



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 07:47 PM GMT

PDB ID : 1GXP
Title : PHOB EFFECTOR DOMAIN IN COMPLEX WITH PHO BOX DNA.
Authors : Blanco, A.G.; Sola, M.; Gomis-Ruth, F.X.; Coll, M.
Deposited on : 2002-04-08
Resolution : 2.50 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

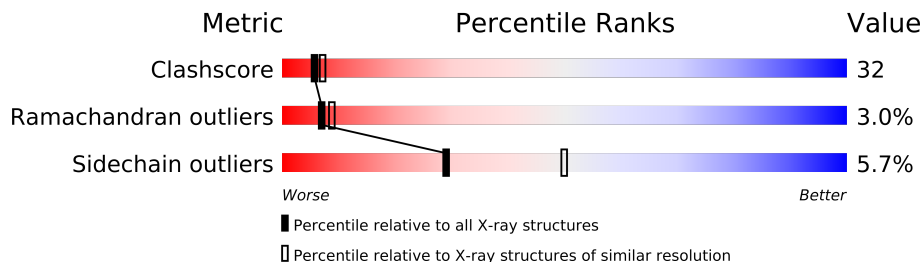
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.50 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	79885	3562 (2.50-2.50)
Ramachandran outliers	78287	3480 (2.50-2.50)
Sidechain outliers	78261	3482 (2.50-2.50)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	106	
1	B	106	
1	E	106	
1	F	106	
2	C	23	
2	G	23	
3	D	23	
3	H	23	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5416 atoms, of which 0 are hydrogen and 0 are deuterium.

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 PHOSPHATE REGULON TRANSCRIPTIONAL REGULATORY PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	103	Total	C	N	O	S	19	0	0
			846	531	156	154	5			
1	B	101	Total	C	N	O	S	20	0	0
			834	523	154	152	5			
1	E	102	Total	C	N	O	S	16	0	0
			841	528	155	153	5			
1	F	102	Total	C	N	O	S	22	0	0
			841	528	155	153	5			

- Molecule 2 is a DNA chain called 5'-D(*GP*AP*GP*CP*TP*GP*TP*CP*AP*TP*AP*AP*AP*GP*TP*TP*GP*TP*CP*AP*CP*GP*G)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	23	Total	C	N	O	P	20	0	0
			473	226	89	136	22			
2	G	23	Total	C	N	O	P	20	0	0
			473	226	89	136	22			

- Molecule 3 is a DNA chain called 5'-D(*CP*CP*CP*GP*TP*GP*AP*CP*AP*AP*CP*TP*TP*TP*AP*TP*GP*AP*CP*AP*GP*CP*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	23	Total	C	N	O	P	5	0	0
			464	223	83	136	22			
3	H	23	Total	C	N	O	P	8	0	0
			464	223	83	136	22			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	38	Total	O	0	0
			38	38		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	31	Total 31	O 31	0	0
4	C	8	Total 8	O 8	0	0
4	D	17	Total 17	O 17	0	0
4	E	41	Total 41	O 41	0	0
4	F	30	Total 30	O 30	0	0
4	G	6	Total 6	O 6	0	0
4	H	9	Total 9	O 9	0	0

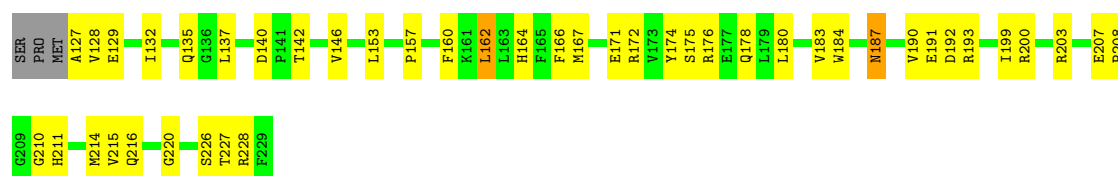
3 Residue-property plots

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.

Note EDS was not executed.

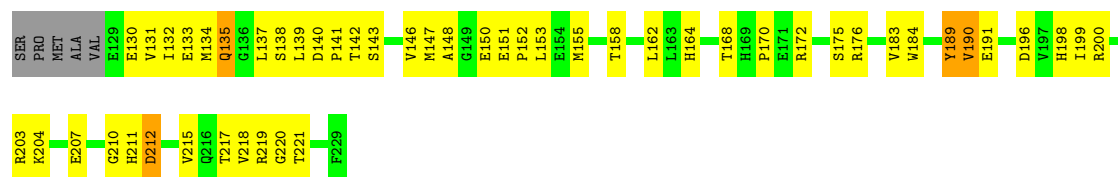
• Molecule 1: PHOSPHATE REGULON TRANSCRIPTIONAL REGULATORY PROTEIN

Chain A: 



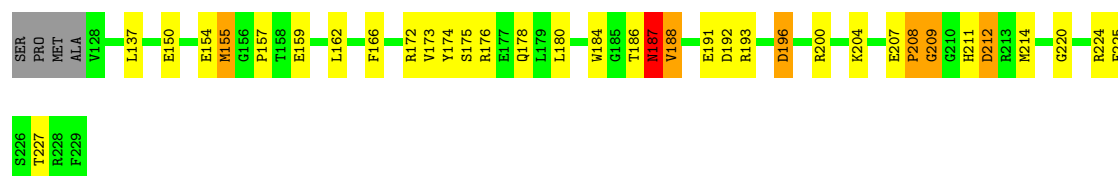
• Molecule 1: PHOSPHATE REGULON TRANSCRIPTIONAL REGULATORY PROTEIN

Chain B: 



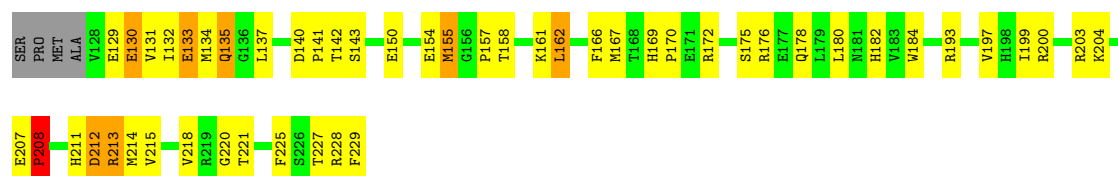
• Molecule 1: PHOSPHATE REGULON TRANSCRIPTIONAL REGULATORY PROTEIN

Chain E: 



• Molecule 1: PHOSPHATE REGULON TRANSCRIPTIONAL REGULATORY PROTEIN

Chain F: 



- Molecule 2: 5'-D(*GP*AP*GP*CP*TP*GP*TP*CP*AP*TP*AP*AP*AP*GP*TP*TP*GP*TP*CP*AP*CP*GP*G)-3'

Chain C: 

G1	A2	G3	C4	T5	G6	T7	C8	A9	T10	A11	A12	A13	G14	T15	T16	G17	T18	C19	A20	C21	G22	G23
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- Molecule 2: 5'-D(*GP*AP*GP*CP*TP*GP*TP*CP*AP*TP*AP*AP*AP*GP*TP*TP*GP*TP*CP*AP*CP*GP*G)-3'

Chain G: 

G1	A2	G3	C4	T5	G6	T7	C8	A9	T10	A11	A12	A13	G14	T15	T16	G17	T18	C19	A20	C21	G22	G23
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- Molecule 3: 5'-D(*CP*CP*CP*GP*TP*GP*AP*CP*AP*AP*CP*TP*TP*TP*AP*TP*GP*AP*CP*AP*GP*CP*T)-3'

Chain D: 

C1	C2	C3	G4	T5	G6	A7	C8	A9	A10	C11	T12	T13	T14	A15	T16	G17	A18	C19	A20	C21	C22	T23
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- Molecule 3: 5'-D(*CP*CP*CP*GP*TP*GP*AP*CP*AP*AP*CP*TP*TP*TP*AP*TP*GP*AP*CP*AP*GP*CP*T)-3'

Chain H: 

C1	C2	C3	G4	T5	G6	A7	C8	A9	A10	C11	T12	T13	T14	A15	T16	G17	A18	C19	A20	C21	C22	T23
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4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	74.11Å 74.11Å 289.68Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.00 – 2.50	Depositor
% Data completeness (in resolution range)	100.0 (40.00-2.50)	Depositor
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC	Depositor
R, R_{free}	0.244 , 0.289	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5416	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

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	0.43	0/866	0.69	0/1168
1	B	0.45	0/854	0.69	0/1151
1	E	0.44	0/861	0.71	0/1161
1	F	0.43	0/861	0.74	1/1161 (0.1%)
2	C	0.41	0/531	0.77	0/819
2	G	0.46	0/531	0.73	0/819
3	D	0.37	0/519	0.71	0/798
3	H	0.45	0/519	0.75	0/798
All	All	0.43	0/5542	0.72	1/7875 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	130	GLU	N-CA-C	6.24	127.85	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	846	0	825	42	0
1	B	834	0	811	46	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	841	0	820	35	0
1	F	841	0	820	49	0
2	C	473	0	261	35	0
2	G	473	0	261	33	0
3	D	464	0	261	38	0
3	H	464	0	261	37	0
4	A	38	0	0	3	0
4	B	31	0	0	4	0
4	C	8	0	0	1	0
4	D	17	0	0	5	0
4	E	41	0	0	4	0
4	F	30	0	0	3	0
4	G	6	0	0	0	0
4	H	9	0	0	2	0
All	All	5416	0	4320	294	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 32.

All (294) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:H:18:DA:H1'	3:H:19:DC:H5'	1.44	0.98
1:F:175:SER:H	1:F:178:GLN:NE2	1.61	0.97
2:C:21:DC:H1'	2:C:22:DG:H5'	1.45	0.94
1:F:175:SER:H	1:F:178:GLN:HE21	0.90	0.90
2:G:22:DG:H2''	2:G:23:DG:H5''	1.52	0.89
3:D:17:DG:H1'	3:D:18:DA:H5''	1.56	0.88
2:G:12:DA:H2''	2:G:13:DA:H5'	1.55	0.87
2:C:12:DA:H2''	2:C:13:DA:H5'	1.57	0.86
1:A:157:PRO:HG2	2:C:5:DT:OP1	1.75	0.86
1:F:175:SER:N	1:F:178:GLN:HE21	1.74	0.84
2:C:11:DA:H1'	2:C:12:DA:H5''	1.60	0.83
3:H:21:DG:H2''	3:H:22:DC:OP2	1.80	0.81
1:A:146:VAL:HG12	1:A:153:LEU:HD12	1.61	0.81
2:C:5:DT:H2''	2:C:6:DG:C8	2.17	0.80
2:G:3:DG:H1'	2:G:4:DC:H5''	1.64	0.79
3:D:7:DA:H3'	4:D:2009:HOH:O	1.82	0.79
2:G:15:DT:H2''	2:G:16:DT:H5'	1.65	0.78
2:C:18:DT:H2''	2:C:19:DC:H5'	1.66	0.77
1:F:197:VAL:HG21	2:G:18:DT:H72	1.66	0.77
3:D:10:DA:H1'	3:D:11:DC:H5'	1.66	0.77
2:G:10:DT:H2''	2:G:11:DA:OP2	1.86	0.76

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:D:18:DA:H1'	3:D:19:DC:H5'	1.67	0.76
1:B:146:VAL:HG12	1:B:153:LEU:HD12	1.65	0.76
2:C:4:DC:H2''	2:C:5:DT:H5'	1.68	0.75
2:C:3:DG:H1'	2:C:4:DC:H5''	1.67	0.75
1:B:175:SER:HB3	4:B:2028:HOH:O	1.88	0.74
2:C:7:DT:H2''	2:C:8:DC:H5'	1.70	0.73
1:F:200:ARG:NH2	3:H:6:DG:OP2	2.22	0.72
1:E:172:ARG:HD3	1:E:174:TYR:OH	1.89	0.72
1:F:131:VAL:HG13	1:F:140:ASP:HA	1.72	0.72
1:A:175:SER:H	1:A:178:GLN:HE21	1.37	0.71
1:A:175:SER:OG	1:A:178:GLN:HG3	1.91	0.70
1:B:158:THR:N	4:B:2012:HOH:O	2.23	0.70
3:H:6:DG:H2''	3:H:7:DA:OP2	1.92	0.70
3:D:15:DA:H1'	3:D:16:DT:H5''	1.75	0.69
1:A:129:GLU:HB3	4:A:2004:HOH:O	1.91	0.69
1:B:135:GLN:HE21	1:B:135:GLN:N	1.90	0.69
3:D:1:DC:H2''	3:D:2:DC:H5''	1.74	0.68
1:A:132:ILE:HG21	1:A:167:MET:HB3	1.76	0.68
2:G:22:DG:C2'	2:G:23:DG:H5''	2.25	0.67
2:G:21:DC:H2''	2:G:22:DG:OP2	1.95	0.67
3:D:18:DA:H2''	3:D:19:DC:O5'	1.95	0.67
1:E:214:MET:CE	1:E:227:THR:HG22	2.25	0.66
3:H:6:DG:H1'	3:H:7:DA:H5''	1.78	0.66
2:C:18:DT:H2''	2:C:19:DC:C5'	2.26	0.66
2:G:12:DA:C2'	2:G:13:DA:H5'	2.24	0.66
3:D:2:DC:O5'	3:D:2:DC:H6	1.79	0.66
3:H:10:DA:H5''	4:H:2004:HOH:O	1.97	0.65
1:F:135:GLN:CD	1:F:227:THR:HB	2.17	0.65
1:B:147:MET:HG2	1:B:152:PRO:HA	1.79	0.65
2:G:7:DT:H2''	2:G:8:DC:H5'	1.78	0.64
3:D:20:DA:H2''	3:D:21:DG:OP2	1.97	0.64
1:F:228:ARG:NH1	4:F:2029:HOH:O	2.30	0.64
2:C:15:DT:H1'	2:C:16:DT:H5'	1.81	0.63
3:H:10:DA:H2''	3:H:11:DC:C5'	2.28	0.63
3:H:13:DT:H1'	3:H:14:DT:H5''	1.79	0.63
2:C:12:DA:H2''	2:C:13:DA:C5'	2.28	0.63
1:A:172:ARG:HG3	1:B:142:THR:HG22	1.81	0.63
1:E:207:GLU:O	1:E:209:GLY:N	2.31	0.63
2:G:18:DT:H4'	2:G:19:DC:OP1	1.99	0.62
1:F:175:SER:OG	1:F:178:GLN:HG3	2.00	0.62
1:A:132:ILE:N	1:A:132:ILE:HD12	2.15	0.62
3:H:10:DA:H1'	3:H:11:DC:H5''	1.82	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:D:2:DC:H2''	3:D:3:DC:H5'	1.82	0.61
2:C:10:DT:H2''	2:C:11:DA:OP2	2.00	0.61
1:E:137:LEU:HB2	1:E:211:HIS:CD2	2.36	0.60
1:E:212:ASP:OD2	1:E:212:ASP:N	2.33	0.60
2:C:22:DG:H1'	2:C:23:DG:H5'	1.82	0.60
3:H:1:DC:H2'	4:H:2001:HOH:O	2.00	0.60
3:D:5:DT:H2''	3:D:6:DG:C8	2.36	0.60
2:C:4:DC:H2'	2:C:5:DT:H71	1.84	0.59
3:D:7:DA:H2''	3:D:8:DC:C5'	2.33	0.58
3:D:6:DG:H2''	3:D:7:DA:OP2	2.03	0.58
2:C:8:DC:H5'	2:C:8:DC:H6	1.69	0.58
2:G:18:DT:H1'	2:G:19:DC:C6	2.38	0.58
1:A:199:ILE:HG23	1:A:215:VAL:HG11	1.84	0.58
2:C:16:DT:H2''	2:C:17:DG:C8	2.39	0.57
1:B:219:ARG:NH1	4:B:2026:HOH:O	2.35	0.57
1:A:183:VAL:HB	1:A:184:TRP:CE3	2.38	0.57
1:F:137:LEU:HB2	1:F:211:HIS:CD2	2.40	0.57
1:E:214:MET:O	1:E:225:PHE:HA	2.05	0.57
1:B:212:ASP:N	1:B:212:ASP:OD2	2.38	0.56
3:D:7:DA:H1'	3:D:8:DC:H5''	1.88	0.56
3:H:7:DA:H5'	3:H:7:DA:H8	1.71	0.56
1:F:207:GLU:N	1:F:208:PRO:HD2	2.20	0.56
3:H:10:DA:H2''	3:H:11:DC:H5''	1.87	0.56
3:H:4:DG:H1'	3:H:5:DT:H5''	1.87	0.55
1:A:146:VAL:CG1	1:A:153:LEU:HD12	2.35	0.55
1:E:173:VAL:HG12	1:F:142:THR:O	2.06	0.55
3:H:10:DA:C2'	3:H:11:DC:H5''	2.37	0.55
3:H:18:DA:H1'	3:H:19:DC:C5'	2.30	0.54
3:D:7:DA:H2''	3:D:8:DC:H5''	1.90	0.54
1:E:175:SER:OG	1:E:178:GLN:HG3	2.08	0.54
2:G:16:DT:H2''	2:G:17:DG:C8	2.42	0.54
1:B:183:VAL:HB	1:B:184:TRP:CE3	2.42	0.54
1:A:207:GLU:O	1:A:210:GLY:N	2.31	0.54
3:D:17:DG:C1'	3:D:18:DA:H5''	2.34	0.54
3:D:17:DG:H2''	3:D:18:DA:OP2	2.07	0.54
3:D:21:DG:H5''	4:D:2017:HOH:O	2.07	0.54
3:D:21:DG:H1'	3:D:22:DC:H5''	1.89	0.54
2:C:7:DT:H2''	2:C:8:DC:C5'	2.36	0.54
2:C:17:DG:H2''	2:C:18:DT:H5''	1.90	0.53
2:C:20:DA:H1'	2:C:21:DC:H5'	1.89	0.53
1:E:172:ARG:HD3	1:E:174:TYR:CZ	2.43	0.53
3:D:7:DA:C2'	3:D:8:DC:H5''	2.39	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:H:5:DT:H2''	3:H:6:DG:C8	2.44	0.53
2:G:20:DA:H1'	2:G:21:DC:H5'	1.90	0.53
1:F:207:GLU:N	1:F:208:PRO:CD	2.72	0.53
1:B:199:ILE:HG23	1:B:215:VAL:CG1	2.38	0.53
1:F:182:HIS:HE1	4:F:2021:HOH:O	1.92	0.53
1:B:151:GLU:OE1	1:F:213:ARG:NH1	2.42	0.52
1:F:150:GLU:HG3	1:F:150:GLU:O	2.09	0.52
1:F:212:ASP:C	1:F:214:MET:H	2.13	0.52
1:B:130:GLU:HG3	1:B:131:VAL:H	1.74	0.52
3:D:1:DC:N4	4:D:2001:HOH:O	2.42	0.52
1:E:186:THR:HG22	1:E:187:ASN:N	2.24	0.52
1:F:193:ARG:NH2	3:H:3:DC:OP2	2.31	0.52
2:C:8:DC:H5'	2:C:8:DC:C6	2.45	0.52
3:H:7:DA:H2''	3:H:8:DC:O5'	2.10	0.52
1:B:204:LYS:HA	1:B:207:GLU:OE2	2.09	0.52
2:G:9:DA:H1'	2:G:10:DT:H5'	1.92	0.52
1:A:175:SER:H	1:A:178:GLN:NE2	2.06	0.52
2:C:19:DC:H2''	2:C:20:DA:C8	2.45	0.51
2:C:8:DC:H2''	2:C:9:DA:C8	2.45	0.51
1:F:162:LEU:HD22	1:F:166:PHE:CE2	2.46	0.51
2:C:3:DG:H8	4:C:2003:HOH:O	1.92	0.51
1:A:208:PRO:HA	4:A:2027:HOH:O	2.09	0.51
2:G:12:DA:H1'	2:G:13:DA:C5'	2.41	0.51
2:C:17:DG:C2'	2:C:18:DT:H5''	2.40	0.50
2:G:10:DT:H1'	2:G:11:DA:C8	2.47	0.50
1:B:218:VAL:HG12	1:B:221:THR:HB	1.94	0.50
1:E:180:LEU:O	1:E:184:TRP:HB2	2.11	0.50
2:G:11:DA:H1'	2:G:12:DA:H5''	1.94	0.50
3:D:21:DG:H2''	3:D:22:DC:C5'	2.41	0.50
1:A:207:GLU:HB2	1:A:208:PRO:HD3	1.94	0.50
1:B:200:ARG:NH2	3:D:6:DG:OP2	2.41	0.50
1:A:214:MET:HE3	1:A:227:THR:HG22	1.94	0.50
1:A:228:ARG:HE	1:B:147:MET:CE	2.25	0.50
1:B:139:LEU:C	1:B:139:LEU:HD23	2.32	0.50
2:C:21:DC:H1'	2:C:22:DG:C5'	2.31	0.50
3:D:17:DG:H1'	3:D:18:DA:C5'	2.35	0.50
1:B:146:VAL:CG1	1:B:153:LEU:HD12	2.38	0.50
1:A:199:ILE:HG23	1:A:215:VAL:CG1	2.41	0.50
1:F:229:PHE:N	1:F:229:PHE:CD2	2.77	0.50
2:C:7:DT:H6	2:C:7:DT:H5'	1.77	0.49
1:B:158:THR:HG21	1:B:198:HIS:CE1	2.48	0.49
1:A:214:MET:CE	1:A:227:THR:HG22	2.42	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:175:SER:N	1:F:178:GLN:NE2	2.45	0.49
2:C:17:DG:H2"	2:C:18:DT:C5'	2.43	0.49
1:A:183:VAL:HB	1:A:184:TRP:HE3	1.76	0.49
3:D:13:DT:H1'	3:D:14:DT:H5"	1.94	0.49
3:D:10:DA:H5"	4:D:2012:HOH:O	2.12	0.49
3:H:7:DA:H2"	3:H:8:DC:C5'	2.43	0.49
3:D:2:DC:H2"	3:D:3:DC:C5'	2.43	0.49
3:H:15:DA:H1'	3:H:16:DT:H5"	1.94	0.49
2:C:2:DA:H2"	2:C:3:DG:OP2	2.13	0.49
1:A:228:ARG:HG3	1:A:228:ARG:HH21	1.78	0.48
3:H:13:DT:H2"	3:H:14:DT:C5'	2.43	0.48
1:B:140:ASP:OD2	1:B:140:ASP:C	2.51	0.48
1:F:129:GLU:OE2	1:F:129:GLU:N	2.46	0.48
1:B:131:VAL:HG22	1:B:140:ASP:HB2	1.94	0.48
1:F:197:VAL:CG2	2:G:18:DT:H72	2.38	0.48
1:F:218:VAL:HG12	1:F:221:THR:HB	1.95	0.48
3:H:17:DG:H2"	3:H:18:DA:OP2	2.12	0.48
1:A:216:GLN:OE1	1:A:226:SER:HB3	2.14	0.48
1:B:147:MET:HG2	1:B:152:PRO:CA	2.44	0.48
3:H:13:DT:H2"	3:H:14:DT:H5'	1.94	0.48
3:D:16:DT:H2"	3:D:17:DG:C8	2.49	0.48
2:G:8:DC:H6	2:G:8:DC:H5'	1.78	0.48
1:A:190:VAL:HG12	1:A:191:GLU:N	2.29	0.47
1:B:175:SER:HA	1:B:221:THR:O	2.14	0.47
1:F:154:GLU:O	1:F:155:MET:HB3	2.14	0.47
1:E:214:MET:HE3	1:E:227:THR:HG22	1.96	0.47
2:G:14:DG:H1'	2:G:15:DT:H5"	1.96	0.47
1:A:216:GLN:HG2	4:A:2031:HOH:O	2.13	0.47
2:C:4:DC:C2'	2:C:5:DT:H71	2.44	0.47
1:E:157:PRO:HA	4:E:2009:HOH:O	2.14	0.47
1:E:155:MET:HA	4:E:2015:HOH:O	2.13	0.47
1:F:175:SER:HB2	4:F:2020:HOH:O	2.15	0.46
2:C:17:DG:H1'	2:C:18:DT:H5"	1.98	0.46
1:F:157:PRO:HG2	2:G:16:DT:OP1	2.15	0.46
3:H:6:DG:H1'	3:H:7:DA:C5'	2.44	0.46
1:E:154:GLU:O	1:E:155:MET:HB3	2.15	0.46
1:B:189:TYR:CD1	1:B:190:VAL:N	2.84	0.46
1:F:204:LYS:O	1:F:207:GLU:HG3	2.15	0.46
1:E:220:GLY:N	3:H:14:DT:OP1	2.48	0.46
1:E:207:GLU:HB2	1:E:208:PRO:HD3	1.95	0.46
1:E:150:GLU:HB3	4:E:2012:HOH:O	2.14	0.46
1:A:190:VAL:HG13	2:C:7:DT:OP2	2.15	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:200:ARG:NH2	1:F:204:LYS:NZ	2.63	0.46
3:H:10:DA:H2''	3:H:11:DC:H5'	1.96	0.46
1:E:186:THR:HG22	1:E:187:ASN:H	1.80	0.46
1:B:135:GLN:HE21	1:B:135:GLN:CA	2.28	0.46
1:E:207:GLU:N	1:E:208:PRO:CD	2.78	0.46
3:H:16:DT:H2''	3:H:17:DG:C8	2.50	0.46
3:H:7:DA:C8	3:H:7:DA:H5'	2.51	0.46
2:G:20:DA:H1'	2:G:21:DC:C5'	2.46	0.45
1:F:166:PHE:HB3	1:F:225:PHE:CB	2.46	0.45
1:A:160:PHE:C	1:A:160:PHE:CD2	2.90	0.45
1:A:228:ARG:HG2	1:B:150:GLU:OE1	2.16	0.45
2:G:4:DC:H2'	2:G:5:DT:H71	1.99	0.45
1:A:171:GLU:O	1:B:143:SER:HB3	2.16	0.45
1:F:203:ARG:NH1	1:F:215:VAL:O	2.45	0.45
1:B:176:ARG:HG3	1:B:176:ARG:HH11	1.82	0.45
1:B:176:ARG:NE	1:B:196:ASP:OD2	2.43	0.45
1:E:187:ASN:O	1:E:188:VAL:C	2.55	0.45
2:G:3:DG:C1'	2:G:4:DC:H5''	2.42	0.45
1:F:212:ASP:C	1:F:214:MET:N	2.70	0.45
1:F:197:VAL:HG21	2:G:18:DT:C7	2.43	0.45
1:B:135:GLN:NE2	1:B:135:GLN:CA	2.80	0.45
3:D:21:DG:H2''	3:D:22:DC:H5'	1.99	0.45
1:F:176:ARG:O	1:F:180:LEU:HG	2.17	0.45
1:A:162:LEU:HD22	1:A:166:PHE:CE2	2.52	0.45
3:D:4:DG:H1'	3:D:5:DT:H5''	1.99	0.44
2:G:12:DA:H1'	2:G:13:DA:H5'	2.00	0.44
1:B:199:ILE:HG23	1:B:215:VAL:HG11	1.99	0.44
1:F:140:ASP:HA	1:F:141:PRO:HD3	1.85	0.44
3:D:12:DT:H2''	3:D:13:DT:OP2	2.18	0.44
1:F:180:LEU:O	1:F:184:TRP:HB2	2.18	0.44
3:H:2:DC:H1'	3:H:3:DC:H5''	1.98	0.44
1:F:169:HIS:N	1:F:170:PRO:CD	2.80	0.44
3:H:15:DA:H2''	3:H:16:DT:C5'	2.48	0.44
1:B:191:GLU:HG3	1:B:191:GLU:O	2.17	0.43
1:F:131:VAL:HG13	1:F:140:ASP:CA	2.44	0.43
1:E:186:THR:O	1:E:188:VAL:N	2.51	0.43
1:A:220:GLY:N	3:D:14:DT:OP1	2.41	0.43
1:F:199:ILE:HG23	1:F:215:VAL:HG11	1.99	0.43
1:E:176:ARG:NE	1:E:196:ASP:OD2	2.51	0.43
1:A:140:ASP:OD2	1:A:142:THR:HB	2.18	0.43
1:F:133:GLU:C	1:F:134:MET:HG3	2.39	0.43
1:A:228:ARG:NH2	1:A:228:ARG:HG3	2.33	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:209:GLY:HA3	1:E:211:HIS:CE1	2.53	0.43
1:B:207:GLU:O	1:B:210:GLY:N	2.41	0.43
2:C:14:DG:H1'	2:C:15:DT:H5''	1.99	0.43
1:B:220:GLY:N	3:D:3:DC:OP1	2.47	0.43
1:A:137:LEU:HB2	1:A:211:HIS:CD2	2.54	0.43
1:B:203:ARG:HD3	4:D:2008:HOH:O	2.17	0.43
1:F:132:ILE:HG21	1:F:167:MET:HB3	2.01	0.43
3:H:11:DC:H2'	3:H:12:DT:H72	2.01	0.43
1:E:176:ARG:HG3	1:E:176:ARG:HH11	1.84	0.43
1:A:127:ALA:HA	1:A:164:HIS:HE1	1.83	0.43
1:F:134:MET:SD	1:F:225:PHE:HE1	2.42	0.43
3:H:15:DA:H2''	3:H:16:DT:H5'	2.01	0.42
2:C:19:DC:H5'	2:C:19:DC:H6	1.83	0.42
2:G:7:DT:H2''	2:G:8:DC:C5'	2.46	0.42
1:E:191:GLU:HB2	4:E:2031:HOH:O	2.18	0.42
1:B:146:VAL:HG11	1:B:155:MET:HE1	2.01	0.42
1:E:172:ARG:NH1	1:E:174:TYR:OH	2.53	0.42
3:D:2:DC:C6	3:D:2:DC:O5'	2.65	0.42
1:A:187:ASN:CG	1:A:187:ASN:O	2.57	0.42
1:A:132:ILE:CD1	1:A:132:ILE:N	2.82	0.42
1:B:183:VAL:HB	1:B:184:TRP:HE3	1.82	0.42
1:B:137:LEU:HB2	1:B:211:HIS:CD2	2.53	0.42
1:B:133:GLU:C	1:B:134:MET:HG3	2.39	0.42
3:D:15:DA:C1'	3:D:16:DT:H5''	2.48	0.42
1:A:172:ARG:HD3	1:A:174:TYR:OH	2.19	0.42
1:B:148:ALA:HB2	1:B:153:LEU:HD11	2.01	0.42
1:B:217:THR:HG22	1:B:218:VAL:N	2.34	0.42
1:B:132:ILE:O	1:B:138:SER:HA	2.20	0.42
1:E:204:LYS:HB2	1:E:204:LYS:HE3	1.73	0.42
2:G:8:DC:H2''	2:G:9:DA:C8	2.55	0.42
1:F:166:PHE:HB3	1:F:225:PHE:HB2	2.02	0.42
1:B:131:VAL:HG22	1:B:140:ASP:CB	2.50	0.42
3:H:10:DA:C1'	3:H:11:DC:H5''	2.47	0.42
3:D:13:DT:H2''	3:D:14:DT:OP2	2.18	0.42
2:G:19:DC:H5'	2:G:19:DC:H6	1.84	0.41
1:F:214:MET:O	1:F:225:PHE:HA	2.20	0.41
1:E:193:ARG:HA	1:E:196:ASP:OD2	2.20	0.41
2:G:20:DA:H2''	2:G:21:DC:OP2	2.20	0.41
1:A:190:VAL:CG1	1:A:191:GLU:N	2.83	0.41
1:E:224:ARG:NH1	1:F:143:SER:O	2.46	0.41
1:A:203:ARG:O	1:A:207:GLU:HG3	2.20	0.41
1:B:140:ASP:HA	1:B:141:PRO:HD3	1.83	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:180:LEU:HD23	1:A:180:LEU:HA	1.74	0.41
2:C:15:DT:H2''	2:C:16:DT:O5'	2.20	0.41
1:B:170:PRO:HA	4:B:2016:HOH:O	2.20	0.41
1:A:200:ARG:NH1	3:D:17:DG:OP2	2.53	0.41
1:F:140:ASP:HB3	1:F:143:SER:OG	2.21	0.41
1:F:220:GLY:N	3:H:3:DC:OP1	2.51	0.41
1:E:166:PHE:HB3	1:E:225:PHE:CG	2.56	0.41
3:H:7:DA:H2''	3:H:8:DC:H5'	2.02	0.41
1:E:207:GLU:C	1:E:209:GLY:H	2.21	0.41
1:F:158:THR:O	1:F:161:LYS:HB2	2.20	0.41
1:B:164:HIS:HD2	1:B:164:HIS:O	2.04	0.41
1:F:207:GLU:HA	1:F:212:ASP:OD2	2.21	0.40
1:E:186:THR:C	1:E:188:VAL:H	2.24	0.40
3:H:22:DC:H2''	3:H:23:DT:OP2	2.22	0.40
3:H:7:DA:H1'	3:H:8:DC:H5'	2.03	0.40
2:G:3:DG:H1'	2:G:4:DC:C5'	2.41	0.40
1:A:176:ARG:NH2	1:A:193:ARG:NH2	2.70	0.40
2:G:21:DC:C2'	2:G:22:DG:OP2	2.68	0.40
1:E:176:ARG:HD3	1:E:192:ASP:HB3	2.03	0.40
3:D:10:DA:H2''	3:D:11:DC:OP2	2.22	0.40
1:E:166:PHE:HB3	1:E:225:PHE:CB	2.51	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	101/106 (95%)	89 (88%)	10 (10%)	2 (2%)	11	17
1	B	99/106 (93%)	90 (91%)	8 (8%)	1 (1%)	22	38
1	E	100/106 (94%)	86 (86%)	9 (9%)	5 (5%)	3	3
1	F	100/106 (94%)	87 (87%)	9 (9%)	4 (4%)	5	5
All	All	400/424 (94%)	352 (88%)	36 (9%)	12 (3%)	7	9

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	130	GLU
1	A	192	ASP
1	B	190	VAL
1	E	187	ASN
1	E	208	PRO
1	F	208	PRO
1	E	155	MET
1	E	209	GLY
1	A	128	VAL
1	F	155	MET
1	F	213	ARG
1	E	188	VAL

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	92/95 (97%)	89 (97%)	3 (3%)	50	76
1	B	91/95 (96%)	85 (93%)	6 (7%)	24	41
1	E	92/95 (97%)	86 (94%)	6 (6%)	24	42
1	F	92/95 (97%)	86 (94%)	6 (6%)	24	42
All	All	367/380 (97%)	346 (94%)	21 (6%)	29	50

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	135	GLN
1	A	162	LEU
1	A	187	ASN
1	B	135	GLN
1	B	162	LEU
1	B	168	THR
1	B	172	ARG
1	B	189	TYR
1	B	212	ASP

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Mol	Chain	Res	Type
1	E	159	GLU
1	E	162	LEU
1	E	187	ASN
1	E	196	ASP
1	E	200	ARG
1	E	212	ASP
1	F	133	GLU
1	F	135	GLN
1	F	162	LEU
1	F	172	ARG
1	F	208	PRO
1	F	212	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	164	HIS
1	A	178	GLN
1	B	135	GLN
1	B	164	HIS
1	B	178	GLN
1	E	164	HIS
1	E	178	GLN
1	F	135	GLN
1	F	178	GLN
1	F	181	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.