



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:57 AM UTC

PDB ID : 2XG8 / pdb_00002xg8
Title : Structural basis of gene regulation by protein PII: The crystal complex of PII and PipX from *Synechococcus elongatus* PCC 7942
Authors : Llacer, J.L.; Rubio, V.
Deposited on : 2010-06-01
Resolution : 3.20 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

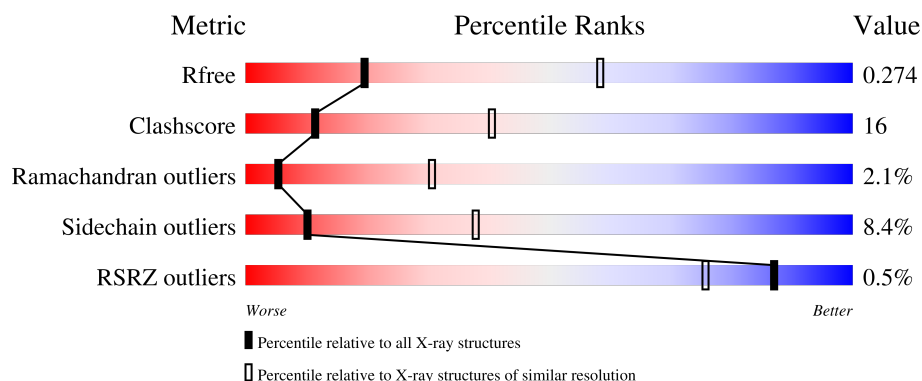
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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}	180053	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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 $\leq 5\%$. 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	112	 75% 17% 5%
1	B	112	 76% 18%
1	C	112	 77% 19%
2	D	89	 61% 31%
2	E	89	 2% 61% 31%

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Mol	Chain	Length	Quality of chain
2	F	89	% 57%22%••12%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4532 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 NITROGEN REGULATORY PROTEIN P-II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	106	Total	C	N	O	S	0	0	0
			824	522	144	156	2			
1	B	108	Total	C	N	O	S	0	0	0
			834	528	145	159	2			
1	C	112	Total	C	N	O	S	0	0	0
			853	540	148	163	2			

- Molecule 2 is a protein called PIPX.

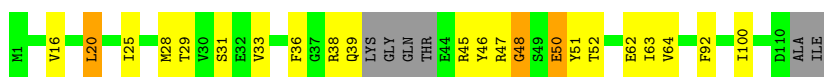
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	87	Total	C	N	O	S	0	0	0
			717	461	125	129	2			
2	E	87	Total	C	N	O	S	0	0	0
			709	453	128	126	2			
2	F	78	Total	C	N	O	S	0	0	0
			592	383	103	104	2			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	O	0	0
			1	1		
3	B	1	Total	O	0	0
			1	1		
3	C	1	Total	O	0	0
			1	1		

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● Molecule 1: NITROGEN REGULATORY PROTEIN P-II

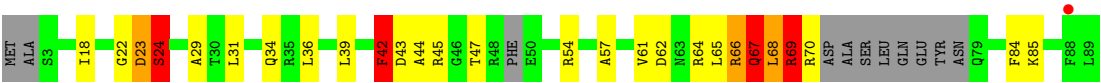


R1
R9
V16
L20
M28
T29
S30
V31
F32
V33
F36
G37
R38
Q39
K40
G41
Q42
Y46
R47
G48
S49
E50
V53
E62
I63
I86
F92
T99
N108
A1A
ASP
A1A
I1E

V11	V16	L20	M28	T29	V30	S31	E32	V33	G37	R38	Q39	K40	G41	G42	T43	E44	R45	Y46	R47	G48	S49	E50	Y51	E54	E62	I63	V64	F92	I100	I112
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Chain F: 57% 22% 12% 9%



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	111.38Å 61.59Å 112.98Å 90.00° 126.44° 90.00°	Depositor
Resolution (Å)	50.00 – 3.20 50.00 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (50.00-3.20) 99.8 (50.00-3.20)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.79 (at 3.19Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.227 , 0.272 0.241 , 0.274	Depositor DCC
R_{free} test set	495 reflections (4.79%)	wwPDB-VP
Wilson B-factor (Å ²)	73.2	Xtriage
Anisotropy	0.166	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 82.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4532	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.72% 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, 0.333 respectively for untwinned datasets, and 0.375, 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	0.54	0/830	0.79	0/1113
1	B	0.55	0/841	0.88	1/1130 (0.1%)
1	C	0.53	0/860	0.96	6/1156 (0.5%)
2	D	0.53	0/732	0.84	2/987 (0.2%)
2	E	0.50	0/724	0.89	2/976 (0.2%)
2	F	0.61	0/603	0.98	3/815 (0.4%)
All	All	0.54	0/4590	0.89	14/6177 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	C	0	1
2	F	0	4
All	All	0	7

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	37	GLY	N-CA-C	9.85	133.87	114.16
2	F	85	LYS	CA-C-N	9.19	134.27	120.31
2	F	85	LYS	C-N-CA	9.19	134.27	120.31
1	B	37	GLY	N-CA-C	7.56	135.23	113.30
2	E	23	ASP	CA-C-N	6.59	133.57	121.70
2	E	23	ASP	C-N-CA	6.59	133.57	121.70
1	C	47	ARG	N-CA-C	-6.54	96.88	110.80
1	C	43	THR	CA-C-N	6.17	132.81	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	43	THR	C-N-CA	6.17	132.81	121.70
1	C	37	GLY	CA-C-N	-5.92	113.66	122.36
1	C	37	GLY	C-N-CA	-5.92	113.66	122.36
2	D	43	ASP	N-CA-C	5.89	118.59	111.40
2	D	47	THR	N-CA-C	5.62	122.77	110.80
2	F	23	ASP	CB-CA-C	-5.45	99.74	110.10

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	46	TYR	Peptide
1	B	46	TYR	Peptide
1	C	37	GLY	Peptide
2	F	42	PHE	Peptide
2	F	67	GLN	Peptide
2	F	68	LEU	Peptide
2	F	69	ARG	Peptide

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	824	0	862	21	0
1	B	834	0	869	20	0
1	C	853	0	884	39	0
2	D	717	0	691	22	0
2	E	709	0	664	28	0
2	F	592	0	534	32	0
3	A	1	0	0	0	0
3	B	1	0	0	1	0
3	C	1	0	0	0	0
All	All	4532	0	4504	145	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 16.

All (145) 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:C:38:ARG:N	1:C:39:GLN:CB	2.08	1.15
2:F:43:ASP:CB	2:F:44:ALA:HB3	1.84	1.07
2:E:43:ASP:N	2:E:44:ALA:HB2	1.73	1.04
2:F:43:ASP:HB3	2:F:44:ALA:HB3	1.36	1.03
2:F:43:ASP:CG	2:F:44:ALA:HB3	1.83	1.03
1:C:39:GLN:N	1:C:40:LYS:HA	1.81	0.95
2:E:73:SER:CB	2:E:74:LEU:CB	2.47	0.93
2:F:69:ARG:HA	2:F:70:ARG:CB	1.98	0.93
1:B:36:PHE:CG	1:B:37:GLY:HA3	2.07	0.88
2:F:43:ASP:HB3	2:F:44:ALA:CB	2.08	0.82
1:C:37:GLY:C	1:C:39:GLN:CB	2.54	0.80
2:E:87:THR:HG23	2:E:88:PHE:CD1	2.17	0.79
2:E:42:PHE:C	2:E:44:ALA:HB2	2.10	0.76
1:C:38:ARG:H	1:C:39:GLN:CB	1.96	0.76
1:C:37:GLY:O	1:C:39:GLN:CB	2.34	0.75
2:F:44:ALA:HA	2:F:45:ARG:CB	2.16	0.74
1:B:36:PHE:CD2	1:B:37:GLY:HA3	2.23	0.73
2:F:42:PHE:HA	2:F:43:ASP:HB2	1.72	0.72
2:F:84:PHE:C	2:F:84:PHE:CD1	2.69	0.69
1:A:38:ARG:HB2	1:A:39:GLN:CB	2.24	0.67
2:D:25:LYS:HD3	2:D:39:LEU:HD22	1.75	0.66
1:C:54:GLU:CG	2:D:34:GLN:HE22	2.09	0.65
2:F:66:ARG:O	2:F:68:LEU:N	2.30	0.64
1:B:39:GLN:N	1:B:40:LYS:HA	2.13	0.64
1:C:37:GLY:CA	1:C:38:ARG:CB	2.75	0.64
1:C:47:ARG:O	1:C:47:ARG:HG3	1.98	0.64
2:F:22:GLY:CA	2:F:23:ASP:HB3	2.28	0.63
1:B:38:ARG:H	1:B:39:GLN:HB2	1.64	0.62
1:C:54:GLU:HG3	2:D:34:GLN:HE22	1.64	0.61
2:F:66:ARG:O	2:F:67:GLN:C	2.44	0.61
2:F:44:ALA:HB1	2:F:45:ARG:C	2.26	0.61
2:E:22:GLY:HA2	2:E:23:ASP:C	2.25	0.59
1:A:31:SER:HB3	1:B:33:VAL:HG12	1.83	0.59
2:F:84:PHE:CD1	2:F:84:PHE:O	2.55	0.59
2:F:84:PHE:C	2:F:84:PHE:HD1	2.11	0.59
1:A:92:PHE:CE1	1:C:64:VAL:HG11	2.38	0.58
2:F:44:ALA:CA	2:F:45:ARG:CB	2.79	0.58
1:C:38:ARG:HA	1:C:39:GLN:C	2.29	0.58
1:C:54:GLU:HG3	2:D:34:GLN:NE2	2.20	0.57
2:E:74:LEU:N	2:E:75:GLN:HB2	2.19	0.57
1:C:28:MET:CB	1:C:63:ILE:HG22	2.35	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:16:VAL:HG12	1:B:20:LEU:HD23	1.87	0.57
1:B:28:MET:CB	1:B:63:ILE:HG22	2.35	0.57
2:D:43:ASP:OD2	2:D:43:ASP:N	2.38	0.57
2:D:45:ARG:N	2:D:46:GLY:HA3	2.20	0.57
2:E:18:ILE:HD11	2:E:29:ALA:HB2	1.87	0.56
2:F:22:GLY:HA3	2:F:23:ASP:HB3	1.88	0.56
2:E:23:ASP:HB2	2:E:24:SER:CB	2.35	0.56
2:F:69:ARG:CA	2:F:70:ARG:CB	2.67	0.56
2:F:43:ASP:CB	2:F:44:ALA:CB	2.71	0.56
1:C:16:VAL:HG12	1:C:20:LEU:HD23	1.88	0.55
2:D:80:LEU:C	2:D:80:LEU:HD23	2.31	0.55
2:E:23:ASP:CB	2:E:24:SER:CB	2.84	0.55
1:A:28:MET:CB	1:A:63:ILE:HG22	2.37	0.55
2:F:18:ILE:HD11	2:F:29:ALA:HB2	1.87	0.55
2:E:44:ALA:HA	2:E:45:ARG:C	2.32	0.55
1:A:16:VAL:HG12	1:A:20:LEU:HD23	1.89	0.54
1:B:38:ARG:CB	1:B:39:GLN:HA	2.37	0.54
2:E:23:ASP:HB2	2:E:24:SER:CA	2.37	0.54
1:B:28:MET:HB2	1:B:63:ILE:HG22	1.88	0.54
1:B:36:PHE:CD1	1:B:37:GLY:HA3	2.42	0.54
2:D:21:PHE:N	2:D:22:GLY:HA3	2.22	0.54
1:B:42:GLN:HG2	1:B:53:VAL:HG13	1.90	0.54
2:D:18:ILE:HD11	2:D:29:ALA:HB2	1.90	0.53
2:E:23:ASP:CB	2:E:24:SER:HB3	2.38	0.53
1:C:38:ARG:N	1:C:39:GLN:CA	2.71	0.53
1:B:29:THR:HG23	1:B:62:GLU:HB2	1.91	0.53
2:D:21:PHE:CE2	2:D:60:LEU:HD22	2.44	0.53
1:A:38:ARG:CB	1:A:39:GLN:CB	2.87	0.53
1:C:37:GLY:HA2	1:C:38:ARG:CB	2.39	0.52
1:B:39:GLN:HG3	1:B:39:GLN:O	2.10	0.52
1:A:29:THR:HG23	1:A:62:GLU:HB2	1.90	0.52
1:A:28:MET:HB2	1:A:63:ILE:HG22	1.92	0.51
1:C:28:MET:HB2	1:C:63:ILE:HG22	1.92	0.51
1:B:38:ARG:N	1:B:39:GLN:HB2	2.25	0.51
2:F:66:ARG:C	2:F:68:LEU:H	2.17	0.51
1:B:39:GLN:O	1:B:39:GLN:CG	2.58	0.51
2:F:66:ARG:C	2:F:68:LEU:N	2.68	0.51
1:B:99:THR:HG22	1:C:92:PHE:HD1	1.75	0.51
2:E:25:LYS:HG3	2:E:41:ALA:HB2	1.93	0.51
1:C:29:THR:HG23	1:C:62:GLU:HB2	1.93	0.50
2:E:65:LEU:HD11	2:E:80:LEU:C	2.36	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:51:TYR:CD2	2:D:36:LEU:HD11	2.47	0.50
1:C:38:ARG:CA	1:C:39:GLN:C	2.85	0.50
2:D:25:LYS:HG2	2:D:41:ALA:HB2	1.94	0.49
1:B:31:SER:HB3	1:C:33:VAL:HG12	1.94	0.49
1:C:39:GLN:H	1:C:40:LYS:HA	1.72	0.49
2:F:43:ASP:HB3	2:F:44:ALA:CA	2.42	0.49
1:A:33:VAL:HG12	1:C:31:SER:HB3	1.94	0.49
1:C:38:ARG:CA	1:C:39:GLN:CB	2.91	0.48
2:E:43:ASP:HB3	2:E:44:ALA:HA	1.95	0.48
2:D:57:ALA:O	2:D:61:VAL:HG23	2.14	0.47
1:C:47:ARG:O	1:C:48:GLY:C	2.56	0.47
1:A:36:PHE:CD2	1:C:28:MET:HE3	2.50	0.47
1:C:46:TYR:O	1:C:47:ARG:HB3	2.15	0.46
2:E:41:ALA:O	2:E:48:ARG:N	2.47	0.46
1:C:41:GLY:HA2	1:C:42:GLN:CB	2.46	0.46
2:E:65:LEU:CD1	2:E:80:LEU:HD22	2.46	0.46
1:C:47:ARG:O	1:C:47:ARG:CG	2.63	0.46
2:D:21:PHE:N	2:D:22:GLY:CA	2.79	0.46
2:E:73:SER:CA	2:E:74:LEU:CB	2.94	0.46
2:E:83:VAL:O	2:E:87:THR:HG22	2.15	0.46
1:A:36:PHE:HB2	1:C:28:MET:HG2	1.98	0.46
1:C:42:GLN:H	1:C:43:THR:HA	1.80	0.46
1:A:64:VAL:HG11	1:B:92:PHE:CE1	2.50	0.46
2:E:34:GLN:OE1	2:E:34:GLN:N	2.49	0.45
1:B:40:LYS:H	1:B:86:ILE:HD11	1.82	0.45
2:E:65:LEU:HD12	2:E:80:LEU:HB3	1.97	0.45
2:F:57:ALA:O	2:F:61:VAL:HG23	2.16	0.45
1:A:38:ARG:H	1:A:39:GLN:C	2.25	0.45
1:B:9:ARG:HG3	3:B:2001:HOH:O	2.17	0.44
1:C:42:GLN:CB	1:C:43:THR:C	2.90	0.44
2:F:43:ASP:CG	2:F:44:ALA:CB	2.72	0.44
1:A:52:THR:HG22	2:F:34:GLN:O	2.17	0.44
2:D:10:PRO:HG2	2:F:47:THR:HG21	1.99	0.44
2:E:57:ALA:O	2:E:61:VAL:HG23	2.17	0.44
2:D:56:GLU:HA	2:D:59:MET:HE3	2.00	0.44
2:E:87:THR:HG23	2:E:88:PHE:CE1	2.53	0.44
2:F:22:GLY:CA	2:F:23:ASP:CB	2.96	0.44
1:A:92:PHE:CD1	1:C:64:VAL:HG11	2.53	0.43
2:E:23:ASP:OD2	2:E:23:ASP:N	2.50	0.43
1:C:43:THR:OG1	1:C:44:GLU:HA	2.18	0.43
2:D:43:ASP:OD2	2:D:47:THR:N	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:47:ARG:O	1:C:48:GLY:O	2.36	0.43
1:A:45:ARG:HB3	1:A:50:GLU:HB3	2.00	0.43
1:A:51:TYR:CD2	2:F:36:LEU:HD11	2.54	0.43
2:D:64:ARG:HA	2:D:67:GLN:HE21	1.85	0.42
2:E:23:ASP:HB2	2:E:24:SER:C	2.45	0.42
2:F:84:PHE:O	2:F:84:PHE:HD1	2.00	0.42
2:D:25:LYS:CD	2:D:39:LEU:HD22	2.45	0.42
2:E:43:ASP:H	2:E:44:ALA:HB2	1.70	0.42
2:F:44:ALA:HB1	2:F:45:ARG:CA	2.49	0.41
1:C:39:GLN:N	1:C:40:LYS:CA	2.64	0.41
2:D:43:ASP:N	2:D:44:ALA:HA	2.35	0.41
2:F:23:ASP:O	2:F:24:SER:CB	2.68	0.41
1:A:25:ILE:HG21	1:A:63:ILE:HD12	2.02	0.41
2:D:89:LEU:O	2:D:89:LEU:HD23	2.21	0.41
1:C:100:ILE:HD12	1:C:100:ILE:N	2.35	0.41
2:D:21:PHE:O	2:D:21:PHE:CG	2.74	0.41
1:A:47:ARG:O	1:A:48:GLY:O	2.39	0.40
1:A:92:PHE:HE1	1:C:64:VAL:HG21	1.86	0.40
1:A:100:ILE:HD12	1:A:100:ILE:N	2.36	0.40
2:E:68:LEU:HD11	2:E:80:LEU:HD12	2.02	0.40
2:F:64:ARG:O	2:F:68:LEU:CB	2.70	0.40
2:E:66:ARG:O	2:E:69:ARG:HB3	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	102/112 (91%)	95 (93%)	6 (6%)	1 (1%)	12	45
1	B	106/112 (95%)	95 (90%)	7 (7%)	4 (4%)	2	18
1	C	110/112 (98%)	101 (92%)	7 (6%)	2 (2%)	6	34

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	85/89 (96%)	77 (91%)	7 (8%)	1 (1%)	10	42
2	E	85/89 (96%)	74 (87%)	9 (11%)	2 (2%)	4	28
2	F	72/89 (81%)	65 (90%)	5 (7%)	2 (3%)	4	25
All	All	560/603 (93%)	507 (90%)	41 (7%)	12 (2%)	5	31

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	24	SER
2	F	67	GLN
1	A	48	GLY
1	B	38	ARG
1	C	48	GLY
2	E	74	LEU
1	B	47	ARG
1	C	47	ARG
2	E	24	SER
1	B	48	GLY
2	D	22	GLY
1	B	37	GLY

5.3.2 Protein sidechains ⓘ

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	88/93 (95%)	86 (98%)	2 (2%)	44	70
1	B	89/93 (96%)	87 (98%)	2 (2%)	45	71
1	C	89/93 (96%)	86 (97%)	3 (3%)	32	64
2	D	74/78 (95%)	62 (84%)	12 (16%)	2	12
2	E	69/78 (88%)	58 (84%)	11 (16%)	2	13
2	F	54/78 (69%)	45 (83%)	9 (17%)	2	11
All	All	463/513 (90%)	424 (92%)	39 (8%)	10	38

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	LEU
1	A	50	GLU
1	B	20	LEU
1	B	50	GLU
1	C	20	LEU
1	C	45	ARG
1	C	50	GLU
2	D	20	SER
2	D	26	GLU
2	D	31	LEU
2	D	34	GLN
2	D	39	LEU
2	D	43	ASP
2	D	54	ARG
2	D	62	ASP
2	D	65	LEU
2	D	70	ARG
2	D	74	LEU
2	D	87	THR
2	E	21	PHE
2	E	23	ASP
2	E	24	SER
2	E	26	GLU
2	E	31	LEU
2	E	39	LEU
2	E	47	THR
2	E	54	ARG
2	E	62	ASP
2	E	65	LEU
2	E	80	LEU
2	F	24	SER
2	F	31	LEU
2	F	39	LEU
2	F	42	PHE
2	F	54	ARG
2	F	62	ASP
2	F	65	LEU
2	F	66	ARG
2	F	69	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	69	GLN
1	B	98	GLN
1	C	69	GLN
1	C	98	GLN
2	D	5	ASN
2	D	34	GLN
2	D	67	GLN
2	E	5	ASN
2	F	5	ASN
2	F	8	ASN
2	F	63	ASN
2	F	67	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 oligosaccharides 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	106/112 (94%)	-0.19	0	100 100	8, 44, 117, 202	0
1	B	108/112 (96%)	-0.07	0	100 100	12, 50, 123, 185	1 (0%)
1	C	112/112 (100%)	-0.11	0	100 100	16, 43, 121, 160	0
2	D	87/89 (97%)	0.02	0	100 100	31, 68, 128, 205	0
2	E	87/89 (97%)	0.30	2 (2%)	61 41	38, 87, 161, 188	0
2	F	78/89 (87%)	0.35	1 (1%)	75 55	60, 102, 143, 176	0
All	All	578/603 (95%)	0.03	3 (0%)	87 76	8, 63, 148, 205	1 (0%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	70	ARG	3.3
2	E	68	LEU	2.9
2	F	88	PHE	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 oligosaccharides 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.