



# Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 07:31 pm BST

PDB ID : 3QSB  
Title : Structure of E. coli polIIIbeta with (Z)-5-(1-((4'-Fluorobiphenyl-4-yl)methoxymino)butyl)-2,2-dimethyl-4,6-dioxocyclohexanecarbonitrile  
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Deposited on : 2011-02-20  
Resolution : 1.90 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

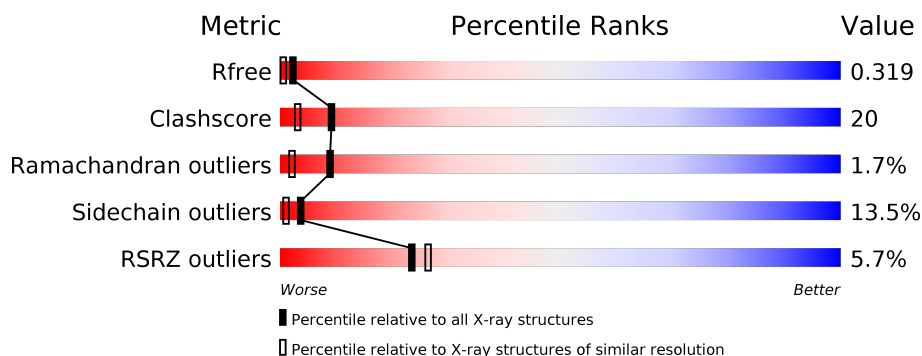
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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	366	<div> <div>5%</div> <div> <div></div> <div>65%</div> <div>26%</div> <div>7%</div> <div>.</div> </div> </div>
2	B	366	<div> <div>6%</div> <div> <div></div> <div>63%</div> <div>30%</div> <div>6%</div> <div>.</div> </div> </div>

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6414 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.

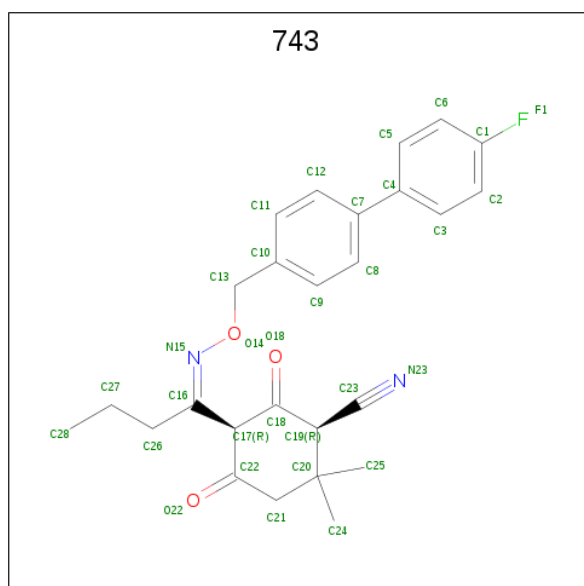
- Molecule 1 is a protein called DNA polymerase III subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	366	Total	C	N	O	S	1	17	0
			3002	1876	534	571	21			

- Molecule 2 is a protein called DNA polymerase III subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	366	Total	C	N	O	S	0	15	0
			2983	1867	532	562	22			

- Molecule 3 is (1R,5R)-5-[(1Z)-N-[(4'-fluorobiphenyl-4-yl)methoxy]butanimidoyl]-2,2-dimethyl-4,6-dioxocyclohexanecarbonitrile (three-letter code: 743) (formula: C<sub>26</sub>H<sub>27</sub>FN<sub>2</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			32	26	1	2	3		
3	B	1	Total	C	F	N	O	0	0
			32	26	1	2	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	180	Total 180	O 180	0	0
4	B	185	Total 185	O 185	0	0

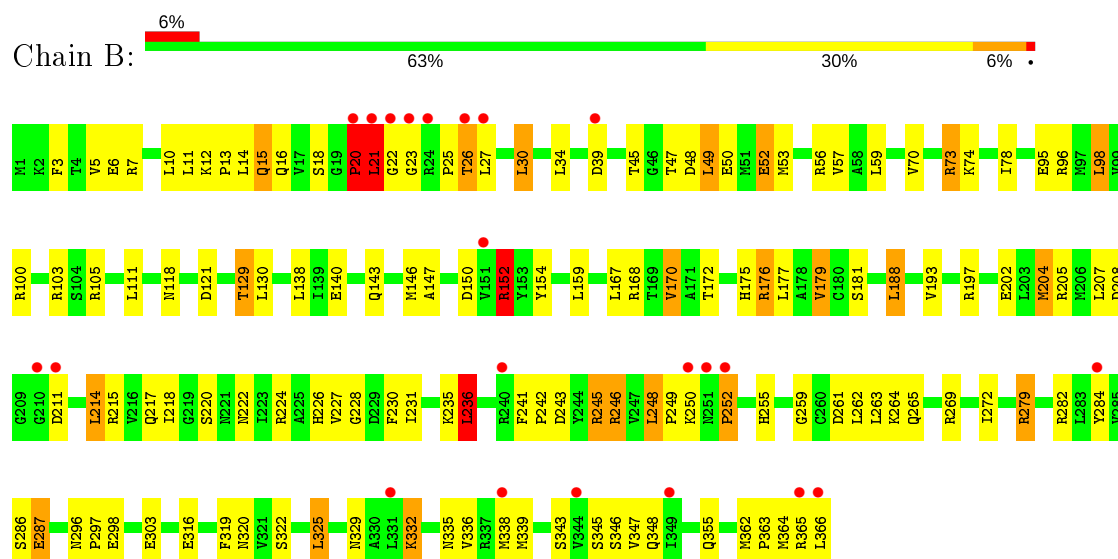
### 3 Residue-property plots

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: DNA polymerase III subunit beta



#### • Molecule 2: DNA polymerase III subunit beta



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	79.18 Å 66.52 Å 83.12 Å 90.00° 115.86° 90.00°	Depositor
Resolution (Å)	24.69 – 1.90 24.35 – 1.90	Depositor EDS
% Data completeness (in resolution range)	93.7 (24.69-1.90) 93.7 (24.35-1.90)	Depositor EDS
$R_{merge}$	0.24	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.23 (at 1.89 Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.242 , 0.317 0.245 , 0.319	Depositor DCC
$R_{free}$ test set	2902 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	18.4	Xtriage
Anisotropy	0.079	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 58.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for h,-k,-h-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6414	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 61.88 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.2260e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CSO, 743

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	1.07	3/3051 (0.1%)	1.17	24/4123 (0.6%)
2	B	1.05	0/3025	1.12	19/4086 (0.5%)
All	All	1.06	3/6076 (0.0%)	1.15	43/8209 (0.5%)

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
1	A	1	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	303	GLU	CD-OE2	5.48	1.31	1.25
1	A	170	VAL	CB-CG2	5.47	1.64	1.52
1	A	173	ASP	CA-CB	5.01	1.65	1.53

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	173	ASP	CB-CG-OD2	10.84	128.05	118.30
1	A	173	ASP	CB-CG-OD1	-10.77	108.60	118.30
1	A	176	ARG	NE-CZ-NH2	-10.61	114.99	120.30
1	A	179	VAL	CB-CA-C	-9.95	92.50	111.40
2	B	245	ARG	NE-CZ-NH1	9.57	125.09	120.30
2	B	245	ARG	NE-CZ-NH2	-9.07	115.77	120.30
2	B	152	ARG	NE-CZ-NH2	-8.63	115.99	120.30
2	B	188	LEU	CA-CB-CG	8.31	134.40	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	188	LEU	CA-CB-CG	8.13	133.99	115.30
1	A	176	ARG	CG-CD-NE	-7.72	95.58	111.80
2	B	56	ARG	NE-CZ-NH2	-7.39	116.61	120.30
2	B	279	ARG	NE-CZ-NH2	-7.31	116.64	120.30
1	A	176	ARG	NE-CZ-NH1	6.94	123.77	120.30
2	B	279	ARG	NE-CZ-NH1	6.84	123.72	120.30
2	B	332	LYS	N-CA-C	-6.80	92.64	111.00
1	A	138	LEU	CA-CB-CG	6.46	130.16	115.30
1	A	98	LEU	CB-CG-CD1	6.40	121.88	111.00
1	A	340	LEU	CA-CB-CG	6.38	129.97	115.30
1	A	197[A]	ARG	NE-CZ-NH2	-6.28	117.16	120.30
1	A	197[B]	ARG	NE-CZ-NH2	-6.28	117.16	120.30
1	A	21	LEU	CA-CB-CG	6.19	129.54	115.30
1	A	279	ARG	NE-CZ-NH1	6.18	123.39	120.30
1	A	366	LEU	CA-CB-CG	6.17	129.50	115.30
2	B	325	LEU	CA-CB-CG	6.17	129.50	115.30
2	B	98	LEU	CB-CG-CD1	6.03	121.26	111.00
1	A	138	LEU	CB-CG-CD1	5.88	121.00	111.00
1	A	179	VAL	CG1-CB-CG2	5.83	120.22	110.90
2	B	208	ASP	CB-CG-OD2	5.66	123.40	118.30
1	A	204	MET	CG-SD-CE	-5.66	91.14	100.20
1	A	137[A]	ARG	NE-CZ-NH1	5.55	123.08	120.30
1	A	137[B]	ARG	NE-CZ-NH1	5.55	123.08	120.30
1	A	188	LEU	CB-CG-CD1	5.37	120.12	111.00
2	B	7	ARG	NE-CZ-NH1	5.37	122.98	120.30
1	A	179	VAL	CA-CB-CG2	5.29	118.83	110.90
2	B	236	LEU	CB-CG-CD1	5.20	119.84	111.00
2	B	7	ARG	NE-CZ-NH2	-5.19	117.71	120.30
2	B	248	LEU	CB-CG-CD1	5.10	119.66	111.00
2	B	176	ARG	NE-CZ-NH2	-5.09	117.75	120.30
2	B	152	ARG	NE-CZ-NH1	5.07	122.83	120.30
2	B	105[A]	ARG	NE-CZ-NH1	5.03	122.82	120.30
2	B	105[B]	ARG	NE-CZ-NH1	5.03	122.82	120.30
1	A	197[A]	ARG	NE-CZ-NH1	5.03	122.81	120.30
1	A	197[B]	ARG	NE-CZ-NH1	5.03	122.81	120.30

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	197[B]	ARG	CA

All (1) planarity outliers are listed below:



Mol	Chain	Res	Type	Group
1	A	23	GLY	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	3002	0	3003	123	7
2	B	2983	0	2994	122	2
3	A	32	0	27	1	0
3	B	32	0	27	2	0
4	A	180	0	0	19	1
4	B	185	0	0	14	6
All	All	6414	0	6051	243	8

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 20.

All (243) 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:340:LEU:HD13	1:A:347:VAL:HG21	1.33	1.10
2:B:16:GLN:HG3	2:B:53:MET:HE2	1.37	1.04
1:A:137[B]:ARG:HG3	1:A:137[B]:ARG:HH11	1.18	1.03
2:B:226[A]:HIS:CD2	2:B:231:ILE:HG12	1.97	1.00
1:A:272:ILE:HG12	4:B:529:HOH:O	1.62	0.99
2:B:16:GLN:HG3	2:B:53:MET:CE	1.95	0.96
1:A:39[A]:ASP:CG	1:A:40:GLY:N	2.17	0.93
1:A:340:LEU:CD1	1:A:347:VAL:HG21	1.99	0.92
2:B:226[A]:HIS:CE1	2:B:228:GLY:O	2.25	0.90
1:A:282[A]:ARG:HH22	1:A:316:GLU:HG2	1.35	0.90
2:B:140[A]:GLU:HG3	2:B:204:MET:CE	2.02	0.90
2:B:246[A]:ARG:HA	2:B:246[A]:ARG:NE	1.87	0.89
1:A:135[B]:MET:HE2	1:A:135[B]:MET:HA	1.53	0.89
1:A:340:LEU:HD13	1:A:347:VAL:CG2	2.02	0.89
1:A:282[A]:ARG:NH2	1:A:316:GLU:HG2	1.88	0.88
2:B:140[A]:GLU:HG3	2:B:204:MET:HE1	1.53	0.88
2:B:47:THR:HB	4:B:368:HOH:O	1.74	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:226[A]:HIS:HD2	2:B:231:ILE:HG12	1.37	0.86
1:A:140[B]:GLU:HG3	1:A:204:MET:HE1	1.58	0.85
2:B:147:ALA:O	2:B:197:ARG:NH2	2.11	0.84
2:B:282:ARG:HD2	2:B:316:GLU:OE1	1.79	0.83
1:A:140[B]:GLU:HG3	1:A:204:MET:CE	2.09	0.82
1:A:77:ASP:OD1	1:A:80:ARG:NH2	2.13	0.81
1:A:137[B]:ARG:HG3	1:A:137[B]:ARG:NH1	1.96	0.79
1:A:226:HIS:HD2	1:A:231:ILE:HG12	1.48	0.79
1:A:170:VAL:HG22	1:A:179:VAL:HG13	1.65	0.78
4:A:519:HOH:O	2:B:103:ARG:HD3	1.84	0.77
2:B:23:GLY:HA2	4:B:549:HOH:O	1.85	0.77
2:B:143:GLN:HG3	2:B:146:MET:CE	2.15	0.77
2:B:118:ASN:OD1	4:B:391:HOH:O	2.03	0.74
1:A:27:LEU:HB3	1:A:30:LEU:HD22	1.69	0.74
1:A:214:LEU:HD13	1:A:227:VAL:CG2	2.19	0.72
1:A:226:HIS:CE1	1:A:228:GLY:O	2.42	0.72
2:B:140[A]:GLU:OE1	4:B:436:HOH:O	2.08	0.72
1:A:250:LYS:HG3	1:A:251:ASN:ND2	2.05	0.72
2:B:243:ASP:HB3	2:B:246[B]:ARG:HG3	1.73	0.71
2:B:346:SER:HA	2:B:363:PRO:HD3	1.71	0.71
2:B:252:PRO:O	2:B:255:HIS:HE1	1.74	0.71
2:B:303[B]:GLU:OE1	4:B:509:HOH:O	2.08	0.71
1:A:91:GLN:OE1	4:A:382:HOH:O	2.08	0.70
1:A:100:ARG:HD3	4:A:426:HOH:O	1.91	0.70
1:A:300:GLU:OE1	2:B:74:LYS:NZ	2.24	0.70
1:A:136:LYS:O	1:A:140[A]:GLU:HG3	1.90	0.70
2:B:319:PHE:CZ	2:B:347:VAL:HG12	2.25	0.70
1:A:137[A]:ARG:NH2	4:A:392:HOH:O	2.24	0.70
2:B:143:GLN:HG3	2:B:146:MET:HE2	1.74	0.69
2:B:159:LEU:HD23	2:B:170[A]:VAL:HG13	1.73	0.69
1:A:123[A]:GLN:HG2	4:A:424:HOH:O	1.92	0.69
1:A:118:ASN:ND2	4:A:381:HOH:O	2.25	0.69
2:B:319:PHE:CE1	2:B:347:VAL:HG12	2.28	0.69
2:B:319:PHE:CZ	2:B:347:VAL:CG1	2.75	0.69
2:B:226[A]:HIS:CD2	2:B:231:ILE:CG1	2.74	0.68
2:B:249:PRO:HD2	2:B:348:GLN:HE21	1.57	0.68
1:A:143:GLN:O	1:A:146:MET:HE2	1.94	0.68
1:A:32:ASN:HB3	1:A:69:THR:HG23	1.76	0.68
1:A:282[A]:ARG:NH1	1:A:316:GLU:OE2	2.29	0.66
1:A:69:THR:HB	1:A:111:LEU:O	1.95	0.66
1:A:135[B]:MET:HA	1:A:135[B]:MET:CE	2.26	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:252:PRO:O	2:B:255:HIS:CE1	2.49	0.66
2:B:252:PRO:HG2	2:B:339[B]:MET:CE	2.26	0.65
1:A:24:ARG:O	1:A:26:THR:N	2.30	0.65
2:B:143:GLN:O	2:B:146:MET:HE2	1.97	0.65
2:B:226[A]:HIS:HE1	2:B:228:GLY:O	1.76	0.65
1:A:39[A]:ASP:OD1	1:A:40:GLY:N	2.26	0.64
1:A:346:SER:HA	1:A:363:PRO:HD3	1.78	0.64
1:A:214:LEU:HD11	1:A:225:ALA:HB1	1.78	0.63
2:B:48:ASP:O	2:B:49:LEU:HB2	1.97	0.63
1:A:16:GLN:HG2	1:A:230:PHE:CD1	2.33	0.63
2:B:320:ASN:HB2	2:B:364:MET:CE	2.29	0.63
2:B:320:ASN:HB2	2:B:364:MET:HE1	1.80	0.63
1:A:299:GLN:HB2	2:B:96:ARG:NH2	2.14	0.63
2:B:226[A]:HIS:HD2	2:B:231:ILE:CG1	2.09	0.62
1:A:226:HIS:HD2	1:A:231:ILE:CG1	2.11	0.62
1:A:250:LYS:HE2	1:A:251:ASN:HD21	1.64	0.62
1:A:287[B]:GLU:OE2	1:A:311:SER:HA	1.99	0.62
2:B:143:GLN:HG3	2:B:146:MET:HE1	1.81	0.61
2:B:296:ASN:HB2	2:B:297:PRO:CD	2.29	0.61
1:A:214:LEU:CD1	1:A:227:VAL:CG2	2.78	0.61
2:B:241:PHE:CD1	2:B:242:PRO:HD2	2.36	0.61
2:B:332:LYS:NZ	2:B:332:LYS:HB2	2.15	0.61
1:A:135[B]:MET:HG3	1:A:214:LEU:CD2	2.31	0.60
1:A:25:PRO:C	1:A:27:LEU:H	2.04	0.60
2:B:16:GLN:HB3	2:B:230:PHE:CE1	2.37	0.60
1:A:340:LEU:HA	1:A:347:VAL:HG22	1.83	0.60
2:B:217:GLN:NE2	2:B:224:ARG:HH11	1.98	0.60
1:A:96:ARG:NH2	1:A:109:SER:HB3	2.17	0.59
1:A:173:ASP:HB3	1:A:175:HIS:H	1.66	0.59
2:B:246[A]:ARG:CA	2:B:246[A]:ARG:NE	2.59	0.58
2:B:129:THR:O	2:B:130:LEU:HB3	2.03	0.58
2:B:50:GLU:HB2	4:B:540:HOH:O	2.03	0.58
2:B:246[A]:ARG:HA	2:B:246[A]:ARG:CZ	2.33	0.58
1:A:214:LEU:CD1	1:A:227:VAL:HG23	2.34	0.58
1:A:296:ASN:HB2	1:A:297:PRO:CD	2.34	0.57
2:B:73[A]:ARG:NH1	4:B:549:HOH:O	2.37	0.57
1:A:201:ILE:HD12	4:A:472:HOH:O	2.03	0.57
1:A:132[A]:GLN:NE2	1:A:212:ASN:O	2.35	0.57
2:B:20:PRO:O	2:B:22:GLY:N	2.38	0.56
2:B:150:ASP:OD1	2:B:152:ARG:CD	2.53	0.56
1:A:53:MET:CE	1:A:206:MET:HG2	2.34	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:140[A]:GLU:HG3	2:B:204:MET:HE3	1.84	0.56
2:B:362[B]:MET:CE	3:B:367:743:H2	2.35	0.56
2:B:224:ARG:HD3	2:B:226[B]:HIS:CE1	2.40	0.56
2:B:286:SER:O	2:B:287:GLU:C	2.42	0.55
1:A:250:LYS:HE2	1:A:251:ASN:ND2	2.21	0.55
1:A:282[A]:ARG:NH2	1:A:316:GLU:CG	2.66	0.55
2:B:150:ASP:OD1	2:B:152:ARG:HD2	2.06	0.55
2:B:25:PRO:O	2:B:26:THR:CB	2.54	0.55
1:A:176:ARG:NH2	1:A:326:ASP:OD2	2.35	0.55
1:A:255:HIS:HE1	4:A:461:HOH:O	1.90	0.54
1:A:168:ARG:HD2	1:A:179:VAL:CG1	2.38	0.54
2:B:74:LYS:O	2:B:78:ILE:HG13	2.07	0.54
1:A:12[B]:LYS:HB3	1:A:13:PRO:HD3	1.89	0.54
2:B:224:ARG:CD	2:B:226[B]:HIS:CE1	2.91	0.54
1:A:24:ARG:C	1:A:26:THR:N	2.61	0.53
2:B:18:SER:O	2:B:20:PRO:HD3	2.08	0.53
2:B:263:LEU:HD21	2:B:338[A]:MET:CE	2.38	0.53
1:A:247:VAL:O	4:A:520:HOH:O	2.19	0.53
1:A:43:SER:OG	1:A:56[A]:ARG:NH1	2.42	0.53
2:B:20:PRO:HD2	2:B:205[A]:ARG:NH1	2.24	0.52
1:A:226:HIS:CD2	1:A:231:ILE:HG12	2.37	0.52
2:B:259:GLY:HA2	2:B:335:ASN:HD22	1.74	0.52
2:B:175:HIS:O	2:B:362[A]:MET:HG2	2.09	0.52
1:A:282[A]:ARG:CZ	1:A:316:GLU:CD	2.79	0.51
2:B:140[A]:GLU:CG	2:B:204:MET:CE	2.85	0.51
1:A:260:CYS:SG	4:A:433:HOH:O	2.29	0.51
1:A:366:LEU:HA	4:A:487:HOH:O	2.10	0.51
2:B:222:ASN:N	2:B:236:LEU:HD22	2.25	0.51
2:B:252:PRO:HG2	2:B:339[B]:MET:HE2	1.92	0.51
2:B:252:PRO:HG2	2:B:339[B]:MET:HE3	1.91	0.51
1:A:317:ILE:HD11	1:A:363:PRO:HB3	1.91	0.51
1:A:135[B]:MET:CA	1:A:135[B]:MET:HE2	2.34	0.51
1:A:25:PRO:HB2	1:A:30:LEU:HB2	1.92	0.51
1:A:53:MET:HE1	1:A:206:MET:HG2	1.92	0.51
2:B:319:PHE:CZ	2:B:347:VAL:HG11	2.46	0.50
2:B:168:ARG:HD2	2:B:179:VAL:HG23	1.93	0.49
1:A:222:ASN:N	1:A:236:LEU:HD22	2.27	0.49
1:A:16:GLN:CB	1:A:230:PHE:CE1	2.95	0.49
1:A:303:GLU:HG2	1:A:304:GLU:N	2.25	0.49
2:B:27:LEU:O	2:B:30:LEU:HB2	2.11	0.49
2:B:12:LYS:HB3	2:B:13:PRO:HD3	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:168:ARG:HD2	2:B:179:VAL:CG2	2.43	0.48
1:A:205[B]:ARG:NH1	4:A:512:HOH:O	2.00	0.48
2:B:282:ARG:NH1	4:B:539:HOH:O	2.32	0.48
2:B:332:LYS:HB2	2:B:332:LYS:HZ3	1.76	0.48
1:A:25:PRO:C	1:A:27:LEU:N	2.65	0.48
1:A:16:GLN:HB3	1:A:230:PHE:CE1	2.49	0.48
1:A:30:LEU:HD11	1:A:49:LEU:HD13	1.95	0.47
2:B:215:ARG:HB3	2:B:226[B]:HIS:HB2	1.96	0.47
1:A:135[B]:MET:HB3	1:A:207:LEU:HD21	1.97	0.47
2:B:279:ARG:HD3	2:B:322:SER:OG	2.15	0.47
1:A:137[B]:ARG:CG	1:A:137[B]:ARG:NH1	2.71	0.47
1:A:226:HIS:CD2	1:A:231:ILE:CG1	2.96	0.47
2:B:143:GLN:CG	2:B:146:MET:HE2	2.45	0.47
2:B:30:LEU:HD11	2:B:49:LEU:HD13	1.97	0.47
1:A:50:GLU:OE1	1:A:50:GLU:HA	2.14	0.47
2:B:20:PRO:HG3	4:B:490:HOH:O	2.15	0.47
2:B:246[A]:ARG:HA	2:B:246[A]:ARG:HE	1.75	0.47
1:A:144:PHE:CD2	1:A:326:ASP:HB3	2.50	0.47
2:B:140[A]:GLU:CG	2:B:204:MET:HE3	2.45	0.46
1:A:132[B]:GLN:NE2	1:A:208:ASP:OD1	2.46	0.46
1:A:249:PRO:HB2	1:A:252:PRO:HG3	1.97	0.46
1:A:289[B]:GLN:CG	4:A:369:HOH:O	2.62	0.46
2:B:202:GLU:OE1	2:B:205[B]:ARG:HD3	2.15	0.46
1:A:32:ASN:CB	1:A:69:THR:HG23	2.45	0.46
2:B:263:LEU:HD21	2:B:338[A]:MET:HE2	1.96	0.46
1:A:147:ALA:O	1:A:197[A]:ARG:NH1	2.49	0.46
2:B:249:PRO:HD2	2:B:348:GLN:NE2	2.28	0.45
2:B:25:PRO:HD3	4:B:541:HOH:O	2.15	0.45
1:A:53:MET:HE3	1:A:206:MET:HG2	1.98	0.45
1:A:48:ASP:O	1:A:49:LEU:HB2	2.17	0.45
2:B:241:PHE:CG	2:B:242:PRO:HD2	2.51	0.45
2:B:362[B]:MET:HE1	3:B:367:743:H2	1.99	0.45
1:A:365:ARG:O	1:A:366:LEU:HB3	2.16	0.45
2:B:243:ASP:OD1	2:B:245:ARG:HB2	2.17	0.45
2:B:224:ARG:HD2	2:B:226[B]:HIS:HE1	1.82	0.44
2:B:3:PHE:HD2	2:B:5:VAL:HG23	1.82	0.44
1:A:280:GLY:HA3	1:A:364:MET:HE1	1.98	0.44
1:A:33:LEU:HG	1:A:72:ALA:CB	2.47	0.44
1:A:48:ASP:OD1	1:A:48:ASP:C	2.55	0.44
2:B:264[A]:LYS:HD2	2:B:329:ASN:OD1	2.17	0.44
1:A:184:ILE:HD11	1:A:188:LEU:HD21	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:269:ARG:HA	2:B:272:ILE:HD12	1.99	0.44
2:B:346:SER:CA	2:B:363:PRO:HD3	2.46	0.44
1:A:202:GLU:OE1	1:A:205[A]:ARG:NH1	2.50	0.43
2:B:261:ASP:O	2:B:265:GLN:HG2	2.18	0.43
1:A:21:LEU:HG	1:A:30:LEU:O	2.18	0.43
1:A:316:GLU:O	1:A:343:SER:HB2	2.18	0.43
1:A:140[B]:GLU:CG	1:A:204:MET:HE1	2.41	0.43
2:B:284:TYR:CZ	2:B:316:GLU:OE1	2.72	0.43
2:B:332:LYS:NZ	4:B:432:HOH:O	2.48	0.43
1:A:282[A]:ARG:NH1	1:A:366:LEU:OXT	2.52	0.43
2:B:217:GLN:HE22	2:B:224:ARG:HH11	1.66	0.43
2:B:263:LEU:HD21	2:B:338[A]:MET:HE3	2.01	0.43
1:A:100:ARG:NH1	4:A:382:HOH:O	2.51	0.42
1:A:26:THR:O	1:A:28:PRO:HD3	2.18	0.42
1:A:302:ALA:HB1	4:A:403:HOH:O	2.18	0.42
1:A:19:GLY:HA3	4:A:512:HOH:O	2.19	0.42
2:B:172:THR:HB	2:B:177:LEU:HD12	2.01	0.42
2:B:52:GLU:HB2	4:B:368:HOH:O	2.19	0.42
2:B:50:GLU:O	2:B:235:LYS:HG2	2.19	0.42
2:B:16:GLN:CG	2:B:53:MET:CE	2.83	0.42
1:A:154:TYR:HB2	1:A:194:ILE:HD13	2.01	0.42
2:B:193:VAL:HG11	2:B:218:ILE:CG2	2.50	0.42
1:A:316:GLU:CD	1:A:316:GLU:C	2.78	0.42
1:A:304:GLU:HG2	4:A:403:HOH:O	2.19	0.42
2:B:21:LEU:HD13	2:B:21:LEU:O	2.20	0.42
2:B:152:ARG:HB3	2:B:154:TYR:CZ	2.55	0.42
2:B:21:LEU:CD1	2:B:21:LEU:O	2.67	0.42
2:B:16:GLN:CB	2:B:230:PHE:CE1	3.03	0.42
1:A:96:ARG:NH2	2:B:298:GLU:O	2.51	0.42
1:A:153:TYR:HD1	1:A:153:TYR:H	1.67	0.42
1:A:297:PRO:C	1:A:299:GLN:H	2.22	0.42
2:B:296:ASN:HB2	2:B:297:PRO:HD3	1.99	0.41
2:B:319:PHE:CE1	2:B:347:VAL:CG1	2.98	0.41
1:A:365:ARG:O	1:A:366:LEU:CB	2.68	0.41
1:A:137[A]:ARG:HD3	1:A:137[A]:ARG:C	2.40	0.41
1:A:202:GLU:OE1	1:A:205[A]:ARG:HD2	2.21	0.41
2:B:34:LEU:HB3	2:B:45:THR:HB	2.02	0.41
2:B:15:GLN:O	2:B:205[B]:ARG:NH2	2.53	0.41
2:B:345:SER:O	2:B:363:PRO:HG3	2.20	0.41
1:A:24:ARG:C	1:A:26:THR:H	2.24	0.41
2:B:146:MET:HE3	2:B:146:MET:HB2	1.90	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:45:THR:HG22	4:B:368:HOH:O	2.20	0.41
1:A:111:LEU:HD12	1:A:111:LEU:HA	1.93	0.41
1:A:255:HIS:CE1	4:A:461:HOH:O	2.71	0.41
1:A:252:PRO:O	1:A:255:HIS:NE2	2.54	0.41
2:B:217:GLN:HE21	2:B:224:ARG:HD2	1.86	0.41
2:B:214:LEU:HD13	2:B:227:VAL:HG22	2.03	0.41
1:A:16:GLN:HB3	1:A:230:PHE:HE1	1.84	0.41
1:A:286:SER:OG	1:A:289[B]:GLN:HG2	2.20	0.41
1:A:331:LEU:HD13	1:A:336[B]:VAL:HG12	2.03	0.41
2:B:202:GLU:OE1	2:B:205[A]:ARG:NH1	2.54	0.41
2:B:48:ASP:OD1	2:B:48:ASP:C	2.59	0.41
1:A:289[B]:GLN:HG2	4:A:369:HOH:O	2.21	0.41
2:B:150:ASP:OD1	2:B:152:ARG:HD3	2.21	0.40
3:A:367:743:H3	3:A:367:743:H8	1.82	0.40
2:B:246[B]:ARG:HH11	2:B:246[B]:ARG:HD3	1.76	0.40
2:B:25:PRO:O	2:B:26:THR:OG1	2.39	0.40
1:A:25:PRO:CB	1:A:30:LEU:HB2	2.51	0.40
1:A:32:ASN:ND2	1:A:71:PRO:HA	2.36	0.40
1:A:252:PRO:O	1:A:255:HIS:CD2	2.74	0.40
1:A:331:LEU:O	1:A:332:LYS:C	2.59	0.40

All (8) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:123[A]:GLN:NE2	4:B:526:HOH:O[1_655]	1.40	0.80
1:A:123[A]:GLN:OE1	4:B:528:HOH:O[1_655]	1.88	0.32
1:A:123[A]:GLN:CD	4:B:526:HOH:O[1_655]	1.98	0.22
1:A:250:LYS:O	4:B:485:HOH:O[1_656]	2.08	0.12
1:A:109:SER:OG	2:B:211:ASP:OD2[2_655]	2.10	0.10
1:A:123[A]:GLN:OE1	4:B:526:HOH:O[1_655]	2.10	0.10
1:A:39[A]:ASP:O	4:B:445:HOH:O[1_655]	2.13	0.07
2:B:39:ASP:OD1	4:A:439:HOH:O[1_455]	2.16	0.04

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	381/366 (104%)	361 (95%)	14 (4%)	6 (2%)	9	2
2	B	378/366 (103%)	356 (94%)	16 (4%)	6 (2%)	9	2
All	All	759/732 (104%)	717 (94%)	30 (4%)	12 (2%)	9	2

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	20	PRO
1	A	24	ARG
1	A	25	PRO
1	A	26	THR
2	B	20	PRO
2	B	21	LEU
2	B	26	THR
2	B	250	LYS
2	B	49	LEU
1	A	21	LEU
1	A	252	PRO
2	B	252	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	330/313 (105%)	283 (86%)	47 (14%)	3	1
2	B	326/312 (104%)	281 (86%)	45 (14%)	3	1
All	All	656/625 (105%)	564 (86%)	92 (14%)	4	1

All (92) residues with a non-rotameric sidechain are listed below:



Mol	Chain	Res	Type
1	A	6[A]	GLU
1	A	6[B]	GLU
1	A	10	LEU
1	A	11	LEU
1	A	15	GLN
1	A	28	PRO
1	A	30	LEU
1	A	32	ASN
1	A	39[A]	ASP
1	A	39[B]	ASP
1	A	42	LEU
1	A	49	LEU
1	A	52	GLU
1	A	57	VAL
1	A	60	VAL
1	A	62	PRO
1	A	69	THR
1	A	70	VAL
1	A	95	GLU
1	A	98	LEU
1	A	107	SER
1	A	111	LEU
1	A	118	ASN
1	A	123[A]	GLN
1	A	123[B]	GLN
1	A	137[A]	ARG
1	A	137[B]	ARG
1	A	138	LEU
1	A	152	ARG
1	A	173	ASP
1	A	176	ARG
1	A	181	SER
1	A	188	LEU
1	A	203	LEU
1	A	207	LEU
1	A	236	LEU
1	A	248	LEU
1	A	287[A]	GLU
1	A	287[B]	GLU
1	A	299	GLN
1	A	303	GLU
1	A	306	LEU
1	A	316	GLU

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Mol	Chain	Res	Type
1	A	343	SER
1	A	344	VAL
1	A	365	ARG
1	A	366	LEU
2	B	6	GLU
2	B	10	LEU
2	B	11	LEU
2	B	14	LEU
2	B	15	GLN
2	B	20	PRO
2	B	21	LEU
2	B	30	LEU
2	B	52	GLU
2	B	57	VAL
2	B	59	LEU
2	B	70	VAL
2	B	73[A]	ARG
2	B	73[B]	ARG
2	B	95	GLU
2	B	98	LEU
2	B	100	ARG
2	B	111	LEU
2	B	121	ASP
2	B	129	THR
2	B	138	LEU
2	B	152	ARG
2	B	167	LEU
2	B	170[A]	VAL
2	B	170[B]	VAL
2	B	176	ARG
2	B	179	VAL
2	B	181	SER
2	B	188	LEU
2	B	204	MET
2	B	207	LEU
2	B	214	LEU
2	B	220	SER
2	B	236	LEU
2	B	246[A]	ARG
2	B	246[B]	ARG
2	B	248	LEU
2	B	262	LEU

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Mol	Chain	Res	Type
2	B	287	GLU
2	B	325	LEU
2	B	336	VAL
2	B	343	SER
2	B	355	GLN
2	B	365	ARG
2	B	366	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	GLN
1	A	91	GLN
1	A	118	ASN
1	A	212	ASN
1	A	226	HIS
1	A	251	ASN
1	A	255	HIS
1	A	288	ASN
1	A	299	GLN
1	A	355	GLN
2	B	36	GLN
2	B	143	GLN
2	B	149	GLN
2	B	212	ASN
2	B	217	GLN
2	B	251	ASN
2	B	255	HIS
2	B	335	ASN
2	B	348	GLN
2	B	355	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	CSO	B	260	2	3,6,7	0.88	0	0,6,8	0.00	-

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	CSO	B	260	2	-	0/1/5/7	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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
3	743	B	367	-	29,34,34	1.24	4 (13%)	35,48,48	1.72	7 (20%)
3	743	A	367	-	29,34,34	1.42	5 (17%)	35,48,48	1.67	4 (11%)

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	743	B	367	-	-	3/13/42/42	0/3/3/3
3	743	A	367	-	-	2/13/42/42	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	367	743	C7-C4	-3.58	1.40	1.49
3	B	367	743	C7-C4	-3.47	1.40	1.49
3	B	367	743	O14-N15	-2.70	1.37	1.42
3	A	367	743	O14-N15	-2.39	1.37	1.42
3	A	367	743	C23-N23	2.37	1.19	1.14
3	B	367	743	C26-C16	2.24	1.52	1.50
3	A	367	743	C6-C1	2.13	1.41	1.37
3	B	367	743	C19-C23	2.07	1.49	1.47
3	A	367	743	C12-C11	2.05	1.42	1.38

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	367	743	C13-O14-N15	6.92	116.60	107.90
3	B	367	743	C13-O14-N15	6.21	115.71	107.90
3	A	367	743	C25-C20-C19	4.84	114.76	109.29
3	B	367	743	C24-C20-C19	-3.87	104.91	109.29
3	B	367	743	C5-C6-C1	2.74	121.19	118.36
3	B	367	743	C25-C20-C19	2.47	112.08	109.29
3	B	367	743	O18-C18-C17	-2.44	118.56	121.96
3	A	367	743	C5-C6-C1	2.24	120.68	118.36
3	B	367	743	C6-C1-C2	-2.05	120.10	122.83
3	B	367	743	C17-C18-C19	2.01	119.70	114.61
3	A	367	743	C24-C20-C19	-2.00	107.02	109.29

There are no chirality outliers.

All (5) torsion outliers are listed below:

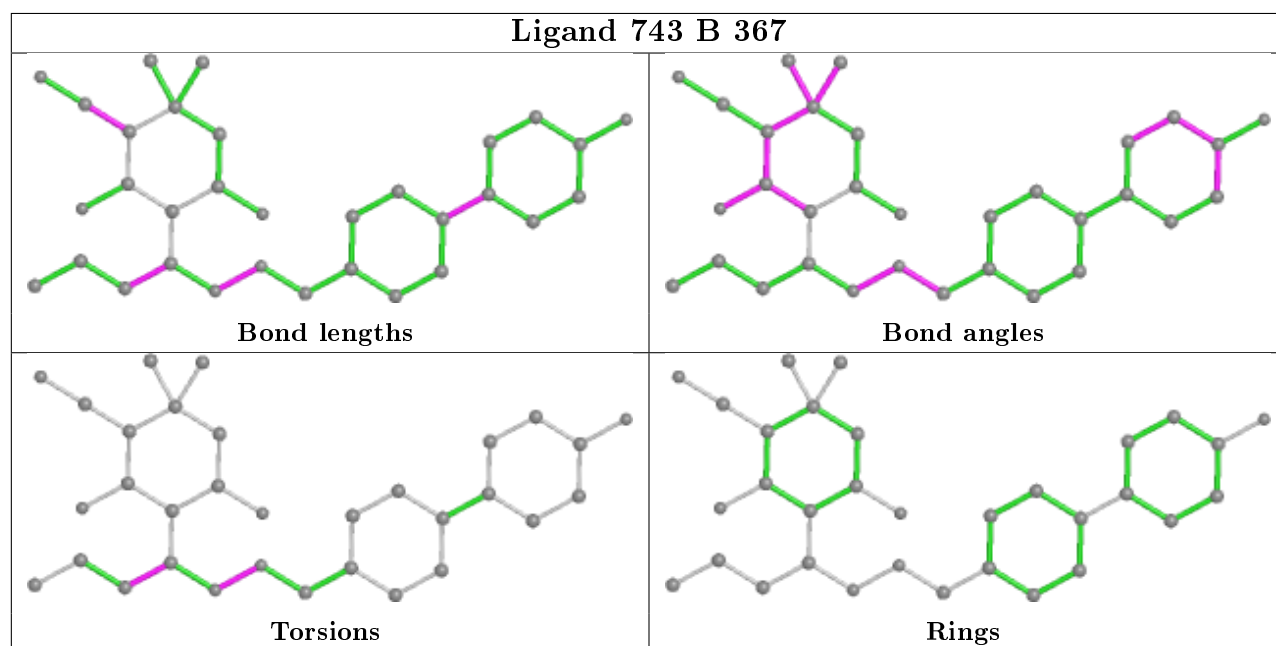
Mol	Chain	Res	Type	Atoms
3	B	367	743	C16-N15-O14-C13
3	A	367	743	C16-N15-O14-C13
3	B	367	743	N15-C16-C26-C27
3	B	367	743	C17-C16-C26-C27
3	A	367	743	C17-C16-C26-C27

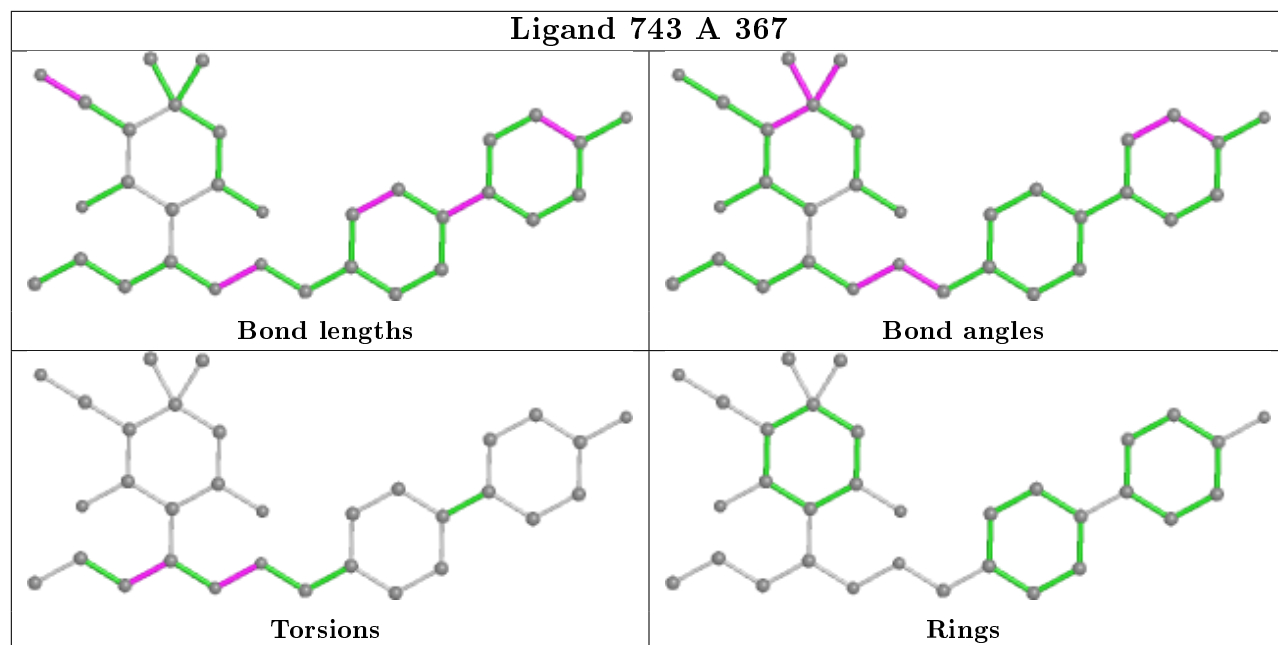
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	367	743	2	0
3	A	367	743	1	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.





## 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	366/366 (100%)	0.39	20 (5%)	25 28	11, 22, 40, 63	1 (0%)
2	B	365/366 (99%)	0.40	22 (6%)	21 24	11, 23, 41, 66	0
All	All	731/732 (99%)	0.40	42 (5%)	23 26	11, 22, 41, 66	1 (0%)

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	22	GLY	9.6
1	A	21	LEU	9.5
2	B	20	PRO	8.2
2	B	22	GLY	7.2
1	A	20	PRO	6.4
2	B	366	LEU	6.4
2	B	24	ARG	5.8
2	B	21	LEU	5.5
1	A	25	PRO	5.1
2	B	210	GLY	4.9
2	B	251	ASN	4.8
1	A	23	GLY	4.7
2	B	23	GLY	4.7
2	B	365	ARG	4.2
2	B	27	LEU	4.0
1	A	148	HIS	3.8
1	A	366	LEU	3.6
2	B	284	TYR	3.5
1	A	24	ARG	3.4
2	B	26	THR	3.3
1	A	284	TYR	3.3
1	A	26	THR	3.3
2	B	151	VAL	3.1
1	A	211	ASP	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	251	ASN	3.0
2	B	331	LEU	3.0
1	A	39[A]	ASP	2.9
2	B	250	LYS	2.8
1	A	365	ARG	2.8
2	B	39	ASP	2.6
2	B	211	ASP	2.6
1	A	60	VAL	2.5
1	A	209	GLY	2.5
1	A	355	GLN	2.4
1	A	298	GLU	2.4
2	B	338[A]	MET	2.3
2	B	240	ARG	2.3
2	B	252	PRO	2.2
1	A	17	VAL	2.1
1	A	344	VAL	2.0
2	B	344	VAL	2.0
2	B	349	ILE	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [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. 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	CSO	B	260	7/8	0.96	0.11	17,20,22,24	1

## 6.3 Carbohydrates [i](#)

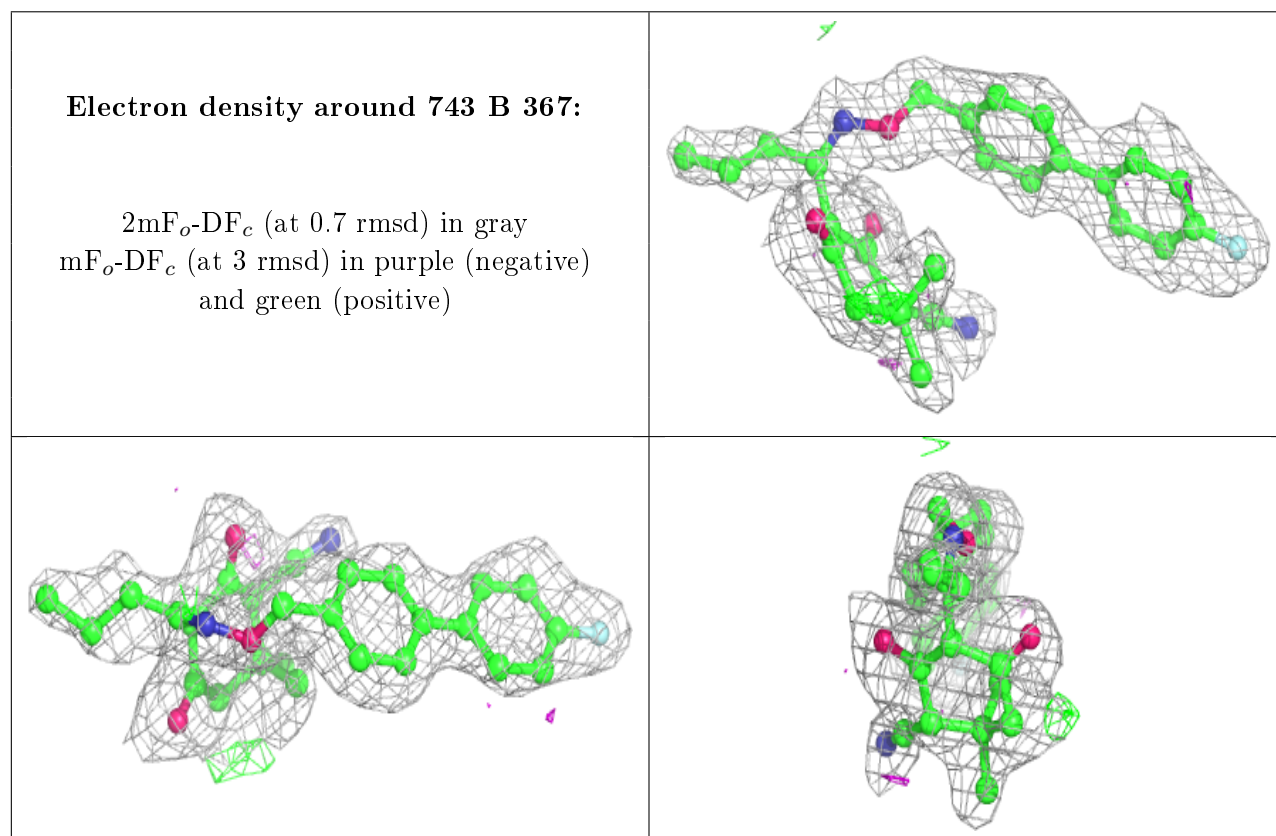
There are no carbohydrates in this entry.

## 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. 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.

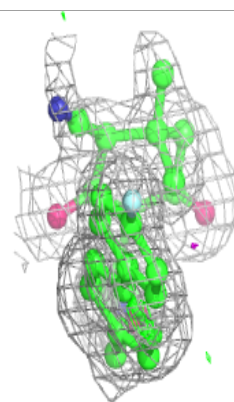
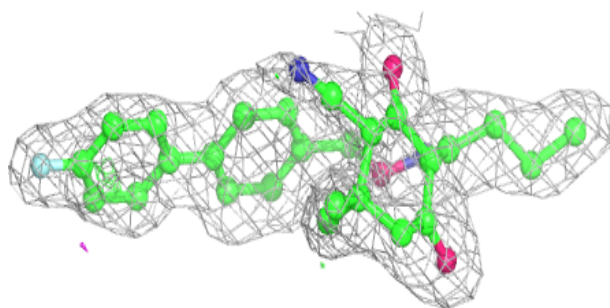
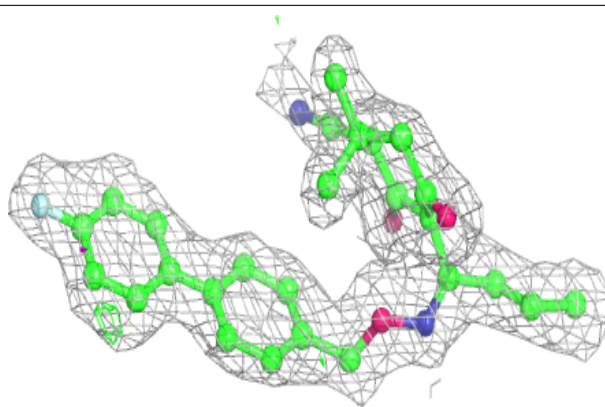
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	743	B	367	32/32	0.82	0.16	21,35,42,43	0
3	743	A	367	32/32	0.87	0.13	19,28,35,41	0

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 743 A 367:**

$2mF_o-DF_c$  (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.