Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:39 PM GMT

PDB ID	:	1BRM
Title	:	ASPARTATE BETA-SEMIALDEHYDE DEHYDROGENASE FROM ESCHERICHIA COLI
Authors	:	Hadfield, A.T.; Kryger, G.; Ouyang, J.; Ringe, D.; Petsko, G.A.; Viola, R.E.
Deposited on	:	1998-08-24
Resolution	:	2.50 Å(reported)

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 (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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.
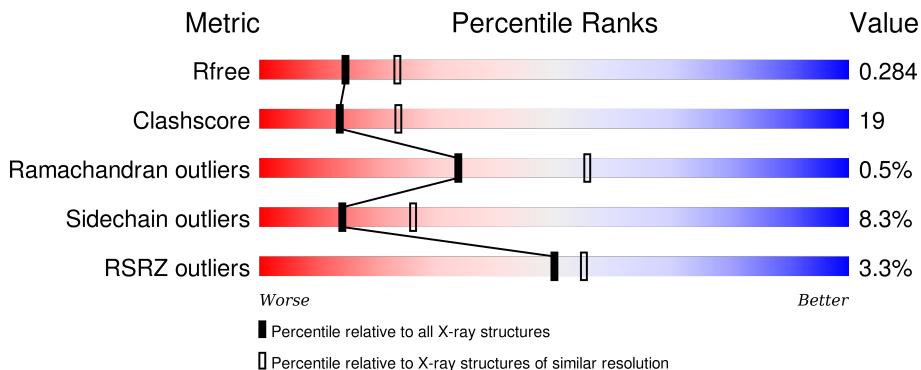
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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. 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 <=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	367	
1	B	367	
1	C	367	
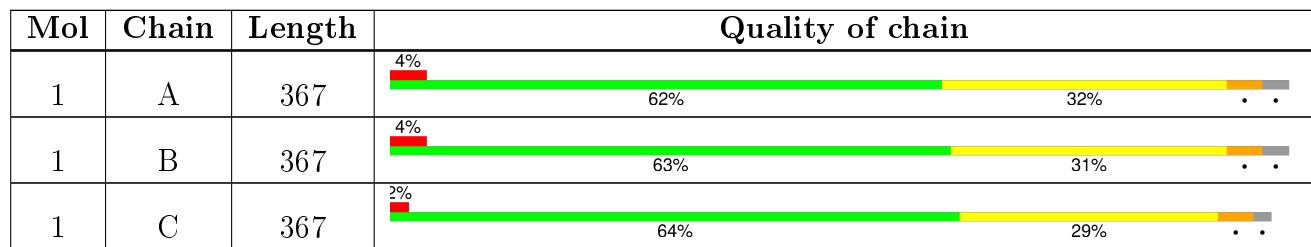
2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 8499 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 ASPARTATE-SEMIALDEHYDE DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	356	Total 2720	C 1722	N 472	O 510	S 16	0	0	0
1	B	357	Total 2729	C 1727	N 473	O 513	S 16	0	0	0
1	C	359	Total 2747	C 1738	N 476	O 517	S 16	0	0	0

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	73	Total 73	O 73	0	0
2	B	113	Total 113	O 113	0	0
2	C	117	Total 117	O 117	0	0

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: ASPARTATE-SEMIALDEHYDE DEHYDROGENASE
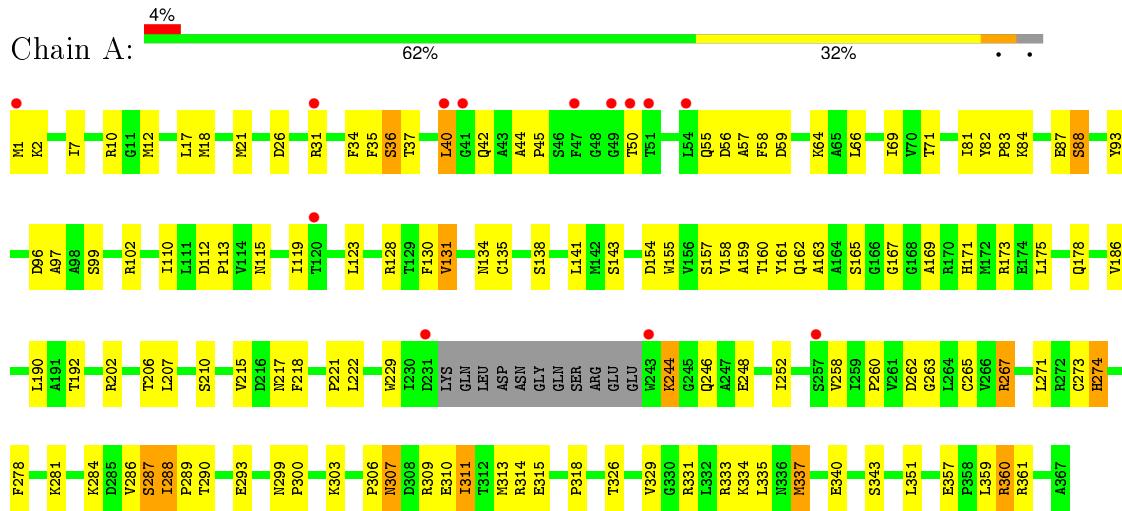
- Molecule 1: ASPARTATE-SEMIALDEHYDE DEHYDROGENASE
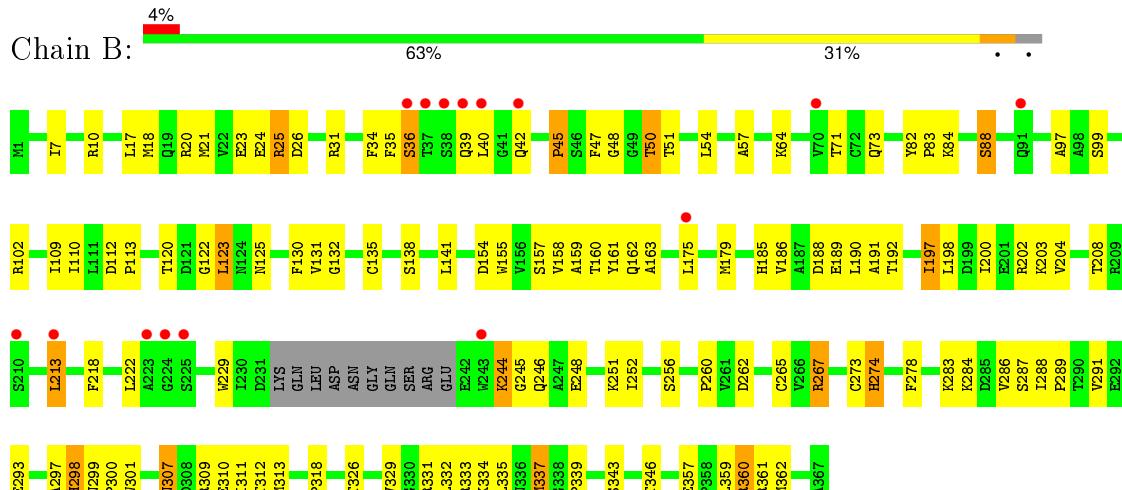
- Molecule 1: ASPARTATE-SEMIALDEHYDE DEHYDROGENASE
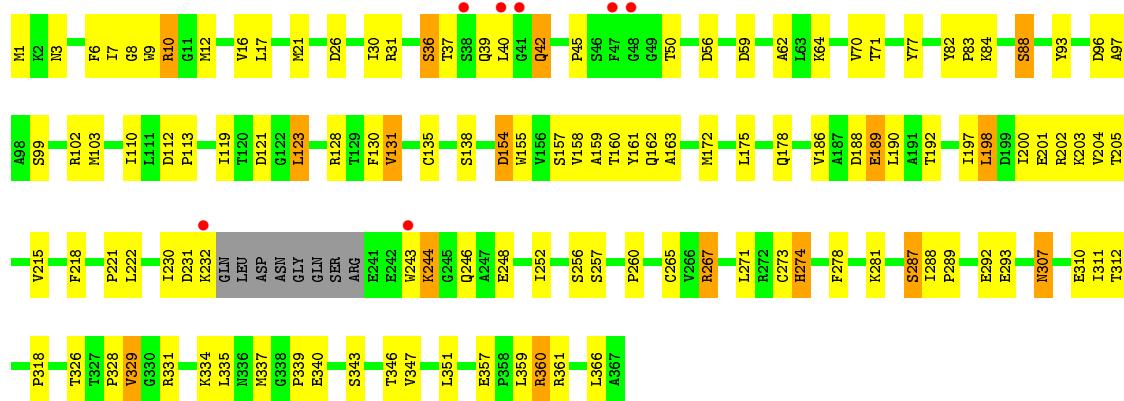
4 Data and refinement statistics (i)

| Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	174.89Å 117.05Å 58.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.00 – 2.50 26.67 – 2.50	Depositor EDS
% Data completeness (in resolution range)	82.6 (28.00-2.50) 82.6 (26.67-2.50)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
$<I/\sigma(I)>$ [1]	4.36 (at 2.50Å)	Xtriage
Refinement program	X-PLOR 3.85	Depositor
R, R_{free}	0.225 , 0.294 0.216 , 0.284	Depositor DCC
R_{free} test set	3466 reflections (9.96%)	DCC
Wilson B-factor (Å^2)	39.9	Xtriage
Anisotropy	0.460	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 57.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$<|L|> = 0.48, <L^2> = 0.31$	Xtriage
Outliers	0 of 34800 reflections	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	8499	wwPDB-VP
Average B, all atoms (Å^2)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.48% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.

[2]Theoretical values of $<|L|>$, $<L^2>$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

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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	0.32	0/2772	0.58	0/3764
1	B	0.32	0/2781	0.59	0/3776
1	C	0.31	0/2799	0.58	0/3799
All	All	0.32	0/8352	0.59	0/11339

There are no bond length outliers.

There are no bond angle outliers.

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	2720	0	2737	104	0
1	B	2729	0	2743	104	0
1	C	2747	0	2762	105	0
2	A	73	0	0	5	0
2	B	113	0	0	6	0
2	C	117	0	0	5	0
All	All	8499	0	8242	305	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 19.

All (305) 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:39:GLN:HB3	1:B:42:GLN:HG3	1.32	1.10
1:C:40:LEU:HD22	1:C:56:ASP:HA	1.44	0.99
1:C:56:ASP:HB3	1:C:59:ASP:HB2	1.44	0.97
1:C:289:PRO:O	1:C:293:GLU:HG3	1.70	0.92
1:C:39:GLN:HB3	1:C:42:GLN:HG3	1.50	0.90
1:B:186:VAL:HG12	1:B:190:LEU:HG	1.54	0.90
1:A:169:ALA:HB1	1:B:197:ILE:HG21	1.52	0.90
1:C:40:LEU:CD2	1:C:56:ASP:HA	2.11	0.80
1:A:96:ASP:HB3	1:A:131:VAL:HG13	1.63	0.78
1:B:208:THR:HA	1:B:213:LEU:HD22	1.66	0.77
1:B:7:ILE:HB	1:B:71:THR:HG22	1.68	0.75
1:C:218:PHE:HE1	1:C:267:ARG:HD3	1.52	0.75
1:B:218:PHE:HE1	1:B:267:ARG:HD3	1.51	0.74
1:A:218:PHE:HE1	1:A:267:ARG:HD3	1.50	0.74
1:B:218:PHE:CE1	1:B:267:ARG:HD3	2.22	0.74
1:A:218:PHE:CE1	1:A:267:ARG:HD3	2.23	0.74
1:A:306:PRO:HD2	1:A:311:ILE:HD13	1.70	0.73
1:B:186:VAL:HG11	1:B:200:ILE:HG23	1.71	0.73
1:C:218:PHE:CE1	1:C:267:ARG:HD3	2.24	0.73
1:C:232:LYS:HB2	1:C:243:TRP:HZ3	1.52	0.72
1:A:289:PRO:O	1:A:293:GLU:HG3	1.90	0.72
1:A:335:LEU:HD21	1:A:343:SER:HB3	1.73	0.71
1:C:232:LYS:HB2	1:C:243:TRP:CZ3	2.26	0.71
1:C:335:LEU:HD21	1:C:343:SER:HB3	1.73	0.70
1:C:175:LEU:HD13	1:C:222:LEU:HD23	1.74	0.70
1:C:7:ILE:HB	1:C:71:THR:HG22	1.74	0.69
1:B:335:LEU:HD21	1:B:343:SER:HB3	1.74	0.69
1:B:24:GLU:O	1:B:25:ARG:HG2	1.92	0.69
1:B:245:GLY:HA2	2:B:804:HOH:O	1.93	0.68
1:A:134:ASN:OD1	2:A:875:HOH:O	2.11	0.68
1:C:40:LEU:HD22	1:C:56:ASP:CA	2.22	0.68
1:A:7:ILE:HB	1:A:71:THR:HG22	1.76	0.67
1:C:288:ILE:O	1:C:292:GLU:HG3	1.94	0.67
1:B:158:VAL:HG22	1:B:278:PHE:CD1	2.30	0.67
1:C:188:ASP:HB2	1:C:189:GLU:OE1	1.94	0.66
1:A:158:VAL:HG22	1:A:278:PHE:CD1	2.30	0.66
1:C:186:VAL:O	1:C:186:VAL:HG12	1.93	0.66
1:B:186:VAL:O	1:B:186:VAL:HG12	1.95	0.66
1:C:158:VAL:HG22	1:C:278:PHE:CD1	2.31	0.65
1:A:44:ALA:HB1	1:A:45:PRO:HD2	1.78	0.65
1:A:56:ASP:HB3	1:A:59:ASP:HB2	1.77	0.65
1:A:202:ARG:O	1:A:206:THR:HG23	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:307:ASN:HD21	1:C:331:ARG:HD2	1.63	0.64
1:C:287:SER:OG	1:C:289:PRO:HD2	1.98	0.64
1:A:340:GLU:HB2	2:A:533:HOH:O	1.96	0.64
1:C:12:MET:HE2	1:C:351:LEU:HD21	1.80	0.64
1:C:112:ASP:CG	1:C:113:PRO:HD3	2.18	0.63
1:C:10:ARG:NH1	1:C:36:SER:HB2	2.12	0.63
1:B:307:ASN:HD21	1:B:331:ARG:HD2	1.64	0.63
1:C:3:ASN:OD1	1:C:31:ARG:HD2	1.99	0.62
1:A:307:ASN:ND2	1:A:331:ARG:HA	2.14	0.62
1:B:39:GLN:HB3	1:B:42:GLN:CG	2.21	0.62
1:B:307:ASN:ND2	1:B:331:ARG:HA	2.14	0.62
1:C:186:VAL:O	1:C:186:VAL:CG1	2.47	0.61
1:C:329:VAL:HG13	1:C:346:THR:HG22	1.80	0.61
1:A:169:ALA:O	1:A:173:ARG:HG3	2.01	0.61
1:C:307:ASN:ND2	1:C:331:ARG:HA	2.14	0.60
1:A:171:HIS:CE1	1:A:217:ASN:HB3	2.37	0.60
1:A:287:SER:OG	1:A:289:PRO:HD2	2.00	0.60
1:C:186:VAL:HG11	1:C:200:ILE:HB	1.84	0.59
1:A:303:LYS:NZ	1:A:315:GLU:HG2	2.17	0.59
1:A:83:PRO:O	1:A:87:GLU:HG2	2.03	0.59
1:C:186:VAL:HG13	1:C:189:GLU:HG2	1.83	0.59
1:B:267:ARG:HD2	2:B:776:HOH:O	2.01	0.59
1:A:307:ASN:HD21	1:A:331:ARG:HD2	1.66	0.59
1:A:1:MET:CE	1:A:31:ARG:HG2	2.33	0.59
1:A:1:MET:HE3	1:A:31:ARG:HG2	1.85	0.59
1:B:329:VAL:HG13	1:B:346:THR:HG22	1.84	0.59
1:B:45:PRO:HB2	1:B:47:PHE:CD2	2.38	0.59
1:B:7:ILE:HG22	1:B:73:GLN:HB2	1.85	0.58
1:C:7:ILE:HD13	1:C:77:TYR:OH	2.04	0.58
1:B:123:LEU:HD13	1:B:130:PHE:CZ	2.38	0.58
1:A:160:THR:OG1	1:A:265:CYS:HA	2.04	0.58
1:A:21:MET:HG2	1:A:26:ASP:HB2	1.86	0.58
1:B:186:VAL:CG1	1:B:186:VAL:O	2.52	0.57
1:C:12:MET:CE	1:C:351:LEU:HD11	2.35	0.57
1:A:99:SER:HA	1:A:102:ARG:HG3	1.86	0.57
1:A:12:MET:SD	1:B:197:ILE:HD13	2.44	0.57
1:B:112:ASP:CG	1:B:113:PRO:HD3	2.25	0.57
1:C:99:SER:HA	1:C:102:ARG:HG3	1.87	0.57
1:C:103:MET:HB2	2:C:851:HOH:O	2.04	0.57
1:B:283:LYS:O	1:B:284:LYS:HG3	2.05	0.56
1:B:99:SER:HA	1:B:102:ARG:HG3	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:132:GLY:HA3	2:B:601:HOH:O	2.05	0.56
1:C:112:ASP:OD1	1:C:113:PRO:HD3	2.05	0.56
1:C:244:LYS:HD3	1:C:265:CYS:SG	2.46	0.56
1:C:160:THR:OG1	1:C:265:CYS:HA	2.05	0.56
1:C:186:VAL:HG13	1:C:189:GLU:CG	2.36	0.55
1:B:112:ASP:OD1	1:B:113:PRO:HD3	2.06	0.55
1:B:197:ILE:HD12	1:B:198:LEU:HD23	1.88	0.55
1:A:96:ASP:HB3	1:A:131:VAL:CG1	2.35	0.55
1:C:288:ILE:HD11	1:C:334:LYS:HG2	1.88	0.55
1:A:309:ARG:HG3	1:B:229:TRP:NE1	2.22	0.55
1:A:248:GLU:HG3	2:A:814:HOH:O	2.07	0.54
1:C:172:MET:HB3	1:C:271:LEU:HD22	1.90	0.54
1:B:309:ARG:HG2	1:B:313:MET:HE3	1.88	0.54
1:A:35:PHE:CZ	1:A:66:LEU:HD11	2.42	0.54
1:C:158:VAL:HG22	1:C:278:PHE:HD1	1.70	0.54
1:A:158:VAL:HG22	1:A:278:PHE:HD1	1.71	0.54
1:B:160:THR:OG1	1:B:265:CYS:HA	2.08	0.54
1:C:7:ILE:HD12	1:C:71:THR:CG2	2.37	0.53
1:B:360:ARG:HH12	1:B:361:ARG:HD2	1.74	0.53
1:C:178:GLN:NE2	1:C:215:VAL:HG12	2.23	0.53
1:A:112:ASP:CG	1:A:113:PRO:HD3	2.29	0.53
1:A:360:ARG:HH12	1:A:361:ARG:HD2	1.73	0.53
1:A:37:THR:HG21	1:A:58:PHE:HE2	1.73	0.53
1:B:39:GLN:CB	1:B:42:GLN:HG3	2.23	0.52
1:C:307:ASN:ND2	1:C:331:ARG:HD2	2.24	0.52
1:B:267:ARG:NH2	2:B:607:HOH:O	2.42	0.52
1:C:7:ILE:HD13	1:C:77:TYR:CZ	2.44	0.52
1:C:12:MET:HE2	1:C:351:LEU:HD11	1.91	0.52
1:C:97:ALA:HB2	1:C:359:LEU:HD11	1.91	0.52
1:A:36:SER:OG	1:A:36:SER:O	2.26	0.52
1:C:360:ARG:HH12	1:C:361:ARG:HD2	1.74	0.52
1:B:123:LEU:HD13	1:B:130:PHE:HZ	1.74	0.52
1:B:162:GLN:HE22	1:B:265:CYS:HB3	1.72	0.52
1:B:97:ALA:HB2	1:B:359:LEU:HD11	1.90	0.52
1:C:288:ILE:CD1	1:C:334:LYS:HG2	2.40	0.51
1:C:232:LYS:CB	1:C:243:TRP:HZ3	2.19	0.51
1:A:229:TRP:NE1	1:B:309:ARG:HG3	2.24	0.51
1:C:39:GLN:CB	1:C:42:GLN:HG3	2.31	0.51
1:C:186:VAL:HG22	1:C:203:LYS:HD3	1.90	0.51
1:B:175:LEU:O	1:B:179:MET:HG3	2.10	0.51
1:B:35:PHE:HB3	1:B:57:ALA:HB2	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:64:LYS:HE2	1:B:88:SER:OG	2.10	0.51
1:B:158:VAL:HG22	1:B:278:PHE:HD1	1.71	0.51
1:A:64:LYS:HE2	1:A:88:SER:OG	2.10	0.51
1:B:10:ARG:HH12	1:B:36:SER:HB2	1.74	0.51
1:A:93:TYR:CE1	1:A:128:ARG:HD3	2.45	0.51
1:C:162:GLN:HE22	1:C:265:CYS:HB3	1.75	0.51
1:B:186:VAL:HG13	1:B:189:GLU:CG	2.41	0.50
1:B:186:VAL:O	1:B:190:LEU:HG	2.11	0.50
1:A:306:PRO:CD	1:A:311:ILE:HD13	2.41	0.50
1:B:185:HIS:O	1:B:203:LYS:HE2	2.10	0.50
1:B:288:ILE:N	1:B:289:PRO:CD	2.75	0.50
1:B:307:ASN:ND2	1:B:331:ARG:HD2	2.26	0.50
1:A:97:ALA:HB2	1:A:359:LEU:HD11	1.92	0.50
1:B:309:ARG:O	1:B:313:MET:HG3	2.11	0.50
1:C:200:ILE:O	1:C:204:VAL:HG13	2.12	0.50
1:A:2:LYS:O	1:A:31:ARG:HG3	2.12	0.50
1:A:17:LEU:HG	1:A:21:MET:HE3	1.93	0.50
1:A:143:SER:HB3	1:A:329:VAL:HG22	1.93	0.50
1:C:6:PHE:CD1	1:C:70:VAL:HB	2.47	0.49
1:A:267:ARG:NH2	2:A:751:HOH:O	2.44	0.49
1:A:167:GLY:HA3	1:A:171:HIS:CD2	2.47	0.49
1:A:84:LYS:O	1:A:88:SER:HB3	2.12	0.49
1:A:307:ASN:ND2	1:A:331:ARG:HD2	2.26	0.49
1:A:21:MET:HG2	1:A:26:ASP:CB	2.42	0.49
1:C:310:GLU:H	1:C:310:GLU:CD	2.16	0.49
1:A:288:ILE:HD11	1:A:334:LYS:HG2	1.94	0.49
1:C:7:ILE:HD12	1:C:71:THR:HG22	1.95	0.49
1:C:288:ILE:N	1:C:289:PRO:CD	2.76	0.49
1:C:158:VAL:HG12	1:C:159:ALA:N	2.28	0.49
1:A:299:ASN:HB2	1:A:300:PRO:HD2	1.94	0.49
1:C:64:LYS:HE2	1:C:88:SER:OG	2.13	0.49
1:C:84:LYS:O	1:C:88:SER:HB3	2.13	0.49
1:B:311:ILE:HG23	1:B:312:THR:N	2.28	0.49
1:A:244:LYS:HE3	1:A:262:ASP:OD1	2.13	0.49
1:C:9:TRP:CH2	1:C:45:PRO:HG3	2.48	0.48
1:A:162:GLN:HE22	1:A:265:CYS:HB3	1.77	0.48
1:A:310:GLU:CD	1:A:310:GLU:H	2.17	0.48
1:C:36:SER:O	1:C:40:LEU:CD2	2.62	0.48
1:C:59:ASP:HB3	1:C:62:ALA:HB3	1.96	0.48
1:B:287:SER:O	1:B:291:VAL:HG23	2.13	0.48
1:A:207:LEU:HD12	1:A:210:SER:OG	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:PHE:CE2	1:A:45:PRO:HG2	2.48	0.48
1:B:244:LYS:HD3	1:B:265:CYS:SG	2.53	0.48
1:C:17:LEU:O	1:C:21:MET:HG3	2.14	0.48
1:B:155:TRP:HZ3	1:B:157:SER:HB2	1.79	0.48
1:C:155:TRP:HZ3	1:C:157:SER:HB2	1.79	0.48
1:B:175:LEU:HD13	1:B:222:LEU:HD23	1.95	0.47
1:B:310:GLU:CD	1:B:310:GLU:H	2.16	0.47
1:A:158:VAL:HG12	1:A:159:ALA:N	2.29	0.47
1:A:163:ALA:HB2	1:A:273:CYS:O	2.14	0.47
1:C:6:PHE:HD1	1:C:70:VAL:HB	1.79	0.47
1:A:155:TRP:HZ3	1:A:157:SER:HB2	1.79	0.47
1:A:309:ARG:O	1:A:313:MET:HG3	2.14	0.47
1:B:18:MET:HE3	1:B:47:PHE:CD2	2.49	0.47
1:B:162:GLN:NE2	1:B:265:CYS:HB3	2.30	0.47
1:C:163:ALA:HB2	1:C:273:CYS:O	2.15	0.47
1:B:158:VAL:HG12	1:B:159:ALA:N	2.29	0.47
1:A:155:TRP:NE1	1:A:281:LYS:HD3	2.30	0.47
1:C:162:GLN:NE2	1:C:265:CYS:HB3	2.30	0.47
1:A:35:PHE:HZ	1:A:66:LEU:HD11	1.78	0.47
1:B:84:LYS:O	1:B:88:SER:HB3	2.14	0.47
1:A:246:GLN:OE1	1:A:260:PRO:HA	2.15	0.47
1:B:186:VAL:HG13	1:B:189:GLU:HG3	1.96	0.47
1:B:248:GLU:O	1:B:252:ILE:HG13	2.15	0.47
1:C:197:ILE:HG23	1:C:198:LEU:HD23	1.97	0.46
1:A:248:GLU:O	1:A:252:ILE:HG13	2.15	0.46
1:B:246:GLN:OE1	1:B:260:PRO:HA	2.15	0.46
1:B:284:LYS:O	1:B:286:VAL:HG23	2.15	0.46
1:C:328:PRO:HD2	1:C:347:VAL:O	2.15	0.46
1:B:189:GLU:H	1:B:189:GLU:CD	2.16	0.46
1:A:34:PHE:CD2	1:A:45:PRO:HD2	2.50	0.46
1:C:123:LEU:HD21	1:C:366:LEU:CD2	2.46	0.46
1:A:178:GLN:NE2	1:A:215:VAL:HG12	2.31	0.46
1:A:12:MET:HE3	1:A:165:SER:HB2	1.98	0.46
1:A:178:GLN:HE22	1:A:221:PRO:HA	1.81	0.46
1:B:300:PRO:HD2	1:B:301:TRP:CE3	2.50	0.46
1:C:96:ASP:HB3	1:C:131:VAL:HG13	1.97	0.46
1:A:186:VAL:HG23	1:A:190:LEU:HD12	1.98	0.46
1:A:58:PHE:HD1	1:A:81:ILE:HD11	1.82	0.45
1:B:109:ILE:HG13	1:B:251:LYS:HE2	1.98	0.45
1:A:82:TYR:HB3	1:A:83:PRO:HD3	1.99	0.45
1:C:246:GLN:OE1	1:C:260:PRO:HA	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:267:ARG:NH2	2:C:508:HOH:O	2.50	0.45
1:A:162:GLN:NE2	1:A:265:CYS:HB3	2.32	0.45
1:B:163:ALA:HB2	1:B:273:CYS:O	2.17	0.45
1:A:262:ASP:OD2	1:B:333:ARG:NH2	2.50	0.45
1:C:334:LYS:HD2	1:C:339:PRO:O	2.17	0.45
1:A:161:TYR:CD1	1:A:161:TYR:N	2.85	0.45
1:C:12:MET:CE	1:C:351:LEU:HD21	2.45	0.44
1:B:334:LYS:HD2	1:B:339:PRO:O	2.17	0.44
1:B:71:THR:HB	2:B:679:HOH:O	2.18	0.44
1:A:7:ILE:CD1	1:A:69:ILE:HG23	2.47	0.44
1:A:87:GLU:OE1	1:A:87:GLU:HA	2.18	0.44
1:B:82:TYR:HB3	1:B:83:PRO:HD3	2.00	0.44
1:C:40:LEU:HD13	1:C:56:ASP:CB	2.48	0.44
1:A:335:LEU:HB3	1:A:337:MET:HE2	2.00	0.44
1:C:12:MET:O	1:C:16:VAL:HG23	2.17	0.44
1:C:288:ILE:HD12	1:C:334:LYS:HE3	2.00	0.44
1:C:230:ILE:HG22	1:C:232:LYS:HG2	2.00	0.44
1:C:198:LEU:O	1:C:202:ARG:HB2	2.17	0.44
1:C:123:LEU:HD21	1:C:366:LEU:HD21	1.99	0.44
1:A:311:ILE:CG2	1:A:314:ARG:HH21	2.30	0.44
1:B:48:GLY:C	1:B:50:THR:N	2.71	0.44
1:C:311:ILE:HG23	1:C:312:THR:N	2.33	0.44
1:C:161:TYR:CD1	1:C:161:TYR:N	2.86	0.43
1:B:161:TYR:CD1	1:B:161:TYR:N	2.85	0.43
1:C:267:ARG:NH1	2:C:541:HOH:O	2.51	0.43
1:C:204:VAL:CG2	1:C:205:THR:N	2.81	0.43
1:C:12:MET:HE3	1:C:351:LEU:HD11	2.00	0.43
1:B:309:ARG:CG	1:B:313:MET:HE3	2.47	0.43
1:C:200:ILE:CG1	1:C:201:GLU:N	2.82	0.43
2:A:842:HOH:O	1:B:313:MET:HE2	2.18	0.43
1:A:35:PHE:CD2	1:A:55:GLN:HB2	2.53	0.43
1:B:273:CYS:HB2	1:B:318:PRO:HB3	2.01	0.43
1:A:36:SER:O	1:A:37:THR:C	2.56	0.43
1:A:273:CYS:HB2	1:A:318:PRO:HB3	2.00	0.43
1:A:333:ARG:NH2	1:B:262:ASP:OD2	2.51	0.43
1:C:248:GLU:O	1:C:252:ILE:HG13	2.18	0.43
1:B:39:GLN:O	1:B:54:LEU:HD12	2.19	0.43
1:B:18:MET:HE1	1:B:34:PHE:CE1	2.54	0.43
1:C:154:ASP:HB2	1:C:281:LYS:O	2.18	0.43
1:B:332:LEU:HB2	2:B:540:HOH:O	2.17	0.43
1:B:110:ILE:HG21	1:B:130:PHE:HB3	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:357:GLU:O	1:C:361:ARG:CG	2.67	0.43
1:C:1:MET:N	2:C:673:HOH:O	2.47	0.43
1:B:357:GLU:O	1:B:361:ARG:CG	2.68	0.42
1:C:119:ILE:HG22	1:C:123:LEU:HD12	2.01	0.42
1:B:50:THR:HG23	1:B:51:THR:N	2.34	0.42
1:A:271:LEU:CD1	1:B:204:VAL:HG21	2.49	0.42
1:A:175:LEU:HD13	1:A:222:LEU:HD23	2.00	0.42
1:B:18:MET:HE1	1:B:34:PHE:HE1	1.84	0.42
1:C:244:LYS:CD	1:C:265:CYS:SG	3.07	0.42
1:A:37:THR:CG2	1:A:58:PHE:HE2	2.31	0.42
1:C:82:TYR:HB3	1:C:83:PRO:HD3	2.00	0.42
1:C:110:ILE:CG2	1:C:130:PHE:HB3	2.49	0.42
1:B:300:PRO:HD2	1:B:301:TRP:CZ3	2.55	0.42
1:A:141:LEU:HA	1:A:141:LEU:HD23	1.90	0.42
1:C:340:GLU:HB2	2:C:795:HOH:O	2.19	0.42
1:A:288:ILE:N	1:A:289:PRO:CD	2.83	0.42
1:A:110:ILE:HG22	1:A:119:ILE:HD11	2.01	0.42
1:A:286:VAL:CG1	1:A:290:THR:HG21	2.50	0.42
1:C:110:ILE:HG21	1:C:130:PHE:HB3	2.02	0.41
1:A:258:VAL:O	1:A:260:PRO:HD3	2.21	0.41
1:B:109:ILE:HA	1:B:131:VAL:O	2.20	0.41
1:C:8:GLY:C	1:C:10:ARG:H	2.24	0.41
1:A:12:MET:SD	1:B:197:ILE:CD1	3.09	0.41
1:B:357:GLU:O	1:B:361:ARG:HG3	2.21	0.41
1:C:21:MET:HE3	1:C:26:ASP:HB3	2.01	0.41
1:A:163:ALA:H	1:A:274:HIS:CE1	2.39	0.41
1:B:122:GLY:HA2	1:B:125:ASN:HD22	1.85	0.41
1:B:188:ASP:O	1:B:191:ALA:HB3	2.21	0.41
1:B:20:ARG:O	1:B:23:GLU:HB3	2.20	0.41
1:A:357:GLU:O	1:A:361:ARG:CG	2.68	0.41
1:C:273:CYS:HB2	1:C:318:PRO:HB3	2.02	0.41
1:B:17:LEU:HD11	1:B:21:MET:HE2	2.03	0.41
1:A:12:MET:HE2	1:A:351:LEU:HD21	2.02	0.41
1:B:289:PRO:O	1:B:293:GLU:HG3	2.21	0.41
1:A:311:ILE:HG23	1:A:314:ARG:HH21	1.85	0.41
1:C:178:GLN:HE22	1:C:221:PRO:HA	1.86	0.41
1:A:37:THR:HG22	1:A:57:ALA:HB3	2.03	0.41
1:B:297:ALA:O	1:B:299:ASN:N	2.54	0.41
1:A:303:LYS:HZ2	1:A:315:GLU:HG2	1.86	0.41
1:B:130:PHE:CE1	1:B:362:MET:HE1	2.56	0.41
1:A:37:THR:HG21	1:A:58:PHE:CE2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:244:LYS:HE3	1:A:263:GLY:H	1.85	0.41
1:C:163:ALA:H	1:C:274:HIS:CE1	2.38	0.41
1:B:20:ARG:NH1	1:B:23:GLU:CD	2.74	0.41
1:A:115:ASN:ND2	1:A:252:ILE:HG23	2.35	0.41
1:C:178:GLN:HE22	1:C:215:VAL:HG12	1.85	0.41
1:B:20:ARG:HD2	1:B:20:ARG:HA	1.78	0.41
1:B:297:ALA:O	1:B:298:HIS:C	2.58	0.41
1:A:40:LEU:HG	1:A:40:LEU:H	1.66	0.41
1:C:93:TYR:CE1	1:C:128:ARG:HD3	2.56	0.41
1:B:110:ILE:CG2	1:B:130:PHE:HB3	2.50	0.40
1:B:163:ALA:H	1:B:274:HIS:CE1	2.39	0.40
1:C:186:VAL:HG12	1:C:190:LEU:HG	2.03	0.40
1:A:357:GLU:O	1:A:361:ARG:HG3	2.21	0.40
1:B:141:LEU:HA	1:B:141:LEU:HD23	1.91	0.40
1:A:12:MET:HE3	1:A:351:LEU:HD11	2.02	0.40
1:A:110:ILE:HG21	1:A:130:PHE:HB3	2.03	0.40
1:B:335:LEU:HB3	1:B:337:MET:HE2	2.04	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	352/367 (96%)	324 (92%)	27 (8%)	1 (0%)	46	68
1	B	353/367 (96%)	320 (91%)	31 (9%)	2 (1%)	30	50
1	C	355/367 (97%)	319 (90%)	34 (10%)	2 (1%)	30	50
All	All	1060/1101 (96%)	963 (91%)	92 (9%)	5 (0%)	34	55

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	42	GLN
1	B	298	HIS
1	C	42	GLN
1	C	231	ASP
1	B	45	PRO

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	292/302 (97%)	269 (92%)	23 (8%)	15	28
1	B	293/302 (97%)	269 (92%)	24 (8%)	14	27
1	C	295/302 (98%)	269 (91%)	26 (9%)	12	23
All	All	880/906 (97%)	807 (92%)	73 (8%)	14	26

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	ARG
1	A	18	MET
1	A	36	SER
1	A	40	LEU
1	A	50	THR
1	A	88	SER
1	A	123	LEU
1	A	131	VAL
1	A	135	CYS
1	A	138	SER
1	A	154	ASP
1	A	192	THR
1	A	244	LYS
1	A	267	ARG
1	A	274	HIS
1	A	284	LYS
1	A	287	SER
1	A	288	ILE

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Mol	Chain	Res	Type
1	A	307	ASN
1	A	311	ILE
1	A	326	THR
1	A	337	MET
1	A	360	ARG
1	B	25	ARG
1	B	26	ASP
1	B	31	ARG
1	B	36	SER
1	B	40	LEU
1	B	50	THR
1	B	88	SER
1	B	120	THR
1	B	123	LEU
1	B	135	CYS
1	B	138	SER
1	B	154	ASP
1	B	192	THR
1	B	197	ILE
1	B	202	ARG
1	B	213	LEU
1	B	244	LYS
1	B	256	SER
1	B	267	ARG
1	B	274	HIS
1	B	307	ASN
1	B	326	THR
1	B	337	MET
1	B	360	ARG
1	C	10	ARG
1	C	30	ILE
1	C	36	SER
1	C	37	THR
1	C	50	THR
1	C	88	SER
1	C	121	ASP
1	C	123	LEU
1	C	131	VAL
1	C	135	CYS
1	C	138	SER
1	C	154	ASP
1	C	189	GLU

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Mol	Chain	Res	Type
1	C	192	THR
1	C	198	LEU
1	C	244	LYS
1	C	256	SER
1	C	257	SER
1	C	267	ARG
1	C	274	HIS
1	C	287	SER
1	C	307	ASN
1	C	326	THR
1	C	329	VAL
1	C	337	MET
1	C	360	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	124	ASN
1	A	150	ASN
1	A	162	GLN
1	A	307	ASN
1	B	3	ASN
1	B	125	ASN
1	B	162	GLN
1	B	307	ASN
1	B	365	GLN
1	C	116	GLN
1	C	150	ASN
1	C	162	GLN
1	C	171	HIS
1	C	217	ASN
1	C	307	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

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	356/367 (97%)	0.21	13 (3%)	45	50	17, 40, 72, 95	0
1	B	357/367 (97%)	0.10	15 (4%)	40	45	13, 34, 64, 90	0
1	C	359/367 (97%)	-0.03	7 (1%)	70	73	13, 34, 69, 93	0
All	All	1072/1101 (97%)	0.09	35 (3%)	50	55	13, 36, 69, 95	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	243	TRP	6.9
1	A	1	MET	4.8
1	A	51	THR	4.7
1	A	243	TRP	4.6
1	C	243	TRP	4.4
1	B	40	LEU	4.2
1	A	40	LEU	3.7
1	A	41	GLY	3.7
1	C	41	GLY	3.6
1	B	38	SER	3.5
1	B	70	VAL	3.3
1	C	40	LEU	3.2
1	C	232	LYS	3.1
1	B	36	SER	3.0
1	A	47	PHE	2.9
1	B	223	ALA	2.8
1	B	224	GLY	2.7
1	A	54	LEU	2.7
1	A	231	ASP	2.6
1	A	50	THR	2.6
1	A	31	ARG	2.6
1	C	47	PHE	2.6
1	A	49	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	91	GLN	2.4
1	C	38	SER	2.4
1	B	225	SER	2.3
1	A	120	THR	2.3
1	C	48	GLY	2.2
1	B	39	GLN	2.2
1	B	37	THR	2.2
1	B	175	LEU	2.1
1	B	42	GLN	2.1
1	A	257	SER	2.1
1	B	213	LEU	2.1
1	B	210	SER	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)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.