



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:47 PM UTC

PDB ID : 3S97 / pdb_00003s97
Title : PTPRZ CNTN1 complex
Authors : Bouyain, S.
Deposited on : 2011-05-31
Resolution : 2.30 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

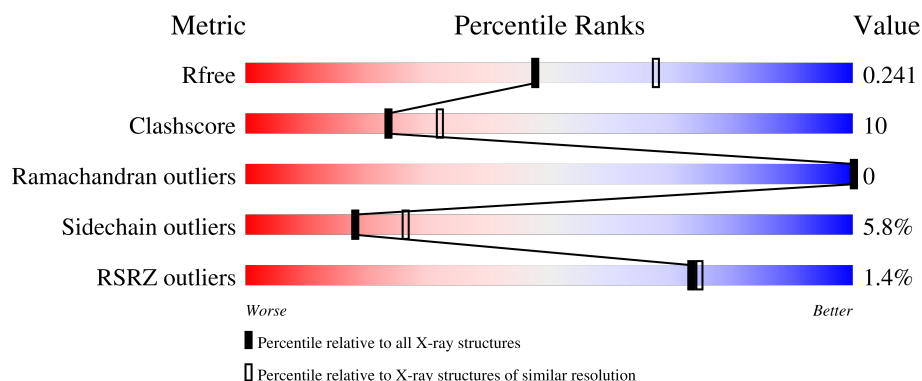
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	273	73% 22% • •
1	B	273	72% 22% • •
2	C	201	76% 21% •
2	D	201	71% 24% • •

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7633 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 zeta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	266	Total	C	N	O	S	0	0	0
			2129	1358	347	413	11			
1	B	265	Total	C	N	O	S	0	0	0
			2121	1352	346	412	11			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	30	GLY	-	expression tag	UNP P23471
A	31	PRO	-	expression tag	UNP P23471
A	32	GLY	-	expression tag	UNP P23471
A	33	SER	-	expression tag	UNP P23471
B	30	GLY	-	expression tag	UNP P23471
B	31	PRO	-	expression tag	UNP P23471
B	32	GLY	-	expression tag	UNP P23471
B	33	SER	-	expression tag	UNP P23471

- Molecule 2 is a protein called Contactin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	197	Total	C	N	O	S	0	0	0
			1580	1019	260	293	8			
2	D	197	Total	C	N	O	S	0	0	0
			1580	1019	260	293	8			

There are 8 discrepancies between the modelled and reference sequences:

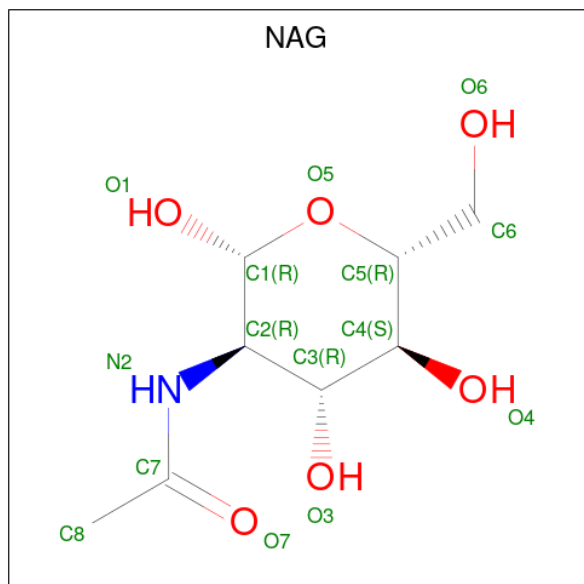
Chain	Residue	Modelled	Actual	Comment	Reference
C	129	GLY	-	expression tag	UNP Q12860
C	130	PRO	-	expression tag	UNP Q12860
C	131	GLY	-	expression tag	UNP Q12860
C	132	SER	-	expression tag	UNP Q12860

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Chain	Residue	Modelled	Actual	Comment	Reference
D	129	GLY	-	expression tag	UNP Q12860
D	130	PRO	-	expression tag	UNP Q12860
D	131	GLY	-	expression tag	UNP Q12860
D	132	SER	-	expression tag	UNP Q12860

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	D	1	Total	C	N	O	0	0
			14	8	1	5		
3	D	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	46	Total	O	0	0
			46	46		
4	B	61	Total	O	0	0
			61	61		
4	C	42	Total	O	0	0
			42	42		

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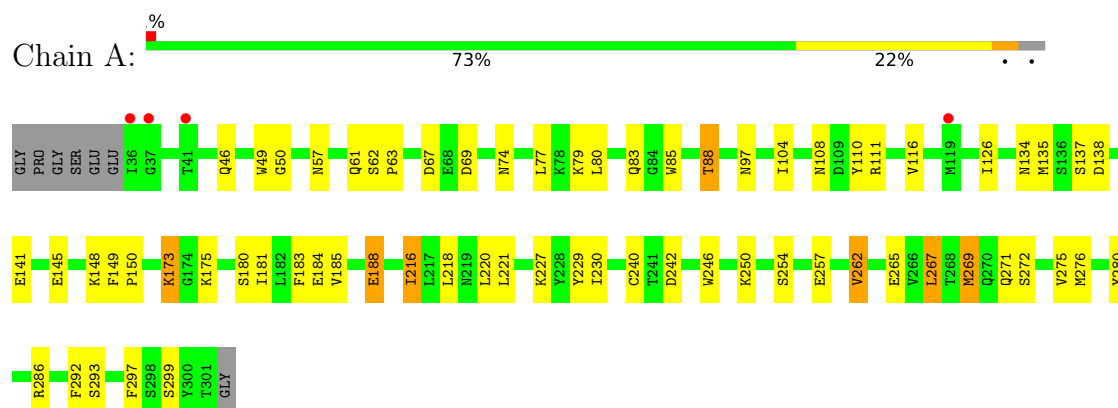
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	18	Total	O	0	0
			18	18		

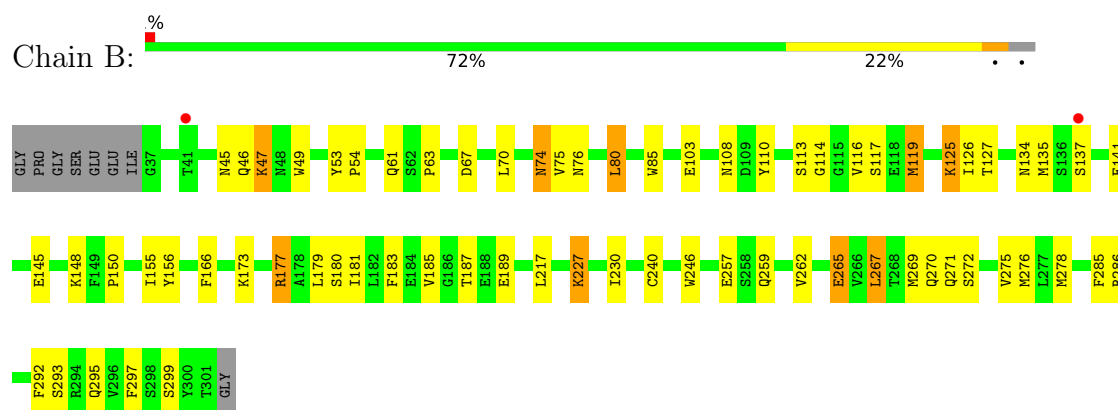
3 Residue-property plots

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.

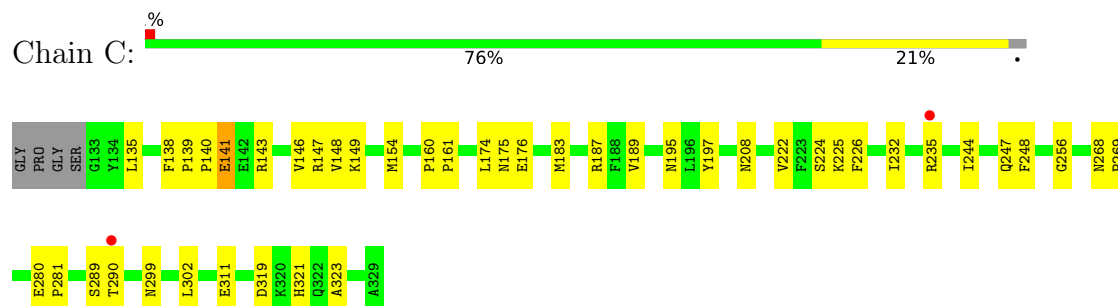
- Molecule 1: Receptor-type tyrosine-protein phosphatase zeta



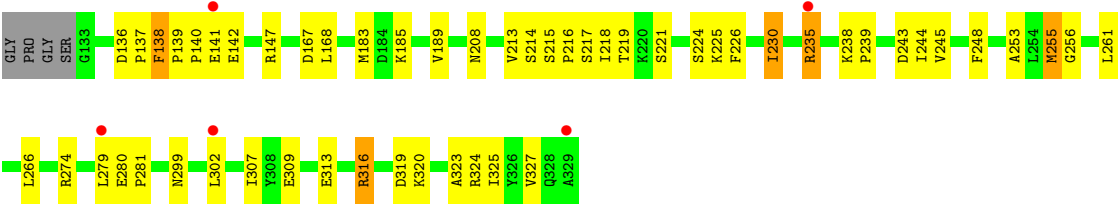
- Molecule 1: Receptor-type tyrosine-protein phosphatase zeta



- Molecule 2: Contactin-1



● Molecule 2: Contactin-1



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	334.11Å 45.96Å 78.95Å 90.00° 101.37° 90.00°	Depositor
Resolution (Å)	33.56 – 2.30 33.56 – 2.30	Depositor EDS
% Data completeness (in resolution range)	95.8 (33.56-2.30) 92.0 (33.56-2.30)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.56 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.7.1_743	Depositor
R, R_{free}	0.201 , 0.255 0.189 , 0.241	Depositor DCC
R_{free} test set	2539 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	34.6	Xtriage
Anisotropy	0.497	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 34.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.025 for -h-2*k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7633	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.42% 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 [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	0.41	0/2177	0.72	0/2942
1	B	0.41	0/2169	0.77	0/2931
2	C	0.46	0/1623	0.83	4/2207 (0.2%)
2	D	0.42	0/1623	0.84	2/2207 (0.1%)
All	All	0.42	0/7592	0.79	6/10287 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	138	PHE	CA-C-N	6.25	124.22	119.66
2	D	138	PHE	C-N-CA	6.25	124.22	119.66
2	C	143	ARG	CA-C-N	5.47	125.41	119.78
2	C	143	ARG	C-N-CA	5.47	125.41	119.78
2	C	139	PRO	CA-C-N	5.00	126.09	119.84
2	C	139	PRO	C-N-CA	5.00	126.09	119.84

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	2129	0	2068	53	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2121	0	2057	47	0
2	C	1580	0	1559	24	0
2	D	1580	0	1559	37	0
3	C	28	0	26	0	0
3	D	28	0	26	0	0
4	A	46	0	0	1	0
4	B	61	0	0	1	0
4	C	42	0	0	2	0
4	D	18	0	0	0	0
All	All	7633	0	7295	150	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 10.

All (150) 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:184:GLU:HG2	2:D:235:ARG:HH12	1.34	0.91
1:A:61:GLN:HE22	1:A:240:CYS:HB3	1.39	0.88
1:B:267:LEU:HG	1:B:278:MET:HE1	1.59	0.84
1:B:63:PRO:HG3	1:B:141:GLU:HB3	1.61	0.83
2:D:183:MET:HE3	2:D:189:VAL:H	1.44	0.81
2:D:248:PHE:CE1	2:D:323:ALA:HB3	2.14	0.81
2:D:139:PRO:HB2	2:D:141:GLU:HG2	1.65	0.79
1:A:63:PRO:HG3	1:A:141:GLU:HB3	1.65	0.77
1:B:74:ASN:HD21	1:B:76:ASN:HB2	1.48	0.76
2:D:253:ALA:O	2:D:327:VAL:HA	1.87	0.74
2:D:307:ILE:HG12	2:D:324:ARG:HG2	1.74	0.68
1:A:272:SER:O	1:A:275:VAL:HG12	1.93	0.68
1:A:61:GLN:HE21	1:A:286:ARG:HH12	1.41	0.68
2:C:140:PRO:O	2:C:225:LYS:HE2	1.94	0.67
1:A:185:VAL:H	2:D:235:ARG:NH2	1.92	0.67
2:D:214:SER:HB3	2:D:221:SER:OG	1.95	0.67
1:A:184:GLU:HG2	2:D:235:ARG:NH1	2.08	0.67
1:B:74:ASN:HD22	1:B:76:ASN:H	1.42	0.67
1:B:278:MET:HG3	2:D:266:LEU:HD12	1.76	0.66
1:B:61:GLN:HE22	1:B:240:CYS:HB3	1.59	0.66
1:B:74:ASN:HD22	1:B:74:ASN:C	2.02	0.65
2:D:280:GLU:HG3	2:D:281:PRO:HD2	1.79	0.65
1:A:185:VAL:H	2:D:235:ARG:HH22	1.45	0.64
1:A:61:GLN:NE2	1:A:286:ARG:HH12	1.96	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:275:VAL:HG21	2:C:154:MET:HE3	1.80	0.64
1:B:67:ASP:OD2	1:B:145:GLU:HA	1.98	0.63
1:B:141:GLU:HG2	1:B:285:PHE:HB2	1.81	0.62
1:B:61:GLN:NE2	1:B:286:ARG:HH12	1.98	0.61
1:B:113:SER:HA	1:B:177:ARG:NH2	2.15	0.61
1:B:148:LYS:HE2	4:B:308:HOH:O	1.99	0.61
1:A:108:ASN:HB2	1:A:110:TYR:CE2	2.36	0.60
2:D:168:LEU:HD13	2:D:213:VAL:CG1	2.31	0.60
1:B:61:GLN:HE21	1:B:286:ARG:HH12	1.49	0.60
1:B:108:ASN:HB2	1:B:110:TYR:CE2	2.37	0.60
1:A:50:GLY:HA3	1:A:57:ASN:ND2	2.17	0.59
1:B:74:ASN:ND2	1:B:76:ASN:H	2.00	0.59
1:B:113:SER:HA	1:B:177:ARG:HH22	1.67	0.59
2:D:261:LEU:HD11	2:D:325:ILE:HD11	1.84	0.59
2:D:309:GLU:OE1	2:D:320:LYS:HD2	2.03	0.58
1:A:46:GLN:HA	1:A:49:TRP:CE2	2.39	0.58
2:D:280:GLU:CG	2:D:281:PRO:HD2	2.35	0.57
2:D:239:PRO:HB2	2:D:316:ARG:HD2	1.87	0.57
2:C:135:LEU:HD21	2:C:222:VAL:HG22	1.86	0.56
2:C:248:PHE:CE1	2:C:323:ALA:HB3	2.41	0.56
2:C:146:VAL:HG12	2:C:148:VAL:HG23	1.88	0.55
1:A:108:ASN:HB2	1:A:110:TYR:CZ	2.41	0.55
1:A:271:GLN:HB2	1:A:276:MET:SD	2.46	0.55
2:D:208:ASN:HB3	2:D:226:PHE:HB3	1.89	0.54
2:D:279:LEU:O	2:D:280:GLU:HB2	2.06	0.54
1:A:135:MET:HE2	1:A:257:GLU:CD	2.32	0.54
1:A:275:VAL:CG2	2:C:154:MET:HE3	2.37	0.53
2:C:208:ASN:HB3	2:C:226:PHE:HB3	1.90	0.53
1:A:85:TRP:HB3	1:A:218:LEU:HB2	1.89	0.53
2:C:141:GLU:H	2:C:141:GLU:CD	2.16	0.53
1:B:45:ASN:CG	1:B:47:LYS:HG2	2.34	0.52
2:C:183:MET:HE2	2:C:189:VAL:HG21	1.92	0.52
1:A:50:GLY:HA3	1:A:57:ASN:HD21	1.75	0.51
1:B:230:ILE:HA	1:B:246:TRP:O	2.11	0.51
2:D:256:GLY:HA2	2:D:299:ASN:OD1	2.10	0.51
1:A:150:PRO:HD2	1:A:183:PHE:O	2.10	0.51
1:B:125:LYS:NZ	1:B:125:LYS:HB2	2.26	0.51
1:B:269:MET:SD	1:B:278:MET:HG2	2.52	0.50
1:B:134:ASN:HB2	1:B:137:SER:HB3	1.92	0.50
1:B:156:TYR:HB3	1:B:166:PHE:CE1	2.46	0.50
2:C:183:MET:HE3	4:C:90:HOH:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:46:GLN:HA	1:B:49:TRP:CE2	2.47	0.50
2:D:168:LEU:HD13	2:D:213:VAL:HG11	1.94	0.50
1:A:62:SER:HB3	1:A:63:PRO:HA	1.94	0.50
1:B:292:PHE:O	1:B:293:SER:HB2	2.10	0.50
1:A:69:ASP:OD1	1:A:69:ASP:N	2.42	0.49
2:C:280:GLU:HB3	2:C:281:PRO:HD2	1.93	0.49
1:B:180:SER:C	1:B:181:ILE:HD12	2.39	0.48
1:A:88:THR:HA	1:A:216:ILE:HD12	1.96	0.48
1:A:46:GLN:HA	1:A:49:TRP:NE1	2.29	0.48
1:B:179:LEU:HD12	1:B:179:LEU:N	2.30	0.47
1:B:259:GLN:O	1:B:262:VAL:HG22	2.14	0.47
1:B:150:PRO:HD2	1:B:183:PHE:O	2.15	0.47
1:A:77:LEU:HD12	1:A:299:SER:HA	1.97	0.47
2:D:302:LEU:H	2:D:302:LEU:HD12	1.80	0.47
2:C:289:SER:HA	4:C:60:HOH:O	2.15	0.46
1:A:173:LYS:N	1:A:173:LYS:HD3	2.31	0.46
1:B:103:GLU:HG3	1:B:127:THR:HG22	1.99	0.45
2:C:195:ASN:HB2	2:C:197:TYR:CE1	2.51	0.45
1:A:181:ILE:HD12	1:A:181:ILE:N	2.31	0.45
1:A:188:GLU:H	1:A:188:GLU:HG2	1.44	0.45
1:B:53:TYR:HA	1:B:54:PRO:HD3	1.84	0.45
2:C:268:ASN:HD22	2:C:269:PRO:HA	1.81	0.45
1:B:269:MET:HG2	1:B:278:MET:HG2	1.99	0.44
1:A:67:ASP:OD2	1:A:145:GLU:HA	2.18	0.44
1:A:184:GLU:CG	2:D:235:ARG:HH12	2.18	0.44
1:B:230:ILE:HG22	1:B:297:PHE:HB2	1.99	0.44
2:D:139:PRO:HA	2:D:140:PRO:HD3	1.84	0.44
1:A:275:VAL:HG21	2:C:154:MET:CE	2.46	0.44
1:B:117:SER:OG	1:B:119:MET:HB2	2.18	0.44
1:A:230:ILE:HG22	1:A:297:PHE:HB2	1.98	0.44
1:A:149:PHE:HB3	1:A:183:PHE:O	2.18	0.43
2:C:149:LYS:HA	2:C:232:ILE:HB	2.00	0.43
2:D:147:ARG:HG2	2:D:230:ILE:HB	2.00	0.43
1:A:184:GLU:HA	2:D:235:ARG:HH22	1.82	0.43
1:A:230:ILE:HA	1:A:246:TRP:O	2.18	0.43
2:D:136:ASP:HB3	2:D:137:PRO:HD2	2.00	0.43
1:A:83:GLN:HB2	1:A:111:ARG:HB2	1.99	0.43
1:A:292:PHE:O	1:A:293:SER:HB2	2.18	0.43
1:B:74:ASN:C	1:B:74:ASN:ND2	2.71	0.43
1:B:80:LEU:HD12	1:B:114:GLY:HA3	2.01	0.43
1:A:269:MET:HE2	1:A:276:MET:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:85:TRP:CD2	1:B:217:LEU:HD23	2.53	0.43
2:D:274:ARG:HH22	2:D:313:GLU:HB3	1.83	0.43
1:A:138:ASP:OD1	1:A:138:ASP:N	2.51	0.43
1:B:108:ASN:HB2	1:B:110:TYR:CZ	2.53	0.43
1:B:181:ILE:HD12	1:B:181:ILE:N	2.33	0.43
2:D:255:MET:HB3	2:D:255:MET:HE2	1.78	0.43
1:A:61:GLN:HE21	1:A:286:ARG:NH1	2.11	0.42
1:A:267:LEU:HD11	1:A:280:TYR:CE1	2.54	0.42
1:A:104:ILE:HB	1:A:126:ILE:HB	2.01	0.42
1:B:227:LYS:HA	1:B:299:SER:OG	2.18	0.42
2:D:218:ILE:O	2:D:219:THR:C	2.62	0.42
1:A:134:ASN:HB2	1:A:137:SER:HB2	2.02	0.42
1:A:180:SER:C	1:A:181:ILE:HD12	2.44	0.42
2:D:243:ASP:OD1	2:D:243:ASP:C	2.62	0.42
1:A:227:LYS:HD2	1:A:227:LYS:HA	1.88	0.42
1:A:148:LYS:HD3	4:A:303:HOH:O	2.19	0.42
1:B:271:GLN:HB2	1:B:276:MET:SD	2.60	0.42
2:C:138:PHE:CD2	2:C:224:SER:HB3	2.55	0.42
1:A:262:VAL:HA	1:A:265:GLU:HG3	2.01	0.42
2:D:244:ILE:HG13	2:D:319:ASP:HB3	2.02	0.42
1:A:61:GLN:NE2	1:A:240:CYS:HB3	2.20	0.41
2:C:174:LEU:O	2:C:175:ASN:HB2	2.18	0.41
1:A:61:GLN:HE21	1:A:286:ARG:HH22	1.67	0.41
1:B:70:LEU:HD23	1:B:70:LEU:HA	1.81	0.41
2:D:225:LYS:HA	2:D:225:LYS:HD3	1.87	0.41
1:B:262:VAL:HA	1:B:265:GLU:HG3	2.01	0.41
1:B:67:ASP:CG	2:C:147:ARG:HH22	2.28	0.41
2:C:247:GLN:HG2	2:C:321:HIS:CG	2.55	0.41
2:C:160:PRO:HA	2:C:161:PRO:HD3	1.86	0.41
1:B:80:LEU:HD12	1:B:80:LEU:HA	1.93	0.41
1:A:134:ASN:O	1:A:137:SER:HB3	2.21	0.41
1:B:126:ILE:HG12	1:B:155:ILE:HG12	2.02	0.41
1:B:135:MET:HB2	1:B:185:VAL:HB	2.02	0.41
2:C:138:PHE:CE2	2:C:224:SER:HB3	2.56	0.41
2:C:244:ILE:HG13	2:C:319:ASP:HB3	2.03	0.41
2:C:256:GLY:O	2:C:299:ASN:HA	2.20	0.41
1:A:74:ASN:C	1:A:74:ASN:HD22	2.28	0.41
1:B:189:GLU:HG2	1:B:259:GLN:NE2	2.36	0.41
2:D:167:ASP:OD2	2:D:216:PRO:HG2	2.21	0.41
2:D:238:LYS:HA	2:D:239:PRO:HD3	1.85	0.41
1:A:97:ASN:OD1	1:A:97:ASN:C	2.63	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:138:PHE:CD2	2:D:224:SER:HB3	2.56	0.40
1:A:229:TYR:OH	1:A:250:LYS:HD2	2.21	0.40
2:D:215:SER:HA	2:D:216:PRO:HD2	1.93	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	264/273 (97%)	255 (97%)	9 (3%)	0	100	100
1	B	263/273 (96%)	255 (97%)	8 (3%)	0	100	100
2	C	195/201 (97%)	186 (95%)	9 (5%)	0	100	100
2	D	195/201 (97%)	189 (97%)	6 (3%)	0	100	100
All	All	917/948 (97%)	885 (96%)	32 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	240/244 (98%)	225 (94%)	15 (6%)	16	23
1	B	239/244 (98%)	221 (92%)	18 (8%)	12	17

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	C	177/179 (99%)	170 (96%)	7 (4%)	28	42
2	D	177/179 (99%)	169 (96%)	8 (4%)	24	37
All	All	833/846 (98%)	785 (94%)	48 (6%)	18	26

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	79	LYS
1	A	80	LEU
1	A	88	THR
1	A	116	VAL
1	A	173	LYS
1	A	175	LYS
1	A	188	GLU
1	A	216	ILE
1	A	220	LEU
1	A	221	LEU
1	A	242	ASP
1	A	254	SER
1	A	262	VAL
1	A	267	LEU
1	A	269	MET
1	B	47	LYS
1	B	74	ASN
1	B	75	VAL
1	B	80	LEU
1	B	116	VAL
1	B	119	MET
1	B	125	LYS
1	B	173	LYS
1	B	177	ARG
1	B	187	THR
1	B	227	LYS
1	B	257	GLU
1	B	265	GLU
1	B	267	LEU
1	B	270	GLN
1	B	272	SER
1	B	275	VAL
1	B	295	GLN
2	C	141	GLU

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Mol	Chain	Res	Type
2	C	176	GLU
2	C	187	ARG
2	C	235	ARG
2	C	290	THR
2	C	302	LEU
2	C	311	GLU
2	D	142	GLU
2	D	185	LYS
2	D	217	SER
2	D	230	ILE
2	D	235	ARG
2	D	245	VAL
2	D	255	MET
2	D	316	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	48	ASN
1	A	57	ASN
1	A	61	GLN
1	A	74	ASN
1	A	76	ASN
1	A	83	GLN
1	A	209	GLN
1	A	219	ASN
1	A	270	GLN
1	A	289	GLN
1	B	61	GLN
1	B	74	ASN
1	B	83	GLN
1	B	219	ASN
1	B	270	GLN
1	B	283	ASN
1	B	295	GLN
2	C	257	GLN
2	C	268	ASN
2	C	322	GLN
2	D	200	ASN
2	D	268	ASN
2	D	322	GLN

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 oligosaccharides 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$
3	NAG	D	208(A)	2	14,14,15	0.56	0	17,19,21	0.94	1 (5%)
3	NAG	D	258(A)	2	14,14,15	1.30	1 (7%)	17,19,21	2.59	9 (52%)
3	NAG	C	208(A)	2	14,14,15	0.48	0	17,19,21	1.09	1 (5%)
3	NAG	C	258(A)	2	14,14,15	0.56	0	17,19,21	1.00	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
3	NAG	D	208(A)	2	-	0/6/23/26	0/1/1/1
3	NAG	D	258(A)	2	-	2/6/23/26	0/1/1/1
3	NAG	C	208(A)	2	-	1/6/23/26	0/1/1/1
3	NAG	C	258(A)	2	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	258(A)	NAG	C1-C2	3.58	1.57	1.52

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	258(A)	NAG	C1-O5-C5	5.60	119.70	112.19
3	D	258(A)	NAG	C1-C2-N2	5.55	119.17	110.43
3	C	208(A)	NAG	C1-O5-C5	3.63	117.05	112.19
3	C	258(A)	NAG	C2-N2-C7	-3.48	118.23	122.90
3	D	258(A)	NAG	O4-C4-C5	3.09	116.93	109.32
3	D	208(A)	NAG	C1-O5-C5	3.04	116.25	112.19
3	D	258(A)	NAG	O5-C1-C2	-2.64	107.21	111.29
3	D	258(A)	NAG	C3-C4-C5	-2.49	105.72	110.23
3	D	258(A)	NAG	O4-C4-C3	-2.38	104.78	110.38
3	D	258(A)	NAG	C8-C7-N2	2.32	119.96	116.12
3	D	258(A)	NAG	O7-C7-C8	-2.30	117.95	122.05
3	D	258(A)	NAG	O3-C3-C4	-2.05	105.55	110.38

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	258(A)	NAG	O5-C5-C6-O6
3	C	258(A)	NAG	C4-C5-C6-O6
3	D	258(A)	NAG	C4-C5-C6-O6
3	C	258(A)	NAG	O5-C5-C6-O6
3	C	208(A)	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	266/273 (97%)	-0.15	4 (1%) 72 73	24, 41, 72, 90	0
1	B	265/273 (97%)	-0.29	2 (0%) 82 83	23, 37, 67, 86	0
2	C	197/201 (98%)	-0.36	2 (1%) 79 80	22, 37, 57, 96	0
2	D	197/201 (98%)	0.18	5 (2%) 58 60	26, 47, 88, 103	0
All	All	925/948 (97%)	-0.16	13 (1%) 73 75	22, 40, 74, 103	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	36	ILE	5.2
2	C	290	THR	3.2
2	D	329	ALA	3.0
1	A	37	GLY	2.7
2	D	235	ARG	2.6
2	D	279	LEU	2.6
2	D	302	LEU	2.4
2	D	141	GLU	2.4
1	B	41	THR	2.3
1	A	119	MET	2.1
1	A	41	THR	2.1
1	B	137	SER	2.0
2	C	235	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 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
3	NAG	D	258(A)	14/15	0.50	0.17	67,103,109,111	0
3	NAG	C	258(A)	14/15	0.80	0.14	65,75,88,94	0
3	NAG	D	208(A)	14/15	0.87	0.10	48,58,63,64	0
3	NAG	C	208(A)	14/15	0.91	0.09	39,54,64,67	0

6.5 Other polymers [i](#)

There are no such residues in this entry.