Full wwPDB X-ray Structure Validation Report (i)

Jun 25, 2024 – 08:17 AM EDT

PDB ID	:	6EKI
Title	:	Structure of a hyperthermostable carbonic anhydrase identified from an active hydrothermal vent chimney
Authors	:	Fredslund, F.; Poulsen, J.-C.N.; Lo Leggio, L.
Deposited on	:	2017-09-26
Resolution	:	2.56 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 (i) 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 (1)) were used in the production of this report:

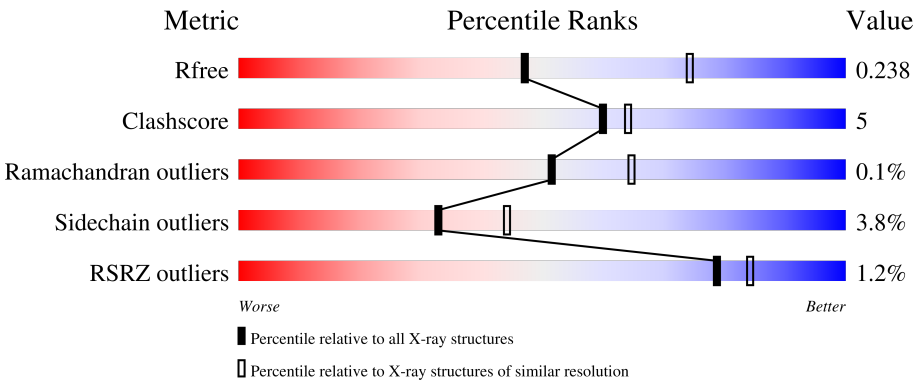
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 2.56 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	1279 (2.58-2.54)
Clashscore	141614	1327 (2.58-2.54)
Ramachandran outliers	138981	1312 (2.58-2.54)
Sidechain outliers	138945	1312 (2.58-2.54)
RSRZ outliers	127900	1269 (2.58-2.54)

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 <=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	226	.% 81% 15% ••
1	B	226	2% 82% 15% •
1	C	226	83% 14% ••
1	D	226	3% 86% 11% ••
1	E	226	81% 15% ••

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Mol	Chain	Length	Quality of chain
1	F	226	% 85% 13%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 21474 atoms, of which 10658 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 Carbonic anhydrase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	222	3568	1144	1779	307	329	9	0	0	0
1	B	222	3567	1144	1778	307	329	9	0	0	0
1	C	222	3568	1144	1779	307	329	9	0	0	0
1	D	222	3567	1144	1778	307	329	9	0	0	0
1	E	222	3561	1144	1772	307	329	9	0	0	0
1	F	222	3561	1144	1772	307	329	9	0	0	0

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
2	A	1	1	1	0	0
2	B	1	1	1	0	0
2	C	1	1	1	0	0
2	D	1	1	1	0	0
2	E	1	1	1	0	0
2	F	1	1	1	0	0

- Molecule 3 is water.

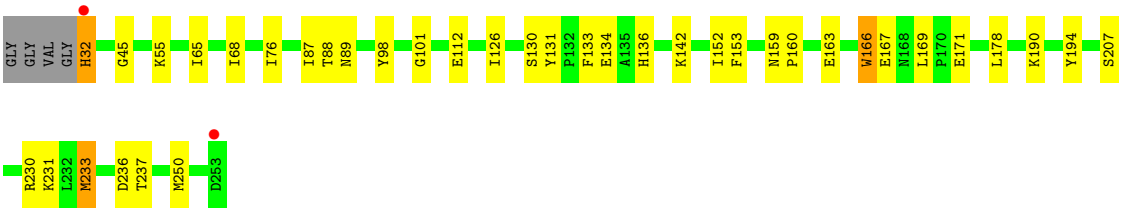
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	11	11	11	0	0
			Total	O		
3	B	14	14	14	0	0
			Total	O		
3	C	11	11	11	0	0
			Total	O		
3	D	16	16	16	0	0
			Total	O		
3	E	12	12	12	0	0
			Total	O		
3	F	12	12	12	0	0

3 Residue-property plots (i)

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.

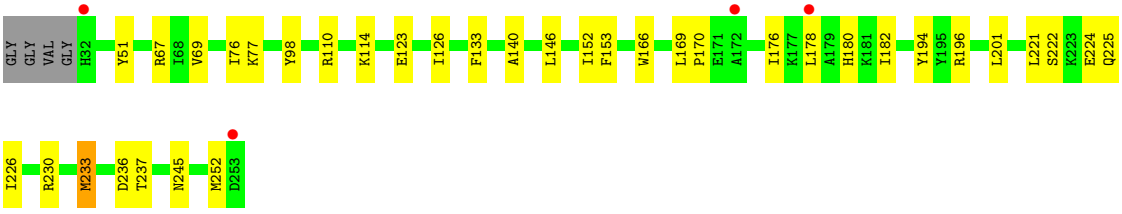
• Molecule 1: Carbonic anhydrase

Chain A:



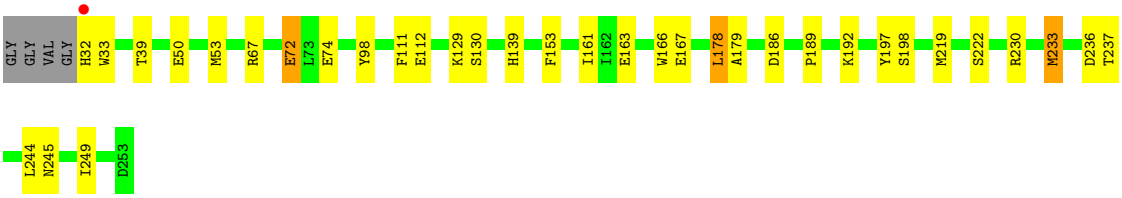
• Molecule 1: Carbonic anhydrase

Chain B:



• Molecule 1: Carbonic anhydrase

Chain C:

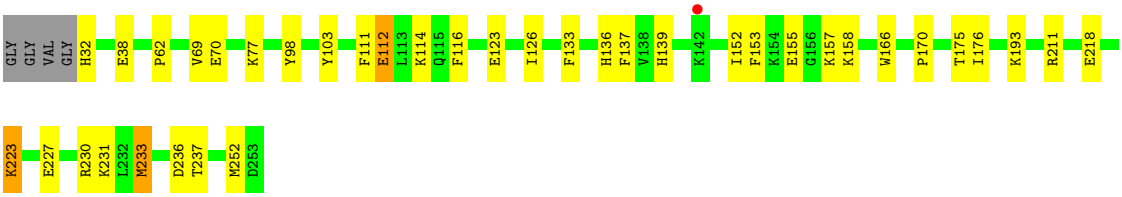


• Molecule 1: Carbonic anhydrase

Chain D:

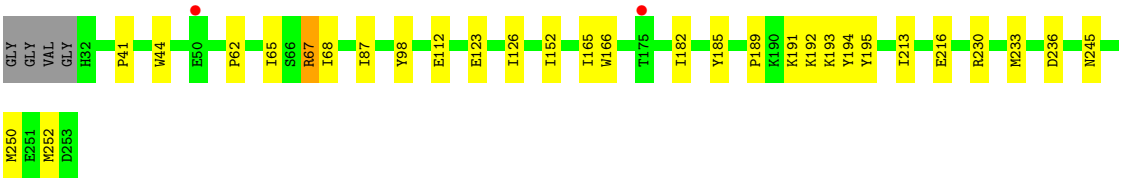
• Molecule 1: Carbonic anhydrase

Chain E:



• Molecule 1: Carbonic anhydrase

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	46.32Å 71.28Å 103.34Å 88.38° 82.84° 84.22°	Depositor
Resolution (Å)	39.99 – 2.56 39.99 – 2.56	Depositor EDS
% Data completeness (in resolution range)	96.3 (39.99-2.56) 96.3 (39.99-2.56)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.23 (at 2.54Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.164 , 0.236 0.169 , 0.238	Depositor DCC
R_{free} test set	2351 reflections (5.77%)	wwPDB-VP
Wilson B-factor (Å^2)	41.0	Xtriage
Anisotropy	0.222	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.40 , 44.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	21474	wwPDB-VP
Average B, all atoms (Å^2)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.63% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: ZN

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	0.74	0/1835	0.79	0/2471
1	B	0.73	0/1835	0.82	1/2471 (0.0%)
1	C	0.76	2/1835 (0.1%)	0.76	0/2471
1	D	0.73	0/1835	0.75	0/2471
1	E	0.75	1/1835 (0.1%)	0.80	1/2471 (0.0%)
1	F	0.68	0/1835	0.75	0/2471
All	All	0.73	3/11010 (0.0%)	0.78	2/14826 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	112	GLU	CG-CD	5.52	1.60	1.51
1	C	197	TYR	CD1-CE1	-5.43	1.31	1.39
1	C	72	GLU	CG-CD	5.39	1.60	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	E	211	ARG	NE-CZ-NH2	-5.36	117.62	120.30
1	B	196	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 (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	1789	1779	1778	19	0
1	B	1789	1778	1778	20	1
1	C	1789	1779	1778	15	0
1	D	1789	1778	1778	14	0
1	E	1789	1772	1778	17	1
1	F	1789	1772	1778	20	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	11	0	0	0	0
3	B	14	0	0	3	0
3	C	11	0	0	0	0
3	D	16	0	0	2	0
3	E	12	0	0	0	0
3	F	12	0	0	0	0
All	All	10816	10658	10668	99	1

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 5.

All (99) 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:D:230:ARG:NH2	1:D:236:ASP:OD1	1.86	1.08
1:B:230:ARG:NH2	1:B:236:ASP:OD1	2.12	0.83
1:F:193:LYS:NZ	1:F:252:MET:O	2.15	0.79
1:B:77:LYS:NZ	3:B:401:HOH:O	2.12	0.78
1:C:230:ARG:NH2	1:C:236:ASP:OD1	2.23	0.72
1:D:233:MET:HG2	1:D:237:THR:HG21	1.71	0.71
1:B:69:VAL:HG11	1:F:250:MET:HE1	1.74	0.69
1:B:230:ARG:NH1	3:B:403:HOH:O	2.25	0.69
1:D:42:GLN:OE1	3:D:401:HOH:O	2.10	0.69
1:F:189:PRO:O	1:F:192:LYS:NZ	2.26	0.69
1:A:230:ARG:NH2	1:A:236:ASP:OD1	2.26	0.68
1:B:222:SER:OG	1:B:224:GLU:O	2.11	0.67
1:D:211:ARG:NH2	3:D:403:HOH:O	2.31	0.64
1:E:112:GLU:OE1	1:E:114:LYS:NZ	2.31	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:67:ARG:NH2	1:C:245:ASN:O	2.33	0.60
1:A:233:MET:HG2	1:A:237:THR:HG21	1.83	0.60
1:B:67:ARG:NH2	1:B:245:ASN:O	2.34	0.60
1:D:193:LYS:HA	1:D:252:MET:HE2	1.83	0.60
1:B:123:GLU:OE2	3:B:402:HOH:O	2.16	0.58
1:B:178:LEU:O	1:B:180:HIS:N	2.35	0.57
1:B:69:VAL:CG1	1:F:250:MET:HE1	2.35	0.57
1:D:195:TYR:CE2	1:D:216:GLU:HG2	2.41	0.56
1:E:230:ARG:NH2	1:E:236:ASP:OD1	2.38	0.56
1:C:189:PRO:O	1:C:192:LYS:NZ	2.37	0.55
1:D:157:LYS:HG2	1:D:158:LYS:N	2.22	0.55
1:E:155:GLU:OE2	1:E:223:LYS:NZ	2.25	0.55
1:F:126:ILE:CD1	1:F:152:ILE:HD13	2.38	0.54
1:A:134:GLU:OE2	1:A:136:HIS:NE2	2.37	0.53
1:D:101:GLY:N	1:D:112:GLU:OE2	2.37	0.53
1:D:107:ASP:OD2	1:D:211:ARG:NH1	2.42	0.52
1:F:67:ARG:NH2	1:F:245:ASN:O	2.43	0.52
1:A:101:GLY:N	1:A:112:GLU:OE2	2.42	0.51
1:A:166:TRP:HA	1:A:169:LEU:HD13	1.92	0.51
1:B:76:ILE:HD13	1:B:194:TYR:CE1	2.46	0.51
1:C:233:MET:HG2	1:C:237:THR:HG21	1.93	0.51
1:C:72:GLU:HG2	1:E:70:GLU:H	1.76	0.50
1:E:77:LYS:NZ	1:E:103:TYR:OH	2.45	0.49
1:A:88:THR:HG22	1:A:89:ASN:N	2.27	0.49
1:D:193:LYS:HA	1:D:252:MET:CE	2.43	0.49
1:A:133:PHE:HB3	1:A:153:PHE:HB2	1.95	0.49
1:C:178:LEU:O	1:C:179:ALA:HB3	2.14	0.48
1:F:230:ARG:NH2	1:F:236:ASP:OD1	2.43	0.48
1:F:41:PRO:HA	1:F:44:TRP:CE2	2.48	0.48
1:C:111:PHE:HB3	1:C:139:HIS:HB3	1.95	0.47
1:E:227:GLU:O	1:E:231:LYS:HB2	2.14	0.47
1:B:233:MET:CG	1:B:237:THR:HG21	2.44	0.47
1:F:126:ILE:HD13	1:F:152:ILE:HD13	1.95	0.47
1:A:163:GLU:O	1:A:167:GLU:HG3	2.15	0.47
1:F:194:TYR:CD2	1:F:213:ILE:HG23	2.50	0.47
1:A:126:ILE:HD12	1:A:152:ILE:HD13	1.97	0.46
1:A:76:ILE:HD12	1:A:194:TYR:CE1	2.50	0.46
1:A:65:ILE:HG22	1:A:68:ILE:HD11	1.97	0.46
1:E:170:PRO:HB3	1:E:176:ILE:HG13	1.98	0.46
1:B:133:PHE:HB3	1:B:153:PHE:HB2	1.97	0.45
1:C:163:GLU:O	1:C:167:GLU:HG3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:ILE:HD11	1:A:178:LEU:HD11	1.99	0.45
1:F:165:ILE:CG1	1:F:182:ILE:HD13	2.47	0.45
1:D:70:GLU:HB3	1:D:253:ASP:OD1	2.16	0.44
1:A:142:LYS:HD3	1:F:191:LYS:HD3	1.98	0.44
1:B:76:ILE:CD1	1:B:194:TYR:CZ	3.00	0.44
1:A:130:SER:HB3	1:C:130:SER:HB3	1.98	0.44
1:D:175:THR:HG22	1:D:176:ILE:N	2.32	0.44
1:F:65:ILE:HG22	1:F:68:ILE:HD11	1.98	0.44
1:F:67:ARG:O	1:F:67:ARG:HG3	2.18	0.44
1:C:50:GLU:HG2	1:C:50:GLU:O	2.18	0.43
1:D:116:PHE:HA	1:D:136:HIS:O	2.18	0.43
1:C:33:TRP:HA	1:C:39:THR:HB	2.01	0.43
1:E:193:LYS:HA	1:E:252:MET:CE	2.49	0.43
1:E:157:LYS:HG2	1:E:158:LYS:O	2.19	0.43
1:A:32:HIS:ND1	1:A:32:HIS:C	2.72	0.43
1:C:244:LEU:HD21	1:C:249:ILE:HD11	2.01	0.43
1:E:152:ILE:HG13	1:E:218:GLU:HG3	2.01	0.43
1:A:45:GLY:HA3	1:A:55:LYS:HB2	2.01	0.42
1:B:221:LEU:HD21	1:B:226:ILE:HD11	2.01	0.42
1:E:116:PHE:HB3	1:E:137:PHE:CD2	2.54	0.42
1:F:62:PRO:HG3	1:F:123:GLU:HB3	2.01	0.42
1:E:133:PHE:HB3	1:E:153:PHE:HB2	2.01	0.42
1:B:146:LEU:CD1	1:B:201:LEU:HD21	2.50	0.42
1:F:87:ILE:HD13	1:F:165:ILE:HG23	2.02	0.42
1:B:114:LYS:HE2	1:B:140:ALA:HB3	2.01	0.42
1:E:111:PHE:HB3	1:E:139:HIS:HB3	2.01	0.42
1:F:126:ILE:HD13	1:F:152:ILE:CD1	2.49	0.42
1:A:159:ASN:HA	1:A:160:PRO:HD3	1.96	0.42
1:E:193:LYS:HE3	1:E:252:MET:O	2.20	0.42
1:A:88:THR:CG2	1:A:89:ASN:N	2.83	0.41
1:B:233:MET:HG2	1:B:237:THR:HG21	2.01	0.41
1:D:37:GLY:O	1:D:43:HIS:CD2	2.73	0.41
1:B:126:ILE:HD13	1:B:152:ILE:HD13	2.03	0.41
1:E:62:PRO:HG3	1:E:123:GLU:HB3	2.03	0.41
1:F:185:TYR:CZ	1:F:192:LYS:HE2	2.55	0.41
1:B:178:LEU:HD13	1:B:182:ILE:HG12	2.02	0.41
1:C:161:ILE:HD11	1:C:186:ASP:HB3	2.03	0.41
1:A:131:TYR:CE2	1:C:129:LYS:HD3	2.55	0.41
1:F:195:TYR:CE2	1:F:216:GLU:HG2	2.55	0.41
1:B:170:PRO:N	1:B:176:ILE:HD12	2.36	0.41
1:C:153:PHE:CD2	1:C:219:MET:HB3	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:152:ILE:O	1:F:152:ILE:HG13	2.22	0.40
1:E:116:PHE:HA	1:E:136:HIS:O	2.21	0.40
1:E:233:MET:CG	1:E:237:THR:HG21	2.51	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:51:TYR:OH	1:E:38:GLU:OE2[1_556]	2.08	0.12

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	220/226 (97%)	213 (97%)	7 (3%)	0	100	100
1	B	220/226 (97%)	210 (96%)	9 (4%)	1 (0%)	29	39
1	C	220/226 (97%)	212 (96%)	8 (4%)	0	100	100
1	D	220/226 (97%)	212 (96%)	8 (4%)	0	100	100
1	E	220/226 (97%)	212 (96%)	8 (4%)	0	100	100
1	F	220/226 (97%)	214 (97%)	6 (3%)	0	100	100
All	All	1320/1356 (97%)	1273 (96%)	46 (4%)	1 (0%)	51	65

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	225	GLN

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	195/196 (100%)	186 (95%)	9 (5%)	27	36
1	B	195/196 (100%)	189 (97%)	6 (3%)	40	52
1	C	195/196 (100%)	185 (95%)	10 (5%)	24	32
1	D	195/196 (100%)	189 (97%)	6 (3%)	40	52
1	E	195/196 (100%)	187 (96%)	8 (4%)	30	41
1	F	195/196 (100%)	190 (97%)	5 (3%)	46	59
All	All	1170/1176 (100%)	1126 (96%)	44 (4%)	33	44

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	HIS
1	A	98	TYR
1	A	166	TRP
1	A	171	GLU
1	A	190	LYS
1	A	207	SER
1	A	231	LYS
1	A	233	MET
1	A	250	MET
1	B	98	TYR
1	B	110	ARG
1	B	166	TRP
1	B	169	LEU
1	B	233	MET
1	B	252	MET
1	C	32	HIS
1	C	53	MET
1	C	74	GLU
1	C	98	TYR
1	C	112	GLU
1	C	166	TRP
1	C	178	LEU

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Mol	Chain	Res	Type
1	C	198	SER
1	C	222	SER
1	C	233	MET
1	D	98	TYR
1	D	166	TRP
1	D	169	LEU
1	D	227	GLU
1	D	233	MET
1	D	253	ASP
1	E	32	HIS
1	E	69	VAL
1	E	98	TYR
1	E	126	ILE
1	E	166	TRP
1	E	175	THR
1	E	223	LYS
1	E	233	MET
1	F	67	ARG
1	F	98	TYR
1	F	112	GLU
1	F	166	TRP
1	F	233	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	92	HIS
1	D	43	HIS
1	E	42	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (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	222/226 (98%)	-0.04	2 (0%)	84	89	29, 45, 65, 99	0
1	B	222/226 (98%)	-0.04	4 (1%)	68	75	24, 46, 74, 89	0
1	C	222/226 (98%)	-0.04	1 (0%)	91	94	25, 43, 65, 79	0
1	D	222/226 (98%)	0.09	6 (2%)	54	63	31, 48, 77, 98	0
1	E	222/226 (98%)	0.01	1 (0%)	91	94	25, 48, 67, 96	0
1	F	222/226 (98%)	0.10	2 (0%)	84	89	30, 52, 73, 87	0
All	All	1332/1356 (98%)	0.01	16 (1%)	79	84	24, 47, 73, 99	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	179	ALA	3.5
1	B	178	LEU	3.3
1	D	180	HIS	3.3
1	B	172	ALA	3.0
1	F	175	THR	2.7
1	A	253	ASP	2.7
1	A	32	HIS	2.6
1	D	253	ASP	2.5
1	D	47	LEU	2.4
1	B	32	HIS	2.3
1	E	142	LYS	2.2
1	D	170	PRO	2.2
1	D	42	GLN	2.2
1	B	253	ASP	2.1
1	C	32	HIS	2.0
1	F	50	GLU	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 monosaccharides 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(Å^2)	Q<0.9
2	ZN	F	301	1/1	0.98	0.13	45,45,45,45	0
2	ZN	B	301	1/1	0.99	0.11	36,36,36,36	0
2	ZN	C	301	1/1	0.99	0.23	50,50,50,50	0
2	ZN	D	301	1/1	0.99	0.12	38,38,38,38	0
2	ZN	E	301	1/1	0.99	0.15	41,41,41,41	0
2	ZN	A	301	1/1	0.99	0.18	47,47,47,47	0

6.5 Other polymers (i)

There are no such residues in this entry.