



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:09 AM GMT

PDB ID : 3HDV
Title : Crystal structure of response regulator receiver protein from *Pseudomonas putida*
Authors : Bagaria, A.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2009-05-07
Resolution : 2.09 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
<http://wwpdb.org/validation/2016/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
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

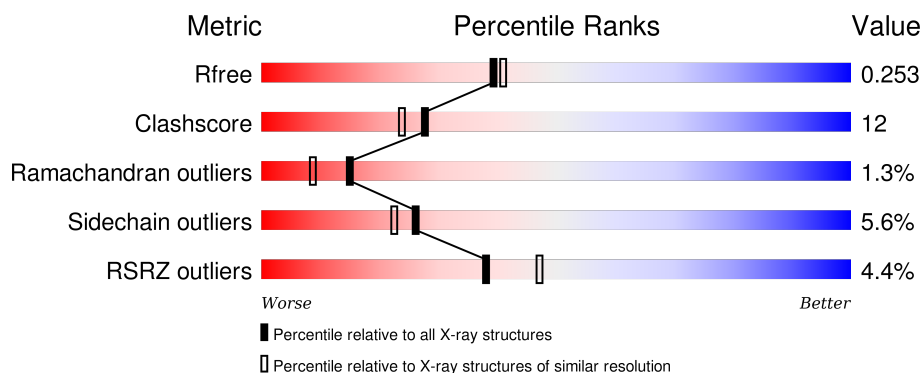
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.09 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	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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 $\leq 5\%$. 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	136	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 99%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 80% 5% •• 12% </div> </div>
1	B	136	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 99%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 84% 9% 7% </div> </div>
1	C	136	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 3%, green 97%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 77% 10% •• 11% </div> </div>
1	D	136	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 10%, green 90%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 60% 15% 10% • 14% </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3978 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 Response regulator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	120	Total	C	N	O	Se	0	0	0
			931	590	162	176	3			
1	B	126	Total	C	N	O	Se	0	0	0
			970	616	168	183	3			
1	C	121	Total	C	N	O	Se	0	0	0
			931	589	162	177	3			
1	D	117	Total	C	N	O	Se	0	0	0
			903	572	155	173	3			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	62	MSE	-	expression tag	UNP Q88QX9
A	63	SER	-	expression tag	UNP Q88QX9
A	64	LEU	-	expression tag	UNP Q88QX9
A	190	GLU	-	expression tag	UNP Q88QX9
A	191	GLY	-	expression tag	UNP Q88QX9
A	192	HIS	-	expression tag	UNP Q88QX9
A	193	HIS	-	expression tag	UNP Q88QX9
A	194	HIS	-	expression tag	UNP Q88QX9
A	195	HIS	-	expression tag	UNP Q88QX9
A	196	HIS	-	expression tag	UNP Q88QX9
A	197	HIS	-	expression tag	UNP Q88QX9
B	62	MSE	-	expression tag	UNP Q88QX9
B	63	SER	-	expression tag	UNP Q88QX9
B	64	LEU	-	expression tag	UNP Q88QX9
B	190	GLU	-	expression tag	UNP Q88QX9
B	191	GLY	-	expression tag	UNP Q88QX9
B	192	HIS	-	expression tag	UNP Q88QX9
B	193	HIS	-	expression tag	UNP Q88QX9
B	194	HIS	-	expression tag	UNP Q88QX9
B	195	HIS	-	expression tag	UNP Q88QX9
B	196	HIS	-	expression tag	UNP Q88QX9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	197	HIS	-	expression tag	UNP Q88QX9
C	62	MSE	-	expression tag	UNP Q88QX9
C	63	SER	-	expression tag	UNP Q88QX9
C	64	LEU	-	expression tag	UNP Q88QX9
C	190	GLU	-	expression tag	UNP Q88QX9
C	191	GLY	-	expression tag	UNP Q88QX9
C	192	HIS	-	expression tag	UNP Q88QX9
C	193	HIS	-	expression tag	UNP Q88QX9
C	194	HIS	-	expression tag	UNP Q88QX9
C	195	HIS	-	expression tag	UNP Q88QX9
C	196	HIS	-	expression tag	UNP Q88QX9
C	197	HIS	-	expression tag	UNP Q88QX9
D	62	MSE	-	expression tag	UNP Q88QX9
D	63	SER	-	expression tag	UNP Q88QX9
D	64	LEU	-	expression tag	UNP Q88QX9
D	190	GLU	-	expression tag	UNP Q88QX9
D	191	GLY	-	expression tag	UNP Q88QX9
D	192	HIS	-	expression tag	UNP Q88QX9
D	193	HIS	-	expression tag	UNP Q88QX9
D	194	HIS	-	expression tag	UNP Q88QX9
D	195	HIS	-	expression tag	UNP Q88QX9
D	196	HIS	-	expression tag	UNP Q88QX9
D	197	HIS	-	expression tag	UNP Q88QX9

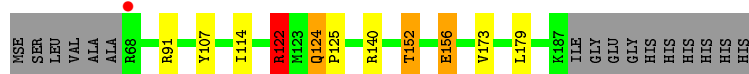
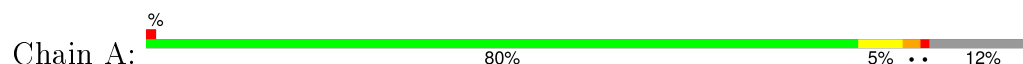
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	58	Total O 58 58	0	0
2	B	81	Total O 81 81	0	0
2	C	68	Total O 68 68	0	0
2	D	36	Total O 36 36	0	0

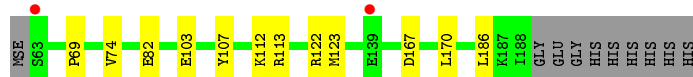
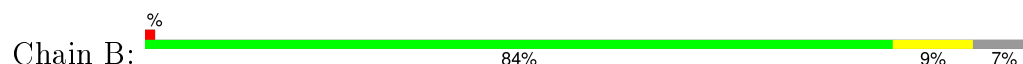
3 Residue-property plots [i](#)

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.

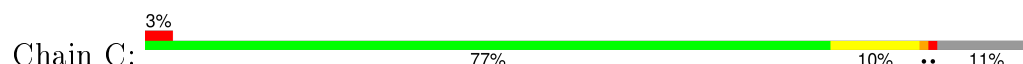
- Molecule 1: Response regulator



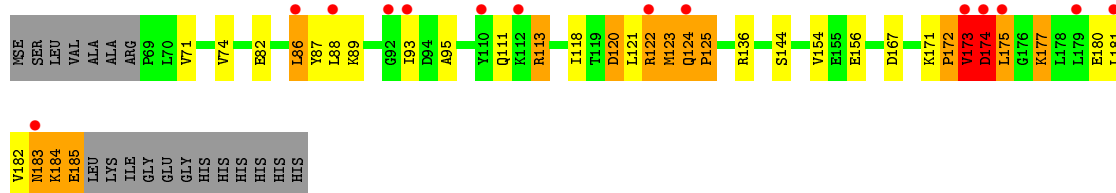
- Molecule 1: Response regulator



- Molecule 1: Response regulator



- Molecule 1: Response regulator



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	110.77Å 140.44Å 48.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.50 – 2.09 44.48 – 2.09	Depositor EDS
% Data completeness (in resolution range)	99.6 (44.50-2.09) 99.6 (44.48-2.09)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.43 (at 2.08Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.205 , 0.245 0.218 , 0.253	Depositor DCC
R_{free} test set	2304 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	38.2	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 45.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 45619 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3978	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 4.45% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	0.91	1/936 (0.1%)	0.87	2/1260 (0.2%)
1	B	0.97	2/975 (0.2%)	0.87	1/1314 (0.1%)
1	C	0.96	1/936 (0.1%)	1.00	5/1262 (0.4%)
1	D	0.92	1/908 (0.1%)	0.91	4/1223 (0.3%)
All	All	0.94	5/3755 (0.1%)	0.91	12/5059 (0.2%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	123	MSE	C-N	9.59	1.56	1.34
1	B	82	GLU	CG-CD	5.30	1.59	1.51
1	B	82	GLU	CD-OE1	5.24	1.31	1.25
1	C	155	GLU	CB-CG	5.11	1.61	1.52
1	A	156	GLU	CD-OE1	-5.08	1.20	1.25

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	122	ARG	NE-CZ-NH2	-12.03	114.28	120.30
1	C	122	ARG	NE-CZ-NH1	8.51	124.56	120.30
1	A	122	ARG	NE-CZ-NH2	-7.69	116.45	120.30
1	D	120	ASP	CB-CG-OD1	5.77	123.49	118.30
1	D	89	LYS	CB-CA-C	-5.75	98.89	110.40
1	C	167	ASP	CB-CG-OD1	5.66	123.39	118.30
1	D	136	ARG	NE-CZ-NH2	-5.42	117.59	120.30
1	D	86	LEU	CB-CG-CD2	5.35	120.10	111.00
1	C	122	ARG	CG-CD-NE	-5.28	100.71	111.80
1	A	152	THR	CB-CA-C	-5.23	97.47	111.60
1	B	167	ASP	CB-CG-OD1	5.04	122.83	118.30
1	C	120	ASP	CB-CG-OD1	5.02	122.82	118.30

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	931	0	978	7	0
1	B	970	0	1024	3	0
1	C	931	0	973	17	0
1	D	903	0	942	64	0
2	A	58	0	0	1	0
2	B	81	0	0	0	0
2	C	68	0	0	5	0
2	D	36	0	0	0	0
All	All	3978	0	3917	90	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 12.

All (90) 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:87:TYR:CZ	1:D:175:LEU:HD22	1.62	1.35
1:D:174:ASP:O	1:D:175:LEU:HG	1.34	1.24
1:D:124:GLN:CB	1:D:125:PRO:HD3	1.73	1.16
1:D:124:GLN:HB2	1:D:125:PRO:HD3	1.14	1.14
1:D:173:VAL:HG13	1:D:174:ASP:H	1.00	1.14
1:D:173:VAL:HG13	1:D:174:ASP:N	1.54	1.07
1:D:124:GLN:HB2	1:D:125:PRO:CD	1.84	1.06
1:D:121:LEU:CB	1:D:122:ARG:HH11	1.71	1.02
1:D:87:TYR:CZ	1:D:175:LEU:CD2	2.45	0.98
1:D:121:LEU:HB3	1:D:122:ARG:NH1	1.82	0.94
1:D:173:VAL:CG1	1:D:174:ASP:N	2.30	0.94
1:D:174:ASP:O	1:D:175:LEU:CG	2.18	0.91
1:D:88:LEU:HD21	1:D:181:LEU:HD22	1.50	0.90
1:D:121:LEU:HB3	1:D:122:ARG:HH11	1.36	0.90
1:D:87:TYR:CE2	1:D:175:LEU:HD22	2.10	0.87
1:D:121:LEU:HB2	1:D:122:ARG:HH11	1.39	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:124:GLN:CB	1:D:125:PRO:CD	2.45	0.86
1:D:82:GLU:O	1:D:86:LEU:HG	1.76	0.85
1:D:173:VAL:CG1	1:D:174:ASP:H	1.83	0.83
1:C:112:LYS:O	1:C:113:ARG:HB2	1.79	0.83
1:D:88:LEU:CD2	1:D:181:LEU:HD22	2.11	0.79
1:D:87:TYR:OH	1:D:175:LEU:HD22	1.81	0.79
1:D:173:VAL:CG2	1:D:177:LYS:HZ1	1.99	0.75
1:C:112:LYS:HG3	1:C:113:ARG:H	1.52	0.75
1:C:77:ASN:HD22	1:C:80:ASN:H	1.35	0.74
1:C:112:LYS:O	1:C:113:ARG:CB	2.37	0.73
1:D:173:VAL:HG22	1:D:177:LYS:HZ1	1.53	0.72
1:D:172:PRO:O	1:D:173:VAL:O	2.09	0.71
1:D:173:VAL:HG22	1:D:177:LYS:NZ	2.06	0.69
1:D:118:ILE:HD11	1:D:181:LEU:HD11	1.74	0.69
1:D:124:GLN:HB3	1:D:125:PRO:HD3	1.69	0.69
1:D:122:ARG:HH12	1:D:156:GLU:CD	1.96	0.68
1:D:177:LYS:HZ3	1:D:177:LYS:CB	2.07	0.68
1:C:112:LYS:CG	1:C:113:ARG:H	2.07	0.68
1:A:124:GLN:HG3	1:A:125:PRO:HA	1.78	0.65
1:D:88:LEU:HD21	1:D:181:LEU:CD2	2.27	0.64
1:D:121:LEU:CB	1:D:122:ARG:NH1	2.44	0.63
1:D:167:ASP:HB2	1:D:180:GLU:OE2	1.99	0.63
1:D:111:GLN:HE21	1:D:113:ARG:HG3	1.66	0.61
1:D:88:LEU:CD2	1:D:181:LEU:CD2	2.79	0.61
1:D:93:ILE:HD11	1:D:182:VAL:CG1	2.30	0.61
1:D:111:GLN:NE2	1:D:113:ARG:HG3	2.16	0.60
1:D:87:TYR:CE2	1:D:175:LEU:CD2	2.79	0.60
1:D:144:SER:HB3	1:D:184:LYS:HD2	1.85	0.59
1:D:173:VAL:CG2	1:D:177:LYS:NZ	2.64	0.58
1:C:122:ARG:NH2	1:C:156:GLU:OE2	2.34	0.57
1:D:74:VAL:HG12	1:D:123:MSE:SE	2.55	0.56
1:C:112:LYS:HG3	1:C:113:ARG:N	2.19	0.55
1:D:122:ARG:NH1	1:D:156:GLU:OE2	2.39	0.55
1:D:177:LYS:HZ3	1:D:177:LYS:HB3	1.70	0.55
1:A:173:VAL:HG12	2:A:204:HOH:O	2.06	0.55
1:C:122:ARG:HH22	1:C:156:GLU:CD	2.10	0.54
1:A:91:ARG:HG3	1:A:179:LEU:HD11	1.91	0.52
1:A:122:ARG:NH2	1:A:156:GLU:OE1	2.38	0.52
1:D:93:ILE:HD11	1:D:182:VAL:HG12	1.91	0.52
1:D:174:ASP:O	1:D:175:LEU:CB	2.56	0.52
1:C:112:LYS:CG	1:C:113:ARG:N	2.73	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:87:TYR:CE1	1:D:175:LEU:CD2	2.92	0.52
1:D:177:LYS:O	1:D:180:GLU:HB2	2.11	0.51
1:D:122:ARG:HE	1:D:122:ARG:N	2.08	0.51
1:D:175:LEU:C	1:D:177:LYS:H	2.15	0.50
1:B:74:VAL:HG12	1:B:123:MSE:SE	2.62	0.50
1:A:122:ARG:HH22	1:A:156:GLU:CD	2.15	0.49
1:D:144:SER:CB	1:D:184:LYS:HD2	2.43	0.48
1:C:136:ARG:NH2	2:C:35:HOH:O	2.42	0.48
1:C:94:ASP:OD2	1:C:113:ARG:NH2	2.47	0.48
1:C:130:ASP:OD2	2:C:222:HOH:O	2.20	0.48
1:D:177:LYS:NZ	1:D:177:LYS:CB	2.77	0.48
1:D:120:ASP:OD1	1:D:121:LEU:N	2.46	0.46
1:D:171:LYS:HB3	1:D:172:PRO:HA	1.98	0.46
1:C:122:ARG:HA	1:C:122:ARG:HD3	1.69	0.45
1:D:122:ARG:HA	1:D:122:ARG:HD3	1.83	0.45
1:C:77:ASN:ND2	1:C:80:ASN:H	2.09	0.44
1:C:122:ARG:HD3	2:C:2:HOH:O	2.17	0.44
1:D:122:ARG:NE	1:D:122:ARG:CA	2.80	0.43
1:D:88:LEU:HD22	1:D:181:LEU:CD2	2.46	0.43
1:D:173:VAL:HG21	1:D:177:LYS:HZ1	1.83	0.42
1:A:114:ILE:O	1:A:140:ARG:NH2	2.48	0.42
1:B:170:LEU:HD21	1:D:154:VAL:HG11	2.01	0.42
1:C:102:GLU:HG2	2:C:214:HOH:O	2.18	0.42
1:D:71:VAL:O	1:D:95:ALA:HA	2.20	0.42
1:D:82:GLU:O	1:D:86:LEU:CG	2.59	0.42
1:D:174:ASP:HB3	1:D:175:LEU:H	1.72	0.41
1:D:185:GLU:HG2	1:D:185:GLU:H	1.54	0.41
1:D:175:LEU:C	1:D:177:LYS:N	2.74	0.41
1:D:86:LEU:HD23	1:D:86:LEU:N	2.35	0.41
1:A:122:ARG:HD3	1:A:122:ARG:HA	1.33	0.41
1:C:155:GLU:HG3	2:C:25:HOH:O	2.19	0.41
1:D:183:ASN:HD22	1:D:183:ASN:C	2.24	0.41
1:B:69:PRO:HB3	1:B:186:LEU:HD13	2.03	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	118/136 (87%)	114 (97%)	4 (3%)	0	100	100
1	B	124/136 (91%)	120 (97%)	4 (3%)	0	100	100
1	C	119/136 (88%)	116 (98%)	2 (2%)	1 (1%)	24	17
1	D	115/136 (85%)	106 (92%)	4 (4%)	5 (4%)	3	1
All	All	476/544 (88%)	456 (96%)	14 (3%)	6 (1%)	15	9

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	124	GLN
1	D	173	VAL
1	D	174	ASP
1	D	175	LEU
1	C	113	ARG
1	D	125	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	102/110 (93%)	98 (96%)	4 (4%)	39	39
1	B	106/110 (96%)	101 (95%)	5 (5%)	32	30
1	C	101/110 (92%)	96 (95%)	5 (5%)	30	27
1	D	99/110 (90%)	90 (91%)	9 (9%)	12	7

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	408/440 (93%)	385 (94%)	23 (6%)	26	22

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	107	TYR
1	A	122	ARG
1	A	124	GLN
1	A	152	THR
1	B	103	GLU
1	B	107	TYR
1	B	112	LYS
1	B	113	ARG
1	B	122	ARG
1	C	65	VAL
1	C	76	ASP
1	C	113	ARG
1	C	122	ARG
1	C	140	ARG
1	D	113	ARG
1	D	122	ARG
1	D	172	PRO
1	D	173	VAL
1	D	174	ASP
1	D	177	LYS
1	D	183	ASN
1	D	184	LYS
1	D	185	GLU

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

Mol	Chain	Res	Type
1	A	124	GLN
1	C	77	ASN
1	D	80	ASN
1	D	109	HIS
1	D	111	GLN
1	D	183	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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	117/136 (86%)	0.25	1 (0%) 85 88	26, 37, 51, 62	0
1	B	123/136 (90%)	0.26	2 (1%) 74 79	25, 35, 52, 61	0
1	C	118/136 (86%)	0.42	4 (3%) 49 58	26, 36, 54, 63	0
1	D	114/136 (83%)	0.82	14 (12%) 5 7	28, 42, 65, 68	0
All	All	472/544 (86%)	0.43	21 (4%) 38 47	25, 37, 58, 68	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	173	VAL	9.8
1	D	175	LEU	6.2
1	D	174	ASP	5.8
1	D	124	GLN	5.0
1	D	92	GLY	3.4
1	D	122	ARG	3.2
1	B	139	GLU	3.2
1	D	93	ILE	2.9
1	D	181	LEU	2.9
1	D	112	LYS	2.5
1	C	67	ALA	2.5
1	D	183	ASN	2.5
1	C	76	ASP	2.5
1	C	139	GLU	2.4
1	D	88	LEU	2.3
1	D	110	TYR	2.3
1	B	63	SER	2.2
1	C	65	VAL	2.2
1	D	86	LEU	2.1
1	D	179	LEU	2.1
1	A	68	ARG	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.