Full wwPDB X-ray Structure Validation Report (i)

Dec 2, 2023 – 03:34 pm GMT

PDB ID	:	2WBL
Title	:	Three-dimensional structure of a binary ROP-PRONE complex
Authors	:	Thomas, C.; Fricke, I.; Weyand, M.; Berken, A.
Deposited on	:	2009-03-02
Resolution	:	2.90 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:

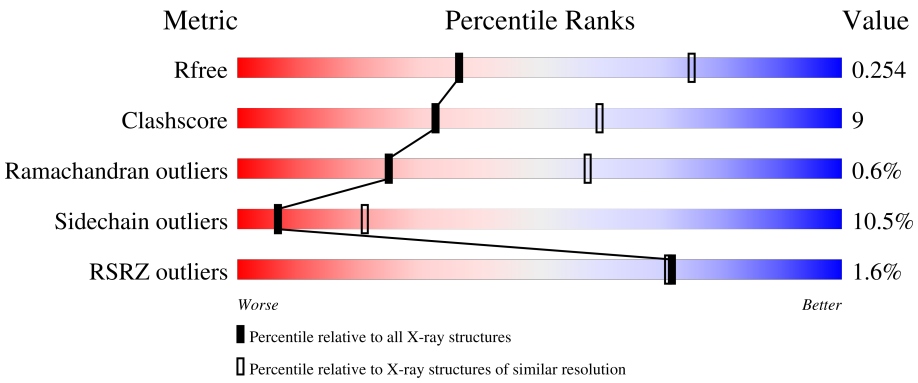
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation 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 2.90 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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	365	.% 71% 16% • 10%
1	B	365	2% 68% 19% • 10%
2	C	180	2% 74% 17% • 7%
2	D	180	.% 72% 18% • 6%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 7497 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 RHO OF PLANTS GUANINE NUCLEOTIDE EXCHANGE FACTOR 8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	329	2493	1583	409	486	15	0	0	0
1	B	329	2489	1587	408	480	14	0	0	1

- Molecule 2 is a protein called RAC-LIKE GTP-BINDING PROTEIN ARAC2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	168	1260	804	215	236	5	0	0	1
2	D	169	1255	802	211	237	5	0	0	1

There are 2 discrepancies between the modelled and reference sequences:

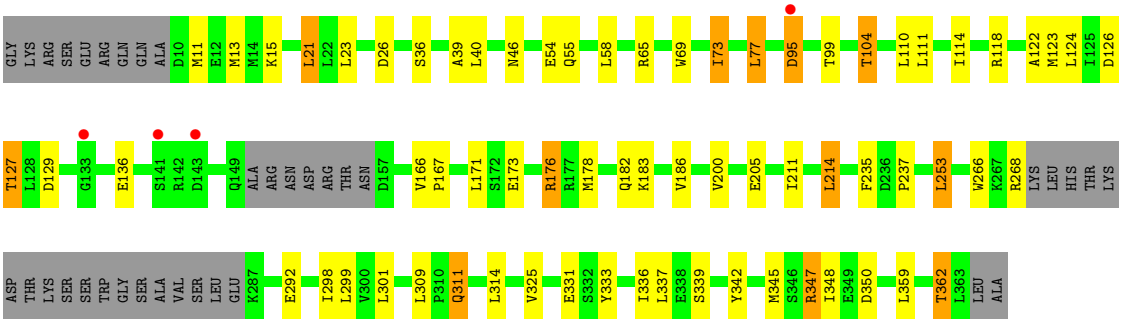
Chain	Residue	Modelled	Actual	Comment	Reference
C	28	GLY	SER	SEE REMARK 999	UNP Q38903
D	28	GLY	SER	SEE REMARK 999	UNP Q38903

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.

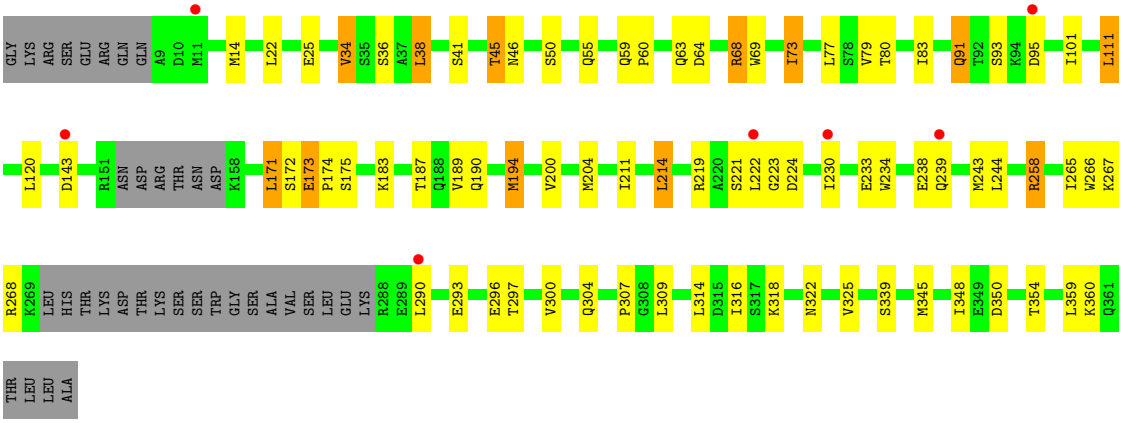
• Molecule 1: RHO OF PLANTS GUANINE NUCLEOTIDE EXCHANGE FACTOR 8

Chain A:



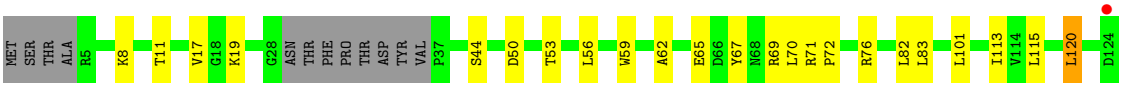
• Molecule 1: RHO OF PLANTS GUANINE NUCLEOTIDE EXCHANGE FACTOR 8

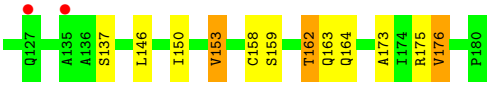
Chain B:



• Molecule 2: RAC-LIKE GTP-BINDING PROTEIN ARAC2

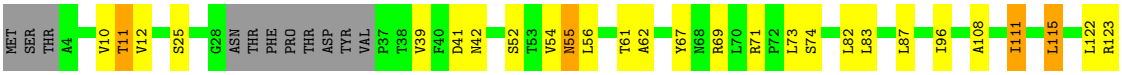
Chain C:

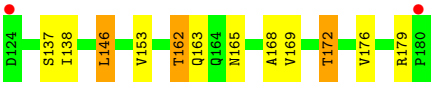




• Molecule 2: RAC-LIKE GTP-BINDING PROTEIN ARAC2

Chain D:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	80.55Å 110.22Å 151.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.41 – 2.90 49.41 – 2.90	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.41-2.90) 99.8 (49.41-2.90)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.84 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.5.0088	Depositor
R, R_{free}	0.215 , 0.261 0.209 , 0.254	Depositor DCC
R_{free} test set	1532 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	51.8	Xtriage
Anisotropy	0.516	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 26.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	7497	wwPDB-VP
Average B, all atoms (Å^2)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.33% 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	0.44	0/2538	0.59	1/3450 (0.0%)
1	B	0.44	0/2535	0.62	1/3448 (0.0%)
2	C	0.50	0/1283	0.63	0/1744
2	D	0.46	1/1278 (0.1%)	0.60	0/1740
All	All	0.45	1/7634 (0.0%)	0.61	2/10382 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	179	ARG	C-N	-5.21	1.24	1.34

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	171	LEU	CA-CB-CG	5.81	128.67	115.30
1	B	111	LEU	CA-CB-CG	5.60	128.18	115.30

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	2493	0	2374	56	0
1	B	2489	0	2371	49	0
2	C	1260	0	1230	23	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	1255	0	1216	23	0
All	All	7497	0	7191	128	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 9.

All (128) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:347:ARG:HG2	1:A:347:ARG:HH11	1.02	1.13
1:A:347:ARG:HG2	1:A:347:ARG:NH1	1.78	0.94
1:A:347:ARG:HH11	1:A:347:ARG:CG	1.83	0.92
1:A:342:TYR:HA	1:A:345:MET:HE3	1.60	0.82
1:B:183:LYS:O	1:B:187:THR:HG23	1.79	0.82
1:A:173:GLU:HA	1:A:176:ARG:HD2	1.64	0.79
1:A:339:SER:HB2	2:C:67:TYR:CD2	2.17	0.79
1:A:46:ASN:OD1	2:C:71:ARG:NH2	2.18	0.76
1:B:187:THR:HG22	1:B:345:MET:SD	2.25	0.76
1:B:350:ASP:O	1:B:354:THR:HG23	1.86	0.76
2:D:123:ARG:HG2	2:D:137:SER:HB2	1.68	0.75
2:C:162:THR:HG22	2:C:164:GLN:H	1.52	0.73
2:D:108:ALA:O	2:D:111:ILE:HD12	1.86	0.73
2:D:168:ALA:O	2:D:172:THR:HG22	1.88	0.73
2:C:62:ALA:O	2:C:71:ARG:NH1	2.20	0.72
1:A:54:GLU:OE1	1:A:65:ARG:HD3	1.89	0.72
2:D:11:THR:HB	2:D:82:LEU:HB2	1.71	0.71
1:B:293:GLU:O	1:B:297:THR:HG23	1.90	0.71
1:A:123:MET:O	1:A:127:THR:HG22	1.91	0.70
1:A:176:ARG:NH2	1:A:359:LEU:HD11	2.06	0.70
1:A:211:ILE:HA	1:A:214:LEU:HD22	1.73	0.69
1:A:339:SER:HB2	2:C:67:TYR:HD2	1.56	0.66
1:A:176:ARG:HH22	1:A:359:LEU:HD11	1.61	0.66
1:B:171:LEU:H	1:B:171:LEU:HD23	1.61	0.66
1:A:124:LEU:HA	1:A:127:THR:HG23	1.79	0.65
2:D:111:ILE:HD13	2:D:111:ILE:O	1.99	0.62
1:A:15:LYS:CG	1:B:14:MET:CE	2.78	0.62
1:A:15:LYS:HG3	1:B:14:MET:HE1	1.82	0.62
1:B:120:LEU:HD22	1:B:189:VAL:HG13	1.82	0.62
1:A:15:LYS:HG2	1:B:14:MET:CE	2.29	0.61
1:B:239:GLN:O	1:B:243:MET:HG2	2.01	0.61
2:D:153:VAL:HG21	2:D:176:VAL:HG21	1.83	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:171:LEU:HD23	1:B:171:LEU:N	2.17	0.59
1:B:204:MET:HE2	1:B:318:LYS:HB3	1.85	0.59
1:A:311:GLN:H	1:A:311:GLN:NE2	2.01	0.58
2:D:83:LEU:HD22	2:D:96:ILE:HD12	1.86	0.58
1:A:339:SER:CB	2:C:67:TYR:CD2	2.87	0.57
1:B:46:ASN:OD1	2:D:71:ARG:NH2	2.38	0.57
1:B:171:LEU:N	1:B:171:LEU:CD2	2.68	0.56
2:C:82:LEU:HD21	2:C:173:ALA:HB2	1.89	0.55
2:D:168:ALA:O	2:D:172:THR:CG2	2.54	0.55
2:C:83:LEU:HD12	2:C:113:ILE:HG21	1.88	0.55
2:D:165:ASN:O	2:D:169:VAL:HG23	2.07	0.55
1:A:73:ILE:HG23	1:A:77:LEU:HD22	1.89	0.55
2:D:62:ALA:O	2:D:71:ARG:NH1	2.38	0.55
1:A:173:GLU:HA	1:A:176:ARG:CD	2.35	0.54
2:C:11:THR:HG22	2:C:19:LYS:HG2	1.89	0.54
1:B:266:TRP:C	1:B:268:ARG:H	2.11	0.54
1:A:15:LYS:HG3	1:B:14:MET:CE	2.37	0.54
1:A:69:TRP:CH2	1:A:73:ILE:HD12	2.43	0.53
1:B:200:VAL:O	1:B:204:MET:HG3	2.09	0.53
1:A:122:ALA:O	1:A:126:ASP:HB2	2.09	0.52
1:B:265:ILE:HD13	1:B:268:ARG:HH21	1.75	0.52
1:B:173:GLU:HG3	1:B:174:PRO:HD3	1.92	0.52
1:A:40:LEU:HD11	1:B:22:LEU:HD13	1.91	0.51
1:B:91:GLN:HG2	1:B:101:ILE:HD13	1.91	0.51
2:C:44:SER:HA	2:C:56:LEU:O	2.10	0.51
1:B:34:VAL:HG21	1:B:38:LEU:HD13	1.92	0.51
2:C:101:LEU:HD13	2:C:150:ILE:HB	1.90	0.51
1:B:219:ARG:O	1:B:223:GLY:HA2	2.11	0.50
1:B:296:GLU:O	1:B:300:VAL:HG23	2.11	0.50
1:B:219:ARG:O	1:B:223:GLY:CA	2.60	0.50
1:B:190:GLN:O	1:B:194:MET:HB2	2.12	0.50
1:A:237:PRO:HB2	1:A:301:LEU:HD12	1.94	0.49
1:A:342:TYR:HA	1:A:345:MET:CE	2.38	0.49
1:A:95:ASP:OD2	1:A:95:ASP:N	2.44	0.49
2:C:162:THR:CG2	2:C:164:GLN:HB2	2.43	0.48
2:D:54:VAL:CG1	2:D:55:ASN:N	2.75	0.48
1:A:114:ILE:O	1:A:118:ARG:HG3	2.12	0.48
1:A:182:GLN:O	1:A:186:VAL:HG23	2.13	0.48
2:D:67:TYR:CE2	2:D:69:ARG:HB2	2.48	0.48
2:D:162:THR:O	2:D:163:GLN:HB2	2.14	0.47
2:D:115:LEU:HD11	2:D:146:LEU:HD13	1.97	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:15:LYS:HG2	1:B:14:MET:HE2	1.96	0.47
1:A:166:VAL:HG13	1:A:167:PRO:HD2	1.96	0.47
1:A:311:GLN:HE22	2:C:69:ARG:HA	1.80	0.47
1:A:15:LYS:CG	1:B:14:MET:HE2	2.45	0.46
1:A:21:LEU:HB3	1:B:22:LEU:HD21	1.96	0.46
1:A:26:ASP:HA	2:C:44:SER:HB3	1.96	0.46
2:D:123:ARG:NH1	2:D:138:ILE:O	2.43	0.45
1:A:253:LEU:HD23	1:A:309:LEU:HD22	1.98	0.45
2:C:153:VAL:HG11	2:C:176:VAL:HG21	1.98	0.45
2:C:162:THR:O	2:C:163:GLN:HB2	2.17	0.45
1:A:55:GLN:O	1:A:347:ARG:NH1	2.38	0.44
1:B:25:GLU:HG2	2:D:42:ASN:O	2.17	0.44
1:A:21:LEU:HD12	1:A:39:ALA:HB1	1.99	0.44
2:C:120:LEU:HB2	2:C:158:CYS:O	2.18	0.44
1:B:221:SER:O	1:B:258:ARG:HG3	2.17	0.44
1:B:80:THR:HA	1:B:83:ILE:HD12	1.99	0.44
2:C:153:VAL:HG21	2:C:176:VAL:HG21	1.99	0.44
1:A:311:GLN:HG3	2:C:72:PRO:HB3	2.00	0.43
1:A:15:LYS:CG	1:B:14:MET:HE1	2.42	0.43
1:A:123:MET:O	1:A:127:THR:CG2	2.64	0.43
1:B:69:TRP:CH2	1:B:73:ILE:HG13	2.53	0.43
1:A:104:THR:O	1:A:104:THR:CG2	2.66	0.43
1:A:266:TRP:C	1:A:268:ARG:H	2.21	0.43
2:D:12:VAL:HG12	2:D:61:THR:HG21	2.01	0.43
2:D:10:VAL:HG21	2:D:74:SER:O	2.19	0.43
1:A:336:ILE:HG13	1:A:337:LEU:N	2.34	0.43
1:B:172:SER:CB	1:B:174:PRO:HD2	2.49	0.42
2:D:41:ASP:O	2:D:42:ASN:C	2.58	0.42
1:A:299:LEU:HD23	1:A:299:LEU:HA	1.83	0.42
1:B:79:VAL:HG11	1:B:325:VAL:HG11	2.00	0.42
1:A:311:GLN:CG	2:C:72:PRO:HB3	2.49	0.42
1:B:183:LYS:HB2	1:B:348:ILE:HG21	2.01	0.42
1:B:204:MET:CE	1:B:322:ASN:HB2	2.49	0.42
2:C:8:LYS:HE3	2:C:59:TRP:CE2	2.54	0.42
1:A:183:LYS:HB2	1:A:348:ILE:HG21	2.01	0.42
1:A:176:ARG:HH22	1:A:359:LEU:CD1	2.31	0.41
1:A:235:PHE:CZ	1:A:298:ILE:HD11	2.55	0.41
1:A:69:TRP:CZ3	1:A:73:ILE:HD12	2.55	0.41
1:B:173:GLU:N	1:B:174:PRO:CD	2.83	0.41
1:B:211:ILE:HA	1:B:214:LEU:HD22	2.02	0.41
1:A:211:ILE:O	1:A:214:LEU:HB2	2.21	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:93:SER:C	1:B:95:ASP:H	2.24	0.41
1:B:316:ILE:HG23	2:D:73:LEU:HD23	2.02	0.41
1:A:69:TRP:HH2	1:A:336:ILE:HD12	1.84	0.41
1:A:339:SER:CB	2:C:67:TYR:HD2	2.25	0.41
1:B:63:GLN:HE21	1:B:64:ASP:H	1.69	0.41
1:B:59:GLN:HB2	1:B:60:PRO:CD	2.51	0.41
2:C:159:SER:OG	2:C:162:THR:HB	2.21	0.41
1:B:304:GLN:O	1:B:307:PRO:HD3	2.21	0.41
1:A:136:GLU:O	1:A:167:PRO:HD3	2.20	0.40
1:B:204:MET:HE3	1:B:322:ASN:HB2	2.02	0.40
1:B:41:SER:O	1:B:45:THR:CG2	2.70	0.40
1:A:77:LEU:HD21	1:A:333:TYR:CE2	2.56	0.40
1:B:68:ARG:HH22	2:D:39:VAL:CG2	2.34	0.40
1:B:339:SER:HB2	2:D:67:TYR:CD2	2.57	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	323/365 (88%)	309 (96%)	13 (4%)	1 (0%)	41	71
1	B	323/365 (88%)	304 (94%)	16 (5%)	3 (1%)	17	48
2	C	164/180 (91%)	156 (95%)	7 (4%)	1 (1%)	25	58
2	D	165/180 (92%)	154 (93%)	10 (6%)	1 (1%)	25	58
All	All	975/1090 (89%)	923 (95%)	46 (5%)	6 (1%)	25	58

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	362	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	224	ASP
1	B	360	LYS
2	C	50	ASP
2	D	87	LEU
1	B	267	LYS

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	260/323 (80%)	231 (89%)	29 (11%)	6 18
1	B	254/323 (79%)	226 (89%)	28 (11%)	6 19
2	C	130/152 (86%)	117 (90%)	13 (10%)	7 23
2	D	128/152 (84%)	117 (91%)	11 (9%)	10 30
All	All	772/950 (81%)	691 (90%)	81 (10%)	7 21

All (81) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	MET
1	A	13	MET
1	A	21	LEU
1	A	23	LEU
1	A	36	SER
1	A	58	LEU
1	A	73	ILE
1	A	77	LEU
1	A	95	ASP
1	A	99	THR
1	A	104	THR
1	A	110	LEU
1	A	111	LEU
1	A	127	THR
1	A	129	ASP
1	A	176	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	178	MET
1	A	200	VAL
1	A	205	GLU
1	A	214	LEU
1	A	253	LEU
1	A	292	GLU
1	A	311	GLN
1	A	314	LEU
1	A	325	VAL
1	A	331	GLU
1	A	347	ARG
1	A	350	ASP
1	A	362	THR
1	B	34	VAL
1	B	36	SER
1	B	38	LEU
1	B	45	THR
1	B	50	SER
1	B	55	GLN
1	B	68	ARG
1	B	73	ILE
1	B	77	LEU
1	B	91	GLN
1	B	111	LEU
1	B	143	ASP
1	B	171	LEU
1	B	173	GLU
1	B	175	SER
1	B	194	MET
1	B	214	LEU
1	B	222	LEU
1	B	230	ILE
1	B	233	GLU
1	B	234	TRP
1	B	238	GLU
1	B	244	LEU
1	B	258	ARG
1	B	290	LEU
1	B	309	LEU
1	B	314	LEU
1	B	359	LEU
2	C	17	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	C	53	THR
2	C	65	GLU
2	C	70	LEU
2	C	76	ARG
2	C	115	LEU
2	C	120	LEU
2	C	137	SER
2	C	146	LEU
2	C	153	VAL
2	C	162	THR
2	C	175	ARG
2	C	176	VAL
2	D	11	THR
2	D	25	SER
2	D	52	SER
2	D	55	ASN
2	D	56	LEU
2	D	111	ILE
2	D	115	LEU
2	D	122	LEU
2	D	146	LEU
2	D	162	THR
2	D	172	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	82	HIS
1	A	91	GLN
1	A	135	ASN
1	A	250	HIS
1	A	304	GLN
1	A	311	GLN
1	B	55	GLN
1	B	63	GLN
1	B	66	GLN
1	B	106	GLN
1	B	130	ASN
1	B	148	GLN
1	B	182	GLN
2	C	42	ASN
2	C	55	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	D	64	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	329/365 (90%)	0.06	4 (1%)	79	79	16, 29, 42, 63	0
1	B	329/365 (90%)	0.10	7 (2%)	63	61	17, 28, 44, 59	0
2	C	168/180 (93%)	-0.12	3 (1%)	68	67	20, 28, 32, 39	0
2	D	169/180 (93%)	-0.16	2 (1%)	79	79	23, 31, 39, 45	0
All	All	995/1090 (91%)	0.00	16 (1%)	72	71	16, 29, 40, 63	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	11	MET	3.5
1	B	95	ASP	3.0
1	A	133	GLY	2.9
2	C	124	ASP	2.8
2	C	135	ALA	2.8
1	B	222	LEU	2.7
1	B	239	GLN	2.4
1	A	141	SER	2.4
1	A	95	ASP	2.3
1	B	230	ILE	2.3
2	D	124	ASP	2.3
1	A	143	ASP	2.3
2	D	180	PRO	2.2
1	B	143	ASP	2.2
1	B	290	LEU	2.1
2	C	127	GLN	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)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.