



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:22 PM UTC

PDB ID : 4XBM / pdb_00004xbm
Title : X-ray crystal structure of Notch ligand Delta-like 1
Authors : Kershaw, N.J.; Burgess, A.W.; Church, N.L.; Luo, C.S.; Adam, T.E.
Deposited on : 2014-12-17
Resolution : 3.20 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

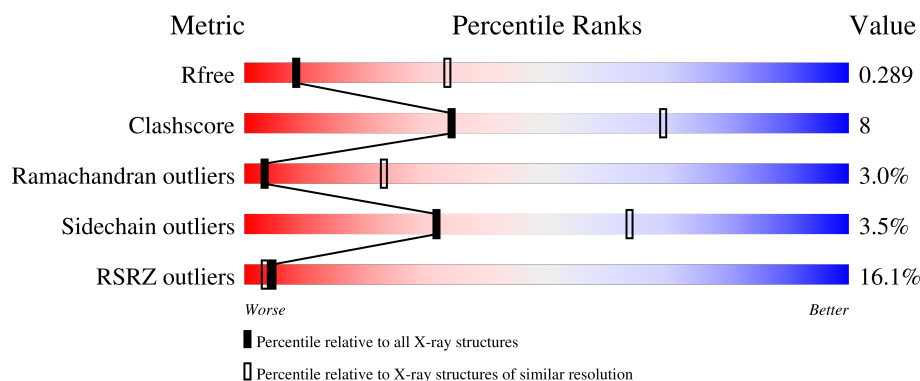
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	531	
1	B	531	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4996 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.

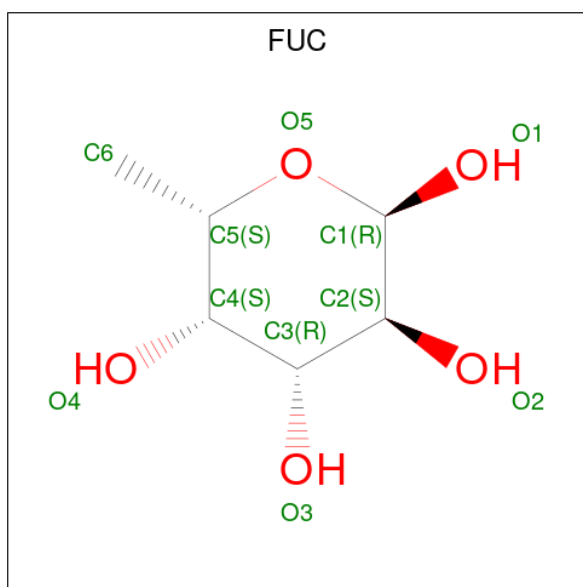
- Molecule 1 is a protein called Delta-like protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	253	Total	C	N	O	S	0	0	0
			1969	1241	347	357	24			
1	B	401	Total	C	N	O	S	0	0	0
			3017	1862	526	580	49			

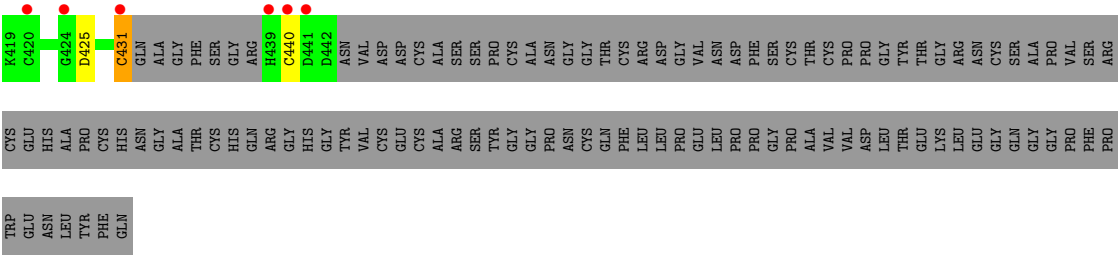
There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	498	GLN	GLU	conflict	UNP O00548
A	502	GLY	ARG	conflict	UNP O00548
A	510	SER	GLY	conflict	UNP O00548
A	546	GLU	-	expression tag	UNP O00548
A	547	ASN	-	expression tag	UNP O00548
A	548	LEU	-	expression tag	UNP O00548
A	549	TYR	-	expression tag	UNP O00548
A	550	PHE	-	expression tag	UNP O00548
A	551	GLN	-	expression tag	UNP O00548
B	498	GLN	GLU	conflict	UNP O00548
B	502	GLY	ARG	conflict	UNP O00548
B	510	SER	GLY	conflict	UNP O00548
B	546	GLU	-	expression tag	UNP O00548
B	547	ASN	-	expression tag	UNP O00548
B	548	LEU	-	expression tag	UNP O00548
B	549	TYR	-	expression tag	UNP O00548
B	550	PHE	-	expression tag	UNP O00548
B	551	GLN	-	expression tag	UNP O00548

- Molecule 2 is alpha-L-fucopyranose (CCD ID: FUC) (formula: C₆H₁₂O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			10	6	4		



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	108.15Å 118.87Å 134.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	67.30 – 3.20 67.30 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (67.30-3.20) 99.8 (67.30-3.20)	Depositor EDS
R_{merge}	0.43	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.21 (at 3.19Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.258 , 0.283 0.263 , 0.289	Depositor DCC
R_{free} test set	1997 reflections (6.82%)	wwPDB-VP
Wilson B-factor (Å ²)	85.6	Xtriage
Anisotropy	0.169	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 60.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	4996	wwPDB-VP
Average B, all atoms (Å ²)	86.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.20% 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: FUC

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	0.32	0/2028	0.81	3/2746 (0.1%)
1	B	0.33	0/3102	0.83	5/4202 (0.1%)
All	All	0.33	0/5130	0.82	8/6948 (0.1%)

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	A	0	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	141	GLU	N-CA-C	-7.21	106.41	114.62
1	A	75	PRO	N-CA-C	6.57	118.71	110.70
1	B	75	PRO	N-CA-C	6.25	118.32	110.70
1	B	51	PRO	CA-C-N	6.13	126.70	120.38
1	B	51	PRO	C-N-CA	6.13	126.70	120.38
1	A	88	GLY	N-CA-C	5.59	118.90	110.18
1	B	88	GLY	N-CA-C	5.07	118.54	111.19
1	B	52	PRO	N-CA-C	5.07	116.88	110.70

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	142	ARG	Peptide
1	B	142	ARG	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	1969	0	1803	34	0
1	B	3017	0	2717	39	0
2	B	10	0	10	0	0
All	All	4996	0	4530	72	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 8.

All (72) 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:A:141:GLU:O	1:A:143:LEU:HB3	1.38	1.23
1:A:141:GLU:O	1:A:143:LEU:CB	1.89	1.20
1:B:384:SER:HB3	1:B:385:PRO:HD2	1.28	1.16
1:B:384:SER:OG	1:B:385:PRO:HD3	1.52	1.10
1:B:384:SER:CB	1:B:385:PRO:CD	2.34	1.06
1:A:141:GLU:O	1:A:142:ARG:C	2.15	0.89
1:A:141:GLU:O	1:A:143:LEU:HB2	1.80	0.80
1:B:384:SER:HB3	1:B:385:PRO:CD	1.98	0.79
1:B:384:SER:CB	1:B:385:PRO:HD3	2.10	0.76
1:B:302:ALA:HB1	1:B:315:SER:O	1.87	0.75
1:A:151:ARG:NH1	1:A:162:ASP:OD2	2.28	0.65
1:B:384:SER:OG	1:B:385:PRO:CD	2.30	0.65
1:B:57:ARG:NH1	1:B:130:ASP:O	2.31	0.64
1:B:195:ARG:NH2	1:B:197:ASP:OD2	2.30	0.64
1:A:43:ASN:HB3	1:A:142:ARG:O	2.00	0.61
1:B:33:VAL:HB	1:B:170:ASP:HB2	1.81	0.61
1:A:33:VAL:HB	1:A:170:ASP:HB3	1.83	0.59
1:A:31:GLU:O	1:A:172:LYS:N	2.35	0.58
1:B:385:PRO:O	1:B:386:ASP:C	2.46	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:141:GLU:O	1:A:143:LEU:N	2.37	0.57
1:B:166:SER:OG	1:B:169:THR:OG1	2.23	0.56
1:A:128:HIS:O	1:A:142:ARG:HB2	2.06	0.56
1:B:31:GLU:O	1:B:172:LYS:N	2.38	0.56
1:A:141:GLU:O	1:A:143:LEU:CA	2.54	0.54
1:B:418:ALA:HB2	1:B:431:CYS:HB3	1.90	0.53
1:A:239:LYS:NZ	1:A:242:GLU:OE1	2.41	0.53
1:A:162:ASP:OD1	1:A:163:LEU:N	2.42	0.52
1:B:143:LEU:O	1:B:144:ILE:HG13	2.10	0.52
1:A:36:LYS:NZ	1:A:56:CYS:SG	2.81	0.52
1:A:127:LEU:HG	1:A:143:LEU:HD12	1.92	0.52
1:B:384:SER:HG	1:B:385:PRO:HD3	1.71	0.51
1:B:44:CYS:O	1:B:54:CYS:HB3	2.10	0.51
1:B:140:PRO:HG2	1:B:142:ARG:HH12	1.76	0.51
1:B:59:PHE:CE1	1:B:129:THR:HG23	2.46	0.51
1:A:119:THR:OG1	1:B:119:THR:OG1	2.25	0.51
1:A:61:ARG:NH2	1:A:75:PRO:HB3	2.25	0.51
1:B:385:PRO:O	1:B:386:ASP:O	2.30	0.50
1:B:65:LYS:HG3	1:B:121:SER:HB2	1.94	0.49
1:B:61:ARG:HH21	1:B:75:PRO:HB3	1.77	0.49
1:B:59:PHE:CE1	1:B:127:LEU:HB2	2.47	0.49
1:A:33:VAL:HA	1:A:91:SER:HB2	1.95	0.48
1:A:65:LYS:HG2	1:A:77:CYS:HA	1.95	0.48
1:B:26:GLU:OE1	1:B:109:ARG:NH1	2.46	0.48
1:B:355:PRO:HB2	1:B:358:PHE:HD1	1.79	0.48
1:B:403:GLU:HG2	1:B:404:LYS:HG3	1.96	0.48
1:B:147:LEU:HB2	1:B:164:HIS:CD2	2.49	0.47
1:A:28:LYS:HB3	1:A:107:PRO:HB3	1.95	0.47
1:A:143:LEU:O	1:A:144:ILE:HG13	2.14	0.47
1:B:36:LYS:HB2	1:B:41:ASN:ND2	2.30	0.47
1:B:127:LEU:HA	1:B:143:LEU:HA	1.97	0.47
1:A:127:LEU:HA	1:A:143:LEU:HA	1.97	0.46
1:B:44:CYS:HA	1:B:128:HIS:CD2	2.50	0.46
1:B:61:ARG:NH2	1:B:75:PRO:HB3	2.30	0.46
1:B:41:ASN:HB3	1:B:44:CYS:SG	2.55	0.46
1:A:124:ILE:HD13	1:A:171:LEU:HD21	1.98	0.45
1:A:124:ILE:HG23	1:A:147:LEU:HB3	1.96	0.45
1:A:140:PRO:HB2	1:A:141:GLU:H	1.62	0.45
1:A:143:LEU:HG	1:A:144:ILE:N	2.32	0.45
1:A:33:VAL:HA	1:A:91:SER:CB	2.47	0.44
1:B:130:ASP:HB3	1:B:132:PRO:HD2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:329:ILE:HG22	1:B:348:ASN:H	1.83	0.44
1:A:42:ARG:H	1:A:42:ARG:HG2	1.62	0.42
1:A:131:SER:CB	1:A:132:PRO:HD2	2.50	0.42
1:A:73:PRO:HA	1:A:123:ILE:HD13	2.02	0.42
1:A:151:ARG:HA	1:A:151:ARG:HD3	1.84	0.42
1:B:25:PHE:CD1	1:B:64:LEU:HD11	2.55	0.42
1:A:61:ARG:NH1	1:A:125:GLU:OE1	2.52	0.41
1:A:26:GLU:OE2	1:A:109:ARG:NH1	2.44	0.41
1:B:193:ARG:HA	1:B:194:PRO:HD3	1.93	0.41
1:A:59:PHE:CE1	1:A:127:LEU:HB2	2.56	0.40
1:B:142:ARG:O	1:B:143:LEU:O	2.39	0.40
1:B:106:ASN:HA	1:B:107:PRO:HD2	1.94	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	243/531 (46%)	213 (88%)	21 (9%)	9 (4%)	2	18
1	B	393/531 (74%)	355 (90%)	28 (7%)	10 (2%)	4	27
All	All	636/1062 (60%)	568 (89%)	49 (8%)	19 (3%)	3	23

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	31	GLU
1	A	142	ARG
1	B	31	GLU
1	B	33	VAL
1	B	143	LEU
1	B	386	ASP

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Mol	Chain	Res	Type
1	B	425	ASP
1	B	144	ILE
1	B	329	ILE
1	B	440	CYS
1	A	33	VAL
1	A	136	ALA
1	A	143	LEU
1	A	241	GLY
1	B	75	PRO
1	A	75	PRO
1	A	167	GLY
1	B	41	ASN
1	A	144	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	212/440 (48%)	207 (98%)	5 (2%)	43	70
1	B	332/440 (76%)	318 (96%)	14 (4%)	26	60
All	All	544/880 (62%)	525 (96%)	19 (4%)	32	64

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	CYS
1	A	54	CYS
1	A	63	CYS
1	A	142	ARG
1	A	197	ASP
1	B	30	GLN
1	B	45	CYS
1	B	46	ARG
1	B	56	CYS
1	B	130	ASP

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Mol	Chain	Res	Type
1	B	144	ILE
1	B	226	CYS
1	B	303	THR
1	B	315	SER
1	B	332	CYS
1	B	361	LYS
1	B	383	ASP
1	B	384	SER
1	B	431	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	213	ASN
1	A	273	GLN
1	A	287	GLN
1	B	128	HIS
1	B	164	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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	FUC	B	1500	1	10,10,11	0.52	0	14,14,16	1.09	1 (7%)

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	FUC	B	1500	1	-	-	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	B	1500	FUC	C1-O5-C5	-2.17	107.83	112.97

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	253/531 (47%)	0.91	38 (15%) 5 4	42, 84, 134, 158	0
1	B	401/531 (75%)	1.02	67 (16%) 4 3	47, 80, 143, 167	0
All	All	654/1062 (61%)	0.98	105 (16%) 4 3	42, 82, 141, 167	0

All (105) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	291	TYR	6.7
1	B	329	ILE	6.4
1	B	327	LEU	6.2
1	B	106	ASN	6.2
1	A	132	PRO	5.9
1	A	170	ASP	5.7
1	B	133	ASP	5.6
1	A	289	LEU	5.4
1	A	137	THR	5.2
1	B	98	GLY	5.1
1	B	143	LEU	5.0
1	B	346	LEU	4.8
1	B	439	HIS	4.7
1	A	52	PRO	4.4
1	B	278	GLU	4.4
1	B	370	CYS	4.4
1	B	140	PRO	4.2
1	B	334	PRO	4.2
1	B	129	THR	4.1
1	B	359	TYR	4.1
1	B	326	GLU	4.1
1	A	290	ASN	4.0
1	A	134	ASP	3.7
1	B	52	PRO	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	31	GLU	3.6
1	A	39	LEU	3.6
1	B	31	GLU	3.5
1	B	196	ASP	3.5
1	A	131	SER	3.4
1	B	323	ALA	3.4
1	B	97	GLY	3.4
1	B	356	PRO	3.4
1	A	40	GLY	3.3
1	A	238	ASP	3.3
1	A	77	CYS	3.3
1	B	293	THR	3.1
1	A	47	GLY	3.1
1	A	141	GLU	3.0
1	B	330	ASP	3.0
1	B	333	ASP	3.0
1	B	441	ASP	3.0
1	B	30	GLN	3.0
1	B	130	ASP	3.0
1	A	186	GLU	2.9
1	B	322	GLY	2.9
1	A	146	ARG	2.9
1	A	264	LEU	2.9
1	B	189	SER	2.9
1	B	385	PRO	2.9
1	B	44	CYS	2.8
1	B	424	GLY	2.8
1	B	49	ALA	2.8
1	B	38	LEU	2.8
1	A	165	SER	2.8
1	B	181	GLU	2.8
1	B	308	GLY	2.8
1	B	431	CYS	2.8
1	A	33	VAL	2.8
1	A	143	LEU	2.8
1	B	32	PHE	2.7
1	A	158	GLU	2.7
1	A	256	GLU	2.6
1	B	387	GLY	2.6
1	B	417	GLY	2.6
1	B	317	ARG	2.6
1	B	335	SER	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	132	PRO	2.6
1	A	90	ASP	2.6
1	A	237	CYS	2.5
1	B	343	CYS	2.5
1	B	420	CYS	2.5
1	A	267	THR	2.5
1	A	283	LEU	2.5
1	B	411	SER	2.4
1	B	344	THR	2.4
1	B	188	CYS	2.4
1	B	50	GLY	2.4
1	B	328	GLY	2.3
1	B	42	ARG	2.3
1	A	240	PRO	2.3
1	B	440	CYS	2.3
1	A	241	GLY	2.3
1	A	284	PHE	2.3
1	B	347	GLU	2.3
1	B	206	GLU	2.3
1	A	136	ALA	2.3
1	B	39	LEU	2.2
1	B	152	HIS	2.2
1	B	351	SER	2.2
1	A	199	PHE	2.2
1	A	168	ARG	2.2
1	B	131	SER	2.2
1	A	41	ASN	2.2
1	B	76	PRO	2.1
1	B	318	PRO	2.1
1	B	398	SER	2.1
1	B	304	CYS	2.1
1	A	167	GLY	2.1
1	B	22	SER	2.0
1	B	342	SER	2.0
1	A	221	CYS	2.0
1	B	381	CYS	2.0
1	B	395	VAL	2.0
1	A	55	ALA	2.0
1	B	297	PRO	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 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	FUC	B	1500	10/11	0.51	0.21	142,162,174,177	0

6.5 Other polymers [i](#)

There are no such residues in this entry.