



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 05:39 AM UTC

PDB ID : 5JNF / pdb_00005jnf
Title : Crystal structure of the LgrA initiation module excluding the Asub domain:
F-A-delta-sub
Authors : Reimer, J.M.; Aloise, M.N.; Schmeing, T.M.
Deposited on : 2016-04-29
Resolution : 2.75 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

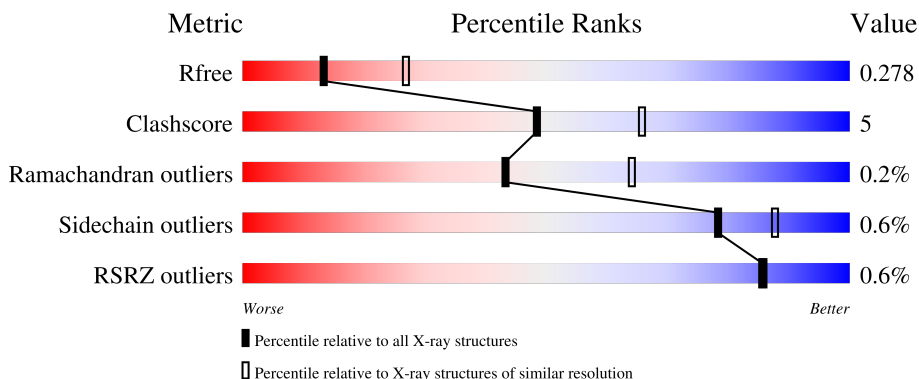
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

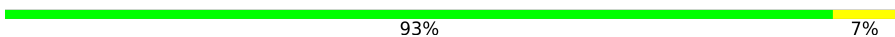
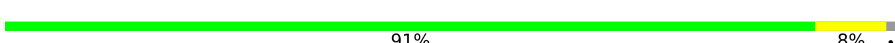
The reported resolution of this entry is 2.75 Å.

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}	180053	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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 of 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	587	
1	B	587	
1	C	587	
1	D	587	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 36748 atoms, of which 18083 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.

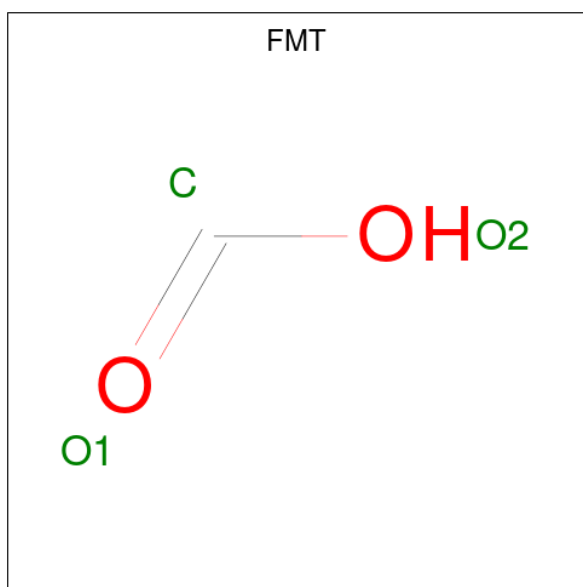
- Molecule 1 is a protein called Linear gramicidin synthase subunit A.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	C	583	Total	C	H	N	O	S	0	0	0
			9221	2957	4577	796	869	22			
1	A	585	Total	C	H	N	O	S	0	0	0
			9239	2962	4585	798	872	22			
1	B	580	Total	C	H	N	O	S	0	0	0
			9033	2918	4452	777	864	22			
1	D	581	Total	C	H	N	O	S	0	0	0
			9050	2920	4459	783	866	22			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-1	GLY	-	expression tag	UNP Q70LM7
C	0	ALA	-	expression tag	UNP Q70LM7
C	1	MET	-	expression tag	UNP Q70LM7
C	2	GLY	-	expression tag	UNP Q70LM7
A	-1	GLY	-	expression tag	UNP Q70LM7
A	0	ALA	-	expression tag	UNP Q70LM7
A	1	MET	-	expression tag	UNP Q70LM7
A	2	GLY	-	expression tag	UNP Q70LM7
B	-1	GLY	-	expression tag	UNP Q70LM7
B	0	ALA	-	expression tag	UNP Q70LM7
B	1	MET	-	expression tag	UNP Q70LM7
B	2	GLY	-	expression tag	UNP Q70LM7
D	-1	GLY	-	expression tag	UNP Q70LM7
D	0	ALA	-	expression tag	UNP Q70LM7
D	1	MET	-	expression tag	UNP Q70LM7
D	2	GLY	-	expression tag	UNP Q70LM7

- Molecule 2 is FORMIC ACID (CCD ID: FMT) (formula: CH₂O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	H	O	0	0
			5	1	2	2		
2	C	1	Total	C	H	O	0	0
			5	1	2	2		
2	C	1	Total	C	H	O	0	0
			5	1	2	2		
2	C	1	Total	C	H	O	0	0
			5	1	2	2		
2	A	1	Total	C	H	O	0	0
			5	1	2	2		

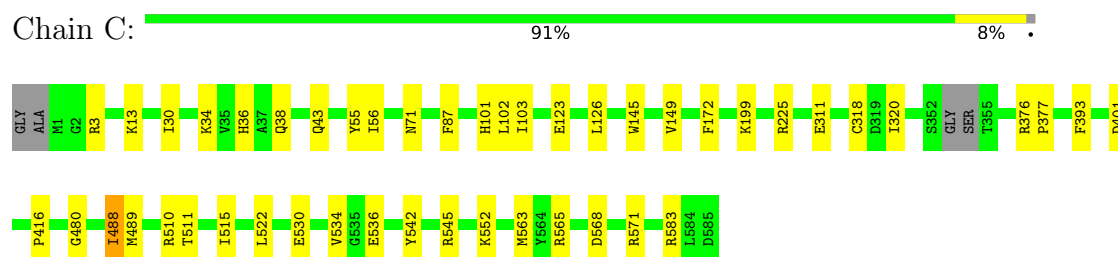
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	83	Total	O	0	0
			83	83		
3	A	73	Total	O	0	0
			73	73		
3	B	10	Total	O	0	0
			10	10		
3	D	14	Total	O	0	0
			14	14		

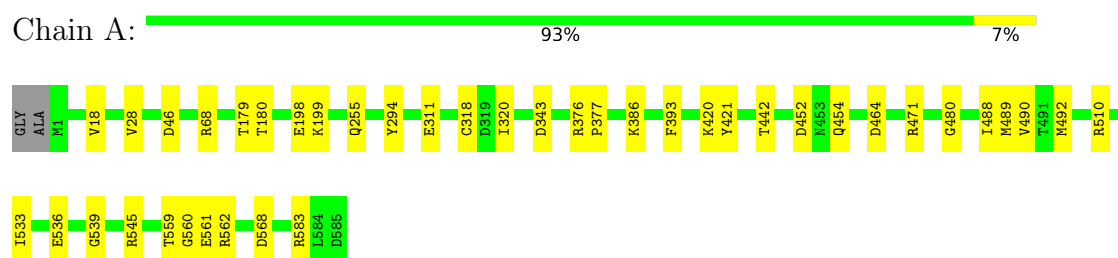
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

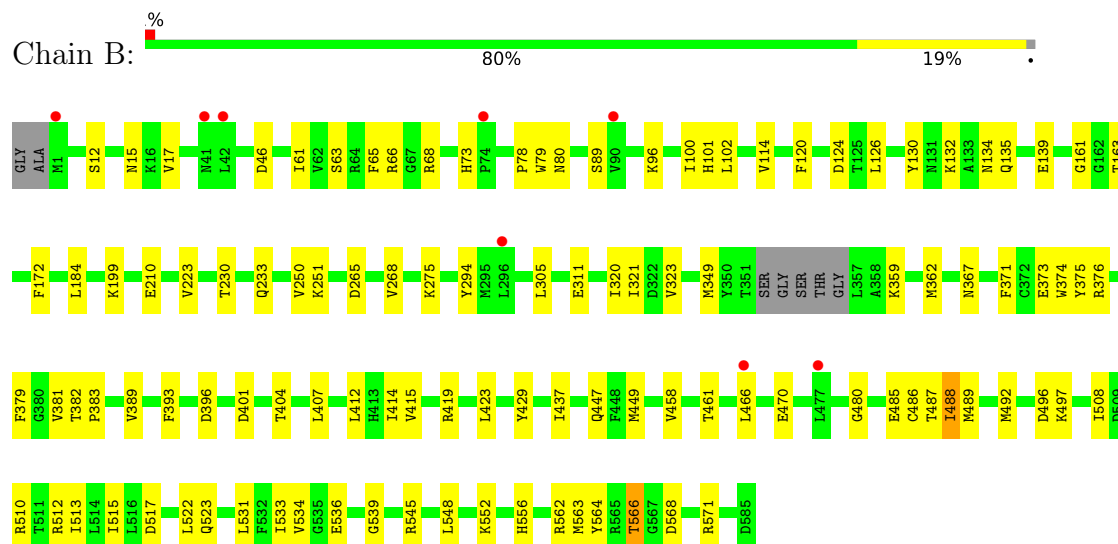
- Molecule 1: Linear gramicidin synthase subunit A



- Molecule 1: Linear gramicidin synthase subunit A

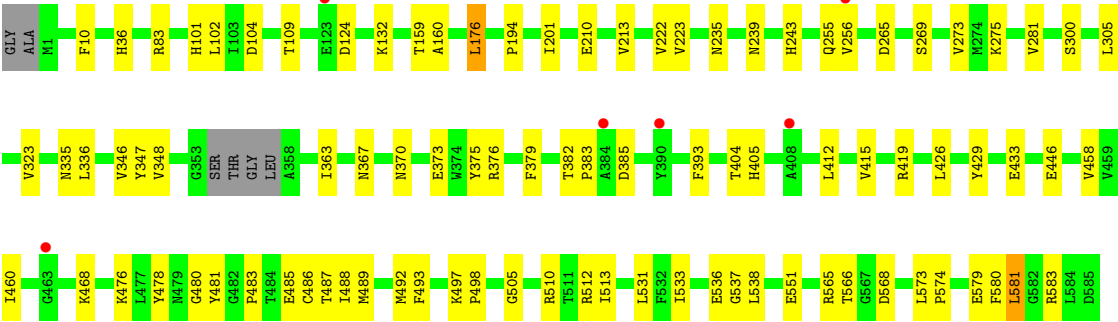


- Molecule 1: Linear gramicidin synthase subunit A



- Molecule 1: Linear gramicidin synthase subunit A

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	161.31Å 161.31Å 139.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.79 – 2.75 44.79 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.8 (44.79-2.75) 99.9 (44.79-2.75)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.57 (at 2.55Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.247 , 0.275 0.252 , 0.278	Depositor DCC
R_{free} test set	5817 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	54.0	Xtriage
Anisotropy	0.284	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 72.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.478 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	36748	wwPDB-VP
Average B, all atoms (Å ²)	106.0	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 37.19 % 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 4.4522e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.16	0/4760	0.36	0/6459
1	B	0.23	1/4686 (0.0%)	0.38	0/6371
1	C	0.15	0/4749	0.35	0/6443
1	D	0.22	1/4696 (0.0%)	0.39	2/6383 (0.0%)
All	All	0.19	2/18891 (0.0%)	0.37	2/25656 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	36	HIS	ND1-CE1	5.26	1.37	1.32
1	B	73	HIS	C-N	5.22	1.46	1.33

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	36	HIS	CB-CG-CD2	-6.82	122.33	131.20
1	D	36	HIS	CB-CG-ND1	5.79	131.38	122.70

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4654	4585	4586	24	0
1	B	4581	4452	4452	88	0
1	C	4644	4577	4577	30	0
1	D	4591	4459	4460	63	0
2	A	3	2	1	0	0
2	C	12	8	4	0	0
3	A	73	0	0	2	0
3	B	10	0	0	1	0
3	C	83	0	0	2	0
3	D	14	0	0	0	0
All	All	18665	18083	18080	200	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 5.

All (200) 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:D:210:GLU:OE2	1:D:275:LYS:NZ	1.64	1.27
1:B:210:GLU:OE2	1:B:275:LYS:NZ	1.74	1.20
1:C:510:ARG:NH1	1:C:536:GLU:OE1	2.09	0.85
1:C:123:GLU:OE2	1:D:574:PRO:O	1.98	0.80
1:A:199:LYS:O	1:A:510:ARG:NH2	2.15	0.79
1:C:199:LYS:O	1:C:510:ARG:NH2	2.18	0.77
1:D:481:TYR:OH	1:D:568:ASP:OD2	2.02	0.76
1:B:531:LEU:O	1:B:566:THR:HG23	1.85	0.76
1:B:367:ASN:HB3	1:B:487:THR:HG21	1.67	0.76
1:D:551:GLU:O	1:D:565:ARG:NH2	2.20	0.75
1:B:486:CYS:O	1:B:487:THR:OG1	2.06	0.73
1:D:375:TYR:CD2	1:D:404:THR:HG21	2.27	0.69
1:D:222:VAL:HG12	1:D:412:LEU:HB2	1.74	0.68
1:D:210:GLU:CD	1:D:275:LYS:NZ	2.51	0.68
1:C:568:ASP:OD1	1:C:583:ARG:NH1	2.27	0.67
1:B:100:ILE:HB	1:B:114:VAL:HG12	1.78	0.66
1:B:79:TRP:HE3	1:B:163:THR:HG23	1.62	0.65
1:B:375:TYR:CD2	1:B:404:THR:CG2	2.81	0.63
1:B:124:ASP:OD1	1:B:132:LYS:NZ	2.31	0.63
1:A:568:ASP:OD1	1:A:583:ARG:NH1	2.29	0.63
1:B:556:HIS:HB2	1:B:563:MET:CE	2.27	0.63
1:B:61:ILE:O	1:B:65:PHE:HD1	1.82	0.63
1:B:80:ASN:H	1:B:163:THR:HG1	1.47	0.62
1:B:311:GLU:HG3	1:B:320:ILE:HD13	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:15:ASN:HB3	1:B:17:VAL:HG12	1.82	0.61
1:B:199:LYS:O	1:B:510:ARG:NH2	2.33	0.61
1:B:359:LYS:HG2	1:B:545:ARG:NH1	2.16	0.60
1:B:359:LYS:CE	1:B:548:LEU:HD21	2.31	0.60
1:B:230:THR:HG23	1:B:233:GLN:H	1.68	0.59
1:C:3:ARG:NH2	1:A:343:ASP:OD2	2.35	0.59
1:D:531:LEU:O	1:D:566:THR:OG1	2.20	0.59
1:B:130:TYR:O	1:B:134:ASN:ND2	2.34	0.59
1:D:235:ASN:ND2	1:D:335:ASN:OD1	2.36	0.59
1:A:452:ASP:OD1	1:A:454:GLN:NE2	2.36	0.58
1:C:34:LYS:NZ	1:C:56:ILE:O	2.36	0.58
1:D:243:HIS:CD2	1:D:336:LEU:HD11	2.38	0.58
1:C:401:ASP:OD1	3:C:701:HOH:O	2.17	0.57
1:A:539:GLY:O	1:A:562:ARG:NH2	2.32	0.57
1:D:210:GLU:CD	1:D:275:LYS:HZ1	2.10	0.57
1:D:375:TYR:CG	1:D:404:THR:HG21	2.40	0.57
1:A:386:LYS:NZ	3:A:704:HOH:O	2.38	0.57
1:B:552:LYS:HB3	1:B:564:TYR:CE1	2.39	0.56
1:B:294:TYR:OH	1:B:545:ARG:NH2	2.38	0.56
1:B:101:HIS:ND1	1:B:102:LEU:O	2.37	0.56
1:B:489:MET:HE1	1:B:492:MET:HE1	1.88	0.56
1:B:311:GLU:CG	1:B:320:ILE:HD13	2.36	0.55
1:B:172:PHE:CZ	1:B:522:LEU:CD1	2.89	0.55
1:D:373:GLU:OE1	1:D:376:ARG:NH2	2.34	0.55
1:D:367:ASN:HB3	1:D:487:THR:HG21	1.89	0.54
1:B:375:TYR:CD2	1:B:404:THR:HG21	2.43	0.54
1:B:250:VAL:C	1:B:251:LYS:HD2	2.33	0.54
1:B:539:GLY:O	1:B:562:ARG:NH2	2.38	0.54
1:B:480:GLY:HA3	1:B:489:MET:CE	2.39	0.53
1:D:347:TYR:O	1:D:363:ILE:HG22	2.09	0.53
1:D:201:ILE:HD11	1:D:536:GLU:HG2	1.91	0.52
1:A:510:ARG:NH1	1:A:536:GLU:OE2	2.41	0.52
1:B:389:VAL:HG11	1:B:414:ILE:HD13	1.92	0.52
1:B:556:HIS:CD2	1:B:563:MET:HE1	2.44	0.52
1:D:255:GLN:OE1	1:D:300:SER:N	2.37	0.52
1:D:281:VAL:HG22	1:D:348:VAL:HB	1.91	0.51
1:D:101:HIS:ND1	1:D:102:LEU:O	2.43	0.51
1:D:458:VAL:HG22	1:D:476:LYS:HD2	1.91	0.51
1:D:480:GLY:HA3	1:D:489:MET:CE	2.40	0.51
1:B:375:TYR:CE2	1:B:381:VAL:HG21	2.46	0.51
1:B:423:LEU:HD12	1:B:447:GLN:HB3	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:492:MET:HE3	1:D:492:MET:HA	1.93	0.51
1:A:559:THR:HG21	1:A:561:GLU:OE2	2.11	0.50
1:B:79:TRP:CE3	1:B:163:THR:HG23	2.42	0.50
1:C:3:ARG:NH1	1:C:43:GLN:O	2.44	0.50
1:C:515:ILE:HD13	1:C:571:ARG:HA	1.94	0.50
1:D:194:PRO:HD3	1:D:512:ARG:NH1	2.27	0.50
1:B:265:ASP:HA	1:B:268:VAL:HG12	1.94	0.50
1:B:265:ASP:HA	1:B:268:VAL:CG1	2.41	0.50
1:A:311:GLU:HG3	1:A:320:ILE:HD13	1.93	0.50
1:D:513:ILE:CD1	1:D:533:ILE:HG12	2.42	0.50
1:C:172:PHE:CZ	1:C:522:LEU:CD1	2.95	0.49
1:B:12:SER:O	1:B:15:ASN:ND2	2.45	0.49
1:B:349:MET:HE1	1:B:396:ASP:N	2.27	0.49
1:B:46:ASP:O	1:B:68:ARG:HD2	2.12	0.49
1:B:375:TYR:CD2	1:B:404:THR:HG22	2.47	0.49
1:B:172:PHE:CZ	1:B:522:LEU:HD11	2.48	0.49
1:B:61:ILE:O	1:B:65:PHE:CD1	2.64	0.49
1:B:552:LYS:O	1:B:564:TYR:HA	2.12	0.49
1:C:311:GLU:HG3	1:C:320:ILE:HD13	1.94	0.49
1:A:18:VAL:HG13	1:A:28:VAL:HG11	1.94	0.48
1:D:446:GLU:HB3	1:D:468:LYS:HD2	1.95	0.48
1:B:487:THR:O	1:B:489:MET:N	2.46	0.48
1:D:478:TYR:HB3	1:D:492:MET:HE2	1.94	0.48
1:B:531:LEU:HD22	1:B:566:THR:HG21	1.96	0.48
1:D:487:THR:HG23	1:D:537:GLY:O	2.13	0.48
1:A:442:THR:HG21	1:A:464:ASP:HB2	1.95	0.48
1:D:486:CYS:O	1:D:487:THR:OG1	2.32	0.47
1:B:415:VAL:HG23	1:B:415:VAL:O	2.15	0.47
1:B:512:ARG:O	1:B:534:VAL:HG12	2.15	0.47
1:C:38:GLN:NE2	1:A:198:GLU:OE1	2.46	0.47
1:D:580:PHE:C	1:D:581:LEU:HD12	2.39	0.47
1:C:534:VAL:HG12	1:C:563:MET:HG2	1.96	0.46
1:D:176:LEU:HD12	1:D:176:LEU:N	2.29	0.46
1:B:96:LYS:HD3	1:B:120:PHE:CD1	2.50	0.46
1:B:371:PHE:CE2	1:B:374:TRP:HZ3	2.34	0.46
1:B:389:VAL:HA	1:B:401:ASP:HB3	1.97	0.46
1:D:375:TYR:CD2	1:D:404:THR:CG2	2.97	0.46
1:D:513:ILE:HD13	1:D:533:ILE:HA	1.97	0.46
1:D:483:PRO:HB2	1:D:485:GLU:OE1	2.15	0.46
1:C:36:HIS:NE2	1:A:561:GLU:OE1	2.47	0.46
1:C:71:ASN:HB2	1:C:103:ILE:HD11	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:123:GLU:HG2	1:C:123:GLU:O	2.15	0.46
1:C:225:ARG:NH1	1:C:416:PRO:CB	2.79	0.46
1:B:63:SER:HA	1:B:66:ARG:HG3	1.96	0.46
1:D:210:GLU:C	1:D:213:VAL:HG22	2.41	0.46
1:D:239:ASN:OD1	1:D:275:LYS:HE2	2.16	0.46
1:B:265:ASP:HB3	1:B:323:VAL:HG11	1.97	0.46
1:B:449:MET:HB3	1:B:470:GLU:O	2.16	0.46
1:A:376:ARG:HB3	1:A:377:PRO:HD3	1.98	0.46
1:D:382:THR:OG1	1:D:383:PRO:HD2	2.15	0.46
1:C:87:PHE:HD1	1:C:126:LEU:HD13	1.80	0.45
1:A:294:TYR:OH	1:A:545:ARG:NH2	2.50	0.45
1:D:458:VAL:HG22	1:D:476:LYS:HB2	1.97	0.45
1:C:13:LYS:HG2	1:C:30:ILE:CD1	2.47	0.45
1:B:359:LYS:HD3	1:B:548:LEU:HD21	1.98	0.45
1:B:496:ASP:OD2	1:B:497:LYS:HE3	2.17	0.45
1:B:359:LYS:HE2	1:B:548:LEU:HD21	1.98	0.45
1:D:159:THR:HG22	1:D:160:ALA:O	2.17	0.45
1:D:370:ASN:OD1	1:D:510:ARG:HD2	2.17	0.45
1:B:515:ILE:HD13	1:B:571:ARG:HA	1.97	0.45
1:D:222:VAL:HG12	1:D:412:LEU:HD12	1.99	0.45
1:D:265:ASP:HB3	1:D:323:VAL:HG11	1.99	0.45
1:B:80:ASN:N	1:B:163:THR:OG1	2.35	0.44
1:D:566:THR:O	1:D:583:ARG:NH1	2.48	0.44
1:B:485:GLU:HG3	1:B:564:TYR:OH	2.16	0.44
1:C:480:GLY:HA3	1:C:489:MET:CE	2.47	0.44
1:B:517:ASP:OD1	1:B:523:GLN:NE2	2.51	0.44
1:C:145:TRP:CE2	1:C:149:VAL:CG2	3.01	0.44
1:B:512:ARG:HB2	1:B:534:VAL:HG13	1.98	0.44
1:D:487:THR:O	1:D:489:MET:N	2.51	0.44
1:C:376:ARG:HB3	1:C:377:PRO:HD3	2.00	0.44
1:B:126:LEU:HD11	1:B:184:LEU:HD22	1.99	0.43
1:B:135:GLN:O	1:B:139:GLU:HG3	2.17	0.43
1:D:573:LEU:HD21	1:D:579:GLU:OE2	2.18	0.43
1:C:552:LYS:HD3	3:C:729:HOH:O	2.18	0.43
1:B:389:VAL:HG11	1:B:414:ILE:CD1	2.48	0.43
1:D:379:PHE:CZ	1:D:460:ILE:HD11	2.53	0.43
1:B:382:THR:OG1	1:B:383:PRO:HD2	2.19	0.43
1:B:480:GLY:HA3	1:B:489:MET:HE1	2.00	0.43
1:D:492:MET:HE3	1:D:493:PHE:H	1.83	0.43
1:B:15:ASN:CB	1:B:17:VAL:HG12	2.47	0.43
1:B:61:ILE:HG23	1:B:65:PHE:CE1	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:243:HIS:NE2	1:D:336:LEU:HD11	2.32	0.43
1:A:46:ASP:O	1:A:68:ARG:HD2	2.19	0.43
1:B:437:ILE:HA	1:B:458:VAL:O	2.17	0.43
1:D:256:VAL:HB	1:D:273:VAL:HG11	2.01	0.43
1:D:269:SER:HB3	1:D:305:LEU:HD13	2.01	0.43
1:B:367:ASN:ND2	1:B:536:GLU:O	2.51	0.42
1:C:55:TYR:CE1	1:A:560:GLY:HA3	2.54	0.42
1:B:223:VAL:N	1:B:412:LEU:O	2.51	0.42
1:D:223:VAL:N	1:D:412:LEU:O	2.47	0.42
1:D:493:PHE:HB2	1:D:505:GLY:CA	2.49	0.42
1:B:373:GLU:OE1	1:B:376:ARG:NH2	2.42	0.42
1:B:461:THR:HG21	1:B:466:LEU:HD13	2.01	0.42
1:B:486:CYS:C	1:B:487:THR:HG1	2.17	0.42
1:C:13:LYS:HG2	1:C:30:ILE:HD12	2.01	0.42
1:A:480:GLY:HA3	1:A:489:MET:HE3	2.01	0.42
1:B:89:SER:HB3	1:B:96:LYS:HG2	2.00	0.42
1:A:490:VAL:HG21	1:A:533:ILE:CD1	2.50	0.42
1:B:512:ARG:HB2	1:B:534:VAL:CG1	2.49	0.42
1:D:124:ASP:OD1	1:D:132:LYS:NZ	2.52	0.42
1:A:420:LYS:HE3	1:A:421:TYR:CE2	2.55	0.42
1:B:419:ARG:HH11	1:B:429:TYR:HB2	1.85	0.42
1:C:542:TYR:HB2	1:C:545:ARG:HB2	2.02	0.41
1:B:78:PRO:CG	1:B:161:GLY:O	2.68	0.41
1:D:10:PHE:CZ	1:D:83:ARG:NH2	2.88	0.41
1:C:145:TRP:CZ2	1:C:149:VAL:HG22	2.55	0.41
1:C:101:HIS:ND1	1:C:102:LEU:O	2.52	0.41
1:B:359:LYS:CD	1:B:548:LEU:HD21	2.50	0.41
1:C:530:GLU:OE2	1:C:565:ARG:HG2	2.20	0.41
1:B:513:ILE:HD13	1:B:533:ILE:HA	2.02	0.41
1:D:104:ASP:OD2	1:D:109:THR:C	2.63	0.41
1:D:429:TYR:CE2	1:D:433:GLU:HG3	2.56	0.41
1:D:367:ASN:ND2	1:D:538:LEU:O	2.51	0.41
1:D:405:HIS:HB2	1:D:412:LEU:HD21	2.03	0.41
1:D:497:LYS:HB3	1:D:498:PRO:HD2	2.02	0.41
1:A:489:MET:CE	1:A:492:MET:HE1	2.51	0.41
1:B:376:ARG:HH11	1:B:407:LEU:HB3	1.86	0.41
1:B:488:ILE:C	1:B:489:MET:HG3	2.45	0.41
1:A:179:THR:HG22	1:A:180:THR:N	2.36	0.41
1:B:379:PHE:HD2	1:B:458:VAL:HG11	1.84	0.41
1:B:362:MET:HE1	3:B:609:HOH:O	2.20	0.41
1:B:371:PHE:CE2	1:B:374:TRP:CZ3	3.08	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:210:GLU:O	1:D:213:VAL:HG22	2.21	0.41
1:A:454:GLN:OE1	1:A:454:GLN:N	2.54	0.41
1:B:566:THR:OG1	1:B:568:ASP:N	2.50	0.41
1:D:419:ARG:NH1	1:D:429:TYR:HB2	2.36	0.41
1:A:471:ARG:O	3:A:701:HOH:O	2.22	0.40
1:B:305:LEU:HD23	1:B:321:ILE:HB	2.02	0.40
1:C:488:ILE:C	1:C:489:MET:HG3	2.46	0.40
1:B:566:THR:HG1	1:B:568:ASP:H	1.67	0.40
1:D:382:THR:HG22	1:D:385:ASP:OD2	2.20	0.40
1:B:531:LEU:CD2	1:B:566:THR:HG21	2.50	0.40
1:D:346:VAL:CG1	1:D:363:ILE:HG23	2.51	0.40
1:D:415:VAL:O	1:D:415:VAL:HG23	2.21	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	583/587 (99%)	565 (97%)	17 (3%)	1 (0%)	43	64
1	B	576/587 (98%)	560 (97%)	14 (2%)	2 (0%)	36	56
1	C	579/587 (99%)	564 (97%)	14 (2%)	1 (0%)	43	64
1	D	577/587 (98%)	560 (97%)	16 (3%)	1 (0%)	43	64
All	All	2315/2348 (99%)	2249 (97%)	61 (3%)	5 (0%)	43	64

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	488	ILE
1	A	488	ILE
1	C	488	ILE

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Mol	Chain	Res	Type
1	D	488	ILE
1	B	508	ILE

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	501/501 (100%)	498 (99%)	3 (1%)	78	88
1	B	488/501 (97%)	486 (100%)	2 (0%)	84	91
1	C	500/501 (100%)	497 (99%)	3 (1%)	78	88
1	D	489/501 (98%)	485 (99%)	4 (1%)	73	84
All	All	1978/2004 (99%)	1966 (99%)	12 (1%)	78	88

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	318	CYS
1	C	393	PHE
1	C	511	THR
1	A	255	GLN
1	A	318	CYS
1	A	393	PHE
1	B	393	PHE
1	B	566	THR
1	D	176	LEU
1	D	393	PHE
1	D	426	LEU
1	D	581	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	212	GLN
1	C	370	ASN

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Mol	Chain	Res	Type
1	C	427	ASN
1	C	454	GLN
1	A	106	HIS
1	A	243	HIS
1	A	324	HIS
1	A	370	ASN
1	B	131	ASN
1	B	324	HIS
1	D	106	HIS
1	D	240	GLN
1	D	307	ASN
1	D	310	HIS
1	D	427	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

5 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	FMT	C	604	-	2,2,2	0.51	0	1,1,1	0.39	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	FMT	C	601	-	2,2,2	0.71	0	1,1,1	0.55	0
2	FMT	C	603	-	2,2,2	0.48	0	1,1,1	0.50	0
2	FMT	C	602	-	2,2,2	0.57	0	1,1,1	0.51	0
2	FMT	A	601	-	2,2,2	0.49	0	1,1,1	0.56	0

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.

5.7 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, 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	585/587 (99%)	-1.01	0 100 100	44, 66, 107, 203	0
1	B	580/587 (98%)	-0.42	8 (1%) 73 73	100, 143, 161, 179	0
1	C	583/587 (99%)	-1.02	0 100 100	42, 66, 103, 126	0
1	D	581/587 (98%)	-0.44	6 (1%) 79 79	106, 142, 162, 176	0
All	All	2329/2348 (99%)	-0.72	14 (0%) 85 86	42, 121, 157, 203	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	41	ASN	2.9
1	D	463	GLY	2.7
1	D	384	ALA	2.7
1	B	74	PRO	2.6
1	D	390	TYR	2.5
1	D	123	GLU	2.5
1	D	256	VAL	2.5
1	B	42	LEU	2.4
1	B	466	LEU	2.4
1	B	296	LEU	2.3
1	B	477	LEU	2.2
1	D	408	ALA	2.1
1	B	1	MET	2.0
1	B	90	VAL	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 [i](#)

There are no oligosaccharides 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($\text{\AA}^2$)	Q<0.9
2	FMT	C	601	3/3	0.98	0.09	76,77,91,96	0
2	FMT	C	602	3/3	0.98	0.12	69,69,82,82	0
2	FMT	C	603	3/3	0.98	0.09	59,65,78,79	0
2	FMT	C	604	3/3	0.98	0.13	64,65,78,84	0
2	FMT	A	601	3/3	0.99	0.10	62,66,74,80	0

6.5 Other polymers [i](#)

There are no such residues in this entry.