



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:57 PM GMT

PDB ID : 4GR4  
Title : Crystal structure of SlgN1deltaAsub  
Authors : Herbst, D.A.; Zocher, G.; Stehle, T.  
Deposited on : 2012-08-24  
Resolution : 2.44 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  
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

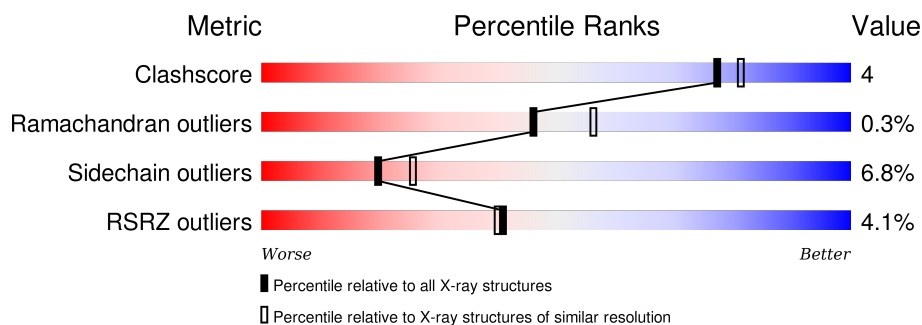
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.44 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	102246	1071 (2.46-2.42)
Ramachandran outliers	100387	1065 (2.46-2.42)
Sidechain outliers	100360	1065 (2.46-2.42)
RSRZ outliers	91569	1005 (2.46-2.42)

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. 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	469	<div> <div>5%</div> <div>84%</div> <div>11%</div> <div>..</div> </div>
1	B	469	<div> <div>%</div> <div>84%</div> <div>10%</div> <div>..</div> </div>
1	C	469	<div> <div>5%</div> <div>86%</div> <div>11%</div> <div>..</div> </div>
1	D	469	<div> <div>5%</div> <div>83%</div> <div>11%</div> <div>.. 5%</div> </div>

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 13783 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 Non-ribosomal peptide synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	454	Total	C	N	O	S	0	0	0
			3357	2122	599	623	13			
1	B	451	Total	C	N	O	S	0	0	0
			3339	2109	595	623	12			
1	C	458	Total	C	N	O	S	0	0	0
			3353	2121	595	625	12			
1	D	444	Total	C	N	O	S	0	0	0
			3291	2080	585	614	12			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP D1GLU5
A	-1	SER	-	EXPRESSION TAG	UNP D1GLU5
A	0	HIS	-	EXPRESSION TAG	UNP D1GLU5
B	-2	GLY	-	EXPRESSION TAG	UNP D1GLU5
B	-1	SER	-	EXPRESSION TAG	UNP D1GLU5
B	0	HIS	-	EXPRESSION TAG	UNP D1GLU5
C	-2	GLY	-	EXPRESSION TAG	UNP D1GLU5
C	-1	SER	-	EXPRESSION TAG	UNP D1GLU5
C	0	HIS	-	EXPRESSION TAG	UNP D1GLU5
D	-2	GLY	-	EXPRESSION TAG	UNP D1GLU5
D	-1	SER	-	EXPRESSION TAG	UNP D1GLU5
D	0	HIS	-	EXPRESSION TAG	UNP D1GLU5

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Cl	0	0
			1	1		

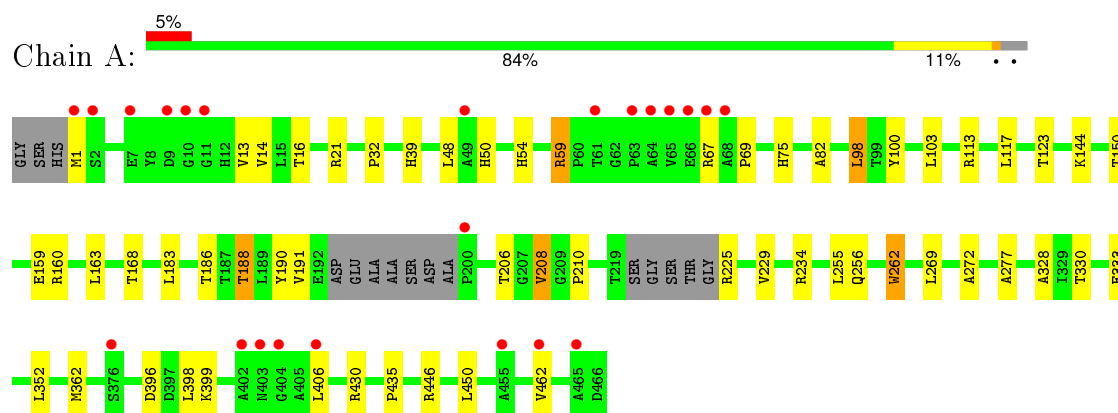
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	139	Total 139	O 139	0	0
3	B	108	Total 108	O 108	0	0
3	C	85	Total 85	O 85	0	0
3	D	110	Total 110	O 110	0	0

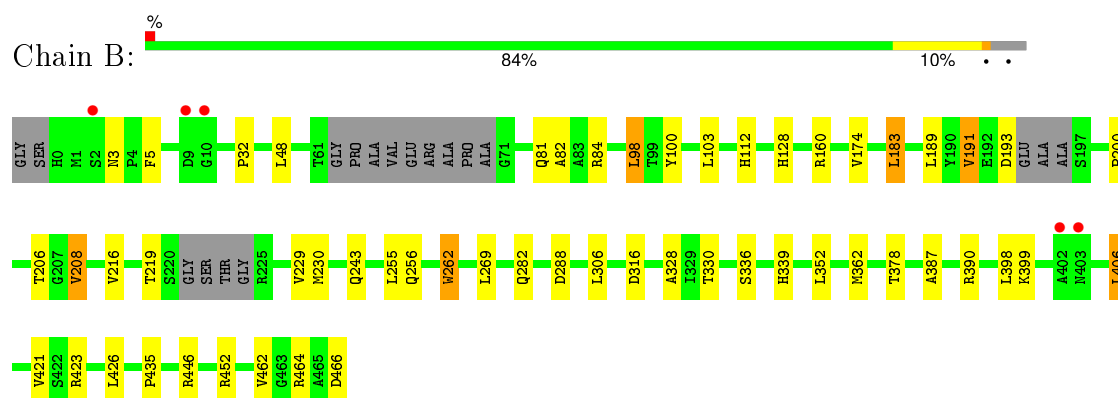
### 3 Residue-property plots [i](#)

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

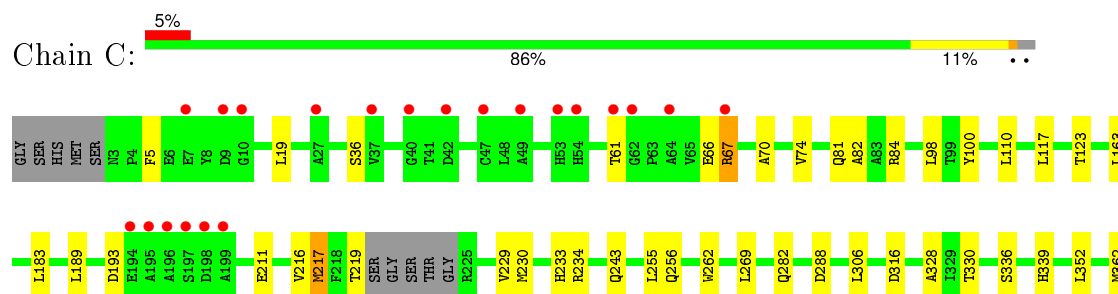
- Molecule 1: Non-ribosomal peptide synthetase

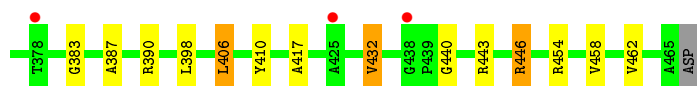


- Molecule 1: Non-ribosomal peptide synthetase

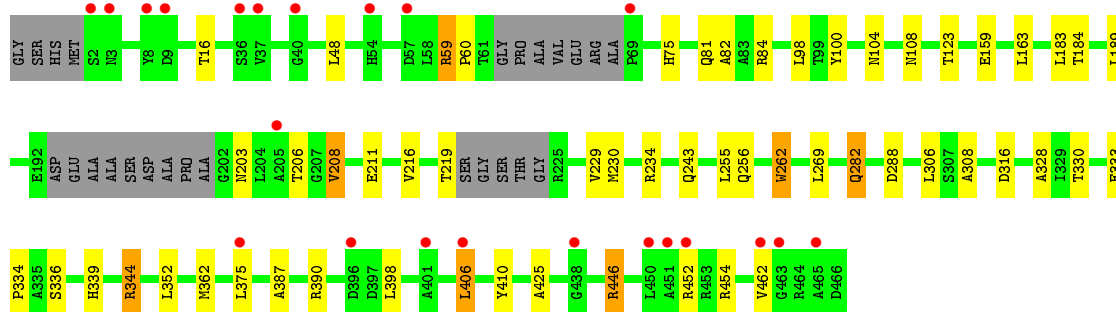
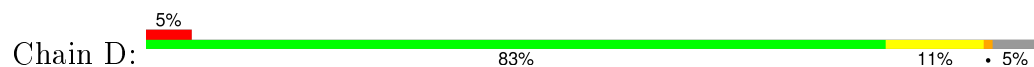


- Molecule 1: Non-ribosomal peptide synthetase





- Molecule 1: Non-ribosomal peptide synthetase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	88.11Å 148.61Å 109.58Å 90.00° 113.34° 90.00°	Depositor
Resolution (Å)	35.72 – 2.44 35.63 – 2.44	Depositor EDS
% Data completeness (in resolution range)	99.6 (35.72-2.44) 99.6 (35.63-2.44)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.12 (at 2.45Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, $R_{free}$	0.180 , 0.213 (Not available) , (Not available)	Depositor DCC
$R_{free}$ test set	NotAvailable	DCC
Wilson B-factor (Å <sup>2</sup> )	36.5	Xtriage
Anisotropy	0.195	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 39.0	EDS
Estimated twinning fraction	0.118 for h,-k,-h-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.44$ , $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	0 of 95737 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	13783	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.54% of the height of the origin peak. No significant pseudotranslation is detected.*

<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, 0.375 respectively for untwinned datasets, and 0.333, 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:  
CL

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	0.52	0/3438	0.75	0/4686
1	B	0.50	0/3419	0.72	0/4659
1	C	0.51	0/3436	0.72	1/4693 (0.0%)
1	D	0.50	0/3370	0.71	0/4591
All	All	0.51	0/13663	0.72	1/18629 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	217	MET	CB-CA-C	5.01	120.42	110.40

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	3357	0	3277	20	0
1	B	3339	0	3243	26	0
1	C	3353	0	3244	25	0
1	D	3291	0	3195	28	0
2	B	1	0	0	0	0
3	A	139	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	108	0	0	0	0
3	C	85	0	0	0	0
3	D	110	0	0	0	0
All	All	13783	0	12959	98	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 4.

All (98) 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:C:110:LEU:HD11	1:C:189:LEU:HD21	1.57	0.84
1:D:104:ASN:HD21	1:D:203:ASN:HD21	1.33	0.76
1:C:406:LEU:HD11	1:C:462:VAL:HG11	1.72	0.71
1:B:406:LEU:HD11	1:B:462:VAL:HG11	1.74	0.70
1:D:406:LEU:HD11	1:D:462:VAL:HG11	1.75	0.68
1:D:108:ASN:HD21	1:D:203:ASN:HD22	1.42	0.67
1:C:282:GLN:HE22	1:C:288:ASP:H	1.42	0.67
1:D:219:THR:HG22	1:D:262:TRP:HH2	1.58	0.66
1:C:243:GLN:HE22	1:C:387:ALA:H	1.43	0.66
1:B:282:GLN:HE22	1:B:288:ASP:H	1.44	0.66
1:C:5:PHE:CZ	1:C:432:VAL:HG11	2.32	0.65
1:D:211:GLU:OE1	1:D:234:ARG:HD3	1.97	0.64
1:D:243:GLN:HE22	1:D:387:ALA:H	1.44	0.64
1:C:211:GLU:OE1	1:C:234:ARG:HD3	1.98	0.63
1:D:282:GLN:HE22	1:D:288:ASP:H	1.46	0.62
1:B:216:VAL:HG22	1:B:230:MET:HB3	1.80	0.61
1:B:219:THR:HG22	1:B:262:TRP:HH2	1.66	0.60
1:C:216:VAL:HG22	1:C:230:MET:HB3	1.84	0.60
1:D:216:VAL:HG22	1:D:230:MET:HB3	1.84	0.59
1:B:243:GLN:HE22	1:B:387:ALA:H	1.47	0.59
1:D:108:ASN:HD21	1:D:203:ASN:ND2	2.01	0.59
1:A:186:THR:O	1:A:188:THR:HG22	2.04	0.57
1:C:410:TYR:CE2	1:C:446:ARG:HB3	2.40	0.55
1:D:219:THR:HG22	1:D:262:TRP:CH2	2.38	0.55
1:D:410:TYR:CE2	1:D:446:ARG:HB3	2.42	0.55
1:C:74:VAL:H	1:C:233:HIS:HD2	1.56	0.54
1:B:219:THR:HG22	1:B:262:TRP:CH2	2.43	0.54
1:A:272:ALA:HA	1:A:277:ALA:HB3	1.90	0.53
1:A:450:LEU:C	1:A:462:VAL:HG22	2.27	0.53
1:A:14:VAL:HB	1:A:39:HIS:HB3	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:19:LEU:HA	1:C:61:THR:HG22	1.90	0.53
1:B:336:SER:OG	1:B:339:HIS:HD2	1.92	0.52
1:C:336:SER:OG	1:C:339:HIS:HD2	1.93	0.52
1:B:306:LEU:O	1:B:330:THR:HA	2.10	0.52
1:D:308:ALA:HB3	1:D:333:GLU:HG2	1.93	0.51
1:D:336:SER:OG	1:D:339:HIS:HD2	1.94	0.51
1:A:450:LEU:HB2	1:A:462:VAL:CG2	2.41	0.51
1:B:328:ALA:HB3	1:B:352:LEU:HD23	1.93	0.51
1:D:59:ARG:HD2	1:D:60:PRO:HD2	1.93	0.51
1:C:74:VAL:H	1:C:233:HIS:CD2	2.29	0.51
1:B:3:ASN:ND2	1:B:446:ARG:HD3	2.26	0.51
1:D:328:ALA:HB3	1:D:352:LEU:HD23	1.93	0.50
1:A:328:ALA:HB3	1:A:352:LEU:HD23	1.93	0.50
1:A:82:ALA:HA	1:A:100:TYR:HB3	1.93	0.50
1:B:316:ASP:OD2	1:B:339:HIS:HE1	1.95	0.50
1:A:117:LEU:HD22	1:C:117:LEU:HD22	1.93	0.50
1:A:21:ARG:NE	3:A:630:HOH:O	2.44	0.50
1:B:174:VAL:CG1	1:B:191:VAL:HG13	2.42	0.49
1:A:150:THR:HG21	1:A:168:THR:HG21	1.95	0.49
1:A:113:ARG:NH1	1:A:190:TYR:O	2.46	0.49
1:C:328:ALA:HB3	1:C:352:LEU:HD23	1.93	0.49
1:C:316:ASP:OD2	1:C:339:HIS:HE1	1.96	0.48
1:B:406:LEU:HD12	1:B:452:ARG:CZ	2.44	0.48
1:D:81:GLN:NE2	1:D:84:ARG:HH21	2.12	0.48
1:D:406:LEU:HB2	1:D:452:ARG:HH21	1.79	0.48
1:C:67:ARG:HH12	1:C:440:GLY:HA3	1.78	0.48
1:C:383:GLY:HA2	1:C:458:VAL:HG13	1.96	0.47
1:C:81:GLN:HE22	1:C:84:ARG:HH21	1.61	0.47
1:D:82:ALA:HA	1:D:100:TYR:HB3	1.97	0.47
1:C:81:GLN:NE2	1:C:84:ARG:HH21	2.12	0.47
1:B:206:THR:OG1	1:B:208:VAL:HG22	2.15	0.47
1:D:333:GLU:HB2	1:D:334:PRO:HD2	1.97	0.47
1:A:98:LEU:HD13	1:A:103:LEU:HB2	1.95	0.47
1:C:5:PHE:CZ	1:C:398:LEU:HD11	2.50	0.47
1:B:81:GLN:NE2	1:B:84:ARG:HH21	2.13	0.46
1:B:32:PRO:HB3	1:B:435:PRO:HA	1.97	0.46
1:D:406:LEU:HD12	1:D:452:ARG:NH2	2.31	0.46
1:D:316:ASP:OD2	1:D:339:HIS:HE1	1.99	0.46
1:A:32:PRO:HB3	1:A:435:PRO:HA	1.98	0.46
1:A:69:PRO:HB2	1:A:210:PRO:HG2	1.98	0.45
1:A:262:TRP:CE3	1:A:262:TRP:HA	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:82:ALA:HA	1:C:100:TYR:HB3	1.97	0.45
1:B:128:HIS:CD2	1:B:183:LEU:HD22	2.52	0.45
1:D:406:LEU:HD12	1:D:452:ARG:HH21	1.82	0.44
1:D:81:GLN:HE22	1:D:84:ARG:HH21	1.64	0.44
1:D:206:THR:OG1	1:D:208:VAL:HG22	2.18	0.44
1:C:306:LEU:O	1:C:330:THR:HA	2.18	0.44
1:B:82:ALA:HA	1:B:100:TYR:HB3	1.99	0.44
1:A:75:HIS:CD2	1:A:144:LYS:HA	2.53	0.44
1:B:5:PHE:HE2	1:B:398:LEU:HD21	1.83	0.44
1:D:344:ARG:HD3	1:D:375:LEU:HD11	1.98	0.44
1:B:262:TRP:HA	1:B:262:TRP:CE3	2.53	0.43
1:A:59:ARG:H	1:A:59:ARG:HG2	1.66	0.43
1:A:1:MET:HG2	1:A:396:ASP:OD1	2.18	0.43
1:B:423:ARG:HH21	1:B:426:LEU:HD22	1.82	0.43
1:A:206:THR:OG1	1:A:208:VAL:HG22	2.18	0.43
1:D:75:HIS:CD2	1:D:75:HIS:H	2.37	0.43
1:B:81:GLN:HE22	1:B:84:ARG:HH21	1.68	0.42
1:C:110:LEU:CD1	1:C:189:LEU:HD21	2.40	0.42
1:D:262:TRP:HA	1:D:262:TRP:CE3	2.55	0.42
1:A:50:HIS:CE1	1:A:54:HIS:CD2	3.08	0.42
1:D:48:LEU:HD22	1:D:425:ALA:HB1	2.01	0.41
1:C:417:ALA:O	1:C:443:ARG:NH2	2.53	0.41
1:B:3:ASN:HD21	1:B:446:ARG:HD3	1.85	0.41
1:B:98:LEU:HD13	1:B:103:LEU:HB2	2.03	0.41
1:C:219:THR:HA	1:C:262:TRP:CZ3	2.56	0.40
1:B:112:HIS:CE1	1:B:200:PRO:HG2	2.56	0.40
1:B:174:VAL:HG12	1:B:191:VAL:HG13	2.04	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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	448/469 (96%)	439 (98%)	8 (2%)	1 (0%)	52	64
1	B	443/469 (94%)	434 (98%)	8 (2%)	1 (0%)	52	64
1	C	454/469 (97%)	441 (97%)	11 (2%)	2 (0%)	39	49
1	D	436/469 (93%)	430 (99%)	5 (1%)	1 (0%)	52	64
All	All	1781/1876 (95%)	1744 (98%)	32 (2%)	5 (0%)	46	56

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	362	MET
1	B	362	MET
1	C	362	MET
1	D	362	MET
1	C	70	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	333/346 (96%)	305 (92%)	28 (8%)	14	16
1	B	332/346 (96%)	312 (94%)	20 (6%)	24	33
1	C	328/346 (95%)	310 (94%)	18 (6%)	27	37
1	D	327/346 (94%)	303 (93%)	24 (7%)	17	23
All	All	1320/1384 (95%)	1230 (93%)	90 (7%)	20	26

All (90) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	VAL
1	A	16	THR
1	A	48	LEU
1	A	59	ARG
1	A	67	ARG
1	A	98	LEU

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Mol	Chain	Res	Type
1	A	123	THR
1	A	159	GLU
1	A	160	ARG
1	A	163	LEU
1	A	183	LEU
1	A	188	THR
1	A	191	VAL
1	A	208	VAL
1	A	225	ARG
1	A	229	VAL
1	A	234	ARG
1	A	255	LEU
1	A	256	GLN
1	A	262	TRP
1	A	269	LEU
1	A	330	THR
1	A	333	GLU
1	A	398	LEU
1	A	399	LYS
1	A	406	LEU
1	A	430	ARG
1	A	446	ARG
1	B	48	LEU
1	B	98	LEU
1	B	160	ARG
1	B	183	LEU
1	B	189	LEU
1	B	191	VAL
1	B	193	ASP
1	B	208	VAL
1	B	229	VAL
1	B	255	LEU
1	B	256	GLN
1	B	262	TRP
1	B	269	LEU
1	B	378	THR
1	B	390	ARG
1	B	399	LYS
1	B	406	LEU
1	B	421	VAL
1	B	464	ARG
1	B	466	ASP

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Mol	Chain	Res	Type
1	C	36	SER
1	C	66	GLU
1	C	67	ARG
1	C	98	LEU
1	C	123	THR
1	C	163	LEU
1	C	183	LEU
1	C	193	ASP
1	C	217	MET
1	C	229	VAL
1	C	255	LEU
1	C	256	GLN
1	C	269	LEU
1	C	390	ARG
1	C	406	LEU
1	C	432	VAL
1	C	446	ARG
1	C	454	ARG
1	D	16	THR
1	D	59	ARG
1	D	98	LEU
1	D	123	THR
1	D	159	GLU
1	D	163	LEU
1	D	183	LEU
1	D	184	THR
1	D	189	LEU
1	D	208	VAL
1	D	229	VAL
1	D	255	LEU
1	D	256	GLN
1	D	262	TRP
1	D	269	LEU
1	D	282	GLN
1	D	306	LEU
1	D	330	THR
1	D	344	ARG
1	D	390	ARG
1	D	398	LEU
1	D	406	LEU
1	D	446	ARG
1	D	454	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50	HIS
1	A	54	HIS
1	B	22	HIS
1	B	75	HIS
1	B	81	GLN
1	B	128	HIS
1	B	243	GLN
1	B	282	GLN
1	B	339	HIS
1	C	75	HIS
1	C	81	GLN
1	C	233	HIS
1	C	243	GLN
1	C	282	GLN
1	C	339	HIS
1	D	75	HIS
1	D	81	GLN
1	D	203	ASN
1	D	243	GLN
1	D	282	GLN
1	D	339	HIS

### 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 carbohydrates in this entry.

### 5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

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

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, 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	454/469 (96%)	-0.13	23 (5%) 32 31	22, 36, 66, 98	0
1	B	451/469 (96%)	-0.24	5 (1%) 82 84	23, 38, 68, 96	0
1	C	458/469 (97%)	-0.08	24 (5%) 31 30	22, 41, 85, 106	0
1	D	444/469 (94%)	0.06	22 (4%) 32 32	24, 38, 79, 109	0
All	All	1807/1876 (96%)	-0.10	74 (4%) 41 40	22, 38, 79, 109	0

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	64	ALA	7.9
1	B	402	ALA	6.2
1	A	65	VAL	5.9
1	A	2	SER	5.1
1	D	2	SER	5.0
1	A	63	PRO	4.9
1	C	198	ASP	4.8
1	A	1	MET	4.5
1	C	197	SER	4.5
1	A	68	ALA	4.1
1	D	69	PRO	4.1
1	A	402	ALA	4.0
1	C	61	THR	3.8
1	B	403	ASN	3.7
1	D	406	LEU	3.7
1	A	200	PRO	3.6
1	D	9	ASP	3.5
1	A	376	SER	3.4
1	A	9	ASP	3.4
1	C	10	GLY	3.3
1	C	194	GLU	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	37	VAL	3.2
1	B	10	GLY	3.2
1	C	53	HIS	3.2
1	D	54	HIS	3.2
1	C	67	ARG	3.2
1	A	462	VAL	3.2
1	C	40	GLY	3.2
1	D	36	SER	3.1
1	D	463	GLY	3.1
1	B	2	SER	3.1
1	C	378	THR	3.0
1	A	61	THR	2.9
1	C	196	ALA	2.9
1	C	64	ALA	2.9
1	A	7	GLU	2.8
1	A	455	ALA	2.8
1	D	205	ALA	2.7
1	C	42	ASP	2.7
1	C	9	ASP	2.7
1	A	66	GLU	2.7
1	D	438	GLY	2.6
1	D	462	VAL	2.6
1	D	401	ALA	2.6
1	C	199	ALA	2.5
1	C	195	ALA	2.5
1	C	425	ALA	2.5
1	A	403	ASN	2.4
1	C	438	GLY	2.4
1	D	375	LEU	2.4
1	A	10	GLY	2.4
1	C	62	GLY	2.4
1	B	9	ASP	2.4
1	D	396	ASP	2.4
1	D	452	ARG	2.4
1	D	450	LEU	2.4
1	D	451	ALA	2.4
1	A	49	ALA	2.3
1	D	57	ASP	2.3
1	C	47	CYS	2.3
1	D	3	ASN	2.2
1	A	11	GLY	2.2
1	A	465	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	37	VAL	2.2
1	C	7	GLU	2.2
1	C	27	ALA	2.1
1	C	54	HIS	2.1
1	A	404	GLY	2.1
1	A	406	LEU	2.1
1	D	465	ALA	2.1
1	D	40	GLY	2.1
1	C	49	ALA	2.0
1	A	67	ARG	2.0
1	D	8	TYR	2.0

## 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. 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, 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	CL	B	501	1/1	0.99	0.08	-	58,58,58,58	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.