



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 12:39 AM GMT

PDB ID : 1DMA  
Title : DOMAIN III OF PSEUDOMONAS AERUGINOSA EXOTOXIN COM-  
PLEXED WITH NICOTINAMIDE AND AMP  
Authors : Li, M.; Dyda, F.; Benhar, I.; Pastan, I.; Davies, D.  
Deposited on : 1995-04-28  
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](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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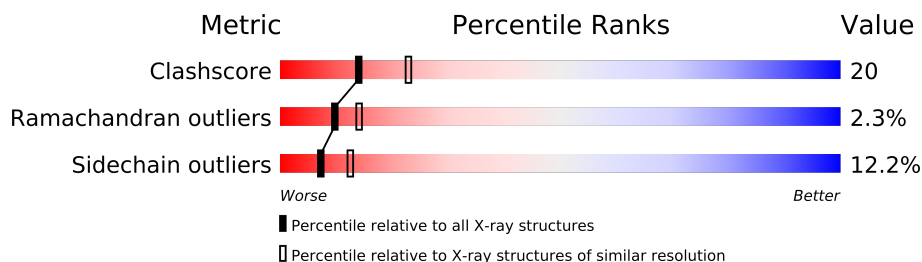
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 21963  
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))
Clashscore	79885	3562 (2.50-2.50)
Ramachandran outliers	78287	3480 (2.50-2.50)
Sidechain outliers	78261	3482 (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  $\geq 3$ , 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	214	
1	B	214	

## 2 Entry composition i

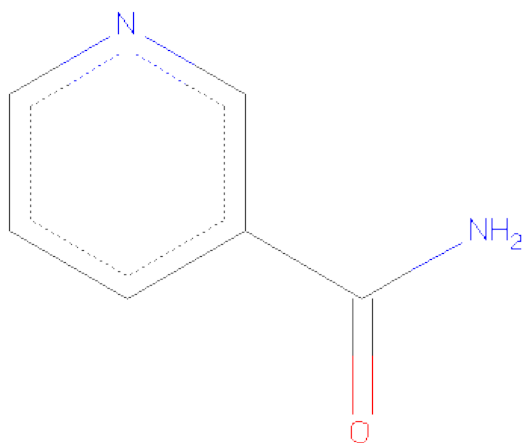
There are 4 unique types of molecules in this entry. The entry contains 3073 atoms, of which 0 are hydrogen and 0 are deuterium.

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 EXOTOXIN A.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	204	Total	C	N	O	5	0	0
			1562	990	280	292			
1	B	189	Total	C	N	O	27	0	0
			1437	914	254	269			

- Molecule 2 is NICOTINAMIDE (three-letter code: NCA) (formula: C<sub>6</sub>H<sub>6</sub>N<sub>2</sub>O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			9	6	2	1		
2	B	1	Total	C	N	O	0	0
			9	6	2	1		

- Molecule 3 is ADENOSINE MONOPHOSPHATE (three-letter code: AMP) (formula: C<sub>10</sub>H<sub>14</sub>N<sub>5</sub>O<sub>7</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
3	B	1	23	10	5	7	1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	18	Total	O	0	0
			18	18		
4	B	15	Total	O	0	0
			15	15		

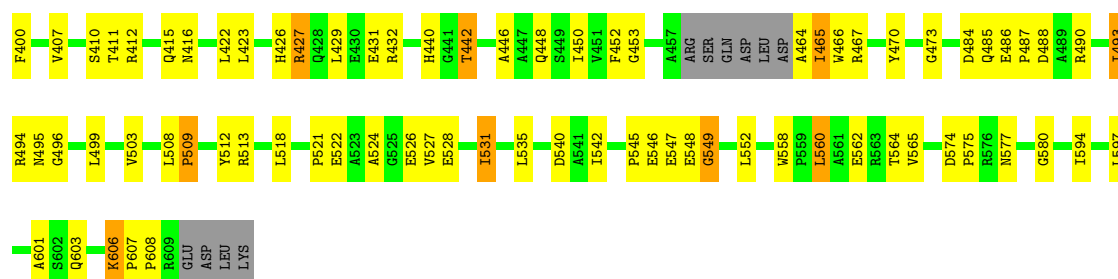
### 3 Residue-property plots

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.

Note EDS was not executed.

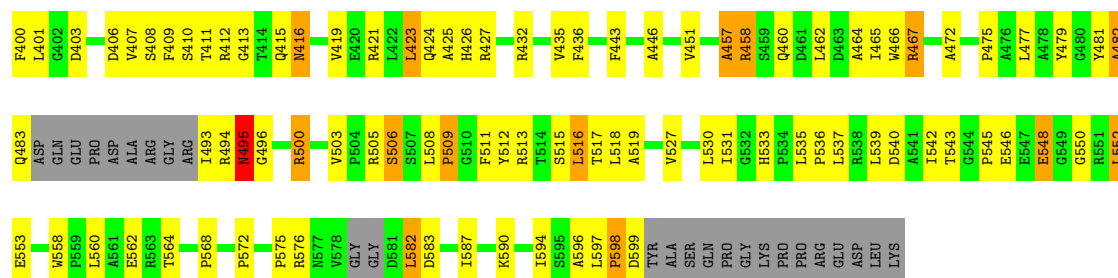
#### • Molecule 1: EXOTOXIN A

Chain A: 



#### • Molecule 1: EXOTOXIN A

Chain B: 



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	87.51Å 87.51Å 134.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.50	Depositor
% Data completeness (in resolution range)	88.9 (8.00-2.50)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, $R_{free}$	0.195 , 0.265	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3073	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	wwPDB-VP

## 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: NCA, AMP

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	0.88	0/1602	1.04	2/2183 (0.1%)
1	B	0.89	1/1471 (0.1%)	1.11	7/2003 (0.3%)
All	All	0.89	1/3073 (0.0%)	1.07	9/4186 (0.2%)

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	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	575	PRO	CA-C	5.05	1.62	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	537	LEU	CA-CB-CG	7.83	133.30	115.30
1	B	516	LEU	CA-CB-CG	7.79	133.22	115.30
1	B	576	ARG	NE-CZ-NH2	7.11	123.86	120.30
1	B	457	ALA	N-CA-C	7.04	130.02	111.00
1	A	606	LYS	CB-CA-C	5.95	122.30	110.40
1	B	494	ARG	N-CA-C	5.59	126.11	111.00
1	B	495	ASN	O-C-N	-5.45	113.93	123.20
1	A	493	ILE	N-CA-C	-5.32	96.63	111.00
1	B	587	ILE	CB-CA-C	-5.04	101.53	111.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	495	ASN	Mainchain

## 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	1562	0	1527	42	0
1	B	1437	0	1402	77	0
2	A	9	0	6	0	0
2	B	9	0	6	2	0
3	B	23	0	12	1	0
4	A	18	0	0	0	0
4	B	15	0	0	0	0
All	All	3073	0	2953	117	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 20.

All (117) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:457:ALA:O	1:B:458:ARG:HB2	1.71	0.89
1:A:400:PHE:HE1	1:B:539:LEU:HD11	1.42	0.82
1:B:482:ALA:HB1	1:B:496:GLY:N	1.93	0.82
1:A:503:VAL:HG12	1:A:564:THR:HG22	1.64	0.79
1:B:598:PRO:HG2	1:B:599:ASP:H	1.47	0.78
1:A:524:ALA:O	1:A:528:GLU:HG3	1.84	0.77
1:B:482:ALA:HB1	1:B:496:GLY:H	1.49	0.76
1:A:429:LEU:HD21	1:A:565:VAL:HG11	1.69	0.73
1:B:558:TRP:O	1:B:562:GLU:HG3	1.89	0.72
1:B:424:GLN:HG3	1:B:427:ARG:HH21	1.53	0.71
1:B:419:VAL:O	1:B:423:LEU:HD22	1.90	0.71
1:A:464:ALA:O	1:A:467:ARG:HG2	1.91	0.69
1:B:436:PHE:HE1	1:B:597:LEU:HD21	1.57	0.68

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:400:PHE:HE1	1:B:539:LEU:CD1	2.06	0.68
1:B:426:HIS:CE1	1:B:594:ILE:HG23	2.28	0.68
1:A:465:ILE:HG23	1:A:535:LEU:HD13	1.78	0.65
1:B:512:TYR:HB2	1:B:542:ILE:HD12	1.80	0.64
1:B:582:LEU:HD22	1:B:583:ASP:N	2.12	0.64
1:B:423:LEU:O	1:B:427:ARG:HG3	1.99	0.63
1:B:512:TYR:O	1:B:543:THR:HG22	1.99	0.62
1:B:435:VAL:HG11	1:B:505:ARG:HE	1.65	0.61
1:B:548:GLU:C	1:B:550:GLY:H	2.02	0.61
1:B:596:ALA:O	1:B:597:LEU:HD23	2.00	0.61
1:A:400:PHE:CE1	1:B:539:LEU:HD11	2.32	0.60
1:B:466:TRP:CD1	1:B:519:ALA:HA	2.36	0.60
1:A:440:HIS:HD2	1:A:470:TYR:O	1.86	0.59
1:A:407:VAL:HA	1:A:415:GLN:O	2.03	0.59
1:B:516:LEU:O	1:B:545:PRO:HD2	2.03	0.58
1:B:466:TRP:HH2	1:B:527:VAL:HG21	1.68	0.58
1:B:503:VAL:HG12	1:B:564:THR:HG22	1.84	0.58
1:B:560:LEU:O	1:B:564:THR:HG23	2.02	0.58
1:B:400:PHE:O	1:B:425:ALA:HB1	2.03	0.58
1:A:540:ASP:O	1:A:560:LEU:HD12	2.03	0.58
1:B:407:VAL:HG11	1:B:568:PRO:HG2	1.84	0.57
1:B:472:ALA:HB2	2:B:700:NCA:H5	1.87	0.57
1:B:506:SER:O	1:B:509:PRO:HD2	2.05	0.57
1:A:546:GLU:CD	1:A:549:GLY:HA3	2.26	0.56
1:B:475:PRO:O	1:B:479:TYR:HB2	2.04	0.56
1:B:443:PHE:CZ	1:B:446:ALA:HB2	2.40	0.56
1:B:513:ARG:HG3	1:B:543:THR:HG23	1.88	0.55
1:A:432:ARG:HH11	1:A:432:ARG:HG2	1.72	0.55
1:A:521:PRO:HG2	1:A:522:GLU:OE1	2.05	0.55
1:A:577:ASN:OD1	1:A:580:GLY:HA3	2.07	0.54
1:B:423:LEU:HD21	1:B:590:LYS:HD3	1.90	0.54
1:A:527:VAL:O	1:A:531:ILE:HB	2.07	0.54
1:B:517:THR:HG22	1:B:545:PRO:HG2	1.91	0.53
1:B:548:GLU:C	1:B:550:GLY:N	2.62	0.53
1:B:419:VAL:HG11	1:B:590:LYS:HB2	1.91	0.52
1:B:598:PRO:HG2	1:B:599:ASP:N	2.22	0.52
1:A:410:SER:OG	1:A:412:ARG:HG2	2.10	0.51
1:B:457:ALA:O	1:B:458:ARG:CB	2.53	0.51
1:B:465:ILE:HG23	1:B:535:LEU:HD13	1.91	0.51
1:A:465:ILE:HG22	1:A:466:TRP:CD1	2.46	0.51
1:B:482:ALA:HB1	1:B:496:GLY:CA	2.40	0.51
1:B:408:SER:HB3	1:B:415:GLN:HB2	1.93	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:426:HIS:NE2	1:A:594:ILE:HG13	2.26	0.50
1:B:481:TYR:CD1	1:B:481:TYR:N	2.80	0.50
1:B:464:ALA:O	1:B:467:ARG:HG2	2.12	0.50
1:B:424:GLN:NE2	1:B:427:ARG:HE	2.10	0.50
1:B:500:ARG:HH22	1:B:594:ILE:HG22	1.77	0.49
1:A:486:GLU:O	1:A:487:PRO:C	2.49	0.49
1:B:401:LEU:HD22	1:B:407:VAL:HG21	1.94	0.49
1:A:547:GLU:O	1:A:548:GLU:HB2	2.10	0.49
1:B:416:ASN:N	1:B:416:ASN:HD22	2.11	0.49
1:A:448:GLN:NE2	1:A:452:PHE:CD2	2.80	0.49
1:B:530:LEU:HD12	1:B:542:ILE:HD11	1.93	0.49
1:B:531:ILE:HG23	1:B:533:HIS:CD2	2.46	0.49
1:A:432:ARG:NH1	1:A:432:ARG:HG2	2.26	0.49
1:B:436:PHE:CE1	1:B:597:LEU:HD21	2.45	0.49
1:B:598:PRO:CG	1:B:599:ASP:H	2.13	0.48
1:A:545:PRO:HA	1:A:552:LEU:HD23	1.96	0.48
1:A:488:ASP:OD1	1:A:488:ASP:C	2.51	0.48
1:B:423:LEU:HD12	1:B:594:ILE:HD11	1.96	0.47
1:B:512:TYR:CD1	1:B:531:ILE:HD12	2.49	0.47
1:B:411:THR:HG23	1:B:495:ASN:OD1	2.15	0.47
1:A:473:GLY:HA3	1:A:597:LEU:HD21	1.96	0.47
1:A:448:GLN:NE2	1:A:452:PHE:HD2	2.13	0.47
1:A:427:ARG:HA	1:A:427:ARG:NE	2.28	0.46
1:B:412:ARG:HD2	1:B:415:GLN:NE2	2.30	0.46
1:B:535:LEU:HB3	1:B:536:PRO:HA	1.98	0.45
1:B:599:ASP:OD1	1:B:599:ASP:O	2.34	0.45
1:A:512:TYR:CD2	1:A:531:ILE:HD12	2.52	0.45
1:A:410:SER:OG	1:A:412:ARG:NE	2.50	0.45
1:B:436:PHE:HE1	1:B:597:LEU:CD2	2.28	0.45
1:A:415:GLN:O	1:A:416:ASN:HB2	2.17	0.44
1:B:413:GLY:O	1:B:415:GLN:NE2	2.50	0.44
1:A:558:TRP:O	1:A:562:GLU:HG3	2.18	0.44
1:B:511:PHE:HD2	1:B:599:ASP:O	2.00	0.44
1:A:442:THR:O	1:A:496:GLY:HA3	2.17	0.44
1:B:435:VAL:HG21	1:B:508:LEU:HD11	1.99	0.43
1:B:553:GLU:HB2	2:B:700:NCA:H6	1.99	0.43
1:A:484:ASP:HB2	1:A:493:ILE:HA	2.01	0.43
1:A:465:ILE:HG23	1:A:535:LEU:CD1	2.45	0.43
1:B:427:ARG:HB2	1:B:427:ARG:CZ	2.49	0.42
1:B:409:PHE:HZ	1:B:451:VAL:HG21	1.83	0.42
1:A:512:TYR:HA	1:A:601:ALA:O	2.19	0.42
1:B:598:PRO:CG	1:B:599:ASP:N	2.77	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:545:PRO:HA	1:B:552:LEU:H	1.84	0.42
1:A:546:GLU:O	1:A:546:GLU:HG3	2.20	0.42
1:B:517:THR:HG22	1:B:545:PRO:O	2.20	0.42
1:B:443:PHE:HB3	1:B:483:GLN:H	1.85	0.42
1:B:512:TYR:HB3	1:B:530:LEU:HD13	2.02	0.41
1:B:424:GLN:N	1:B:424:GLN:OE1	2.53	0.41
1:A:607:PRO:HA	1:A:608:PRO:HD2	1.70	0.41
3:B:701:AMP:H2'	3:B:701:AMP:N3	2.36	0.41
1:B:465:ILE:HD13	1:B:465:ILE:HA	1.92	0.41
1:B:458:ARG:HH21	1:B:460:GLN:HB3	1.84	0.41
1:B:515:SER:O	1:B:545:PRO:HG3	2.21	0.41
1:B:443:PHE:CE2	1:B:446:ALA:HB2	2.56	0.41
1:B:482:ALA:HB1	1:B:496:GLY:HA3	2.03	0.40
1:B:432:ARG:HA	1:B:432:ARG:HD3	1.75	0.40
1:B:410:SER:HB3	1:B:415:GLN:NE2	2.36	0.40
1:A:574:ASP:OD1	1:A:575:PRO:HD2	2.21	0.40
1:A:526:GLU:HA	1:A:526:GLU:OE1	2.21	0.40
1:A:508:LEU:N	1:A:509:PRO:CD	2.84	0.40
1:B:458:ARG:NH2	1:B:460:GLN:HB3	2.37	0.40
1:A:446:ALA:O	1:A:450:ILE:HG13	2.22	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	200/214 (94%)	183 (92%)	14 (7%)	3 (2%)	15	25
1	B	183/214 (86%)	165 (90%)	12 (7%)	6 (3%)	6	7
All	All	383/428 (90%)	348 (91%)	26 (7%)	9 (2%)	10	14

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	546	GLU
1	B	548	GLU
1	B	598	PRO
1	B	458	ARG
1	B	482	ALA
1	A	453	GLY
1	A	495	ASN
1	B	495	ASN
1	A	549	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	158/169 (94%)	139 (88%)	19 (12%)	7	13
1	B	146/169 (86%)	128 (88%)	18 (12%)	7	12
All	All	304/338 (90%)	267 (88%)	37 (12%)	7	12

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	411	THR
1	A	422	LEU
1	A	423	LEU
1	A	427	ARG
1	A	431	GLU
1	A	442	THR
1	A	465	ILE
1	A	485	GLN
1	A	490	ARG
1	A	494	ARG
1	A	499	LEU
1	A	509	PRO
1	A	513	ARG
1	A	518	LEU
1	A	531	ILE
1	A	542	ILE
1	A	560	LEU

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Mol	Chain	Res	Type
1	A	603	GLN
1	A	606	LYS
1	B	403	ASP
1	B	406	ASP
1	B	416	ASN
1	B	421	ARG
1	B	423	LEU
1	B	462	LEU
1	B	467	ARG
1	B	477	LEU
1	B	493	ILE
1	B	495	ASN
1	B	500	ARG
1	B	506	SER
1	B	509	PRO
1	B	518	LEU
1	B	540	ASP
1	B	552	LEU
1	B	572	PRO
1	B	582	LEU

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	428	GLN
1	A	440	HIS
1	B	415	GLN
1	B	416	ASN
1	B	428	GLN
1	B	448	GLN
1	B	577	ASN

### 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 ⓘ

3 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	NCA	A	700	-	9,9,9	1.82	2 (22%)	11,11,11	1.74	3 (27%)
2	NCA	B	700	-	9,9,9	2.16	3 (33%)	11,11,11	2.63	6 (54%)
3	AMP	B	701	-	25,25,25	1.35	4 (16%)	38,38,38	2.88	16 (42%)

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	NCA	A	700	-	-	0/4/4/4	0/1/1/1
2	NCA	B	700	-	-	0/4/4/4	0/1/1/1
3	AMP	B	701	-	-	0/10/26/26	0/1/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	700	NCA	C2-N1	4.76	1.45	1.34
2	A	700	NCA	C3-C7	3.36	1.56	1.50
2	A	700	NCA	C2-N1	3.30	1.41	1.34
3	B	701	AMP	C2'-C1'	-3.29	1.48	1.53
3	B	701	AMP	P-O3P	2.99	1.65	1.54
2	B	700	NCA	C4-C3	-2.47	1.35	1.39
3	B	701	AMP	C5-C4	-2.17	1.35	1.40
2	B	700	NCA	C6-N1	2.09	1.40	1.33
3	B	701	AMP	C8-N9	-2.09	1.33	1.36

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	701	AMP	N3-C2-N1	-9.03	121.16	128.71
3	B	701	AMP	N3-C4-N9	6.96	138.00	125.43
3	B	701	AMP	C8-N9-C4	5.55	111.14	106.90
2	B	700	NCA	C4-C5-C6	5.30	126.80	118.95
3	B	701	AMP	C4'-O4'-C1'	5.23	115.43	109.75
3	B	701	AMP	O4'-C1'-C2'	-3.89	100.80	106.77
3	B	701	AMP	C5-C4-N3	-3.85	117.32	125.70
3	B	701	AMP	O3P-P-O5'	-3.73	96.36	106.65
2	B	700	NCA	O7-C7-N7	3.57	127.75	122.59
3	B	701	AMP	O2P-P-O5'	3.21	115.49	106.65
2	A	700	NCA	C3-C7-N7	3.14	121.35	117.77
2	B	700	NCA	C5-C4-C3	-3.01	116.42	120.32
3	B	701	AMP	O4'-C4'-C5'	-2.87	99.09	109.36
2	A	700	NCA	C4-C3-C2	-2.74	114.28	117.64
2	B	700	NCA	C3-C7-N7	-2.66	114.74	117.77
3	B	701	AMP	N7-C8-N9	-2.59	107.04	114.36
3	B	701	AMP	C2-N3-C4	2.56	121.29	114.01
3	B	701	AMP	O4'-C4'-C3'	-2.44	100.22	105.17
2	B	700	NCA	C5-C6-N1	-2.44	114.91	122.48
2	B	700	NCA	C4-C3-C2	2.40	120.58	117.64
3	B	701	AMP	C5-C4-N9	-2.39	103.72	107.16
2	A	700	NCA	C5-C4-C3	2.32	123.33	120.32
3	B	701	AMP	C6-C5-C4	2.26	121.39	117.25
3	B	701	AMP	O3P-P-O2P	2.08	115.70	107.61
3	B	701	AMP	C1'-N9-C4	-2.06	123.08	126.64

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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.