Full wwPDB X-ray Structure Validation Report (i)

Dec 12, 2023 – 09:27 pm GMT

PDB ID	:	4B3Z
Title	:	Structure of the human collapsin response mediator protein-1, a lung cancer suppressor
Authors	:	Liu, S.H.; Lin, Y.H.; Huang, S.F.; Niou, Y.K.; Huang, L.L.; Chen, Y.J.
Deposited on	:	2012-07-27
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 (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:

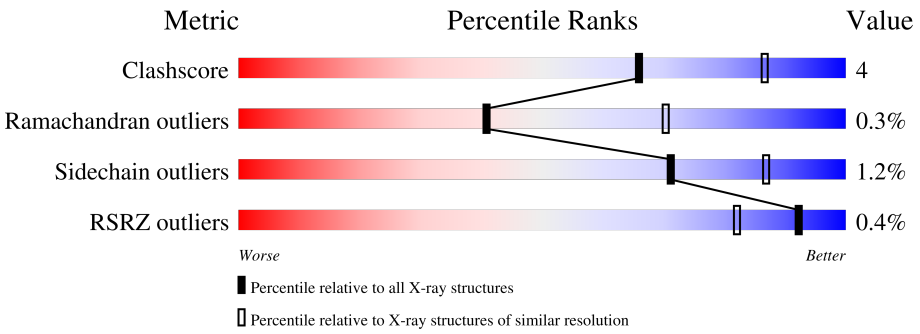
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 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))
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 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	572	71% 12% 17%
1	B	572	73% 9% 18%
1	C	572	% 71% 10% • 18%
1	D	572	75% 8% 17%

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 14530 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 DIHYDROPYRIMIDINASE-RELATED PROTEIN 1.

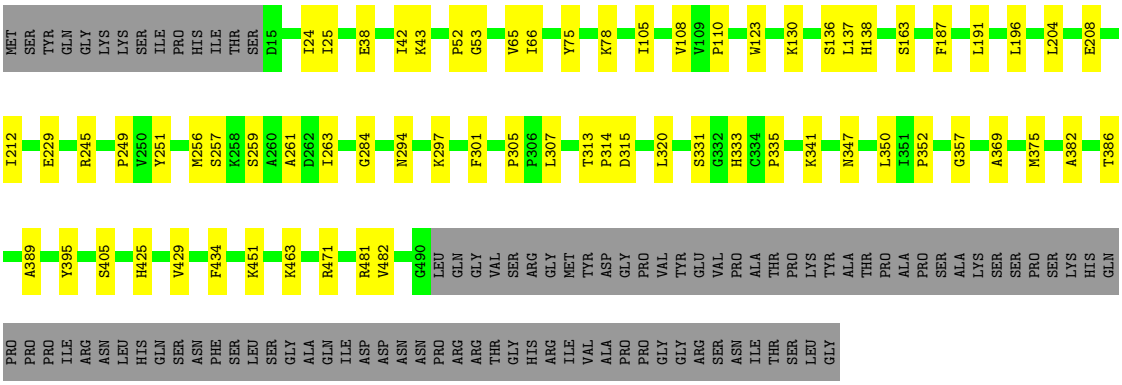
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	476	3654	2322	618	697	17	0	0	0
1	B	469	3608	2295	610	687	16	0	0	0
1	C	468	3603	2292	609	686	16	0	0	0
1	D	477	3665	2329	622	698	16	0	0	0

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.

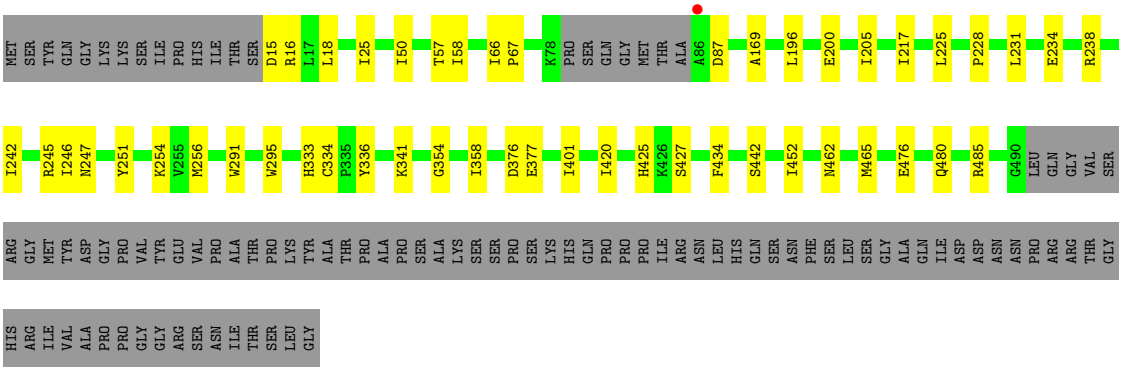
• Molecule 1: DIHYDROPYRIMIDINASE-RELATED PROTEIN 1

Chain A:



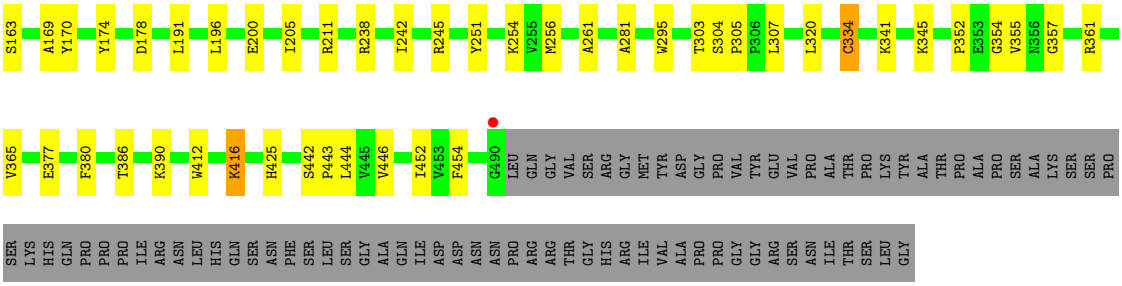
• Molecule 1: DIHYDROPYRIMIDINASE-RELATED PROTEIN 1

Chain B:



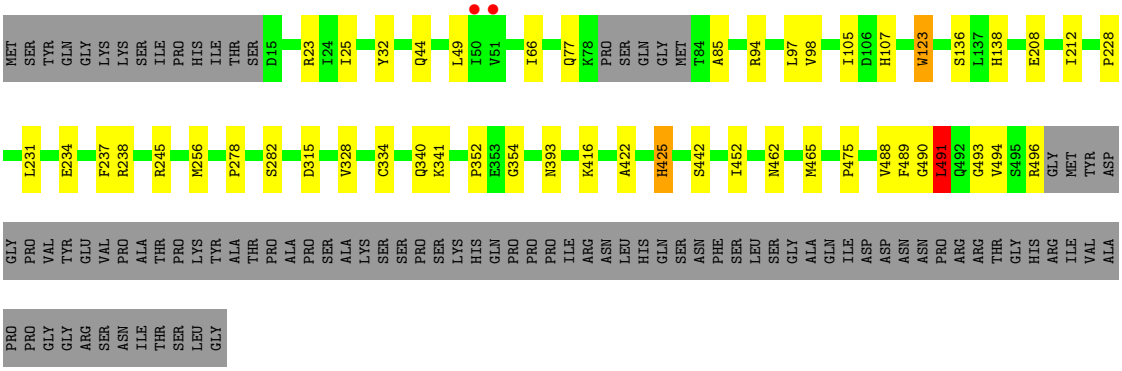
• Molecule 1: DIHYDROPYRIMIDINASE-RELATED PROTEIN 1

Chain C:



• Molecule 1: DIHYDROPYRIMIDINASE-RELATED PROTEIN 1

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	84.81Å 147.85Å 86.02Å 90.00° 94.61° 90.00°	Depositor
Resolution (Å)	39.18 – 3.05 29.62 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.6 (39.18-3.05) 96.7 (29.62-3.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.93 (at 3.00Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.174 , 0.227 0.174 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	56.9	Xtriage
Anisotropy	0.156	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.27 , 20.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.025 for l,-k,h	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	14530	wwPDB-VP
Average B, all atoms (Å^2)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.50% 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)

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	0.42	0/3730	0.58	0/5059
1	B	0.41	0/3682	0.56	0/4992
1	C	0.44	1/3677 (0.0%)	0.58	0/4985
1	D	0.44	0/3739	0.58	0/5069
All	All	0.43	1/14828 (0.0%)	0.57	0/20105

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	334	CYS	CB-SG	-5.17	1.73	1.81

There are no bond angle outliers.

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	3654	0	3641	32	0
1	B	3608	0	3596	29	0
1	C	3603	0	3591	33	0
1	D	3665	0	3657	28	0
All	All	14530	0	14485	116	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 4.

All (116) 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:416:LYS:HD3	1:D:442:SER:HB3	1.72	0.72
1:B:25:ILE:HB	1:B:66:ILE:HG22	1.71	0.71
1:B:295:TRP:HB2	1:D:496:ARG:HD3	1.73	0.70
1:C:25:ILE:HB	1:C:66:ILE:HG22	1.76	0.67
1:C:386:THR:HG22	1:C:390:LYS:HE3	1.78	0.64
1:B:58:ILE:HD11	1:B:452:ILE:HD11	1.81	0.63
1:A:331:SER:OG	1:A:333:HIS:HB2	2.00	0.62
1:B:462:ASN:HB2	1:B:465:MET:HB2	1.82	0.60
1:A:25:ILE:HB	1:A:66:ILE:HG22	1.83	0.59
1:D:488:VAL:O	1:D:490:GLY:N	2.37	0.58
1:A:382:ALA:HA	1:A:386:THR:HB	1.86	0.58
1:C:23:ARG:HD2	1:C:32:TYR:CZ	2.38	0.58
1:D:452:ILE:H	1:D:452:ILE:HD12	1.69	0.58
1:C:412:TRP:CZ3	1:C:443:PRO:HD3	2.38	0.57
1:B:67:PRO:HG3	1:B:401:ILE:HG12	1.84	0.57
1:C:138:HIS:ND1	1:C:163:SER:OG	2.31	0.57
1:D:422:ALA:HA	1:D:425:HIS:CD2	2.38	0.57
1:B:247:ASN:HD21	1:B:485:ARG:HB2	1.70	0.56
1:A:187:PHE:CE2	1:A:191:LEU:HD11	2.41	0.56
1:A:137:LEU:O	1:A:471:ARG:NH2	2.38	0.56
1:A:75:TYR:HB3	1:A:78:LYS:HB2	1.87	0.56
1:A:341:LYS:HA	1:A:352:PRO:HD2	1.87	0.55
1:A:335:PRO:HB2	1:A:429:VAL:HG11	1.88	0.55
1:C:23:ARG:HB2	1:C:32:TYR:CE1	2.40	0.55
1:A:369:ALA:HB1	1:A:375:MET:HG2	1.89	0.54
1:B:334:CYS:HB3	1:B:354:GLY:HA3	1.89	0.54
1:A:130:LYS:HG2	1:A:463:LYS:HE3	1.90	0.54
1:C:196:LEU:HG	1:C:251:TYR:CD1	2.44	0.53
1:C:261:ALA:HB1	1:C:320:LEU:HD12	1.90	0.53
1:C:200:GLU:HB3	1:C:205:ILE:HD11	1.90	0.53
1:A:196:LEU:HG	1:A:251:TYR:CD1	2.44	0.52
1:A:204:LEU:HD21	1:C:242:ILE:HG12	1.91	0.52
1:A:105:ILE:HG23	1:A:136:SER:HB2	1.93	0.51
1:A:313:THR:HB	1:A:314:PRO:HD3	1.93	0.51
1:D:25:ILE:HB	1:D:66:ILE:HG22	1.92	0.50
1:D:278:PRO:HD2	1:D:328:VAL:O	2.11	0.50
1:C:303:THR:HB	1:C:355:VAL:HG13	1.93	0.49
1:D:494:VAL:HG13	1:D:496:ARG:HB2	1.94	0.49
1:B:25:ILE:HD12	1:B:377:GLU:HG3	1.95	0.49
1:B:336:TYR:HB2	1:B:341:LYS:HE2	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:347:ASN:HB3	1:A:350:LEU:HD12	1.94	0.49
1:D:23:ARG:HD2	1:D:32:TYR:CZ	2.48	0.48
1:C:44:GLN:HB3	1:C:49:LEU:HD22	1.96	0.48
1:B:242:ILE:O	1:B:246:ILE:HG13	2.14	0.48
1:A:256:MET:HG3	1:A:305:PRO:HB2	1.96	0.48
1:B:234:GLU:OE1	1:B:238:ARG:NH1	2.47	0.48
1:B:169:ALA:HB2	1:B:200:GLU:HB2	1.96	0.47
1:A:38:GLU:HB3	1:A:43:LYS:HG3	1.95	0.47
1:C:416:LYS:HG2	1:C:442:SER:HB3	1.96	0.47
1:B:476:GLU:HG3	1:B:480:GLN:HG3	1.96	0.47
1:C:341:LYS:HA	1:C:352:PRO:HD2	1.96	0.47
1:B:228:PRO:HD2	1:B:231:LEU:HD12	1.97	0.46
1:A:249:PRO:HD3	1:A:482:VAL:HG21	1.96	0.46
1:B:333:HIS:CG	1:B:358:ILE:HB	2.50	0.46
1:B:225:LEU:HD21	1:B:291:TRP:HZ2	1.81	0.46
1:A:261:ALA:HB1	1:A:320:LEU:HD12	1.98	0.46
1:C:254:LYS:O	1:C:256:MET:HG2	2.16	0.45
1:B:15:ASP:OD1	1:B:16:ARG:N	2.48	0.45
1:D:334:CYS:HB3	1:D:354:GLY:HA3	1.97	0.45
1:B:234:GLU:HA	1:D:237:PHE:HE2	1.82	0.45
1:B:225:LEU:HD21	1:D:491:LEU:HD21	1.99	0.45
1:A:42:ILE:HD12	1:A:405:SER:HB2	2.00	0.44
1:D:208:GLU:O	1:D:212:ILE:HG12	2.17	0.44
1:D:234:GLU:OE1	1:D:238:ARG:NH1	2.51	0.44
1:D:462:ASN:HB2	1:D:465:MET:HB2	1.98	0.44
1:C:334:CYS:HB3	1:C:354:GLY:HA3	2.00	0.44
1:B:217:ILE:HG12	1:D:493:GLY:HA3	1.99	0.44
1:C:178:ASP:OD1	1:C:238:ARG:HD2	2.17	0.44
1:D:44:GLN:HB3	1:D:49:LEU:HD22	1.98	0.44
1:D:105:ILE:HA	1:D:136:SER:O	2.18	0.44
1:D:341:LYS:HA	1:D:352:PRO:HD2	2.00	0.44
1:C:111:GLU:C	1:C:113:GLY:H	2.21	0.44
1:C:153:GLU:HG2	1:C:191:LEU:HD21	1.99	0.43
1:C:169:ALA:HB2	1:C:200:GLU:HB2	1.98	0.43
1:C:304:SER:HA	1:C:305:PRO:C	2.38	0.43
1:A:259:SER:O	1:A:263:ILE:HG12	2.18	0.43
1:A:229:GLU:HB2	1:A:257:SER:CB	2.49	0.43
1:A:481:ARG:HG3	1:B:376:ASP:HB3	2.01	0.43
1:C:108:VAL:HG22	1:C:123:TRP:CG	2.54	0.43
1:B:200:GLU:HB3	1:B:205:ILE:HD11	2.00	0.42
1:C:58:ILE:HD11	1:C:452:ILE:HD11	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:107:HIS:HA	1:D:138:HIS:HB2	2.00	0.42
1:A:108:VAL:O	1:A:110:PRO:HD3	2.19	0.42
1:A:138:HIS:ND1	1:A:163:SER:OG	2.42	0.42
1:C:97:LEU:HD23	1:C:97:LEU:HA	1.95	0.42
1:D:393:ASN:HB2	1:D:475:PRO:HB3	2.00	0.42
1:B:254:LYS:O	1:B:256:MET:HG2	2.19	0.42
1:C:361:ARG:O	1:C:365:VAL:HG23	2.20	0.42
1:A:24:ILE:HA	1:A:65:VAL:HB	2.01	0.42
1:A:229:GLU:HB2	1:A:257:SER:HB2	2.01	0.42
1:C:23:ARG:HG2	1:C:25:ILE:HD11	2.00	0.42
1:A:389:ALA:HB1	1:A:395:TYR:HA	2.02	0.42
1:A:208:GLU:O	1:A:212:ILE:HG12	2.20	0.41
1:B:87:ASP:OD2	1:B:427:SER:OG	2.36	0.41
1:D:85:ALA:HA	1:D:340:GLN:HE22	1.85	0.41
1:D:77:GLN:HB2	1:D:123:TRP:NE1	2.35	0.41
1:D:228:PRO:HD2	1:D:231:LEU:HD12	2.02	0.41
1:D:491:LEU:C	1:D:491:LEU:HD22	2.40	0.41
1:C:446:VAL:HB	1:C:454:PHE:HB3	2.03	0.41
1:C:295:TRP:CZ2	1:C:345:LYS:HA	2.56	0.41
1:C:16:ARG:HB3	1:C:55:VAL:HG12	2.02	0.41
1:C:281:ALA:HB1	1:C:307:LEU:HD21	2.02	0.41
1:A:301:PHE:HD2	1:A:307:LEU:HD13	1.85	0.41
1:B:420:ILE:HG21	1:B:434:PHE:HB2	2.03	0.41
1:B:50:ILE:H	1:B:50:ILE:HG13	1.62	0.41
1:B:196:LEU:HD22	1:B:251:TYR:CD1	2.56	0.41
1:C:107:HIS:HA	1:C:138:HIS:HB2	2.03	0.41
1:C:170:TYR:HB2	1:C:174:TYR:HB2	2.02	0.41
1:C:377:GLU:O	1:C:380:PHE:HB3	2.21	0.41
1:D:256:MET:HE3	1:D:282:SER:HB3	2.02	0.41
1:D:97:LEU:HD23	1:D:97:LEU:HA	1.88	0.41
1:B:238:ARG:O	1:B:242:ILE:HG13	2.20	0.40
1:B:18:LEU:HD23	1:B:57:THR:HG23	2.04	0.40
1:A:284:GLY:HA3	1:A:434:PHE:HE1	1.87	0.40
1:A:294:ASN:ND2	1:A:297:LYS:HB2	2.37	0.40
1:D:94:ARG:O	1:D:98:VAL:HG23	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	474/572 (83%)	445 (94%)	26 (6%)	3 (1%)	25	55
1	B	465/572 (81%)	446 (96%)	19 (4%)	0	100	100
1	C	464/572 (81%)	447 (96%)	16 (3%)	1 (0%)	47	77
1	D	473/572 (83%)	451 (95%)	20 (4%)	2 (0%)	34	64
All	All	1876/2288 (82%)	1789 (95%)	81 (4%)	6 (0%)	41	70

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	489	PHE
1	D	491	LEU
1	A	52	PRO
1	A	53	GLY
1	C	357	GLY
1	A	357	GLY

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	392/473 (83%)	387 (99%)	5 (1%)	69	86
1	B	387/473 (82%)	384 (99%)	3 (1%)	81	91
1	C	387/473 (82%)	381 (98%)	6 (2%)	62	83
1	D	393/473 (83%)	388 (99%)	5 (1%)	69	86

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1559/1892 (82%)	1540 (99%)	19 (1%)	71 87

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	123	TRP
1	A	245	ARG
1	A	315	ASP
1	A	425	HIS
1	A	451	LYS
1	B	245	ARG
1	B	425	HIS
1	B	442	SER
1	C	123	TRP
1	C	211	ARG
1	C	245	ARG
1	C	416	LYS
1	C	425	HIS
1	C	444	LEU
1	D	123	TRP
1	D	245	ARG
1	D	315	ASP
1	D	425	HIS
1	D	491	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	294	ASN
1	B	247	ASN
1	B	425	HIS
1	D	61	ASN
1	D	425	HIS
1	D	449	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)

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	476/572 (83%)	-0.58	0	100	100	30, 53, 91, 168	0
1	B	469/572 (81%)	-0.54	1 (0%)	95	89	38, 61, 89, 134	0
1	C	468/572 (81%)	-0.58	4 (0%)	84	66	28, 52, 89, 165	0
1	D	477/572 (83%)	-0.52	2 (0%)	92	82	34, 59, 98, 165	0
All	All	1890/2288 (82%)	-0.55	7 (0%)	92	82	28, 56, 92, 168	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	87	ASP	3.3
1	D	50	ILE	3.2
1	C	112	PRO	3.0
1	C	490	GLY	2.6
1	B	86	ALA	2.4
1	D	51	VAL	2.2
1	C	54	GLY	2.1

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)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.