90	111.54
3	A	602	CEL	C2-C3-C5	-2.84	123.40	128.10
3	A	602	CEL	C2-C1-N1	-2.75	107.20	111.54
2	A	601	COH	CBD-CAD-C3D	-2.72	107.70	112.49
2	A	601	COH	CHC-C4B-C3B	-2.72	126.24	129.66
2	A	601	COH	CBC-CAC-C3C	-2.38	122.87	126.34
3	B	602	CEL	C2-C3-C5	-2.34	124.24	128.10
2	B	601	COH	CBC-CAC-C3C	-2.26	123.05	126.34
3	A	602	CEL	F2-C4-C1	-2.15	108.62	112.54
4	A	603	BOG	C6-C5-C4	-2.00	107.97	112.99
2	A	601	COH	CMB-C2B-C3B	2.01	129.01	125.09
3	B	602	CEL	C17-C12-N2	2.29	122.95	119.55
2	A	601	COH	C4B-C3B-C2B	2.45	107.54	104.44
3	A	602	CEL	O2-S1-N3	2.74	110.65	107.25
2	B	601	COH	C4B-C3B-C2B	2.81	107.98	104.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602	CEL	C17-C12-N2	2.85	123.79	119.55
3	B	602	CEL	O2-S1-N3	2.86	110.80	107.25
3	B	602	CEL	C3-C2-C1	2.97	108.80	104.76
3	A	602	CEL	C3-C2-C1	3.02	108.88	104.76
3	B	602	CEL	C5-C3-N2	3.06	128.29	123.40
3	B	602	CEL	O1-S1-N3	3.07	111.07	107.25
3	B	602	CEL	O2-S1-C15	3.18	111.00	107.41
3	A	602	CEL	O2-S1-C15	3.23	111.05	107.41
3	A	602	CEL	O1-S1-N3	3.46	111.55	107.25
3	A	602	CEL	C5-C3-N2	3.69	129.29	123.40
3	B	602	CEL	O1-S1-C15	3.76	111.65	107.41
3	A	602	CEL	O1-S1-C15	4.21	112.15	107.41
3	B	602	CEL	C4-C1-N1	5.04	126.00	119.69
3	A	602	CEL	C4-C1-N1	5.23	126.23	119.69
2	B	601	COH	CHB-C4A-NA	9.22	124.25	110.12
2	B	601	COH	CHD-C4C-NC	9.32	124.41	110.12
2	A	601	COH	CHD-C4C-NC	9.45	124.61	110.12
2	A	601	COH	CHB-C4A-NA	9.75	125.06	110.12
2	B	601	COH	CHA-C1A-NA	10.12	125.62	110.12
2	A	601	COH	CHA-C1A-NA	10.20	125.76	110.12
2	B	601	COH	CHC-C1C-NC	10.24	125.82	110.12
2	A	601	COH	CHC-C1C-NC	10.65	126.44	110.12

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	COH	4	0
3	A	602	CEL	1	0
4	A	603	BOG	1	0
5	A	607	NAG	2	0
2	B	601	COH	3	0
5	B	603	NAG	1	0
5	B	606	NAG	1	0

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	552/552 (100%)	-0.32	6 (1%)	82 75	29, 40, 58, 81	0
1	B	551/552 (99%)	-0.36	3 (0%)	91 88	31, 40, 60, 95	0
All	All	1103/1104 (99%)	-0.34	9 (0%)	87 81	29, 40, 59, 95	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	74	PHE	3.7
1	B	74	PHE	3.0
1	B	82	LEU	2.9
1	A	75	LEU	2.8
1	A	79	LYS	2.4
1	A	584	GLN	2.2
1	A	81	LEU	2.1
1	A	240	ASP	2.1
1	B	240	ASP	2.0

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

### 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron

density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	COH	B	601	43/43	0.84	0.32	4.02	48,78,100,104	0
5	NAG	A	604	14/15	0.91	0.30	3.68	50,53,64,70	0
5	NAG	A	607	14/15	0.88	0.29	2.93	56,66,79,92	0
2	COH	A	601	43/43	0.82	0.32	2.88	50,87,105,109	0
5	NAG	B	606	14/15	0.86	0.35	2.03	56,66,82,82	0
5	NAG	B	603	14/15	0.88	0.29	1.81	59,66,77,83	0
4	BOG	A	603	20/20	0.95	0.22	1.28	43,49,66,71	0
5	NAG	B	604	14/15	0.95	0.21	0.61	35,38,44,48	0
5	NAG	A	605	14/15	0.97	0.20	-0.01	31,34,41,49	0
3	CEL	B	602	26/26	0.96	0.15	-0.17	37,43,49,67	0
3	CEL	A	602	26/26	0.96	0.14	-0.65	36,38,48,49	0
5	NAG	A	606	14/15	0.89	0.24	-	41,46,57,62	0
5	NAG	B	605	14/15	0.84	0.32	-	47,57,69,69	0

## 6.5 Other polymers

There are no such residues in this entry.