



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 06:55 am BST

PDB ID : 2ZII
Title : Crystal Structure of Yeast Vps74-N-term Truncation Variant
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Deposited on : 2008-02-18
Resolution : 3.05 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

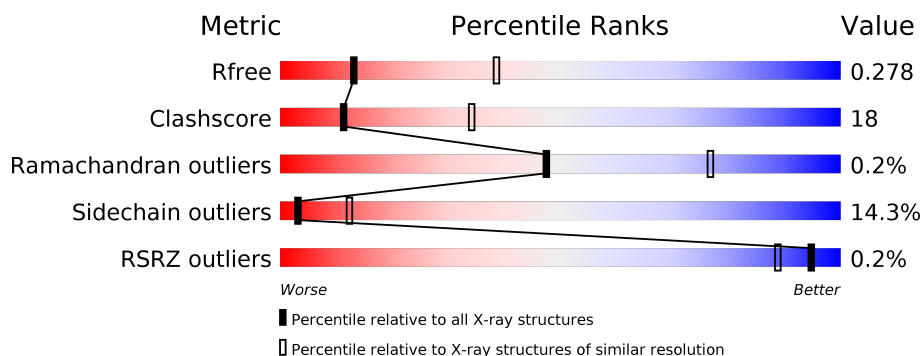
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.05 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}	130704	1754 (3.10-3.02)
Clashscore	141614	1864 (3.10-3.02)
Ramachandran outliers	138981	1794 (3.10-3.02)
Sidechain outliers	138945	1793 (3.10-3.02)
RSRZ outliers	127900	1713 (3.10-3.02)

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 ≥ 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	288	
1	B	288	
1	C	288	
1	D	288	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8590 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 Vacuolar protein sorting-associated protein 74.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	281	Total	C	N	O	S	0	0	0
			2149	1365	356	416	12			
1	B	281	Total	C	N	O	S	0	0	0
			2126	1348	356	411	11			
1	C	279	Total	C	N	O	S	0	0	0
			2125	1351	355	408	11			
1	D	283	Total	C	N	O	S	0	0	0
			2188	1391	371	414	12			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	58	GLY	-	EXPRESSION TAG	UNP Q06385
A	59	SER	-	EXPRESSION TAG	UNP Q06385
B	58	GLY	-	EXPRESSION TAG	UNP Q06385
B	59	SER	-	EXPRESSION TAG	UNP Q06385
C	58	GLY	-	EXPRESSION TAG	UNP Q06385
C	59	SER	-	EXPRESSION TAG	UNP Q06385
D	58	GLY	-	EXPRESSION TAG	UNP Q06385
D	59	SER	-	EXPRESSION TAG	UNP Q06385

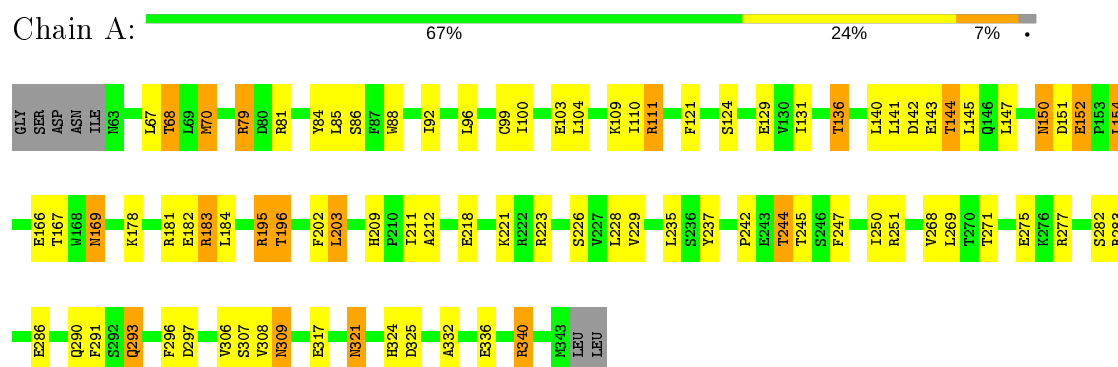
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	O	0	0
			1	1		
2	D	1	Total	O	0	0
			1	1		

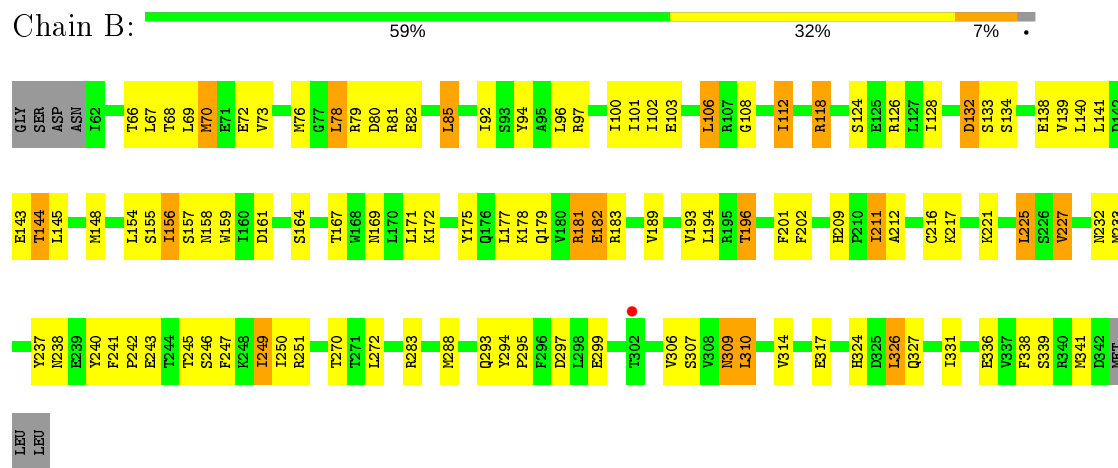
3 Residue-property plots

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

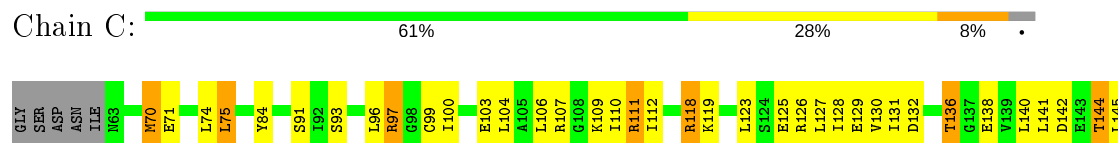
- Molecule 1: Vacuolar protein sorting-associated protein 74

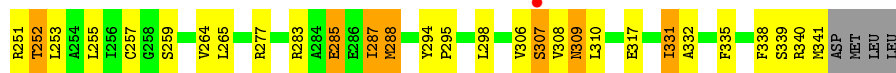


- Molecule 1: Vacuolar protein sorting-associated protein 74



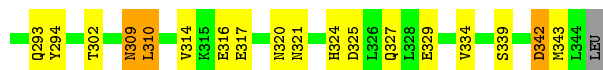
- Molecule 1: Vacuolar protein sorting-associated protein 74





- Molecule 1: Vacuolar protein sorting-associated protein 74

Chain D: 68% 24% 7%



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	104.08Å 104.08Å 292.76Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.10 – 3.05 49.10 – 3.05	Depositor EDS
% Data completeness (in resolution range)	99.7 (49.10-3.05) 99.8 (49.10-3.05)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.25 (at 3.07Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.220 , 0.281 0.219 , 0.278	Depositor DCC
R_{free} test set	1790 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	57.8	Xtriage
Anisotropy	0.283	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 59.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.013 for -h,-k,l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8590	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 22.79 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.3113e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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

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	0.67	0/2180	0.82	2/2959 (0.1%)
1	B	0.58	0/2157	0.76	0/2934
1	C	0.63	0/2156	0.85	6/2929 (0.2%)
1	D	0.68	1/2218 (0.0%)	0.82	2/3005 (0.1%)
All	All	0.64	1/8711 (0.0%)	0.81	10/11827 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	99	CYS	CB-SG	-5.99	1.72	1.81

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	228	LEU	CA-CB-CG	8.54	134.93	115.30
1	A	235	LEU	CA-CB-CG	5.83	128.70	115.30
1	A	154	LEU	CB-CG-CD2	-5.82	101.10	111.00
1	C	111	ARG	NE-CZ-NH1	5.71	123.16	120.30
1	C	170	LEU	CB-CA-C	-5.59	99.59	110.20
1	C	172	LYS	N-CA-C	-5.27	96.77	111.00
1	C	97	ARG	NE-CZ-NH1	5.25	122.92	120.30
1	C	175	TYR	CA-CB-CG	5.13	123.15	113.40
1	D	147	LEU	CA-CB-CG	5.08	126.97	115.30
1	D	97	ARG	NE-CZ-NH1	5.07	122.83	120.30

There are no chirality outliers.

There are no planarity outliers.

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	2149	0	2079	73	0
1	B	2126	0	2026	96	0
1	C	2125	0	2049	79	0
1	D	2188	0	2162	62	0
2	A	1	0	0	0	0
2	D	1	0	0	0	0
All	All	8590	0	8316	300	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 18.

All (300) 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:B:106:LEU:HD13	1:B:331:ILE:HD11	1.35	1.05
1:A:70:MET:H	1:A:70:MET:HE2	1.18	1.05
1:A:242:PRO:HD2	1:A:245:THR:CG2	1.87	1.04
1:C:233:MET:HE1	1:C:310:LEU:HD13	1.45	0.97
1:C:198:MET:HE1	1:C:205:ASP:HB3	1.50	0.93
1:C:100:ILE:HG21	1:C:144:THR:HG21	1.50	0.92
1:D:309:ASN:H	1:D:309:ASN:HD22	1.18	0.91
1:C:233:MET:CE	1:C:310:LEU:HD13	2.07	0.84
1:D:324:HIS:HD2	1:D:327:GLN:OE1	1.62	0.83
1:A:242:PRO:HD2	1:A:245:THR:HG21	1.60	0.82
1:B:194:LEU:HD13	1:B:209:HIS:HB3	1.59	0.81
1:B:106:LEU:CD1	1:B:331:ILE:HD11	2.12	0.80
1:D:233:MET:H	1:D:309:ASN:HD21	1.28	0.80
1:B:155:SER:OG	1:B:158:ASN:HB2	1.83	0.79
1:D:101:ILE:HD11	1:D:334:VAL:HG22	1.65	0.79
1:D:101:ILE:HD11	1:D:334:VAL:CG2	2.13	0.78
1:A:70:MET:N	1:A:70:MET:HE2	1.97	0.76
1:C:309:ASN:HD22	1:C:309:ASN:H	1.33	0.76
1:C:136:THR:HG23	1:C:138:GLU:H	1.51	0.75
1:D:79:ARG:HG3	1:D:86:SER:OG	1.87	0.74
1:A:136:THR:HG21	1:A:141:LEU:HB2	1.70	0.74
1:B:242:PRO:HD2	1:B:245:THR:CG2	2.19	0.73
1:A:79:ARG:HG3	1:A:86:SER:OG	1.89	0.72
1:A:178:LYS:O	1:A:183:ARG:NH1	2.22	0.72
1:C:309:ASN:H	1:C:309:ASN:ND2	1.87	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:189:VAL:HG23	1:B:194:LEU:HD12	1.72	0.72
1:A:242:PRO:HB2	1:A:244:THR:HB	1.72	0.72
1:D:100:ILE:HG21	1:D:144:THR:HG21	1.73	0.71
1:B:81:ARG:NH1	1:D:167:THR:O	2.24	0.71
1:B:309:ASN:HD22	1:B:309:ASN:H	1.36	0.70
1:C:283:ARG:HE	1:C:307:SER:HB2	1.56	0.70
1:B:94:TYR:HD2	1:B:341:MET:CE	2.05	0.70
1:B:97:ARG:HG3	1:B:164:SER:HB3	1.74	0.70
1:C:283:ARG:HE	1:C:307:SER:CB	2.06	0.69
1:D:270:THR:HA	1:D:277:ARG:HD3	1.74	0.69
1:C:100:ILE:CG2	1:C:144:THR:HG21	2.20	0.69
1:A:100:ILE:HG21	1:A:144:THR:HG21	1.74	0.69
1:B:118:ARG:HD2	1:B:126:ARG:HD3	1.73	0.69
1:B:310:LEU:O	1:B:314:VAL:HG23	1.92	0.69
1:C:208:THR:HG23	1:C:210:PRO:HD3	1.76	0.68
1:B:309:ASN:H	1:B:309:ASN:ND2	1.91	0.68
1:D:257:CYS:SG	1:D:288:MET:HG2	2.34	0.68
1:B:108:GLY:HA2	1:B:327:GLN:HE21	1.60	0.67
1:C:197:GLU:HB3	1:C:208:THR:HG22	1.75	0.67
1:B:242:PRO:HD2	1:B:245:THR:HG21	1.75	0.66
1:B:288:MET:HE1	1:B:338:PHE:HB2	1.77	0.66
1:B:238:ASN:OD1	1:B:241:PHE:N	2.26	0.66
1:D:197:GLU:O	1:D:208:THR:HB	1.96	0.65
1:C:220:ILE:HG23	1:C:255:LEU:HD21	1.79	0.65
1:B:194:LEU:HA	1:B:212:ALA:H	1.62	0.65
1:C:198:MET:CE	1:C:205:ASP:HB3	2.24	0.65
1:A:196:THR:HB	1:A:209:HIS:HD2	1.62	0.64
1:B:237:TYR:CD2	1:B:243:GLU:HA	2.32	0.64
1:C:147:LEU:HD22	1:C:175:TYR:CD1	2.32	0.64
1:A:218:GLU:HA	1:A:218:GLU:OE2	1.97	0.64
1:B:69:LEU:O	1:B:73:VAL:HG23	1.98	0.64
1:D:288:MET:HE1	1:D:339:SER:HA	1.80	0.64
1:A:242:PRO:O	1:A:245:THR:HG23	1.98	0.63
1:A:111:ARG:NH1	1:A:129:GLU:OE1	2.32	0.62
1:D:188:LEU:HB3	1:D:194:LEU:HD22	1.81	0.62
1:B:227:VAL:O	1:B:307:SER:HB2	2.00	0.62
1:C:283:ARG:HH21	1:C:307:SER:HB3	1.65	0.62
1:C:233:MET:HE1	1:C:310:LEU:CD1	2.26	0.62
1:D:140:LEU:O	1:D:144:THR:HG23	1.99	0.61
1:D:247:PHE:HB3	1:D:250:ILE:HB	1.82	0.61
1:B:189:VAL:HG23	1:B:194:LEU:CD1	2.31	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:94:TYR:HD2	1:B:341:MET:HE2	1.65	0.61
1:C:233:MET:H	1:C:309:ASN:HD21	1.49	0.61
1:A:103:GLU:OE2	1:A:251:ARG:NH2	2.29	0.60
1:A:247:PHE:O	1:A:251:ARG:HG3	2.00	0.60
1:B:297:ASP:OD1	1:B:299:GLU:HB2	2.01	0.60
1:A:151:ASP:OD1	1:A:152:GLU:N	2.31	0.60
1:C:140:LEU:O	1:C:144:THR:HG22	2.01	0.60
1:C:169:ASN:HD22	1:C:169:ASN:C	2.04	0.60
1:B:233:MET:H	1:B:309:ASN:HD21	1.50	0.60
1:D:324:HIS:CD2	1:D:327:GLN:OE1	2.50	0.60
1:A:309:ASN:H	1:A:309:ASN:HD22	1.49	0.60
1:A:136:THR:HG21	1:A:141:LEU:CB	2.32	0.60
1:C:70:MET:H	1:C:70:MET:HE3	1.66	0.59
1:B:100:ILE:HG21	1:B:144:THR:HG21	1.84	0.59
1:B:169:ASN:ND2	1:B:172:LYS:HE2	2.18	0.59
1:C:250:ILE:HG12	1:C:317:GLU:OE1	2.02	0.59
1:C:112:ILE:HD11	1:C:126:ARG:HB3	1.84	0.58
1:C:136:THR:CG2	1:C:138:GLU:H	2.16	0.58
1:B:80:ASP:O	1:B:211:ILE:HD13	2.04	0.58
1:C:96:LEU:O	1:C:99:CYS:HB2	2.03	0.58
1:D:112:ILE:HD12	1:D:329:GLU:HB3	1.84	0.57
1:B:238:ASN:HD21	1:B:240:TYR:HB2	1.67	0.57
1:A:84:TYR:CE1	1:A:85:LEU:HG	2.38	0.57
1:D:108:GLY:HA2	1:D:327:GLN:NE2	2.18	0.57
1:C:224:VAL:O	1:C:228:LEU:HD23	2.04	0.57
1:B:178:LYS:O	1:B:183:ARG:NH1	2.38	0.56
1:B:140:LEU:O	1:B:144:THR:HG23	2.04	0.56
1:B:294:TYR:CD1	1:B:295:PRO:HA	2.40	0.56
1:A:121:PHE:CE2	1:C:277:ARG:HD2	2.40	0.56
1:A:202:PHE:CD2	1:A:203:LEU:HD23	2.41	0.56
1:B:179:GLN:O	1:B:183:ARG:HD3	2.05	0.56
1:A:181:ARG:NH1	1:B:201:PHE:O	2.39	0.56
1:B:283:ARG:HE	1:B:307:SER:HB3	1.71	0.56
1:B:189:VAL:CG2	1:B:194:LEU:HD12	2.33	0.56
1:B:126:ARG:O	1:B:155:SER:HA	2.06	0.56
1:B:309:ASN:HD22	1:B:309:ASN:N	2.02	0.56
1:A:144:THR:HA	1:A:147:LEU:HD12	1.88	0.55
1:A:195:ARG:HG2	1:A:212:ALA:HB2	1.86	0.55
1:C:103:GLU:HG3	1:C:107:ARG:HE	1.70	0.55
1:D:309:ASN:H	1:D:309:ASN:ND2	1.98	0.55
1:B:100:ILE:HG23	1:B:141:LEU:HD23	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:123:LEU:HA	1:C:126:ARG:HD2	1.88	0.55
1:B:154:LEU:HB3	1:B:158:ASN:HD22	1.70	0.55
1:D:151:ASP:OD1	1:D:152:GLU:N	2.39	0.55
1:D:195:ARG:HG3	1:D:212:ALA:HB2	1.86	0.55
1:C:294:TYR:HA	1:C:295:PRO:C	2.27	0.55
1:B:324:HIS:HD2	1:B:327:GLN:OE1	1.90	0.55
1:D:171:LEU:HD12	1:D:171:LEU:H	1.71	0.55
1:B:194:LEU:HB3	1:B:211:ILE:HA	1.89	0.55
1:B:124:SER:HB2	1:B:157:SER:HB2	1.89	0.55
1:B:161:ASP:HB3	1:B:167:THR:HG23	1.89	0.54
1:B:103:GLU:OE2	1:B:251:ARG:NH2	2.41	0.54
1:B:82:GLU:OE1	1:B:82:GLU:HA	2.06	0.54
1:B:182:GLU:CD	1:B:182:GLU:H	2.10	0.54
1:C:178:LYS:O	1:C:183:ARG:NH1	2.40	0.54
1:C:202:PHE:CG	1:D:86:SER:HB2	2.43	0.54
1:B:233:MET:N	1:B:309:ASN:HD21	2.05	0.54
1:A:136:THR:N	1:A:142:ASP:OD1	2.36	0.54
1:C:179:GLN:O	1:C:183:ARG:HB2	2.09	0.53
1:B:67:LEU:HD21	1:B:193:VAL:HG21	1.90	0.53
1:C:136:THR:HG21	1:C:141:LEU:HB2	1.91	0.53
1:C:285:GLU:HA	1:C:288:MET:HB2	1.91	0.53
1:C:309:ASN:HD22	1:C:309:ASN:N	2.03	0.53
1:D:154:LEU:HB3	1:D:158:ASN:ND2	2.24	0.53
1:D:178:LYS:O	1:D:183:ARG:NH1	2.41	0.53
1:D:194:LEU:HD12	1:D:211:ILE:HA	1.90	0.53
1:B:194:LEU:HA	1:B:212:ALA:N	2.23	0.53
1:B:94:TYR:HD2	1:B:341:MET:HE1	1.74	0.53
1:D:233:MET:N	1:D:309:ASN:HD21	2.03	0.53
1:D:158:ASN:O	1:D:162:LEU:HG	2.09	0.52
1:B:94:TYR:CD2	1:B:341:MET:HE2	2.44	0.52
1:B:112:ILE:CD1	1:B:326:LEU:HD22	2.40	0.52
1:C:283:ARG:O	1:C:287:ILE:HG23	2.09	0.52
1:A:104:LEU:HB3	1:A:110:ILE:HG12	1.92	0.52
1:B:216:CYS:HB2	1:B:240:TYR:CE1	2.44	0.52
1:B:171:LEU:C	1:B:172:LYS:HG2	2.29	0.52
1:A:218:GLU:OE1	1:D:273:GLU:HA	2.09	0.52
1:A:136:THR:HG22	1:A:142:ASP:CG	2.30	0.51
1:A:196:THR:HB	1:A:209:HIS:CD2	2.44	0.51
1:A:321:ASN:ND2	1:A:324:HIS:CE1	2.79	0.51
1:C:172:LYS:HG2	1:C:175:TYR:OH	2.10	0.51
1:D:250:ILE:HG12	1:D:317:GLU:OE1	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:271:THR:HB	1:D:271:THR:HB	1.91	0.51
1:B:154:LEU:HD11	1:B:172:LYS:NZ	2.26	0.51
1:B:247:PHE:O	1:B:251:ARG:HG3	2.09	0.51
1:B:112:ILE:HD11	1:B:326:LEU:HD22	1.92	0.51
1:A:169:ASN:C	1:A:169:ASN:ND2	2.65	0.51
1:D:102:ILE:HG13	1:D:334:VAL:HG11	1.91	0.51
1:A:290:GLN:HB3	1:A:291:PHE:CD1	2.47	0.50
1:D:195:ARG:NH1	1:D:197:GLU:OE2	2.44	0.50
1:A:79:ARG:NH1	1:A:81:ARG:O	2.44	0.50
1:B:156:ILE:HG13	1:B:157:SER:N	2.26	0.50
1:B:144:THR:HG22	1:B:177:LEU:HD21	1.94	0.50
1:D:96:LEU:O	1:D:99:CYS:HB2	2.12	0.50
1:C:288:MET:HE1	1:C:335:PHE:O	2.12	0.50
1:D:111:ARG:HD3	1:D:131:ILE:HD11	1.93	0.50
1:B:106:LEU:CD1	1:B:331:ILE:CD1	2.89	0.50
1:B:247:PHE:HB3	1:B:250:ILE:HB	1.94	0.49
1:C:112:ILE:HD11	1:C:126:ARG:CB	2.42	0.49
1:B:68:THR:OG1	1:B:70:MET:HG2	2.13	0.49
1:C:169:ASN:ND2	1:C:172:LYS:H	2.10	0.49
1:B:132:ASP:OD1	1:B:133:SER:N	2.45	0.49
1:A:84:TYR:O	1:B:201:PHE:HB3	2.12	0.49
1:D:148:MET:HG2	1:D:159:TRP:CZ2	2.48	0.49
1:B:288:MET:HE1	1:B:338:PHE:CB	2.42	0.49
1:B:249:ILE:HD11	1:B:314:VAL:HG22	1.93	0.49
1:C:181:ARG:HD2	1:D:202:PHE:O	2.13	0.49
1:C:242:PRO:HD2	1:C:245:THR:CG2	2.43	0.49
1:A:140:LEU:O	1:A:144:THR:CG2	2.61	0.49
1:A:242:PRO:HD2	1:A:245:THR:HG23	1.89	0.49
1:B:297:ASP:C	1:B:299:GLU:H	2.17	0.49
1:C:288:MET:SD	1:C:339:SER:HB3	2.53	0.49
1:A:111:ARG:HD2	1:A:131:ILE:HG21	1.95	0.48
1:B:101:ILE:O	1:B:102:ILE:C	2.48	0.48
1:D:171:LEU:N	1:D:171:LEU:HD12	2.28	0.48
1:B:143:GLU:CD	1:B:183:ARG:NH2	2.67	0.48
1:C:71:GLU:O	1:C:75:LEU:HD22	2.13	0.48
1:B:249:ILE:HG23	1:B:317:GLU:OE2	2.13	0.48
1:A:169:ASN:HD22	1:A:169:ASN:C	2.17	0.48
1:D:101:ILE:CD1	1:D:334:VAL:CG2	2.89	0.48
1:A:150:ASN:CG	1:A:150:ASN:O	2.51	0.48
1:B:181:ARG:HB2	1:B:182:GLU:OE2	2.14	0.48
1:C:259:SER:HB2	1:C:264:VAL:HG23	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:136:THR:OG1	1:D:142:ASP:OD2	2.25	0.48
1:A:247:PHE:HB3	1:A:250:ILE:HB	1.95	0.47
1:C:136:THR:HG21	1:C:141:LEU:CB	2.44	0.47
1:B:242:PRO:O	1:B:245:THR:HG23	2.14	0.47
1:C:75:LEU:HD21	1:C:188:LEU:HD21	1.95	0.47
1:D:221:LYS:HG3	1:D:268:VAL:HG22	1.97	0.47
1:A:250:ILE:HG12	1:A:317:GLU:OE1	2.15	0.47
1:B:233:MET:H	1:B:309:ASN:ND2	2.11	0.47
1:A:242:PRO:C	1:A:244:THR:H	2.18	0.47
1:A:86:SER:HB2	1:B:202:PHE:CG	2.49	0.47
1:D:255:LEU:HD22	1:D:255:LEU:O	2.15	0.46
1:A:182:GLU:OE2	1:A:182:GLU:N	2.45	0.46
1:D:316:GLU:O	1:D:320:ASN:ND2	2.48	0.46
1:D:277:ARG:O	1:D:281:ILE:HG12	2.15	0.46
1:A:143:GLU:OE2	1:A:183:ARG:NH2	2.49	0.46
1:A:96:LEU:O	1:A:99:CYS:HB2	2.15	0.46
1:D:171:LEU:CD1	1:D:171:LEU:H	2.29	0.46
1:B:196:THR:HB	1:B:209:HIS:CD2	2.50	0.46
1:C:70:MET:HE3	1:C:70:MET:N	2.30	0.46
1:A:283:ARG:NH1	1:A:286:GLU:OE1	2.48	0.46
1:B:108:GLY:HA2	1:B:327:GLN:NE2	2.28	0.46
1:A:242:PRO:HD2	1:A:245:THR:HG22	1.91	0.46
1:C:84:TYR:CG	1:D:202:PHE:HB2	2.51	0.46
1:B:96:LEU:O	1:B:100:ILE:HG13	2.16	0.45
1:D:139:VAL:O	1:D:143:GLU:HG3	2.16	0.45
1:A:269:LEU:O	1:A:277:ARG:HD3	2.17	0.45
1:A:124:SER:HB3	1:A:336:GLU:OE1	2.16	0.45
1:D:103:GLU:HG2	1:D:141:LEU:HD11	1.97	0.45
1:D:310:LEU:O	1:D:314:VAL:HG23	2.16	0.45
1:B:196:THR:HB	1:B:209:HIS:HD2	1.80	0.45
1:B:237:TYR:N	1:B:237:TYR:CD1	2.85	0.45
1:B:324:HIS:CD2	1:B:327:GLN:OE1	2.70	0.45
1:C:147:LEU:HD22	1:C:175:TYR:CG	2.52	0.45
1:D:233:MET:HG2	1:D:235:LEU:HD13	1.99	0.45
1:C:136:THR:N	1:C:142:ASP:OD1	2.47	0.45
1:B:78:LEU:O	1:B:85:LEU:HD12	2.16	0.44
1:C:332:ALA:O	1:C:335:PHE:HB2	2.18	0.44
1:A:306:VAL:O	1:A:307:SER:C	2.55	0.44
1:B:138:GLU:HB2	1:B:141:LEU:HD12	2.00	0.44
1:C:247:PHE:O	1:C:251:ARG:HG3	2.17	0.44
1:A:145:LEU:HD23	1:A:145:LEU:HA	1.76	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:321:ASN:ND2	1:A:324:HIS:HE1	2.16	0.44
1:C:171:LEU:O	1:C:173:ILE:N	2.49	0.44
1:C:169:ASN:CG	1:C:172:LYS:HB2	2.38	0.44
1:C:253:LEU:HD21	1:C:310:LEU:HD11	2.00	0.44
1:A:321:ASN:HD21	1:A:324:HIS:CE1	2.35	0.44
1:D:294:TYR:CD2	1:D:329:GLU:HG3	2.53	0.44
1:A:336:GLU:O	1:A:340:ARG:HG2	2.17	0.44
1:D:162:LEU:CD2	1:D:167:THR:HG21	2.47	0.44
1:C:97:ARG:HA	1:C:100:ILE:HD12	1.99	0.43
1:B:247:PHE:HA	1:B:317:GLU:CD	2.38	0.43
1:D:253:LEU:HD21	1:D:310:LEU:HD21	1.99	0.43
1:D:79:ARG:NH1	1:D:81:ARG:O	2.51	0.43
1:C:195:ARG:O	1:C:209:HIS:HA	2.19	0.43
1:A:70:MET:HG3	1:A:184:LEU:HD21	2.00	0.43
1:B:97:ARG:CG	1:B:164:SER:HB3	2.47	0.43
1:A:221:LYS:HG3	1:A:268:VAL:HB	2.00	0.43
1:C:118:ARG:HB2	1:C:126:ARG:HG2	2.00	0.43
1:A:103:GLU:CG	1:A:141:LEU:HD11	2.48	0.43
1:B:154:LEU:HD11	1:B:172:LYS:HZ1	1.83	0.43
1:D:342:ASP:OD1	1:D:342:ASP:N	2.41	0.43
1:A:88:TRP:CD2	1:A:181:ARG:HD3	2.54	0.43
1:C:257:CYS:SG	1:C:288:MET:HG2	2.59	0.43
1:B:238:ASN:ND2	1:B:240:TYR:HB2	2.34	0.42
1:C:242:PRO:O	1:C:245:THR:HG23	2.19	0.42
1:C:233:MET:CE	1:C:310:LEU:CD1	2.86	0.42
1:D:309:ASN:N	1:D:309:ASN:ND2	2.64	0.42
1:A:296:PHE:HE2	1:A:332:ALA:HB2	1.82	0.42
1:C:283:ARG:NH2	1:C:307:SER:HB3	2.34	0.42
1:B:148:MET:HG2	1:B:159:TRP:CZ2	2.54	0.42
1:A:109:LYS:HB3	1:A:145:LEU:HD11	2.01	0.42
1:C:109:LYS:HA	1:C:132:ASP:HB3	2.01	0.42
1:C:127:LEU:HA	1:C:155:SER:HA	2.02	0.42
1:B:79:ARG:HB2	1:B:82:GLU:O	2.19	0.42
1:D:100:ILE:HG23	1:D:141:LEU:HD23	2.01	0.42
1:A:136:THR:HG22	1:A:142:ASP:OD2	2.19	0.42
1:A:68:THR:HB	1:A:70:MET:CE	2.49	0.42
1:B:112:ILE:HG13	1:B:326:LEU:O	2.20	0.42
1:C:294:TYR:CG	1:C:295:PRO:HA	2.54	0.42
1:C:70:MET:H	1:C:70:MET:CE	2.33	0.42
1:D:247:PHE:HA	1:D:317:GLU:OE1	2.20	0.42
1:A:111:ARG:HD2	1:A:131:ILE:CG2	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:288:MET:SD	1:B:339:SER:HB3	2.60	0.42
1:B:94:TYR:CD2	1:B:341:MET:CE	2.94	0.41
1:C:119:LYS:HA	1:C:126:ARG:NH2	2.35	0.41
1:C:130:VAL:HG21	1:C:149:LYS:HA	2.02	0.41
1:C:249:ILE:O	1:C:253:LEU:HG	2.20	0.41
1:C:248:LYS:O	1:C:252:THR:HG23	2.20	0.41
1:A:293:GLN:OE1	1:A:297:ASP:HB2	2.21	0.41
1:A:296:PHE:CE2	1:A:332:ALA:HB2	2.55	0.41
1:A:152:GLU:O	1:A:154:LEU:HD13	2.20	0.41
1:A:140:LEU:O	1:A:144:THR:HG22	2.20	0.41
1:A:166:GLU:HA	1:A:166:GLU:OE1	2.20	0.41
1:B:128:ILE:HD13	1:B:156:ILE:HG22	2.01	0.41
1:B:306:VAL:O	1:B:307:SER:HB3	2.21	0.41
1:D:73:VAL:HG12	1:D:96:LEU:HD21	2.02	0.41
1:A:223:ARG:HG3	1:A:223:ARG:HH11	1.85	0.41
1:A:242:PRO:C	1:A:244:THR:N	2.74	0.41
1:C:331:ILE:HD12	1:C:335:PHE:CE2	2.55	0.41
1:D:233:MET:CG	1:D:235:LEU:HD13	2.51	0.41
1:C:141:LEU:O	1:C:144:THR:HG23	2.21	0.41
1:C:338:PHE:O	1:C:341:MET:HG2	2.21	0.41
1:B:221:LYS:O	1:B:225:LEU:HB2	2.20	0.41
1:B:72:GLU:O	1:B:76:MET:HG2	2.21	0.41
1:C:125:GLU:HA	1:C:125:GLU:OE1	2.21	0.41
1:D:127:LEU:HA	1:D:155:SER:HA	2.02	0.41
1:C:182:GLU:O	1:C:185:ALA:HB3	2.21	0.41
1:C:104:LEU:HB3	1:C:110:ILE:HG12	2.03	0.40
1:C:233:MET:HE3	1:C:310:LEU:HD13	1.97	0.40
1:D:309:ASN:N	1:D:309:ASN:HD22	1.94	0.40
1:A:237:TYR:N	1:A:237:TYR:CD1	2.90	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	279/288 (97%)	258 (92%)	21 (8%)	0	100	100
1	B	279/288 (97%)	250 (90%)	28 (10%)	1 (0%)	34	64
1	C	277/288 (96%)	253 (91%)	23 (8%)	1 (0%)	34	64
1	D	281/288 (98%)	256 (91%)	25 (9%)	0	100	100
All	All	1116/1152 (97%)	1017 (91%)	97 (9%)	2 (0%)	47	77

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	270	THR
1	C	173	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	226/260 (87%)	197 (87%)	29 (13%)	4	16
1	B	219/260 (84%)	188 (86%)	31 (14%)	3	13
1	C	221/260 (85%)	186 (84%)	35 (16%)	2	9
1	D	233/260 (90%)	199 (85%)	34 (15%)	3	12
All	All	899/1040 (86%)	770 (86%)	129 (14%)	3	12

All (129) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	LEU
1	A	68	THR
1	A	70	MET
1	A	79	ARG
1	A	92	ILE
1	A	111	ARG
1	A	136	THR
1	A	144	THR
1	A	150	ASN

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Mol	Chain	Res	Type
1	A	152	GLU
1	A	167	THR
1	A	169	ASN
1	A	183	ARG
1	A	195	ARG
1	A	196	THR
1	A	203	LEU
1	A	211	ILE
1	A	226	SER
1	A	228	LEU
1	A	229	VAL
1	A	244	THR
1	A	275	GLU
1	A	282	SER
1	A	293	GLN
1	A	308	VAL
1	A	309	ASN
1	A	321	ASN
1	A	325	ASP
1	A	340	ARG
1	B	66	THR
1	B	70	MET
1	B	78	LEU
1	B	85	LEU
1	B	92	ILE
1	B	106	LEU
1	B	112	ILE
1	B	118	ARG
1	B	132	ASP
1	B	134	SER
1	B	139	VAL
1	B	144	THR
1	B	145	LEU
1	B	156	ILE
1	B	175	TYR
1	B	181	ARG
1	B	182	GLU
1	B	196	THR
1	B	211	ILE
1	B	217	LYS
1	B	225	LEU
1	B	227	VAL

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Mol	Chain	Res	Type
1	B	232	ASN
1	B	246	SER
1	B	249	ILE
1	B	272	LEU
1	B	293	GLN
1	B	309	ASN
1	B	310	LEU
1	B	326	LEU
1	B	336	GLU
1	C	70	MET
1	C	74	LEU
1	C	75	LEU
1	C	91	SER
1	C	93	SER
1	C	106	LEU
1	C	111	ARG
1	C	118	ARG
1	C	128	ILE
1	C	129	GLU
1	C	131	ILE
1	C	136	THR
1	C	144	THR
1	C	145	LEU
1	C	158	ASN
1	C	166	GLU
1	C	169	ASN
1	C	170	LEU
1	C	181	ARG
1	C	183	ARG
1	C	208	THR
1	C	211	ILE
1	C	228	LEU
1	C	252	THR
1	C	265	LEU
1	C	285	GLU
1	C	287	ILE
1	C	288	MET
1	C	298	LEU
1	C	306	VAL
1	C	307	SER
1	C	308	VAL
1	C	309	ASN

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Mol	Chain	Res	Type
1	C	331	ILE
1	C	340	ARG
1	D	65	PRO
1	D	73	VAL
1	D	75	LEU
1	D	81	ARG
1	D	101	ILE
1	D	112	ILE
1	D	139	VAL
1	D	144	THR
1	D	147	LEU
1	D	154	LEU
1	D	167	THR
1	D	169	ASN
1	D	170	LEU
1	D	174	ASN
1	D	183	ARG
1	D	194	LEU
1	D	208	THR
1	D	210	PRO
1	D	228	LEU
1	D	229	VAL
1	D	231	ARG
1	D	232	ASN
1	D	236	SER
1	D	255	LEU
1	D	268	VAL
1	D	288	MET
1	D	293	GLN
1	D	302	THR
1	D	309	ASN
1	D	310	LEU
1	D	321	ASN
1	D	325	ASP
1	D	342	ASP
1	D	343	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	169	ASN
1	A	176	GLN

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Mol	Chain	Res	Type
1	A	209	HIS
1	A	309	ASN
1	A	321	ASN
1	A	324	HIS
1	B	158	ASN
1	B	176	GLN
1	B	209	HIS
1	B	267	ASN
1	B	309	ASN
1	B	324	HIS
1	B	327	GLN
1	C	146	GLN
1	C	158	ASN
1	C	169	ASN
1	C	176	GLN
1	C	309	ASN
1	C	320	ASN
1	C	324	HIS
1	D	158	ASN
1	D	174	ASN
1	D	176	GLN
1	D	309	ASN
1	D	324	HIS

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 ⓘ

There are no ligands in this entry.

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 [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	281/288 (97%)	-0.47	0 100 100	25, 38, 53, 68	0
1	B	281/288 (97%)	-0.25	1 (0%) 92 82	32, 49, 65, 70	0
1	C	279/288 (96%)	-0.30	1 (0%) 92 82	31, 46, 63, 66	0
1	D	283/288 (98%)	-0.37	0 100 100	22, 40, 56, 66	0
All	All	1124/1152 (97%)	-0.35	2 (0%) 95 89	22, 44, 62, 70	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	302	THR	2.4
1	C	307	SER	2.2

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.