



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:36 PM GMT

PDB ID : 4YFG  
Title : Crystal structure of PTP delta meA3/meB minus variant Ig1-Fn1  
Authors : Yamagata, A.; Fukai, S.  
Deposited on : 2015-02-25  
Resolution : 3.49 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

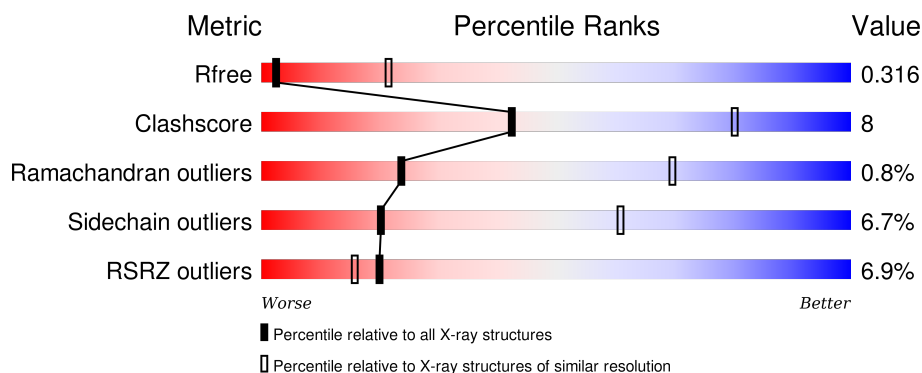
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.49 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))
$R_{free}$	91344	1051 (3.60-3.40)
Clashscore	102246	1157 (3.60-3.40)
Ramachandran outliers	100387	1120 (3.60-3.40)
Sidechain outliers	100360	1121 (3.60-3.40)
RSRZ outliers	91569	1058 (3.60-3.40)

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	488	<div> <div>7%</div> <div>73%</div> <div>23%</div> <div>..</div> </div>
1	B	488	<div> <div>7%</div> <div>76%</div> <div>19%</div> <div>..</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7468 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.

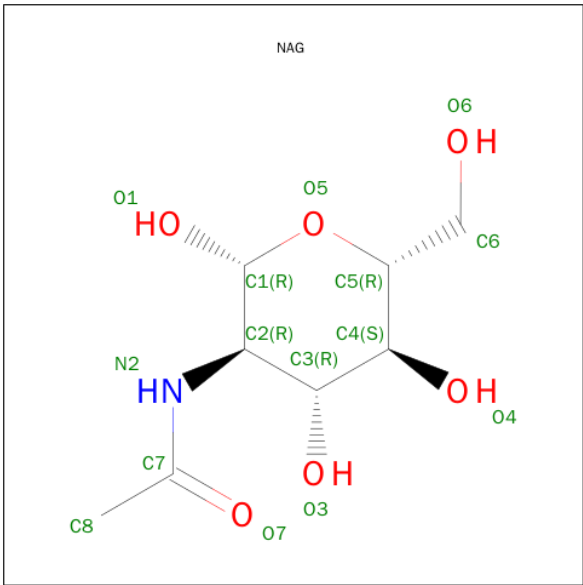
- Molecule 1 is a protein called Receptor-type tyrosine-protein phosphatase delta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	481	Total	C	N	O	S	0	0	0
			3706	2317	642	731	16			
1	A	481	Total	C	N	O	S	0	0	0
			3706	2317	642	731	16			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	509	LYS	-	expression tag	UNP Q64487
B	510	HIS	-	expression tag	UNP Q64487
B	511	HIS	-	expression tag	UNP Q64487
B	512	HIS	-	expression tag	UNP Q64487
B	513	HIS	-	expression tag	UNP Q64487
B	514	HIS	-	expression tag	UNP Q64487
B	515	HIS	-	expression tag	UNP Q64487
A	509	LYS	-	expression tag	UNP Q64487
A	510	HIS	-	expression tag	UNP Q64487
A	511	HIS	-	expression tag	UNP Q64487
A	512	HIS	-	expression tag	UNP Q64487
A	513	HIS	-	expression tag	UNP Q64487
A	514	HIS	-	expression tag	UNP Q64487
A	515	HIS	-	expression tag	UNP Q64487

- Molecule 2 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).

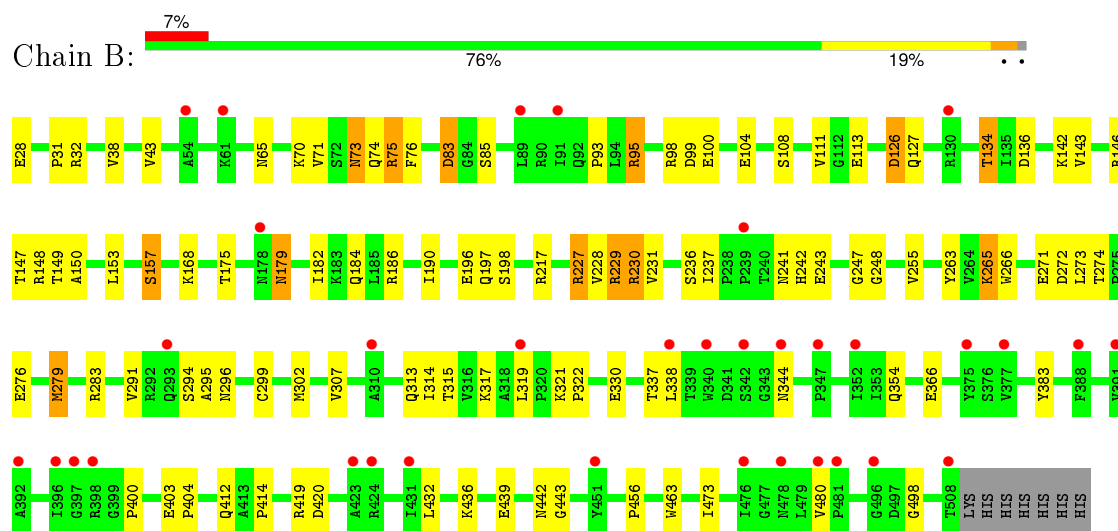


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		

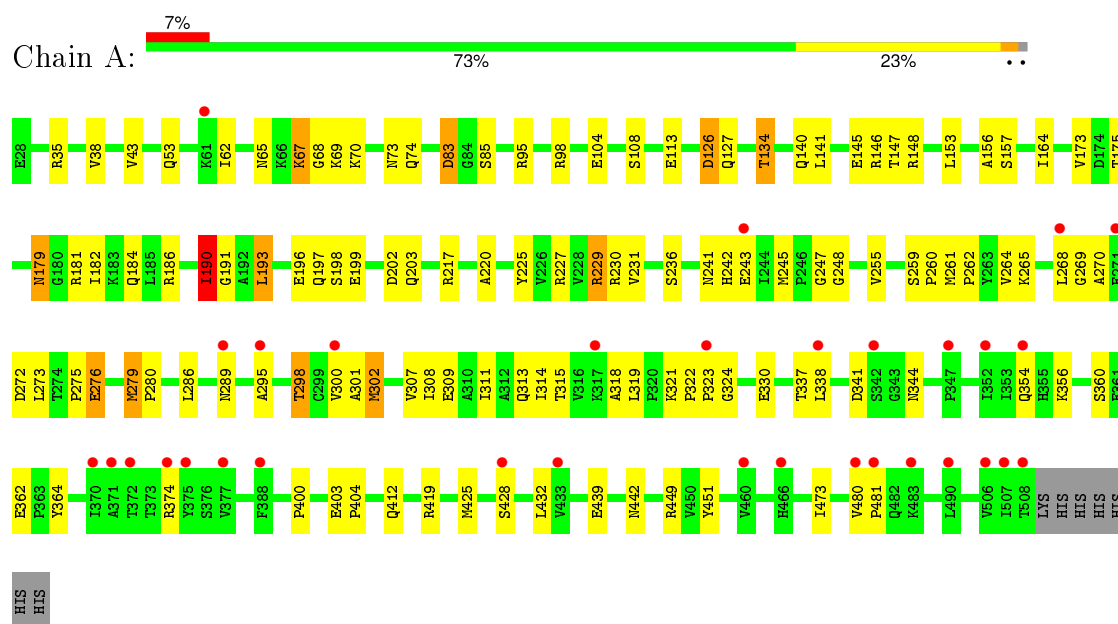
### 3 Residue-property plots

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: Receptor-type tyrosine-protein phosphatase delta



- Molecule 1: Receptor-type tyrosine-protein phosphatase delta



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	70.49Å 72.63Å 94.41Å 107.65° 94.44° 108.47°	Depositor
Resolution (Å)	48.41 – 3.49 48.40 – 3.49	Depositor EDS
% Data completeness (in resolution range)	88.4 (48.41-3.49) 78.4 (48.40-3.49)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.19	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.47 (at 3.48Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, $R_{free}$	0.291 , 0.315 0.291 , 0.316	Depositor DCC
$R_{free}$ test set	969 reflections (5.20%)	DCC
Wilson B-factor (Å <sup>2</sup> )	68.4	Xtriage
Anisotropy	0.443	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 5.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.38$ , $\langle L^2 \rangle = 0.20$	Xtriage
Outliers	0 of 18642 reflections	Xtriage
$F_o, F_c$ correlation	0.82	EDS
Total number of atoms	7468	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	84.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.10% 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: NAG

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	0.26	0/3788	0.52	0/5172
1	B	0.25	0/3788	0.50	0/5172
All	All	0.25	0/7576	0.51	0/10344

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	190	ILE	Peptide

### 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	3706	0	3658	69	0
1	B	3706	0	3658	63	0
2	A	28	0	26	0	0
2	B	28	0	26	0	0
All	All	7468	0	7368	123	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 (123) 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:164:ILE:HD12	1:A:191:GLY:H	1.48	0.78
1:A:156:ALA:H	1:A:190:ILE:HG13	1.48	0.78
1:B:95:ARG:NH2	1:A:302:MET:SD	2.59	0.75
1:A:181:ARG:NH1	1:A:197:GLN:O	2.17	0.74
1:B:108:SER:HB3	1:B:113:GLU:HG3	1.69	0.74
1:A:146:ARG:HG2	1:A:147:THR:HG23	1.69	0.73
1:A:175:THR:HG21	1:A:184:GLN:HG3	1.72	0.72
1:B:98:ARG:NH1	1:A:272:ASP:OD1	2.23	0.71
1:B:272:ASP:OD1	1:A:98:ARG:NH1	2.25	0.70
1:B:265:LYS:NZ	1:B:276:GLU:O	2.26	0.69
1:B:98:ARG:HH21	1:A:276:GLU:HG3	1.60	0.67
1:B:266:TRP:HB3	1:B:273:LEU:HD12	1.76	0.67
1:B:330:GLU:HB2	1:B:337:THR:HB	1.77	0.67
1:B:136:ASP:OD2	1:B:157:SER:OG	2.13	0.67
1:A:330:GLU:HB2	1:A:337:THR:HB	1.77	0.66
1:B:179:ASN:OD1	1:B:179:ASN:N	2.30	0.65
1:A:83:ASP:OD2	1:A:83:ASP:N	2.31	0.64
1:A:428:SER:HB2	1:A:481:PRO:HB3	1.78	0.64
1:B:75:ARG:NH1	1:B:93:PRO:O	2.31	0.64
1:B:83:ASP:N	1:B:83:ASP:OD2	2.30	0.64
1:B:302:MET:HB3	1:B:307:VAL:HG22	1.81	0.63
1:A:145:GLU:OE1	1:A:148:ARG:NH1	2.33	0.62
1:A:126:ASP:OD1	1:A:126:ASP:N	2.31	0.62
1:A:321:LYS:N	1:A:344:ASN:OD1	2.32	0.62
1:A:231:VAL:N	1:A:259:SER:O	2.24	0.61
1:A:175:THR:HG22	1:A:182:ILE:HG22	1.83	0.61
1:B:230:ARG:HH11	1:A:220:ALA:HB3	1.65	0.60
1:A:356:LYS:HE3	1:A:364:TYR:CE1	2.37	0.59
1:A:356:LYS:HE3	1:A:364:TYR:HE1	1.66	0.59
1:A:190:ILE:HD13	1:A:190:ILE:H	1.66	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:SER:HB3	1:A:113:GLU:HG3	1.85	0.59
1:B:175:THR:HG22	1:B:182:ILE:HG22	1.84	0.59
1:A:247:GLY:N	1:A:248:GLY:HA3	2.16	0.59
1:B:295:ALA:N	1:B:314:ILE:O	2.24	0.58
1:A:268:LEU:O	1:A:270:ALA:N	2.37	0.57
1:B:383:TYR:CD2	1:B:443:GLY:HA2	2.39	0.57
1:A:241:ASN:O	1:A:242:HIS:ND1	2.39	0.56
1:B:65:ASN:HB2	1:B:104:GLU:HG3	1.88	0.56
1:A:179:ASN:OD1	1:A:179:ASN:N	2.34	0.56
1:B:142:LYS:HD3	1:B:150:ALA:HB1	1.87	0.56
1:A:403:GLU:HG3	1:A:404:PRO:HD2	1.88	0.55
1:B:263:TYR:HB3	1:B:279:MET:HG2	1.87	0.55
1:A:243:GLU:HA	1:A:315:THR:O	2.06	0.55
1:A:300:VAL:HG13	1:A:309:GLU:HG2	1.89	0.55
1:A:302:MET:HB3	1:A:307:VAL:HG22	1.89	0.54
1:B:247:GLY:N	1:B:291:VAL:O	2.37	0.54
1:B:126:ASP:OD1	1:B:126:ASP:N	2.40	0.54
1:B:419:ARG:NH1	1:B:439:GLU:OE2	2.42	0.53
1:A:295:ALA:N	1:A:314:ILE:O	2.25	0.53
1:A:330:GLU:OE2	1:A:374:ARG:NH2	2.42	0.52
1:B:247:GLY:N	1:B:248:GLY:HA3	2.23	0.52
1:A:229:ARG:O	1:A:231:VAL:HG23	2.09	0.52
1:B:302:MET:CB	1:B:307:VAL:HG22	2.38	0.52
1:A:203:GLN:HG3	1:A:225:TYR:HA	1.92	0.52
1:A:140:GLN:N	1:A:140:GLN:OE1	2.43	0.52
1:B:294:SER:HB3	1:B:315:THR:HG22	1.91	0.52
1:A:156:ALA:O	1:A:190:ILE:HG21	2.10	0.52
1:B:241:ASN:HB3	1:B:313:GLN:HB2	1.92	0.52
1:A:419:ARG:NH1	1:A:439:GLU:OE2	2.43	0.51
1:B:73:ASN:HA	1:B:76:PHE:HB2	1.91	0.51
1:B:241:ASN:O	1:B:242:HIS:ND1	2.45	0.50
1:B:403:GLU:HG3	1:B:404:PRO:HD2	1.92	0.50
1:A:83:ASP:OD1	1:A:85:SER:OG	2.25	0.50
1:B:28:GLU:HB3	1:B:111:VAL:HG21	1.94	0.50
1:A:322:PRO:HD3	1:A:400:PRO:HG2	1.93	0.49
1:B:420:ASP:HB2	1:B:436:LYS:HG3	1.95	0.49
1:A:412:GLN:HB2	1:A:442:ASN:HB3	1.93	0.49
1:B:255:VAL:HG22	1:B:283:ARG:HG3	1.94	0.49
1:B:321:LYS:N	1:B:344:ASN:OD1	2.44	0.49
1:A:273:LEU:HD12	1:A:286:LEU:HD11	1.96	0.48
1:A:73:ASN:OD1	1:A:74:GLN:N	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:83:ASP:OD1	1:B:85:SER:OG	2.26	0.48
1:B:322:PRO:HD3	1:B:400:PRO:HG2	1.96	0.47
1:A:134:THR:HA	1:A:217:ARG:HD2	1.96	0.47
1:A:67:LYS:HA	1:A:68:GLY:HA2	1.62	0.47
1:A:145:GLU:CD	1:A:229:ARG:HH22	2.18	0.47
1:B:296:ASN:OD1	1:B:313:GLN:HG2	2.14	0.47
1:B:95:ARG:HD3	1:A:265:LYS:NZ	2.30	0.47
1:B:243:GLU:HA	1:B:315:THR:O	2.15	0.47
1:A:35:ARG:HB3	1:A:53:GLN:HB2	1.97	0.47
1:B:274:THR:HA	1:B:276:GLU:N	2.30	0.46
1:B:143:VAL:HG11	1:B:227:ARG:HH21	1.81	0.46
1:B:147:THR:N	1:B:198:SER:OG	2.48	0.45
1:A:65:ASN:HB2	1:A:104:GLU:HG3	1.98	0.45
1:A:69:LYS:HG2	1:A:70:LYS:O	2.17	0.45
1:A:301:ALA:O	1:A:308:ILE:HG12	2.17	0.45
1:A:95:ARG:HD2	1:A:95:ARG:HA	1.62	0.45
1:A:230:ARG:HD2	1:A:260:PRO:HD3	1.99	0.45
1:B:354:GLN:HG2	1:B:366:GLU:HG2	1.99	0.45
1:B:175:THR:HG21	1:B:184:GLN:HG3	1.99	0.45
1:B:412:GLN:HB2	1:B:442:ASN:HB3	1.97	0.45
1:B:432:LEU:HD21	1:B:473:ILE:HD11	1.99	0.44
1:B:31:PRO:O	1:B:32:ARG:NH1	2.40	0.44
1:B:146:ARG:HE	1:B:228:VAL:HG12	1.82	0.44
1:B:98:ARG:NH2	1:A:276:GLU:HG3	2.29	0.44
1:B:237:ILE:HB	1:B:255:VAL:HB	1.98	0.44
1:B:134:THR:HA	1:B:217:ARG:HD2	1.98	0.44
1:B:236:SER:HB3	1:B:255:VAL:HG12	2.00	0.43
1:A:324:GLY:HA3	1:A:341:ASP:HB3	1.99	0.43
1:B:419:ARG:HD3	1:B:439:GLU:HG2	2.00	0.43
1:B:265:LYS:HZ3	1:B:272:ASP:HB3	1.83	0.43
1:A:354:GLN:HB3	1:A:364:TYR:HB3	2.00	0.43
1:A:245:MET:HG2	1:A:318:ALA:HA	2.01	0.43
1:B:95:ARG:HD3	1:A:265:LYS:HZ3	1.83	0.43
1:A:322:PRO:HA	1:A:323:PRO:HD3	1.85	0.43
1:A:264:VAL:O	1:A:280:PRO:HD2	2.19	0.43
1:A:298:THR:HG23	1:A:311:ILE:HG13	2.01	0.43
1:A:356:LYS:HD2	1:A:360:SER:O	2.19	0.43
1:A:261:MET:HG3	1:A:262:PRO:HD2	2.01	0.42
1:A:173:VAL:HG11	1:A:193:LEU:HD11	2.01	0.42
1:B:95:ARG:HG2	1:B:98:ARG:NH1	2.35	0.42
1:B:229:ARG:O	1:B:231:VAL:HG23	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:199:GLU:N	1:A:202:ASP:OD2	2.50	0.42
1:B:75:ARG:NH2	1:B:99:ASP:OD1	2.53	0.41
1:A:432:LEU:HD21	1:A:473:ILE:HD11	2.01	0.41
1:B:149:THR:HG22	1:B:196:GLU:HA	2.01	0.41
1:B:148:ARG:O	1:B:198:SER:HB3	2.21	0.41
1:B:95:ARG:HH12	1:A:279:MET:HG2	1.85	0.41
1:B:414:PRO:HB2	1:B:498:GLY:HA3	2.03	0.41
1:A:236:SER:HB3	1:A:255:VAL:HG12	2.03	0.41
1:B:456:PRO:HA	1:B:463:TRP:CZ2	2.56	0.40
1:A:419:ARG:HD3	1:A:439:GLU:HG2	2.04	0.40
1:A:449:ARG:HD2	1:A:451:TYR:OH	2.21	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	479/488 (98%)	460 (96%)	15 (3%)	4 (1%)	24	70
1	B	479/488 (98%)	461 (96%)	14 (3%)	4 (1%)	24	70
All	All	958/976 (98%)	921 (96%)	29 (3%)	8 (1%)	24	70

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	190	ILE
1	A	269	GLY
1	A	276	GLU
1	B	73	ASN
1	B	197	GLN
1	B	74	GLN
1	A	275	PRO

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Mol	Chain	Res	Type
1	B	100	GLU

### 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	420/427 (98%)	391 (93%)	29 (7%)	19	59
1	B	420/427 (98%)	393 (94%)	27 (6%)	22	62
All	All	840/854 (98%)	784 (93%)	56 (7%)	20	61

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	38	VAL
1	B	43	VAL
1	B	70	LYS
1	B	71	VAL
1	B	75	ARG
1	B	83	ASP
1	B	95	ARG
1	B	126	ASP
1	B	127	GLN
1	B	134	THR
1	B	153	LEU
1	B	157	SER
1	B	168	LYS
1	B	179	ASN
1	B	186	ARG
1	B	190	ILE
1	B	227	ARG
1	B	229	ARG
1	B	230	ARG
1	B	265	LYS
1	B	271	GLU
1	B	279	MET

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Mol	Chain	Res	Type
1	B	299	CYS
1	B	317	LYS
1	B	319	LEU
1	B	338	LEU
1	B	480	VAL
1	A	38	VAL
1	A	43	VAL
1	A	62	ILE
1	A	67	LYS
1	A	83	ASP
1	A	126	ASP
1	A	127	GLN
1	A	134	THR
1	A	141	LEU
1	A	153	LEU
1	A	157	SER
1	A	179	ASN
1	A	186	ARG
1	A	190	ILE
1	A	193	LEU
1	A	196	GLU
1	A	198	SER
1	A	227	ARG
1	A	229	ARG
1	A	279	MET
1	A	289	ASN
1	A	298	THR
1	A	302	MET
1	A	313	GLN
1	A	319	LEU
1	A	338	LEU
1	A	362	GLU
1	A	425	MET
1	A	480	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA molecules 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 ⓘ

4 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	NAG	A	601	1	14,14,15	0.59	0	15,19,21	0.56	0
2	NAG	A	602	1	14,14,15	0.42	0	15,19,21	0.62	1 (6%)
2	NAG	B	601	1	14,14,15	0.30	0	15,19,21	0.62	1 (6%)
2	NAG	B	602	1	14,14,15	0.45	0	15,19,21	0.61	1 (6%)

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	601	1	-	0/6/23/26	0/1/1/1
2	NAG	A	602	1	-	0/6/23/26	0/1/1/1
2	NAG	B	601	1	-	0/6/23/26	0/1/1/1
2	NAG	B	602	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	602	NAG	C1-O5-C5	2.13	114.95	112.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	602	NAG	C1-O5-C5	2.18	115.02	112.25
2	B	601	NAG	C1-O5-C5	2.21	115.05	112.25

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, 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	481/488 (98%)	0.64	32 (6%)	21 16	43, 83, 120, 151	0
1	B	481/488 (98%)	0.56	34 (7%)	19 15	42, 84, 121, 152	0
All	All	962/976 (98%)	0.60	66 (6%)	20 16	42, 84, 120, 152	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	371	ALA	5.5
1	A	508	THR	4.8
1	A	352	ILE	4.7
1	B	396	ILE	4.6
1	A	375	TYR	4.4
1	A	338	LEU	3.8
1	B	496	GLY	3.8
1	B	338	LEU	3.6
1	B	54	ALA	3.6
1	A	507	ILE	3.4
1	B	130	ARG	3.3
1	A	243	GLU	3.3
1	B	61	LYS	3.3
1	A	61	LYS	3.3
1	B	293	GLN	3.2
1	A	354	GLN	3.2
1	A	295	ALA	3.1
1	A	377	VAL	3.1
1	A	271	GLU	3.1
1	B	377	VAL	3.0
1	A	480	VAL	2.9
1	B	352	ILE	2.9
1	B	239	PRO	2.9
1	A	506	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	340	TRP	2.8
1	A	374	ARG	2.8
1	A	490	LEU	2.8
1	A	460	VAL	2.7
1	A	433	VAL	2.7
1	A	372	THR	2.7
1	B	431	ILE	2.7
1	B	178	ASN	2.6
1	A	388	PHE	2.6
1	B	89	LEU	2.6
1	B	476	ILE	2.5
1	B	392	ALA	2.5
1	B	91	ILE	2.5
1	B	310	ALA	2.4
1	A	481	PRO	2.4
1	B	423	ALA	2.3
1	A	347	PRO	2.3
1	B	481	PRO	2.3
1	B	388	PHE	2.3
1	A	466	HIS	2.3
1	A	300	VAL	2.3
1	A	268	LEU	2.3
1	B	319	LEU	2.2
1	A	428	SER	2.2
1	A	323	PRO	2.2
1	B	391	VAL	2.2
1	B	508	THR	2.2
1	B	480	VAL	2.2
1	B	478	ASN	2.1
1	B	375	TYR	2.1
1	A	483	LYS	2.1
1	B	342	SER	2.1
1	B	397	GLY	2.1
1	A	289	ASN	2.1
1	A	317	LYS	2.1
1	B	451	TYR	2.1
1	A	342	SER	2.1
1	B	398	ARG	2.0
1	A	370	ILE	2.0
1	B	344	ASN	2.0
1	B	347	PRO	2.0
1	B	424	ARG	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( $\text{\AA}^2$ )	Q<0.9
2	NAG	B	601	14/15	0.84	0.17	-	74,95,103,104	0
2	NAG	B	602	14/15	0.69	0.21	-	94,107,112,119	0
2	NAG	A	602	14/15	0.78	0.17	-	74,87,95,96	0
2	NAG	A	601	14/15	0.85	0.17	-	65,83,114,120	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.