



# Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 10:00 am BST

PDB ID : 6I1G  
Title : Crystal structure of TP domain from Chlamydia trachomatis Penicillin-Binding Protein 3 in complex with piperacillin  
Authors : Bellini, D.; Koekemoer, L.; Newman, H.; Dowson, C.G.  
Deposited on : 2018-10-28  
Resolution : 2.13 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

<https://www.wwpdb.org/validation/2017/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	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

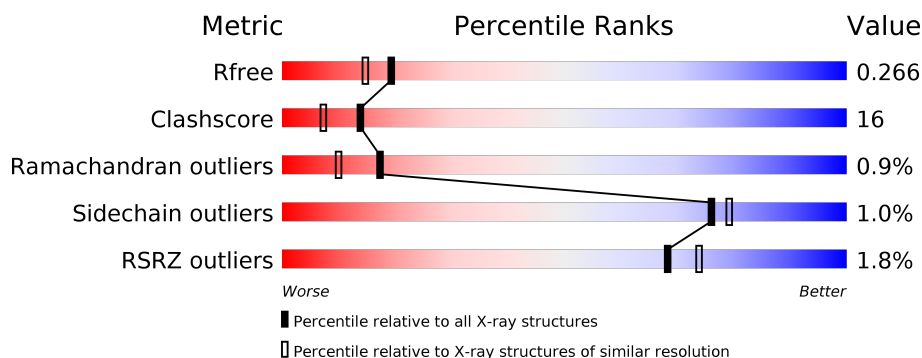
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.13 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))
$R_{free}$	130704	2523 (2.16-2.12)
Clashscore	141614	2653 (2.16-2.12)
Ramachandran outliers	138981	2618 (2.16-2.12)
Sidechain outliers	138945	2617 (2.16-2.12)
RSRZ outliers	127900	2485 (2.16-2.12)

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 respectively. 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	346	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="display: flex; justify-content: space-between; align-items: center;"> <span>%</span> <span style="flex-grow: 1; border-bottom: 1px solid black; position: relative;"> <span style="position: absolute; left: 0; top: -5px; width: 100%; height: 10px; background-color: red;"></span> <span style="position: absolute; left: 0; top: 0; width: 65%; height: 10px; background-color: green;"></span> <span style="position: absolute; left: 65%; top: 0; width: 29%; height: 10px; background-color: yellow;"></span> <span style="position: absolute; left: 95%; top: 0; width: 6%; height: 10px; background-color: grey;"></span> </span></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>65%</span> <span>29%</span> <span>• 6%</span> </div> </div>
1	B	346	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="display: flex; justify-content: space-between; align-items: center;"> <span>3%</span> <span style="flex-grow: 1; border-bottom: 1px solid black; position: relative;"> <span style="position: absolute; left: 0; top: -5px; width: 100%; height: 10px; background-color: red;"></span> <span style="position: absolute; left: 0; top: 0; width: 73%; height: 10px; background-color: green;"></span> <span style="position: absolute; left: 73%; top: 0; width: 25%; height: 10px; background-color: yellow;"></span> <span style="position: absolute; left: 98%; top: 0; width: 2%; height: 10px; background-color: grey;"></span> </span></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>73%</span> <span>25%</span> <span>••</span> </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	JPP	A	701	X	-	-	-
2	JPP	B	701	X	-	-	-

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5023 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.

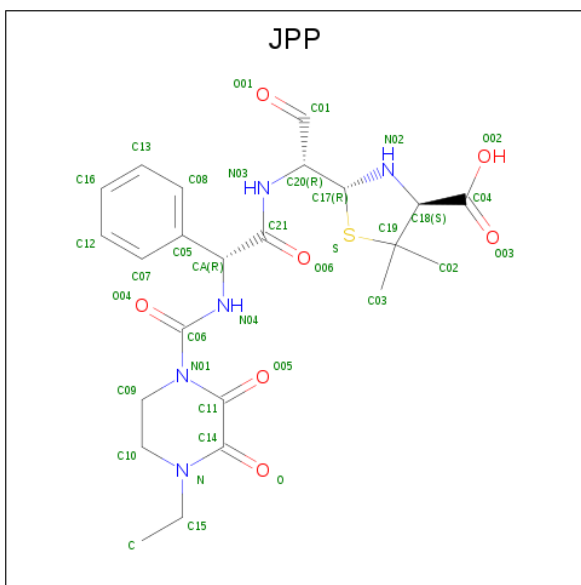
- Molecule 1 is a protein called Penicillin-binding protein, Penicillin-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	325	Total	C	N	O	S	0	0	0
			2407	1512	410	477	8			
1	B	341	Total	C	N	O	S	0	0	0
			2527	1585	427	506	9			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	256	GLY	-	expression tag	UNP A0A0E9FXJ8
A	257	PRO	-	expression tag	UNP A0A0E9FXJ8
A	258	GLY	-	expression tag	UNP A0A0E9FXJ8
A	259	TYR	-	expression tag	UNP A0A0E9FXJ8
A	260	GLN	-	expression tag	UNP A0A0E9FXJ8
A	261	ASP	-	expression tag	UNP A0A0E9FXJ8
A	262	PRO	-	expression tag	UNP A0A0E9FXJ8
A	306	PRO	-	linker	UNP A0A0E9FXJ8
A	307	GLY	-	linker	UNP A0A0E9FXJ8
B	256	GLY	-	expression tag	UNP A0A0E9FXJ8
B	257	PRO	-	expression tag	UNP A0A0E9FXJ8
B	258	GLY	-	expression tag	UNP A0A0E9FXJ8
B	259	TYR	-	expression tag	UNP A0A0E9FXJ8
B	260	GLN	-	expression tag	UNP A0A0E9FXJ8
B	261	ASP	-	expression tag	UNP A0A0E9FXJ8
B	262	PRO	-	expression tag	UNP A0A0E9FXJ8
B	306	PRO	-	linker	UNP A0A0E9FXJ8
B	307	GLY	-	linker	UNP A0A0E9FXJ8

- Molecule 2 is Piperacillin (Open Form) (three-letter code: JPP) (formula:  $C_{23}H_{29}N_5O_7S$ ) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			36	23	5	7	1		
2	B	1	Total	C	N	O	S	0	0
			36	23	5	7	1		

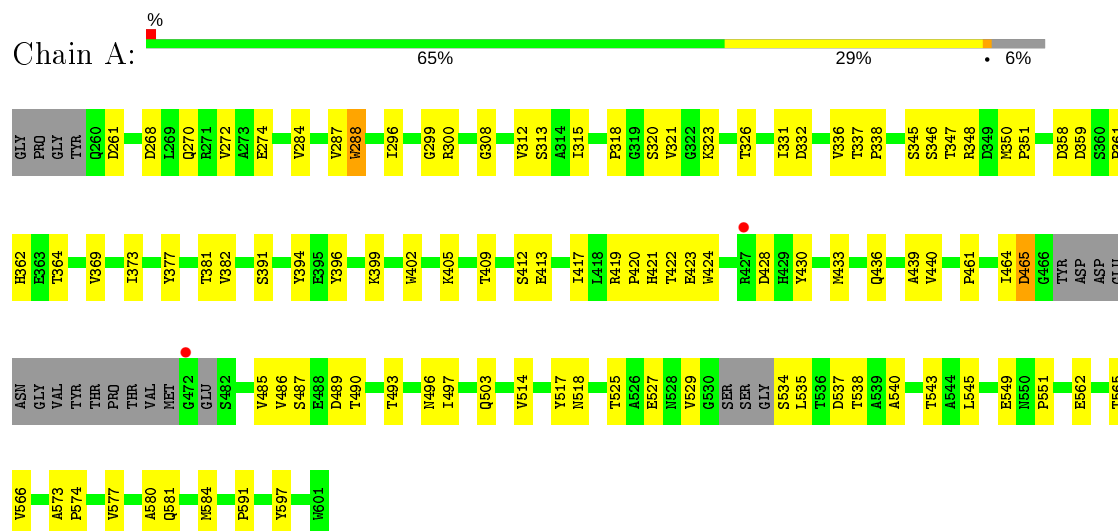
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	11	Total	O	0	0
			11	11		
3	B	6	Total	O	0	0
			6	6		

### 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 the various outlier classes 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: Penicillin-binding protein, Penicillin-binding protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	107.80Å 107.80Å 117.73Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	93.36 – 2.13 93.36 – 2.13	Depositor EDS
% Data completeness (in resolution range)	48.5 (93.36-2.13) 48.5 (93.36-2.13)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.39 (at 2.12Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, $R_{free}$	0.240 , 0.284 0.226 , 0.266	Depositor DCC
$R_{free}$ test set	1118 reflections (5.14%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	16.0	Xtriage
Anisotropy	0.757	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 24.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.038 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	5023	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.99% 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: JPP

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	0.57	2/2450 (0.1%)	0.68	0/3336
1	B	0.51	1/2575 (0.0%)	0.64	0/3511
All	All	0.54	3/5025 (0.1%)	0.66	0/6847

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	402	TRP	CD2-CE2	5.68	1.48	1.41
1	A	288	TRP	CD2-CE2	5.65	1.48	1.41
1	B	402	TRP	CD2-CE2	5.16	1.47	1.41

There are no bond angle outliers.

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	2407	0	2395	88	0
1	B	2527	0	2499	68	0
2	A	36	0	26	9	0
2	B	36	0	27	3	0
3	A	11	0	0	0	0
3	B	6	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5023	0	4947	157	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 16.

All (157) 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:320:SER:OG	2:B:701:JPP:C01	1.67	1.42
1:A:486:VAL:HG13	1:A:490:THR:OG1	1.70	0.90
1:B:449:VAL:HG11	1:B:543:THR:HB	1.56	0.87
1:B:332:ASP:HA	1:B:490:THR:HG21	1.56	0.85
1:A:486:VAL:CG1	1:A:490:THR:OG1	2.28	0.82
1:A:373:ILE:HD11	1:A:497:ILE:HG22	1.64	0.80
1:B:382:VAL:HG13	1:B:430:TYR:CD1	2.16	0.80
1:A:296:ILE:HD13	1:A:551:PRO:HB2	1.63	0.79
1:B:344:VAL:HG12	1:B:362:HIS:HE1	1.48	0.79
1:A:332:ASP:HA	1:A:490:THR:HG21	1.64	0.79
1:A:525:THR:HG22	2:A:701:JPP:H7	1.67	0.77
1:B:271:ARG:O	1:B:275:GLU:HG2	1.87	0.74
1:A:312:VAL:HG12	1:A:413:GLU:HA	1.68	0.74
1:A:405:LYS:HG2	1:A:413:GLU:HG2	1.70	0.74
1:A:323:LYS:HE2	1:A:381:THR:HG21	1.71	0.72
1:A:525:THR:HG21	2:A:701:JPP:H11	1.74	0.69
1:A:348:ARG:HG2	1:A:350:MET:HE1	1.73	0.69
1:B:287:VAL:HG23	1:B:288:TRP:CD1	2.28	0.68
1:A:525:THR:CG2	2:A:701:JPP:H11	2.23	0.68
1:B:327:LEU:O	1:B:331:ILE:HG12	1.95	0.67
1:A:382:VAL:HG13	1:A:430:TYR:CD1	2.30	0.67
1:B:381:THR:HA	1:B:384:ILE:HD12	1.77	0.66
1:B:312:VAL:HG12	1:B:413:GLU:HA	1.78	0.66
1:A:529:VAL:HG22	1:A:535:LEU:HD22	1.78	0.66
1:A:261:ASP:OD1	1:B:283:ARG:NH2	2.29	0.66
1:A:525:THR:O	2:A:701:JPP:H18	1.96	0.66
1:B:465:ASP:O	1:B:477:THR:HG23	1.97	0.65
1:B:320:SER:CB	2:B:701:JPP:C01	2.74	0.65
1:A:270:GLN:O	1:A:274:GLU:HG3	1.97	0.64
1:A:321:VAL:HG11	1:A:543:THR:CG2	2.27	0.64
1:B:486:VAL:HG22	1:B:490:THR:OG1	1.97	0.64
1:A:421:HIS:C	1:A:423:GLU:H	2.01	0.64
1:A:326:THR:HG21	1:A:433:MET:HA	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:514:VAL:HG13	1:A:581:GLN:HG2	1.81	0.63
1:B:408:ILE:HD11	1:B:463:ILE:N	2.14	0.63
1:B:336:VAL:CG1	1:B:369:VAL:HG21	2.29	0.62
1:B:344:VAL:HG12	1:B:362:HIS:CE1	2.31	0.62
1:A:318:PRO:HB3	1:A:440:VAL:HG11	1.80	0.61
1:A:537:ASP:OD1	1:A:562:GLU:O	2.19	0.61
1:B:323:LYS:HE2	1:B:381:THR:HG21	1.83	0.61
1:B:503:GLN:OE1	1:B:598:LYS:HD3	2.01	0.61
1:A:538:THR:OG1	1:A:566:VAL:HG12	2.00	0.61
1:A:321:VAL:HG11	1:A:543:THR:HG23	1.84	0.60
1:B:348:ARG:CG	1:B:350:MET:HE1	2.32	0.60
1:B:347:THR:HG22	1:B:357:SER:HB3	1.83	0.60
1:A:424:TRP:HB3	1:A:428:ASP:HB2	1.84	0.60
1:B:529:VAL:HA	1:B:535:LEU:HA	1.82	0.60
1:A:490:THR:HA	1:A:493:THR:HG22	1.83	0.60
1:A:486:VAL:HG13	1:A:490:THR:HG1	1.67	0.59
1:B:344:VAL:CG1	1:B:362:HIS:HE1	2.16	0.59
1:B:517:TYR:CE2	1:B:591:PRO:HA	2.37	0.59
1:B:264:THR:HG22	1:B:300:ARG:HE	1.66	0.59
1:B:323:LYS:CE	1:B:381:THR:HG21	2.32	0.59
1:A:296:ILE:HD11	1:A:545:LEU:HD22	1.85	0.58
1:A:296:ILE:CD1	1:A:551:PRO:HB2	2.31	0.58
1:A:268:ASP:O	1:A:272:VAL:HG23	2.05	0.57
1:A:382:VAL:HG13	1:A:430:TYR:CE1	2.40	0.57
1:A:284:VAL:HG11	1:A:566:VAL:HG21	1.87	0.57
1:A:377:TYR:O	1:A:381:THR:HG23	2.04	0.57
1:B:399:LYS:HB3	1:B:485:VAL:HG13	1.87	0.57
1:A:487:SER:OG	1:A:490:THR:HG23	2.05	0.56
1:B:385:GLY:HA3	1:B:433:MET:SD	2.45	0.56
1:A:336:VAL:CG1	1:A:369:VAL:HG21	2.36	0.56
1:B:395:GLU:O	1:B:399:LYS:HG2	2.06	0.55
2:B:701:JPP:C01	2:B:701:JPP:O06	2.45	0.55
1:B:288:TRP:CE3	1:B:310:ARG:HD2	2.42	0.55
1:B:503:GLN:HG3	1:B:596:LEU:HD21	1.89	0.54
1:B:305:ALA:HB1	1:B:306:PRO:HA	1.88	0.54
1:B:503:GLN:HG3	1:B:596:LEU:CD2	2.38	0.53
1:B:394:TYR:CZ	1:B:420:PRO:HA	2.44	0.53
1:A:337:THR:HB	1:A:338:PRO:HD2	1.89	0.53
1:B:464:ILE:HG22	1:B:477:THR:HG21	1.91	0.53
1:A:345:SER:HA	1:A:364:THR:HG22	1.92	0.52
1:A:320:SER:HB3	1:A:323:LYS:HD2	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:527:GLU:HB3	1:A:535:LEU:HD12	1.90	0.52
1:A:490:THR:HA	1:A:493:THR:CG2	2.39	0.52
2:A:701:JPP:O06	2:A:701:JPP:C01	2.58	0.52
1:A:394:TYR:CE1	1:A:420:PRO:HA	2.45	0.51
1:A:300:ARG:HD2	1:A:465:ASP:OD1	2.10	0.51
1:A:359:ASP:HB3	2:A:701:JPP:S	2.51	0.51
1:B:287:VAL:HG11	1:B:562:GLU:CD	2.31	0.51
1:A:346:SER:HB3	1:A:361:PRO:HA	1.92	0.51
1:A:345:SER:HB2	1:A:347:THR:O	2.11	0.50
1:A:496:ASN:HB3	1:A:597:TYR:CD1	2.47	0.50
1:B:356:ILE:HG13	1:B:382:VAL:HG11	1.94	0.50
1:A:346:SER:CB	1:A:361:PRO:HA	2.42	0.50
1:A:421:HIS:HA	1:A:424:TRP:CE2	2.46	0.50
1:B:394:TYR:CE1	1:B:420:PRO:HA	2.46	0.50
1:A:284:VAL:HG11	1:A:566:VAL:CG2	2.42	0.50
1:A:287:VAL:HG23	1:A:288:TRP:CD1	2.46	0.50
1:A:573:ALA:N	1:A:574:PRO:CD	2.74	0.50
1:B:469:ASP:O	1:B:471:ASN:HB2	2.12	0.50
1:A:358:ASP:OD2	1:A:362:HIS:NE2	2.40	0.49
1:A:538:THR:HG21	1:A:565:THR:HA	1.94	0.49
1:B:266:ASN:HB3	1:B:269:LEU:HB3	1.93	0.49
1:B:337:THR:HB	1:B:338:PRO:HD2	1.94	0.49
1:A:348:ARG:CG	1:A:350:MET:HE1	2.41	0.48
1:A:331:ILE:HG22	1:A:490:THR:HB	1.93	0.48
1:B:299:GLY:O	1:B:461:PRO:HA	2.12	0.48
1:B:348:ARG:CD	1:B:350:MET:HE1	2.43	0.48
1:A:348:ARG:HG2	1:A:350:MET:CE	2.41	0.48
1:B:598:LYS:HD2	1:B:599:TYR:O	2.14	0.48
1:B:348:ARG:HG2	1:B:350:MET:HE1	1.94	0.47
1:A:409:THR:OG1	1:A:464:ILE:HD12	2.13	0.47
1:A:421:HIS:C	1:A:423:GLU:N	2.66	0.47
1:A:430:TYR:O	1:A:433:MET:HB2	2.14	0.47
1:B:523:THR:HG22	1:B:542:PHE:HD1	1.80	0.47
1:A:399:LYS:HE2	1:A:485:VAL:HA	1.96	0.47
1:A:419:ARG:HD2	1:A:423:GLU:OE1	2.14	0.47
1:B:336:VAL:HG13	1:B:369:VAL:HG21	1.97	0.46
1:B:348:ARG:HG2	1:B:350:MET:CE	2.45	0.46
1:A:313:SER:HB3	1:A:412:SER:O	2.16	0.46
1:A:296:ILE:CD1	1:A:551:PRO:CB	2.93	0.46
1:B:284:VAL:HG11	1:B:571:ALA:HA	1.97	0.46
1:B:523:THR:HG22	1:B:542:PHE:CD1	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:421:HIS:O	1:A:423:GLU:N	2.49	0.45
1:B:318:PRO:HB3	1:B:440:VAL:HG11	1.99	0.45
1:B:546:ILE:HG13	1:B:547:PRO:HA	1.98	0.45
1:A:573:ALA:HB3	1:A:574:PRO:HD3	1.99	0.45
1:B:399:LYS:CB	1:B:485:VAL:HG13	2.47	0.45
1:A:350:MET:HG3	1:A:430:TYR:CE2	2.52	0.44
1:A:321:VAL:HG11	1:A:543:THR:HG21	1.98	0.44
1:A:336:VAL:HG12	1:A:337:THR:N	2.33	0.44
1:B:580:ALA:O	1:B:584:MET:HG2	2.17	0.44
1:A:391:SER:HB3	1:A:421:HIS:CE1	2.52	0.44
1:B:504:PRO:HA	1:B:509:GLU:CD	2.39	0.43
1:A:331:ILE:HG22	1:A:490:THR:CG2	2.48	0.43
1:A:525:THR:HG23	2:A:701:JPP:H18	2.00	0.43
1:A:573:ALA:O	1:A:577:VAL:HG23	2.19	0.43
1:A:424:TRP:HB3	1:A:428:ASP:CB	2.48	0.42
1:A:525:THR:HG21	2:A:701:JPP:C02	2.46	0.42
1:B:377:TYR:O	1:B:381:THR:HG23	2.19	0.42
1:A:527:GLU:HB2	2:A:701:JPP:H19	2.00	0.42
1:A:517:TYR:CE1	1:A:591:PRO:HA	2.54	0.42
1:B:327:LEU:HD22	1:B:373:ILE:HD12	2.01	0.42
1:A:299:GLY:O	1:A:461:PRO:HA	2.20	0.42
1:A:332:ASP:HB3	1:A:396:TYR:OH	2.20	0.42
1:B:409:THR:OG1	1:B:464:ILE:HD12	2.19	0.42
1:B:528:ASN:C	1:B:535:LEU:HD23	2.40	0.42
1:A:489:ASP:O	1:A:493:THR:HG22	2.19	0.42
1:A:529:VAL:HG13	1:A:534:SER:HA	2.02	0.42
1:B:348:ARG:HB2	1:B:379:THR:HG22	2.02	0.41
1:B:527:GLU:HB3	1:B:535:LEU:HD22	2.01	0.41
1:A:405:LYS:HE2	1:A:413:GLU:O	2.20	0.41
1:A:417:ILE:HB	1:A:439:ALA:HB3	2.01	0.41
1:B:337:THR:HB	1:B:338:PRO:CD	2.50	0.41
1:A:315:ILE:HA	1:A:440:VAL:O	2.21	0.41
1:B:394:TYR:O	1:B:398:GLN:HG3	2.20	0.41
1:A:296:ILE:HD11	1:A:545:LEU:CD2	2.50	0.41
1:B:327:LEU:HD22	1:B:373:ILE:CD1	2.50	0.41
1:A:580:ALA:O	1:A:584:MET:HG2	2.21	0.41
1:A:518:ASN:HB2	1:A:549:GLU:OE2	2.19	0.41
1:B:449:VAL:HG23	1:B:554:ALA:HB1	2.01	0.41
1:B:382:VAL:HG13	1:B:430:TYR:CE1	2.56	0.41
1:A:540:ALA:HB2	1:A:566:VAL:HG13	2.03	0.40
1:B:280:ALA:O	1:B:284:VAL:HG22	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:344:VAL:HG11	1:B:377:TYR:CD2	2.56	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	318/346 (92%)	295 (93%)	19 (6%)	4 (1%)	12	6
1	B	337/346 (97%)	313 (93%)	22 (6%)	2 (1%)	25	17
All	All	655/692 (95%)	608 (93%)	41 (6%)	6 (1%)	17	10

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	422	THR
1	B	320	SER
1	A	308	GLY
1	A	465	ASP
1	B	329	ALA
1	A	351	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.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	256/273 (94%)	254 (99%)	2 (1%)	81	85
1	B	270/273 (99%)	267 (99%)	3 (1%)	73	76
All	All	526/546 (96%)	521 (99%)	5 (1%)	76	79

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

Mol	Chain	Res	Type
1	A	436	GLN
1	A	503	GLN
1	B	283	ARG
1	B	311	SER
1	B	531	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	436	GLN
1	A	447	GLN
1	B	421	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

2 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	JPP	A	701	-	31,38,38	2.03	8 (25%)	39,55,55	3.02	8 (20%)
2	JPP	B	701	-	31,38,38	2.10	9 (29%)	39,55,55	4.03	19 (48%)

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	JPP	A	701	-	1/1/12/16	8/22/64/64	0/3/3/3
2	JPP	B	701	-	1/1/12/16	5/22/64/64	0/3/3/3

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	701	JPP	C14-C11	-5.12	1.45	1.53
2	A	701	JPP	C06-N01	-5.07	1.32	1.42
2	B	701	JPP	C06-N01	-4.95	1.32	1.42
2	A	701	JPP	C19-S	-4.66	1.75	1.85
2	B	701	JPP	C19-S	-4.56	1.76	1.85
2	A	701	JPP	C14-C11	-4.54	1.46	1.53
2	B	701	JPP	C17-S	-4.22	1.74	1.84
2	A	701	JPP	C17-S	-4.09	1.74	1.84
2	A	701	JPP	C11-N01	-3.47	1.33	1.40
2	B	701	JPP	C11-N01	-3.42	1.33	1.40
2	A	701	JPP	C05-CA	-2.81	1.48	1.52
2	B	701	JPP	CA-C21	-2.70	1.47	1.54
2	B	701	JPP	C05-CA	-2.54	1.48	1.52
2	A	701	JPP	C20-N03	-2.52	1.43	1.46
2	B	701	JPP	C18-N02	2.34	1.49	1.46
2	B	701	JPP	C14-N	-2.14	1.32	1.35
2	A	701	JPP	CA-N04	-2.09	1.42	1.46

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	701	JPP	C20-N03-C21	-18.63	102.77	123.12
2	A	701	JPP	C20-N03-C21	-11.24	110.83	123.12
2	A	701	JPP	C05-CA-C21	9.05	128.50	108.16
2	A	701	JPP	CA-N04-C06	-7.85	106.33	121.17
2	B	701	JPP	CA-N04-C06	-6.20	109.45	121.17
2	B	701	JPP	O06-C21-CA	-5.78	110.61	120.64
2	B	701	JPP	C05-CA-C21	5.43	120.36	108.16
2	B	701	JPP	O-C14-N	5.02	128.88	123.66
2	B	701	JPP	CA-C21-N03	4.85	129.37	116.82
2	B	701	JPP	O05-C11-N01	4.22	127.14	119.98
2	A	701	JPP	N04-C06-N01	3.83	119.61	114.04
2	A	701	JPP	O01-C01-C20	-3.77	114.33	124.83
2	B	701	JPP	C08-C05-CA	-3.73	114.68	120.80
2	A	701	JPP	O05-C11-N01	3.53	125.98	119.98
2	B	701	JPP	N04-C06-N01	3.52	119.16	114.04
2	B	701	JPP	O01-C01-C20	-3.39	115.39	124.83
2	B	701	JPP	C17-C20-N03	3.24	117.04	109.98
2	B	701	JPP	O-C14-C11	3.16	124.36	118.83
2	B	701	JPP	C19-S-C17	3.03	100.48	93.99
2	B	701	JPP	C21-CA-N04	2.61	116.44	108.44
2	A	701	JPP	O06-C21-CA	-2.43	116.42	120.64
2	B	701	JPP	C08-C05-C07	2.43	121.32	118.29
2	B	701	JPP	C15-N-C14	2.20	122.27	119.64
2	B	701	JPP	C10-N-C15	2.19	120.87	116.79
2	B	701	JPP	O05-C11-C14	2.17	122.63	118.83
2	B	701	JPP	C10-N-C14	-2.11	116.66	120.69
2	A	701	JPP	O05-C11-C14	2.01	122.36	118.83

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	701	JPP	C20
2	B	701	JPP	C20

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701	JPP	C17-C20-N03-C21
2	A	701	JPP	C01-C20-N03-C21
2	A	701	JPP	O06-C21-N03-C20
2	A	701	JPP	CA-C21-N03-C20
2	B	701	JPP	C-C15-N-C10
2	B	701	JPP	CA-C21-N03-C20

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Mol	Chain	Res	Type	Atoms
2	B	701	JPP	O06-C21-N03-C20
2	B	701	JPP	N03-C21-CA-N04
2	A	701	JPP	O06-C21-CA-C05
2	A	701	JPP	N03-C21-CA-C05
2	B	701	JPP	C-C15-N-C14
2	A	701	JPP	O06-C21-CA-N04
2	A	701	JPP	N03-C21-CA-N04

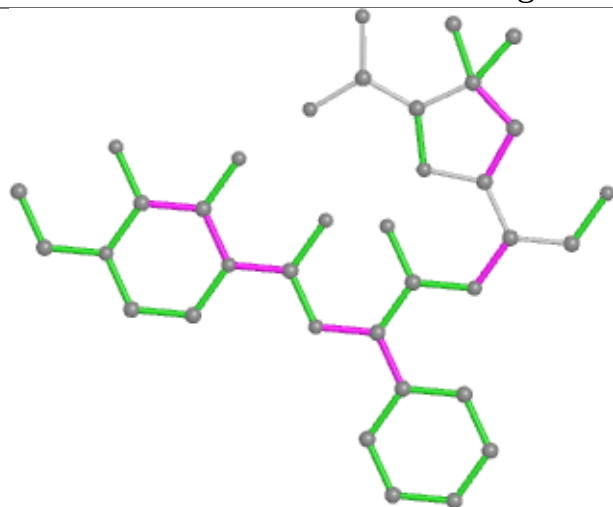
There are no ring outliers.

2 monomers are involved in 12 short contacts:

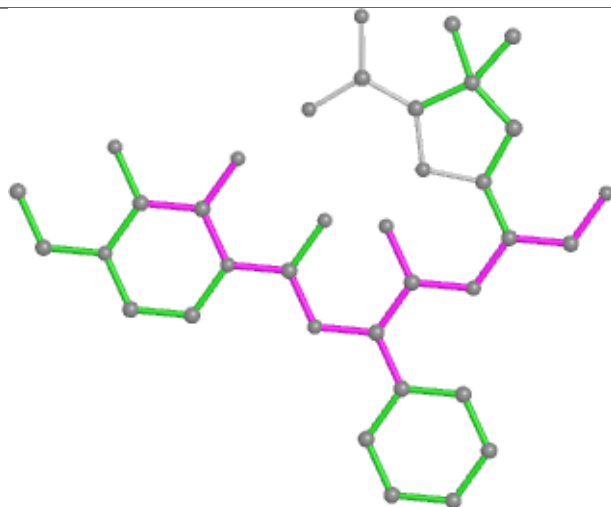
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	701	JPP	9	0
2	B	701	JPP	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

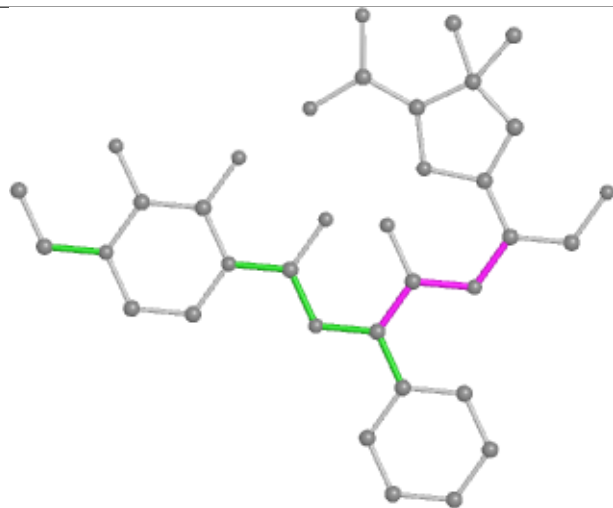
## Ligand JPP A 701



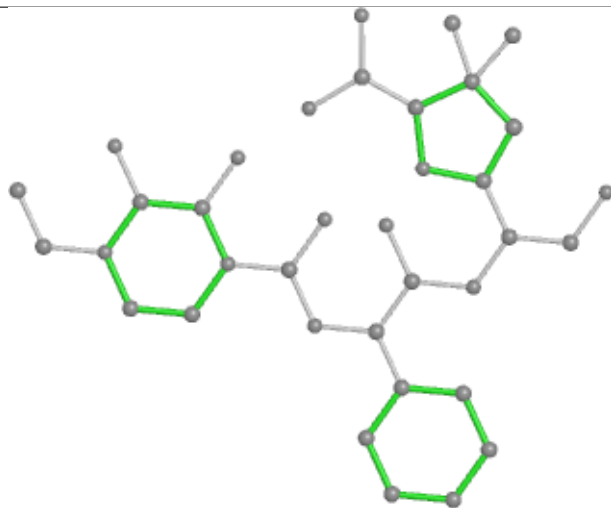
Bond lengths



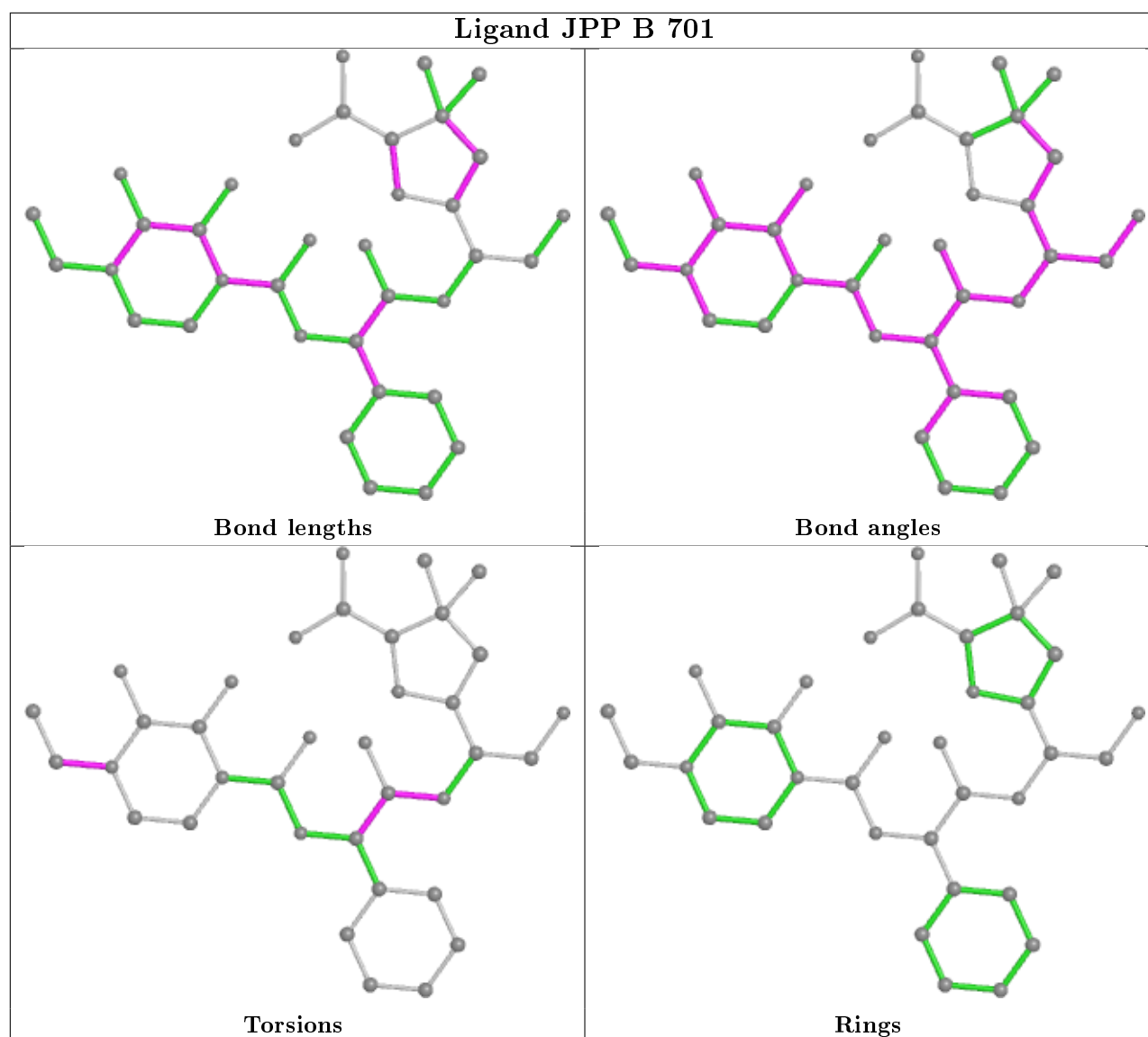
Bond angles



Torsions



Rings



## 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 [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	325/346 (93%)	-0.24	2 (0%) 89 91	3, 9, 21, 112	0
1	B	341/346 (98%)	-0.01	10 (2%) 51 59	7, 17, 41, 85	0
All	All	666/692 (96%)	-0.12	12 (1%) 68 74	3, 13, 37, 112	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	472	GLY	12.1
1	B	531	SER	5.1
1	B	530	GLY	3.5
1	B	532	SER	3.3
1	B	534	SER	3.1
1	B	479	MET	3.0
1	B	427	ARG	2.8
1	B	467	TYR	2.6
1	B	529	VAL	2.6
1	B	478	VAL	2.4
1	B	533	GLY	2.4
1	A	427	ARG	2.1

### 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 ⓘ

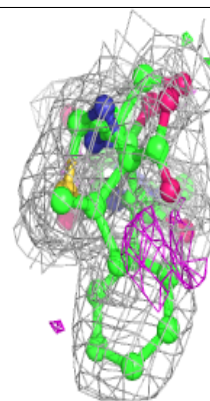
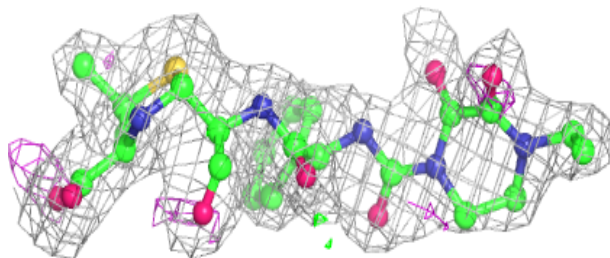
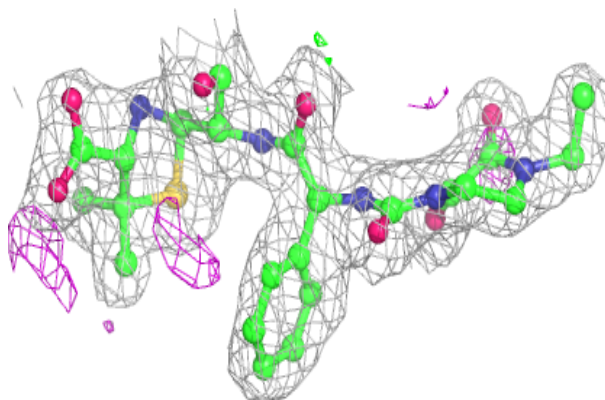
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	JPP	A	701	36/36	0.92	0.15	13,17,19,20	0
2	JPP	B	701	36/36	0.93	0.14	13,17,22,23	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

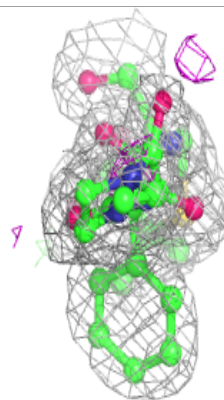
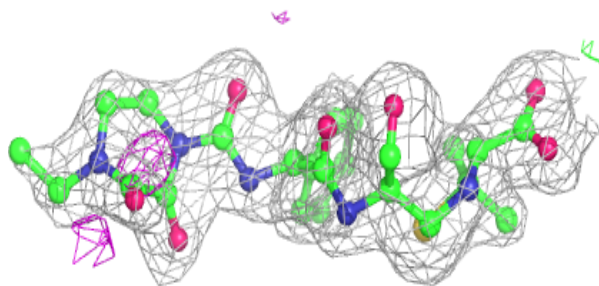
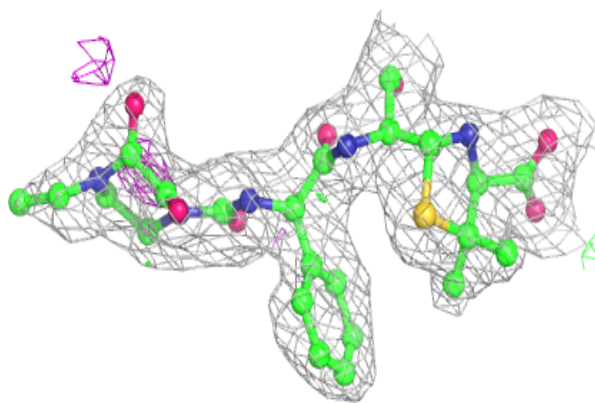
### Electron density around JPP A 701:

2mF<sub>o</sub>-DF<sub>c</sub> (at 0.7 rmsd) in gray  
mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around JPP B 701:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



## 6.5 Other polymers [i](#)

There are no such residues in this entry.