



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 08:07 am BST

PDB ID : 4EMN
Title : Crystal structure of RpfB catalytic domain in complex with benzamidine
Authors : Ruggiero, A.; Marchant, J.; Squeglia, F.; Makarov, V.; De Simone, A.; Berisio, R.
Deposited on : 2012-04-12
Resolution : 1.17 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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
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

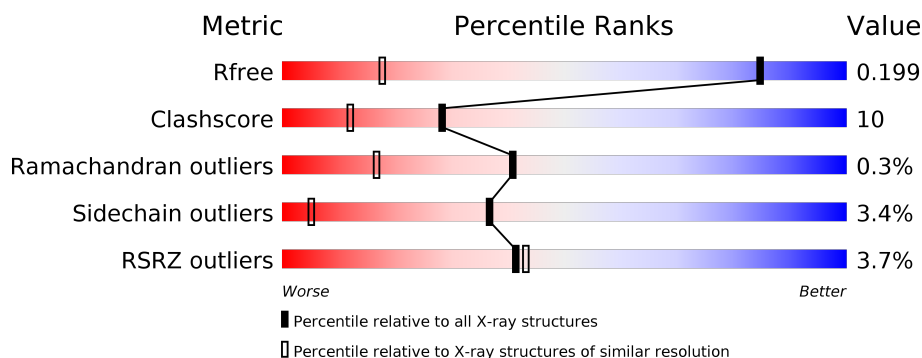
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.17 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	1123 (1.20-1.16)
Clashscore	141614	1182 (1.20-1.16)
Ramachandran outliers	138981	1134 (1.20-1.16)
Sidechain outliers	138945	1134 (1.20-1.16)
RSRZ outliers	127900	1102 (1.20-1.16)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria 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	81	<div> <div>%</div> <div> <div></div> <div>83%</div> <div>12%</div> <div>• •</div> </div> </div>
1	B	81	<div> <div>4%</div> <div> <div></div> <div>78%</div> <div>11%</div> <div>10%</div> <div>•</div> </div> </div>
1	C	81	<div> <div>4%</div> <div> <div></div> <div>89%</div> <div>6%</div> <div>• •</div> </div> </div>
1	D	81	<div> <div>6%</div> <div> <div></div> <div>84%</div> <div>12%</div> <div>• •</div> </div> </div>

2 Entry composition [i](#)

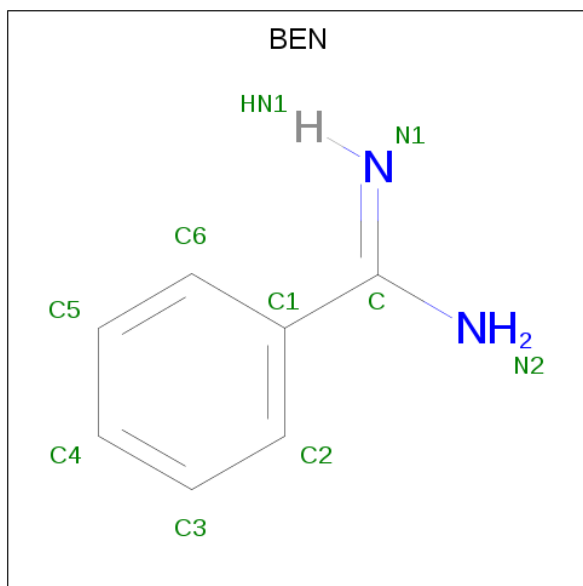
There are 4 unique types of molecules in this entry. The entry contains 3033 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Probable resuscitation-promoting factor rpfb.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	81	Total	C	N	O	S	0	0	0
			600	373	112	113	2			
1	B	80	Total	C	N	O	S	0	0	0
			602	374	114	112	2			
1	C	81	Total	C	N	O	S	0	0	0
			602	375	112	113	2			
1	D	80	Total	C	N	O	S	0	0	0
			598	373	111	112	2			

- Molecule 2 is BENZAMIDINE (three-letter code: BEN) (formula: C₇H₈N₂).



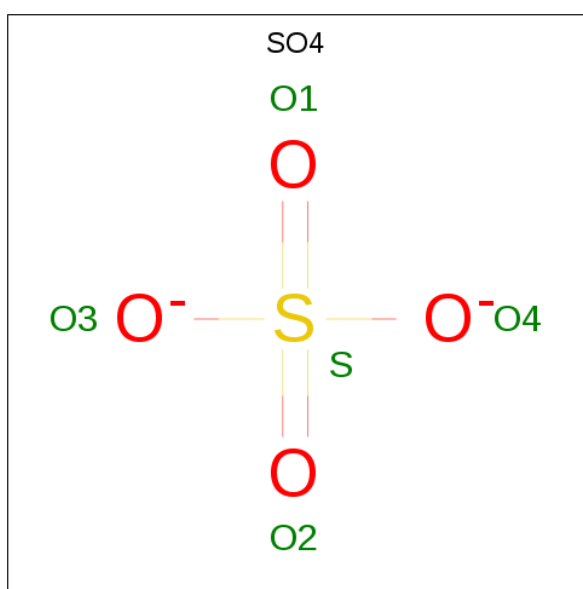
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	0	0
			9	7	2		
2	A	1	Total	C	N	0	0
			9	7	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	N	0	0
			9	7	2		
2	B	1	Total	C	N	0	0
			9	7	2		
2	C	1	Total	C	N	0	0
			9	7	2		
2	C	1	Total	C	N	0	0
			9	7	2		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	153	Total	O	0	0
			153	153		
4	B	158	Total	O	0	0
			158	158		
4	C	127	Total	O	0	0
			127	127		

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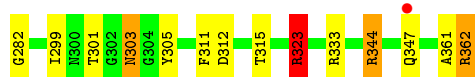
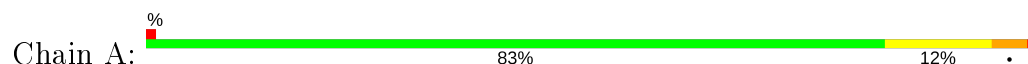
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	129	Total	O	0	0
			129	129		

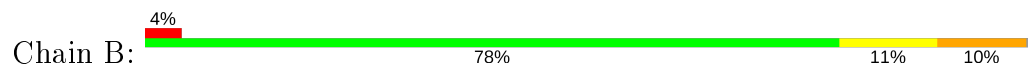
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

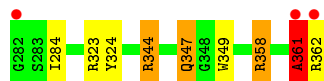
- Molecule 1: Probable resuscitation-promoting factor rpfb



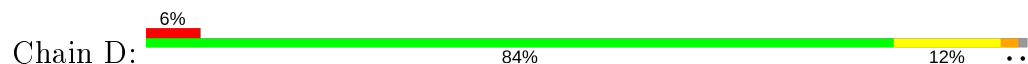
- Molecule 1: Probable resuscitation-promoting factor rpfb



- Molecule 1: Probable resuscitation-promoting factor rpfb



- Molecule 1: Probable resuscitation-promoting factor rpfb



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	42.55Å 50.73Å 66.42Å 90.00° 104.09° 90.00°	Depositor
Resolution (Å)	15.00 – 1.17 11.81 – 1.17	Depositor EDS
% Data completeness (in resolution range)	(Not available) (15.00-1.17) 87.6 (11.81-1.17)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	-0.12 (at 1.17Å)	Xtriage
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.127 , 0.188 0.144 , 0.199	Depositor DCC
R_{free} test set	4265 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	9.0	Xtriage
Anisotropy	0.656	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.43 , 99.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	3033	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.71% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: BEN, SO4

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	0.69	0/615	1.43	12/838 (1.4%)
1	B	0.73	0/617	1.93	17/840 (2.0%)
1	C	0.68	0/617	1.47	7/840 (0.8%)
1	D	0.62	0/613	1.27	5/835 (0.6%)
All	All	0.68	0/2462	1.54	41/3353 (1.2%)

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	D	0	1

There are no bond length outliers.

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	362	ARG	NE-CZ-NH2	-27.24	106.68	120.30
1	B	346	ARG	NE-CZ-NH2	-19.29	110.66	120.30
1	C	358	ARG	NE-CZ-NH2	-16.16	112.22	120.30
1	C	358	ARG	NE-CZ-NH1	11.11	125.85	120.30
1	B	362	ARG	NH1-CZ-NH2	10.51	130.96	119.40
1	C	344	ARG	NE-CZ-NH2	-10.41	115.10	120.30
1	B	358	ARG	NE-CZ-NH2	9.28	124.94	120.30
1	D	358	ARG	CD-NE-CZ	8.81	135.94	123.60
1	B	360	GLY	C-N-CA	8.81	143.73	121.70
1	A	323	ARG	NE-CZ-NH1	-8.55	116.02	120.30
1	A	311	PHE	CB-CG-CD2	-8.01	115.20	120.80
1	A	362	ARG	CA-CB-CG	-7.90	96.03	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	305	TYR	CB-CG-CD1	-7.86	116.28	121.00
1	A	361	ALA	C-N-CA	-7.75	102.33	121.70
1	B	346	ARG	NH1-CZ-NH2	7.63	127.79	119.40
1	B	362	ARG	CD-NE-CZ	-7.61	112.94	123.60
1	B	362	ARG	N-CA-CB	7.55	124.19	110.60
1	C	358	ARG	CD-NE-CZ	7.43	134.00	123.60
1	B	358	ARG	NE-CZ-NH1	-7.28	116.66	120.30
1	B	344	ARG	NE-CZ-NH1	6.99	123.80	120.30
1	B	361	ALA	N-CA-CB	-6.86	100.50	110.10
1	A	344	ARG	NE-CZ-NH1	6.60	123.60	120.30
1	B	360	GLY	N-CA-C	6.58	129.55	113.10
1	A	312	ASP	CB-CG-OD1	6.55	124.20	118.30
1	B	346	ARG	CD-NE-CZ	6.46	132.64	123.60
1	D	361	ALA	C-N-CA	6.33	137.53	121.70
1	B	358	ARG	CD-NE-CZ	6.28	132.39	123.60
1	D	286	ASP	CB-CG-OD2	6.05	123.75	118.30
1	B	334	GLU	OE1-CD-OE2	-5.87	116.26	123.30
1	A	305	TYR	CB-CG-CD2	5.85	124.51	121.00
1	B	324	TYR	CB-CG-CD1	-5.67	117.60	121.00
1	B	327	ARG	NE-CZ-NH1	5.54	123.07	120.30
1	C	361	ALA	O-C-N	5.50	131.50	122.70
1	A	362	ARG	CB-CA-C	5.49	121.37	110.40
1	A	333	ARG	NE-CZ-NH2	-5.48	117.56	120.30
1	D	346	ARG	NE-CZ-NH2	-5.44	117.58	120.30
1	C	347	GLN	CA-CB-CG	5.42	125.32	113.40
1	A	362	ARG	NE-CZ-NH2	-5.20	117.70	120.30
1	A	344	ARG	NE-CZ-NH2	-5.19	117.70	120.30
1	D	342	VAL	CG1-CB-CG2	-5.13	102.69	110.90
1	C	324	TYR	CB-CG-CD2	-5.03	117.98	121.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	361	ALA	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	600	0	548	10	0
1	B	602	0	555	23	0
1	C	602	0	552	4	0
1	D	598	0	549	9	0
2	A	18	0	14	1	0
2	B	18	0	14	0	0
2	C	18	0	14	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	153	0	0	5	0
4	B	158	0	0	15	0
4	C	127	0	0	3	0
4	D	129	0	0	5	0
All	All	3033	0	2246	46	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 10.

All (46) 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:362:ARG:HG2	4:B:7139:HOH:O	1.91	0.69
1:D:358:ARG:HD2	4:D:437:HOH:O	1.91	0.68
1:A:323:ARG:HD3	4:A:568:HOH:O	1.95	0.67
1:C:358:ARG:HD2	4:C:564:HOH:O	1.98	0.63
1:A:323:ARG:HD3	4:A:623:HOH:O	1.98	0.62
1:B:349:TRP:HE1	1:B:362:ARG:H	1.52	0.58
1:B:351:ALA:HB2	4:B:7137:HOH:O	2.06	0.56
1:B:334:GLU:HG2	4:B:7054:HOH:O	2.06	0.54
1:B:347:GLN:HG2	4:B:7061:HOH:O	2.07	0.54
1:B:362:ARG:HD3	4:B:7128:HOH:O	2.09	0.51
1:B:346:ARG:NE	4:B:7017:HOH:O	2.43	0.51
1:B:283:SER:HA	4:B:7036:HOH:O	2.11	0.51
1:B:323:ARG:HD2	4:B:7143:HOH:O	2.10	0.50
1:D:284:ILE:H	1:D:284:ILE:HD12	1.76	0.50
1:B:349:TRP:HE1	1:B:362:ARG:N	2.09	0.49
1:A:282:GLY:N	4:A:638:HOH:O	2.45	0.48
1:D:323:ARG:HD2	4:D:417:HOH:O	2.13	0.48
1:A:362:ARG:HB2	4:C:545:HOH:O	2.15	0.47
1:B:349:TRP:NE1	1:B:361:ALA:HA	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:349:TRP:HE1	1:C:362:ARG:C	2.18	0.47
1:A:362:ARG:NH1	4:A:535:HOH:O	2.48	0.46
1:B:349:TRP:CE2	1:B:361:ALA:HA	2.50	0.46
1:A:315:THR:HG21	2:A:401:BEN:H4	1.98	0.46
1:B:319:ASN:HD21	1:B:347:GLN:NE2	2.13	0.45
1:D:283:SER:N	4:D:519:HOH:O	2.48	0.45
1:A:347:GLN:O	1:A:347:GLN:HG3	2.16	0.45
1:B:362:ARG:NH2	4:B:7001:HOH:O	2.49	0.45
1:C:323:ARG:NH1	4:C:548:HOH:O	2.50	0.45
1:B:344:ARG:NH1	4:B:7041:HOH:O	2.49	0.45
1:B:362:ARG:NE	4:B:7001:HOH:O	2.50	0.45
1:D:303:ASN:HD21	1:D:310:GLN:HE22	1.65	0.45
1:C:284:ILE:CD1	1:C:361:ALA:HA	2.46	0.44
1:D:349:TRP:N	4:D:471:HOH:O	2.50	0.44
1:B:303:ASN:ND2	1:B:305:TYR:H	2.16	0.44
1:B:349:TRP:CD1	1:B:361:ALA:HB1	2.54	0.43
1:D:358:ARG:NH2	4:D:450:HOH:O	2.50	0.43
1:D:349:TRP:HD1	1:D:362:ARG:HG3	1.83	0.43
1:B:361:ALA:N	4:B:7071:HOH:O	2.52	0.42
1:A:299:ILE:HG12	1:A:301:THR:HG23	2.00	0.42
1:A:362:ARG:N	4:A:583:HOH:O	2.49	0.42
1:B:349:TRP:NE1	1:B:362:ARG:N	2.65	0.42
1:B:360:GLY:HA2	4:B:7147:HOH:O	2.20	0.41
1:B:360:GLY:HA2	4:B:7117:HOH:O	2.21	0.41
1:B:362:ARG:HB3	4:B:7039:HOH:O	2.21	0.40
1:D:284:ILE:H	1:D:284:ILE:CD1	2.35	0.40
1:A:303:ASN:N	1:A:303:ASN:HD22	2.20	0.40

There are no symmetry-related clashes.

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	79/81 (98%)	77 (98%)	2 (2%)	0	100	100
1	B	78/81 (96%)	75 (96%)	3 (4%)	0	100	100
1	C	79/81 (98%)	76 (96%)	2 (2%)	1 (1%)	12	1
1	D	78/81 (96%)	76 (97%)	2 (3%)	0	100	100
All	All	314/324 (97%)	304 (97%)	9 (3%)	1 (0%)	41	17

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	361	ALA

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	51/52 (98%)	48 (94%)	3 (6%)	19	1
1	B	52/52 (100%)	50 (96%)	2 (4%)	33	4
1	C	51/52 (98%)	49 (96%)	2 (4%)	32	4
1	D	51/52 (98%)	51 (100%)	0	100	100
All	All	205/208 (99%)	198 (97%)	7 (3%)	37	5

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	303	ASN
1	A	323	ARG
1	A	344	ARG
1	B	283	SER
1	B	303	ASN
1	C	344	ARG
1	C	347	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	303	ASN
1	A	310	GLN
1	B	303	ASN
1	B	310	GLN
1	B	347	GLN
1	D	310	GLN

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 ⓘ

8 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	BEN	C	401	-	9,9,9	0.89	1 (11%)	7,11,11	1.12	1 (14%)
2	BEN	B	402	-	9,9,9	0.81	1 (11%)	7,11,11	1.06	0
3	SO4	B	403	-	4,4,4	0.50	0	6,6,6	0.67	0
3	SO4	A	403	-	4,4,4	0.35	0	6,6,6	0.38	0
2	BEN	A	401	-	9,9,9	0.92	0	7,11,11	1.88	2 (28%)
2	BEN	B	401	-	9,9,9	0.74	0	7,11,11	0.89	0
2	BEN	A	402	-	9,9,9	0.83	0	7,11,11	1.36	1 (14%)
2	BEN	C	402	-	9,9,9	1.08	1 (11%)	7,11,11	0.70	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	BEN	C	401	-	-	2/4/4/4	0/1/1/1
2	BEN	B	402	-	-	1/4/4/4	0/1/1/1
2	BEN	A	401	-	-	0/4/4/4	0/1/1/1
2	BEN	B	401	-	-	0/4/4/4	0/1/1/1
2	BEN	A	402	-	-	0/4/4/4	0/1/1/1
2	BEN	C	402	-	-	1/4/4/4	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	401	BEN	C1-C	2.14	1.51	1.47
2	C	402	BEN	C1-C	2.14	1.51	1.47
2	B	402	BEN	C1-C	2.05	1.51	1.47

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	BEN	C3-C2-C1	-3.10	116.67	120.34
2	A	401	BEN	C4-C5-C6	-2.47	116.43	120.19
2	A	402	BEN	C4-C3-C2	-2.02	117.12	120.19
2	C	401	BEN	C1-C-N2	-2.00	115.02	118.05

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	401	BEN	N2-C-C1-C6
2	B	402	BEN	N2-C-C1-C6
2	C	402	BEN	N2-C-C1-C6
2	C	401	BEN	N1-C-C1-C6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	BEN	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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	81/81 (100%)	-0.07	1 (1%) 79 79	8, 11, 21, 37	0
1	B	80/81 (98%)	0.10	3 (3%) 40 42	8, 11, 30, 44	0
1	C	81/81 (100%)	0.07	3 (3%) 41 43	7, 12, 28, 64	0
1	D	80/81 (98%)	0.17	5 (6%) 20 21	9, 14, 31, 50	0
All	All	322/324 (99%)	0.07	12 (3%) 41 43	7, 12, 29, 64	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	361	ALA	6.8
1	C	362	ARG	5.6
1	B	362	ARG	4.3
1	D	362	ARG	4.1
1	C	361	ALA	3.9
1	D	360	GLY	3.5
1	B	360	GLY	3.4
1	C	282	GLY	3.3
1	A	347	GLN	2.7
1	D	361	ALA	2.7
1	D	284	ILE	2.6
1	D	358	ARG	2.2

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. The B-factors column lists the minimum, median, 95th 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 ²)	Q<0.9
2	BEN	C	402	9/9	0.82	0.17	18,20,23,29	0
2	BEN	B	401	9/9	0.92	0.10	17,22,28,30	0
2	BEN	C	401	9/9	0.92	0.09	18,21,23,24	0
2	BEN	B	402	9/9	0.94	0.10	15,18,23,23	0
3	SO4	B	403	5/5	0.96	0.12	16,18,23,23	0
2	BEN	A	401	9/9	0.96	0.07	10,12,19,21	0
2	BEN	A	402	9/9	0.98	0.05	11,12,13,13	0
3	SO4	A	403	5/5	0.99	0.06	11,12,14,15	0

6.5 Other polymers [i](#)

There are no such residues in this entry.