



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:47 AM UTC

PDB ID : 4YFG / pdb_00004yfg
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

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

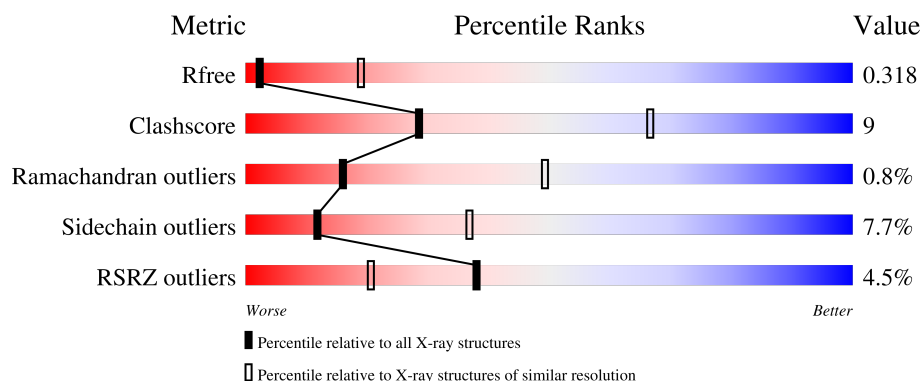
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.49 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}	180053	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)

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	488	5% 72% 24% ..
1	B	488	4% 74% 22% ..

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 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).

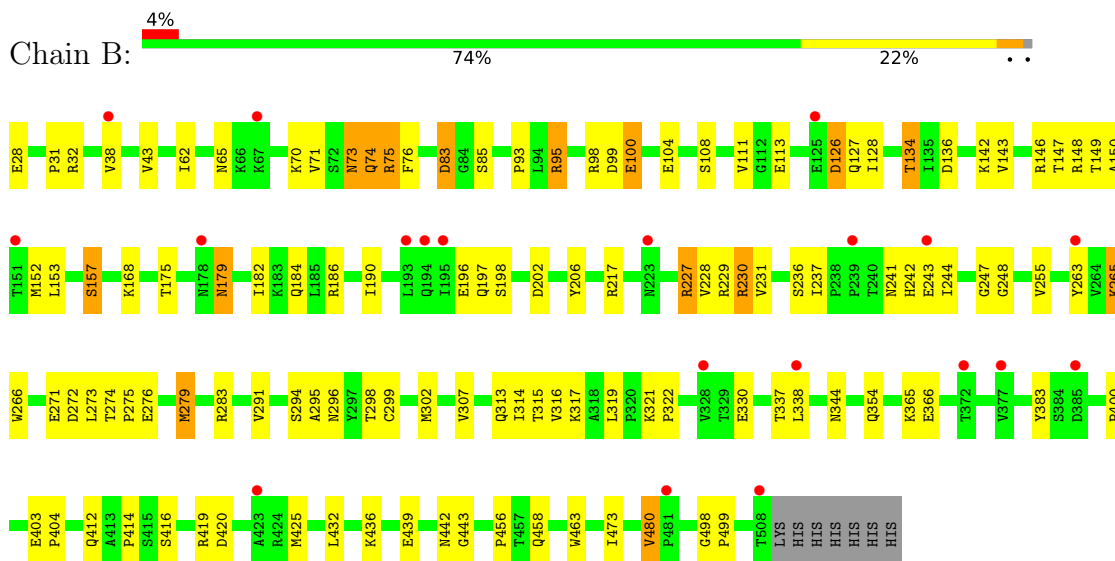


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		

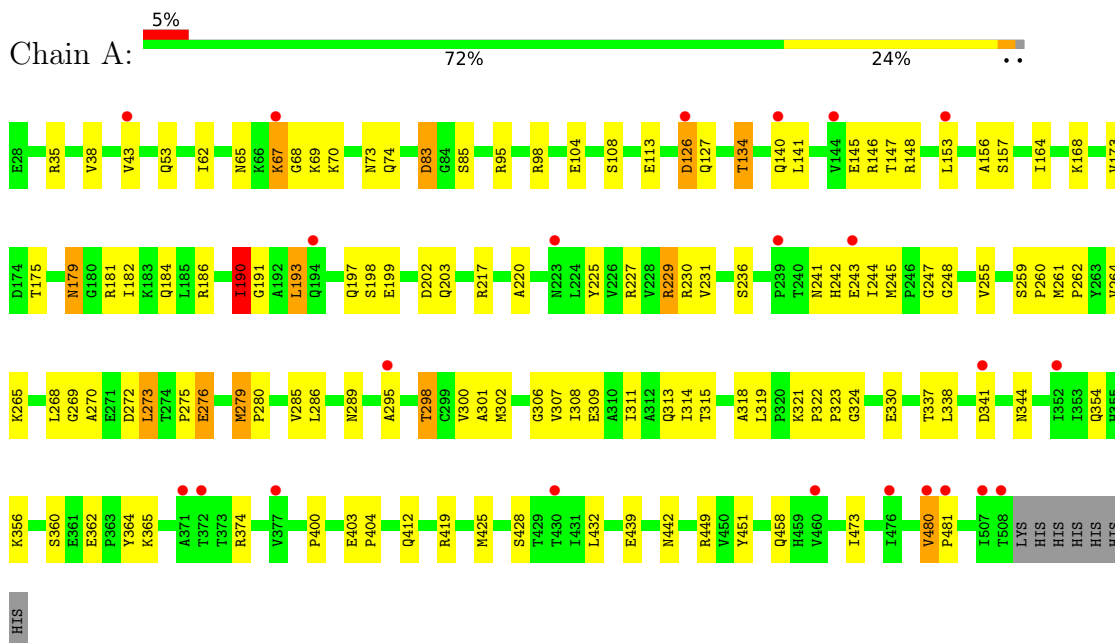
3 Residue-property plots

These plots are drawn for all protein, RNA, 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: 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, α , β , γ	70.49Å 72.63Å 94.41Å 107.65° 94.44° 108.47°	Depositor
Resolution (Å)	48.41 – 3.49 48.41 – 3.49	Depositor EDS
% Data completeness (in resolution range)	88.4 (48.41-3.49) 88.4 (48.41-3.49)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.19	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.47 (at 3.48Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.291 , 0.315 0.291 , 0.318	Depositor DCC
R_{free} test set	969 reflections (4.59%)	wwPDB-VP
Wilson B-factor (Å ²)	68.4	Xtriage
Anisotropy	0.443	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 17.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.38$, $\langle L^2 \rangle = 0.20$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.82	EDS
Total number of atoms	7468	wwPDB-VP
Average B, all atoms (Å ²)	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.*

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: 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	0.33	0/3788	0.83	3/5172 (0.1%)
1	B	0.33	0/3788	0.83	4/5172 (0.1%)
All	All	0.33	0/7576	0.83	7/10344 (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

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	316	VAL	N-CA-C	8.17	119.65	107.80
1	A	306	GLY	N-CA-C	6.58	119.86	111.37
1	B	480	VAL	CA-C-N	6.10	126.06	119.78
1	B	480	VAL	C-N-CA	6.10	126.06	119.78
1	B	74	GLN	N-CA-C	-5.29	99.54	110.80
1	A	480	VAL	CA-C-N	5.23	125.17	119.78
1	A	480	VAL	C-N-CA	5.23	125.17	119.78

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

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	73	0
1	B	3706	0	3658	66	0
2	A	28	0	26	0	0
2	B	28	0	26	0	0
All	All	7468	0	7368	131	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 9.

All (131) 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:OD2	1:B:83:ASP:N	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:108:SER:HB3	1:A:113:GLU:HG3	1.85	0.59
1:A:190:ILE:H	1:A:190:ILE:HD13	1.66	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:A:276:GLU:HA	1:A:279:MET:HE1	1.87	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:B:247:GLY:N	1:B:248:GLY:HA3	2.23	0.52
1:A:330:GLU:OE2	1:A:374:ARG:NH2	2.42	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:140:GLN:N	1:A:140:GLN:OE1	2.43	0.52
1:A:203:GLN:HG3	1:A:225:TYR:HA	1.92	0.52
1:B:294:SER:HB3	1:B:315:THR:HG22	1.91	0.52
1:A:145:GLU:CD	1:A:229:ARG:HH22	2.18	0.52
1:B:241:ASN:HB3	1:B:313:GLN:HB2	1.92	0.52
1:A:156:ALA:O	1:A:190:ILE:HG21	2.10	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:420:ASP:HB2	1:B:436:LYS:HG3	1.95	0.49
1:A:322:PRO:HD3	1:A:400:PRO:HG2	1.93	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
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: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:265:LYS:NZ	1:A:272:ASP:OD1	2.48	0.45
1:A:95:ARG:HD2	1:A:95:ARG:HA	1.62	0.45
1:A:301:ALA:O	1:A:308:ILE:HG12	2.17	0.45
1:B:354:GLN:HG2	1:B:366:GLU:HG2	1.99	0.45
1:A:230:ARG:HD2	1:A:260:PRO:HD3	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:274:THR:HA	1:B:275:PRO:C	2.43	0.44
1:B:98:ARG:NH2	1:A:276:GLU:HG3	2.29	0.44
1:B:134:THR:HA	1:B:217:ARG:HD2	1.98	0.44
1:B:237:ILE:HB	1:B:255:VAL:HB	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:202:ASP:O	1:B:206:TYR:OH	2.34	0.43
1:B:419:ARG:HD3	1:B:439:GLU:HG2	2.00	0.43
1:A:354:GLN:HB3	1:A:364:TYR:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:245:MET:HG2	1:A:318:ALA:HA	2.01	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:322:PRO:HA	1:A:323:PRO:HD3	1.85	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:416:SER:C	1:B:499:PRO:HG2	2.45	0.42
1:B:229:ARG:O	1:B:231:VAL:HG23	2.20	0.42
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:425:MET:HE3	1:B:425:MET:HB3	1.81	0.41
1:A:259:SER:HA	1:A:260:PRO:C	2.45	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:A:425:MET:HE2	1:A:425:MET:HB3	1.96	0.41
1:B:456:PRO:HA	1:B:463:TRP:CZ2	2.56	0.40
1:A:265:LYS:CD	1:A:279:MET:HE2	2.51	0.40
1:B:152:MET:HE3	1:B:152:MET:HB3	1.92	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%)	16	49
1	B	479/488 (98%)	461 (96%)	14 (3%)	4 (1%)	16	49
All	All	958/976 (98%)	921 (96%)	29 (3%)	8 (1%)	16	49

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
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%)	388 (92%)	32 (8%)	12	37
1	B	420/427 (98%)	387 (92%)	33 (8%)	11	36
All	All	840/854 (98%)	775 (92%)	65 (8%)	12	37

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	38	VAL
1	B	43	VAL
1	B	62	ILE
1	B	70	LYS
1	B	71	VAL
1	B	75	ARG
1	B	83	ASP
1	B	95	ARG

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Mol	Chain	Res	Type
1	B	100	GLU
1	B	126	ASP
1	B	127	GLN
1	B	128	ILE
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	230	ARG
1	B	244	ILE
1	B	265	LYS
1	B	271	GLU
1	B	279	MET
1	B	298	THR
1	B	299	CYS
1	B	317	LYS
1	B	319	LEU
1	B	338	LEU
1	B	365	LYS
1	B	458	GLN
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	168	LYS
1	A	179	ASN
1	A	186	ARG
1	A	190	ILE
1	A	193	LEU
1	A	198	SER

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Mol	Chain	Res	Type
1	A	227	ARG
1	A	229	ARG
1	A	244	ILE
1	A	273	LEU
1	A	279	MET
1	A	285	VAL
1	A	289	ASN
1	A	298	THR
1	A	313	GLN
1	A	319	LEU
1	A	338	LEU
1	A	362	GLU
1	A	365	LYS
1	A	458	GLN
1	A	480	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	472	GLN
1	A	472	GLN

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](#)

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	602	1	14,14,15	0.44	0	17,19,21	0.65	1 (5%)
2	NAG	B	601	1	14,14,15	0.32	0	17,19,21	0.64	1 (5%)
2	NAG	A	601	1	14,14,15	0.61	1 (7%)	17,19,21	0.63	0
2	NAG	B	602	1	14,14,15	0.47	0	17,19,21	0.69	1 (5%)

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

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	NAG	C1-C2	2.11	1.55	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	NAG	C1-O5-C5	2.14	115.05	112.19
2	B	602	NAG	C1-O5-C5	2.11	115.02	112.19
2	A	602	NAG	C1-O5-C5	2.06	114.95	112.19

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	602	NAG	O5-C5-C6-O6
2	A	602	NAG	C4-C5-C6-O6
2	B	601	NAG	O5-C5-C6-O6
2	A	601	NAG	O5-C5-C6-O6
2	B	601	NAG	C4-C5-C6-O6
2	B	602	NAG	C4-C5-C6-O6

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	481/488 (98%)	0.53	23 (4%) 35 19	43, 83, 120, 151	0
1	B	481/488 (98%)	0.35	20 (4%) 40 22	42, 84, 121, 152	0
All	All	962/976 (98%)	0.44	43 (4%) 38 20	42, 84, 120, 152	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	481	PRO	3.9
1	A	140	GLN	3.8
1	A	341	ASP	3.7
1	A	377	VAL	3.7
1	A	480	VAL	3.3
1	A	153	LEU	3.1
1	A	481	PRO	3.1
1	A	126	ASP	3.1
1	A	243	GLU	3.0
1	A	194	GLN	3.0
1	B	223	ASN	3.0
1	B	263	TYR	2.8
1	B	178	ASN	2.7
1	A	371	ALA	2.7
1	A	223	ASN	2.7
1	B	338	LEU	2.7
1	A	352	ILE	2.6
1	A	460	VAL	2.6
1	B	377	VAL	2.5
1	A	508	THR	2.5
1	B	193	LEU	2.5
1	B	194	GLN	2.4
1	B	243	GLU	2.4
1	A	144	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	372	THR	2.3
1	B	151	THR	2.3
1	A	239	PRO	2.3
1	B	423	ALA	2.2
1	A	476	ILE	2.2
1	A	507	ILE	2.2
1	B	372	THR	2.2
1	B	38	VAL	2.2
1	B	125	GLU	2.2
1	B	239	PRO	2.2
1	A	67	LYS	2.2
1	A	295	ALA	2.2
1	B	508	THR	2.1
1	B	328	VAL	2.1
1	B	67	LYS	2.1
1	B	195	ILE	2.1
1	B	385	ASP	2.1
1	A	430	THR	2.1
1	A	43	VAL	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	NAG	B	602	14/15	0.65	0.12	94,107,112,119	0
2	NAG	B	601	14/15	0.68	0.12	74,95,103,104	0
2	NAG	A	602	14/15	0.73	0.11	74,87,95,96	0
2	NAG	A	601	14/15	0.79	0.10	65,83,114,120	0

6.5 Other polymers [i](#)

There are no such residues in this entry.