



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 10:55 PM GMT

PDB ID : 3A3K
Title : Reversibly bound chloride in the atrial natriuretic peptide receptor hormone-binding domain
Authors : Ogawa, H.; Qiu, Y.; Ogata, C.M.; Misono, K.S.
Deposited on : 2009-06-14
Resolution : 2.50 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

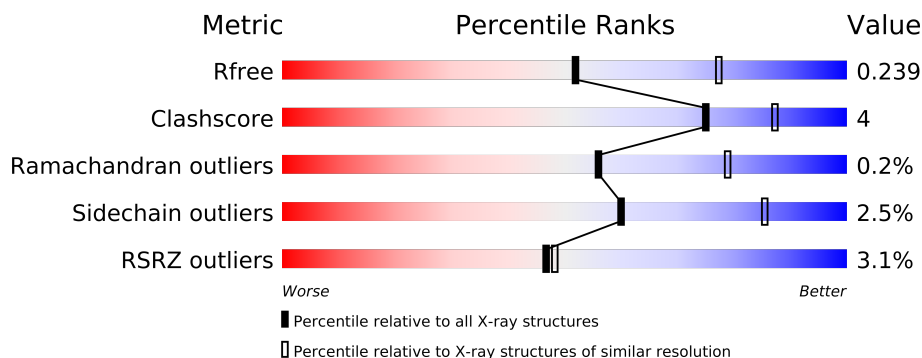
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.50 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}	66092	2784 (2.50-2.50)
Clashscore	79885	3562 (2.50-2.50)
Ramachandran outliers	78287	3480 (2.50-2.50)
Sidechain outliers	78261	3482 (2.50-2.50)
RSRZ outliers	66119	2785 (2.50-2.50)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	435	
1	B	435	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	NAG	B	451	-	X
2	NAG	B	471	-	X
3	SO4	A	504	-	X
3	SO4	B	507	-	X
3	SO4	B	509	-	X

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 7189 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Atrial natriuretic peptide receptor A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	425	Total	C	N	O	S	0	1	0
			3351	2150	583	607	11			
1	B	429	Total	C	N	O	S	0	1	0
			3389	2167	589	621	12			

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



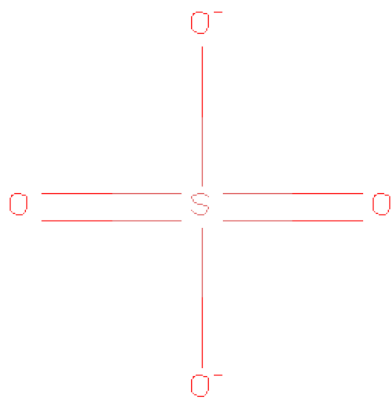
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		

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

- 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	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

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

- Molecule 4 is BROMIDE ION (three-letter code: BR) (formula: Br).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Br	0	0
			1	1		
4	A	1	Total	Br	0	0
			1	1		

- Molecule 5 is water.

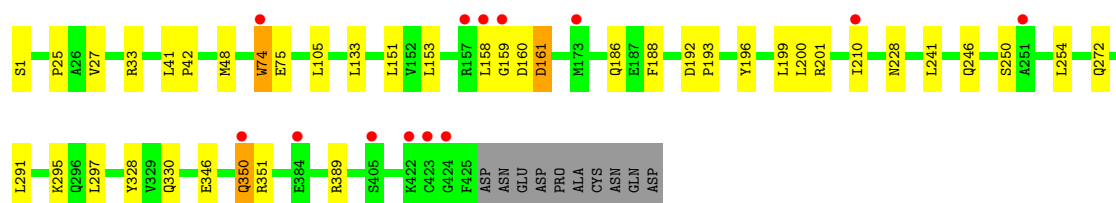
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	137	Total	O	0	0
			137	137		
5	B	125	Total	O	0	0
			125	125		

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 ($RSRZ > 2$). 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.

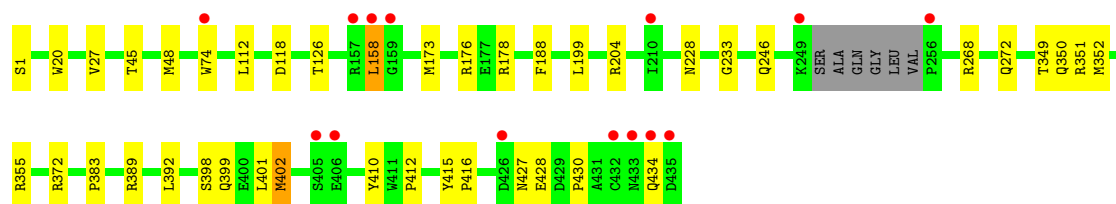
- Molecule 1: Atrial natriuretic peptide receptor A

Chain A: 



- Molecule 1: Atrial natriuretic peptide receptor A

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	120.52Å 120.52Å 160.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.92 – 2.50 19.93 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.92-2.50) 92.5 (19.93-2.50)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.37 (at 2.50Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.205 , 0.243 0.203 , 0.239	Depositor DCC
R_{free} test set	1936 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	26.2	Xtriage
Anisotropy	0.063	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 11.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 38394 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7189	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.63% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, BR, 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	0.86	10/3447 (0.3%)	0.78	12/4697 (0.3%)
1	B	0.76	10/3485 (0.3%)	0.81	8/4746 (0.2%)
All	All	0.81	20/6932 (0.3%)	0.80	20/9443 (0.2%)

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	74[A]	TRP	CG-CD1	-19.25	1.09	1.36
1	A	74[B]	TRP	CG-CD1	-19.25	1.09	1.36
1	B	74[A]	TRP	CG-CD1	-17.14	1.12	1.36
1	B	74[B]	TRP	CG-CD1	-17.14	1.12	1.36
1	A	74[A]	TRP	CD2-CE2	-16.75	1.21	1.41
1	A	74[B]	TRP	CD2-CE2	-16.75	1.21	1.41
1	B	74[A]	TRP	CD2-CE2	-12.20	1.26	1.41
1	B	74[B]	TRP	CD2-CE2	-12.20	1.26	1.41
1	A	74[A]	TRP	CD1-NE1	11.99	1.58	1.38
1	A	74[B]	TRP	CD1-NE1	11.99	1.58	1.38
1	B	74[A]	TRP	CD1-NE1	-10.86	1.19	1.38
1	B	74[B]	TRP	CD1-NE1	-10.86	1.19	1.38
1	B	74[A]	TRP	NE1-CE2	-6.96	1.28	1.37
1	B	74[B]	TRP	NE1-CE2	-6.96	1.28	1.37
1	A	74[A]	TRP	CG-CD2	6.64	1.54	1.43
1	A	74[B]	TRP	CG-CD2	6.64	1.54	1.43
1	A	74[A]	TRP	CD2-CE3	-5.62	1.31	1.40
1	A	74[B]	TRP	CD2-CE3	-5.62	1.31	1.40
1	B	74[A]	TRP	CG-CD2	5.06	1.52	1.43
1	B	74[B]	TRP	CG-CD2	5.06	1.52	1.43

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	74[A]	TRP	CD1-NE1-CE2	21.88	128.69	109.00
1	B	74[B]	TRP	CD1-NE1-CE2	21.88	128.69	109.00
1	A	74[A]	TRP	NE1-CE2-CZ2	-16.24	112.53	130.40
1	A	74[B]	TRP	NE1-CE2-CZ2	-16.24	112.53	130.40
1	A	74[A]	TRP	CD2-CE2-CZ2	14.09	139.20	122.30
1	A	74[B]	TRP	CD2-CE2-CZ2	14.09	139.20	122.30
1	B	74[A]	TRP	NE1-CE2-CD2	-13.72	93.58	107.30
1	B	74[B]	TRP	NE1-CE2-CD2	-13.72	93.58	107.30
1	A	74[A]	TRP	CD1-NE1-CE2	-8.67	101.19	109.00
1	A	74[B]	TRP	CD1-NE1-CE2	-8.67	101.19	109.00
1	B	74[A]	TRP	CG-CD1-NE1	-6.93	103.17	110.10
1	B	74[B]	TRP	CG-CD1-NE1	-6.93	103.17	110.10
1	B	74[A]	TRP	CD2-CE2-CZ2	6.90	130.58	122.30
1	B	74[B]	TRP	CD2-CE2-CZ2	6.90	130.58	122.30
1	A	74[A]	TRP	CE2-CD2-CG	6.75	112.70	107.30
1	A	74[B]	TRP	CE2-CD2-CG	6.75	112.70	107.30
1	A	74[A]	TRP	CH2-CZ2-CE2	-6.47	110.93	117.40
1	A	74[B]	TRP	CH2-CZ2-CE2	-6.47	110.93	117.40
1	A	74[A]	TRP	CE2-CD2-CE3	-5.08	112.61	118.70
1	A	74[B]	TRP	CE2-CD2-CE3	-5.08	112.61	118.70

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3351	0	3301	24	0
1	B	3389	0	3312	25	0
2	A	56	0	52	3	0
2	B	84	0	78	4	0
3	A	20	0	0	0	0
3	B	25	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	137	0	0	7	0
5	B	125	0	0	2	0
All	All	7189	0	6743	51	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 4.

All (51) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:272:GLN:HG3	5:A:579:HOH:O	1.49	1.13
2:A:461:NAG:O4	2:A:462:NAG:C1	2.20	0.89
1:B:272:GLN:HG3	5:B:537:HOH:O	1.75	0.87
2:B:471:NAG:HO4	2:B:472:NAG:C1	1.97	0.77
1:B:233:GLY:HA3	1:B:412:PRO:HG2	1.65	0.77
1:B:178:ARG:HH22	1:B:399:GLN:HE22	1.31	0.75
1:A:161:ASP:HB2	5:A:553:HOH:O	1.87	0.75
1:A:330:GLN:HG2	5:A:490:HOH:O	1.87	0.74
1:B:415:TYR:HB2	1:B:416:PRO:HD2	1.72	0.71
2:B:461:NAG:O4	2:B:462:NAG:C1	2.44	0.65
1:B:27:VAL:HG11	1:B:48:MET:CE	2.29	0.62
1:A:160:ASP:HB2	5:A:553:HOH:O	2.01	0.61
1:A:74[B]:TRP:HD1	1:A:75:GLU:HG3	1.70	0.56
1:A:25:PRO:HB2	1:A:297:LEU:HD12	1.88	0.56
1:B:27:VAL:HG11	1:B:48:MET:HE1	1.86	0.56
1:A:228:ASN:ND2	1:B:204:ARG:HD3	2.22	0.55
1:A:196:TYR:CZ	5:A:436:HOH:O	2.60	0.54
1:B:427:ASN:HD22	1:B:434:GLN:NE2	2.05	0.54
1:A:228:ASN:HD21	1:B:204:ARG:HD3	1.72	0.53
1:A:27:VAL:HG11	1:A:48:MET:CE	2.38	0.53
1:A:200:LEU:HB3	1:B:228:ASN:OD1	2.10	0.52
1:B:392:LEU:HD13	1:B:401:LEU:HD22	1.92	0.51
1:B:427:ASN:HD22	1:B:434:GLN:HE21	1.57	0.51
1:A:291:LEU:O	1:A:295:LYS:HG3	2.10	0.51
1:B:349:THR:HA	1:B:352:MET:HE2	1.92	0.51
1:A:153:LEU:HD23	1:A:186:GLN:HB3	1.93	0.50
1:A:133:LEU:HD21	1:A:241:LEU:HD22	1.93	0.50
1:A:151:LEU:HD23	1:A:210:ILE:HD12	1.93	0.50
1:B:428:GLU:O	1:B:430:PRO:HD3	2.12	0.49
1:B:178:ARG:HH22	1:B:399:GLN:NE2	2.07	0.49
1:B:402:MET:HG3	5:B:565:HOH:O	2.12	0.49
1:A:105:LEU:HD11	1:A:328:TYR:HB2	1.94	0.49
1:B:178:ARG:NH2	1:B:399:GLN:HE22	2.07	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:346:GLU:HB2	5:A:485:HOH:O	2.12	0.48
2:B:451:NAG:C4	2:B:452:NAG:C1	2.91	0.48
1:A:188:PHE:CG	1:A:199:LEU:HD11	2.50	0.46
1:B:351:ARG:O	1:B:355:ARG:HD3	2.15	0.46
1:B:20:TRP:HH2	1:B:48:MET:HE1	1.82	0.44
1:A:350:GLN:HG2	1:A:351:ARG:N	2.33	0.44
2:A:451:NAG:C4	2:A:452:NAG:C1	2.95	0.43
1:A:196:TYR:CE1	5:A:436:HOH:O	2.57	0.43
1:A:105:LEU:HD11	1:A:328:TYR:CB	2.50	0.42
1:B:126:THR:O	1:B:372:ARG:HD2	2.20	0.42
1:A:41:LEU:N	1:A:42:PRO:HD3	2.35	0.42
1:B:398:SER:HB3	2:B:461:NAG:H61	2.01	0.42
1:B:188:PHE:CG	1:B:199:LEU:HD11	2.55	0.42
1:A:192:ASP:HA	1:A:193:PRO:HD3	1.89	0.41
2:A:461:NAG:C4	2:A:462:NAG:C1	2.98	0.41
1:A:201:ARG:NH2	1:B:268:ARG:HH22	2.18	0.41
1:B:112:LEU:HD22	1:B:173:MET:HG3	2.02	0.40
1:B:383:PRO:HG3	1:B:410:TYR:CD1	2.56	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	424/435 (98%)	417 (98%)	6 (1%)	1 (0%)	56	79
1	B	426/435 (98%)	416 (98%)	9 (2%)	1 (0%)	56	79
All	All	850/870 (98%)	833 (98%)	15 (2%)	2 (0%)	56	79

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	158	LEU
1	A	159	GLY

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	353/361 (98%)	344 (98%)	9 (2%)	60	85
1	B	358/361 (99%)	349 (98%)	9 (2%)	60	85
All	All	711/722 (98%)	693 (98%)	18 (2%)	60	85

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	SER
1	A	33	ARG
1	A	158	LEU
1	A	161	ASP
1	A	246	GLN
1	A	250	SER
1	A	254	LEU
1	A	350	GLN
1	A	389	ARG
1	B	1	SER
1	B	45	THR
1	B	118	ASP
1	B	158	LEU
1	B	176	ARG
1	B	246	GLN
1	B	350	GLN
1	B	389	ARG
1	B	402	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	228	ASN
1	A	246	GLN
1	A	272	GLN
1	B	185	HIS
1	B	296	GLN
1	B	399	GLN

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Mol	Chain	Res	Type
1	B	434	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

Of 21 ligands modelled in this entry, 2 are monoatomic - leaving 19 for Mogul analysis.

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	NAG	A	451	1,2	12,14,15	0.77	0	15,19,21	1.14	1 (6%)
2	NAG	A	452	2	12,14,15	0.63	0	15,19,21	1.01	1 (6%)
2	NAG	A	461	1	12,14,15	0.86	1 (8%)	15,19,21	1.41	3 (20%)
2	NAG	A	462	-	12,14,15	0.60	0	15,19,21	1.29	1 (6%)
3	SO4	A	501	-	4,4,4	0.27	0	6,6,6	0.34	0
3	SO4	A	503	-	4,4,4	0.22	0	6,6,6	0.14	0
3	SO4	A	504	-	4,4,4	0.20	0	6,6,6	0.09	0
3	SO4	A	505	-	4,4,4	0.18	0	6,6,6	0.14	0
2	NAG	B	451	1,2	12,14,15	0.81	1 (8%)	15,19,21	1.61	4 (26%)
2	NAG	B	452	2	12,14,15	0.72	0	15,19,21	1.01	1 (6%)
2	NAG	B	461	1	12,14,15	0.73	0	15,19,21	1.37	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	B	462	-	12,14,15	0.52	0	15,19,21	1.15	1 (6%)
2	NAG	B	471	1,2	12,14,15	0.68	0	15,19,21	2.02	2 (13%)
2	NAG	B	472	2	12,14,15	0.67	0	15,19,21	0.78	0
3	SO4	B	502	-	4,4,4	0.24	0	6,6,6	0.09	0
3	SO4	B	506	-	4,4,4	0.20	0	6,6,6	0.19	0
3	SO4	B	507	-	4,4,4	0.24	0	6,6,6	0.21	0
3	SO4	B	508	-	4,4,4	0.23	0	6,6,6	0.27	0
3	SO4	B	509	-	4,4,4	0.17	0	6,6,6	0.17	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	NAG	A	451	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	452	2	-	0/6/23/26	0/1/1/1
2	NAG	A	461	1	-	0/6/23/26	0/1/1/1
2	NAG	A	462	-	-	0/6/23/26	0/1/1/1
3	SO4	A	501	-	-	0/0/0/0	0/0/0/0
3	SO4	A	503	-	-	0/0/0/0	0/0/0/0
3	SO4	A	504	-	-	0/0/0/0	0/0/0/0
3	SO4	A	505	-	-	0/0/0/0	0/0/0/0
2	NAG	B	451	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	452	2	-	0/6/23/26	0/1/1/1
2	NAG	B	461	1	-	0/6/23/26	0/1/1/1
2	NAG	B	462	-	-	0/6/23/26	0/1/1/1
2	NAG	B	471	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	472	2	-	0/6/23/26	0/1/1/1
3	SO4	B	502	-	-	0/0/0/0	0/0/0/0
3	SO4	B	506	-	-	0/0/0/0	0/0/0/0
3	SO4	B	507	-	-	0/0/0/0	0/0/0/0
3	SO4	B	508	-	-	0/0/0/0	0/0/0/0
3	SO4	B	509	-	-	0/0/0/0	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	451	NAG	O5-C5	-2.28	1.41	1.45
2	A	461	NAG	O5-C5	-2.16	1.41	1.45

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	471	NAG	O5-C5-C6	6.19	113.48	106.98
2	A	462	NAG	O5-C5-C6	4.12	111.30	106.98
2	B	471	NAG	O5-C5-C4	-3.80	105.83	110.65
2	B	451	NAG	C3-C2-N2	-3.73	106.09	111.76
2	A	461	NAG	O5-C5-C6	3.27	110.41	106.98
2	B	462	NAG	O5-C5-C6	3.25	110.39	106.98
2	B	461	NAG	C3-C4-C5	2.97	115.51	110.20
2	A	451	NAG	C3-C2-N2	-2.88	107.37	111.76
2	A	461	NAG	O4-C4-C3	-2.64	104.43	110.35
2	B	451	NAG	O5-C5-C6	2.55	109.66	106.98
2	B	452	NAG	O5-C5-C6	2.43	109.53	106.98
2	B	451	NAG	O4-C4-C3	-2.40	104.96	110.35
2	B	451	NAG	C3-C4-C5	2.32	114.35	110.20
2	A	461	NAG	C3-C2-N2	-2.29	108.27	111.76
2	A	452	NAG	C2-N2-C7	2.09	126.59	123.09

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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 ⓘ

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, 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	425/435 (97%)	-0.16	13 (3%)	47 48	10, 21, 36, 49	2 (0%)
1	B	429/435 (98%)	-0.15	14 (3%)	44 45	12, 23, 39, 61	2 (0%)
All	All	854/870 (98%)	-0.15	27 (3%)	47 47	10, 22, 37, 61	4 (0%)

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	158	LEU	7.9
1	B	158	LEU	6.2
1	A	423	CYS	6.0
1	B	435	ASP	5.1
1	B	434	GLN	3.9
1	B	157	ARG	3.5
1	B	159	GLY	3.1
1	B	433	ASN	3.0
1	A	251	ALA	3.0
1	A	157	ARG	3.0
1	A	159	GLY	2.9
1	A	74[A]	TRP	2.7
1	A	210	ILE	2.7
1	B	405	SER	2.7
1	A	350	GLN	2.6
1	B	432	CYS	2.6
1	A	422	LYS	2.4
1	A	384	GLU	2.3
1	A	405	SER	2.3
1	B	74[A]	TRP	2.3
1	B	406	GLU	2.2
1	A	424	GLY	2.2
1	B	210	ILE	2.1
1	B	256	PRO	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	426	ASP	2.1
1	B	249	LYS	2.0
1	A	173	MET	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	SO4	A	504	5/5	0.22	9.82	41,42,42,43	0
2	NAG	B	471	14/15	0.33	8.76	47,51,53,54	0
3	SO4	B	507	5/5	0.23	5.55	45,46,46,46	0
3	SO4	B	509	5/5	0.29	3.58	88,88,89,89	0
2	NAG	B	451	14/15	0.18	2.71	31,35,36,39	0
2	NAG	B	462	14/15	0.24	0.91	46,48,49,50	0
3	SO4	B	508	5/5	0.13	0.75	40,41,41,42	0
2	NAG	B	461	14/15	0.19	0.73	30,33,35,36	0
3	SO4	B	506	5/5	0.20	0.65	45,45,46,46	0
2	NAG	A	451	14/15	0.13	0.50	22,24,28,29	0
3	SO4	A	505	5/5	0.19	0.47	42,42,43,43	0
2	NAG	B	452	14/15	0.31	0.38	58,59,60,60	0
2	NAG	A	462	14/15	0.19	0.31	40,43,44,45	0
2	NAG	A	461	14/15	0.13	0.19	24,25,26,27	0
3	SO4	B	502	5/5	0.16	0.13	39,39,40,40	5
2	NAG	A	452	14/15	0.25	0.03	44,45,45,45	0
3	SO4	A	501	5/5	0.10	-0.78	20,22,22,23	0
3	SO4	A	503	5/5	0.08	-1.42	26,26,26,26	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	BR	B	601	1/1	0.02	-3.58	25,25,25,25	0
4	BR	A	600	1/1	0.02	-4.89	21,21,21,21	0
2	NAG	B	472	14/15	0.45	-	64,66,66,66	0

6.5 Other polymers ⓘ

There are no such residues in this entry.