Full wwPDB X-ray Structure Validation Report ⓘ

Sep 17, 2023 – 10:30 AM EDT

PDB ID	:	4YX7
Title	:	Complex of SpaO(SPOA1,2) and OrgB(APAR)::T4lysozyme fusion protein
Authors	:	Notti, R.Q.; Stebbins, C.E.
Deposited on	:	2015-03-22
Resolution	:	2.00 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:

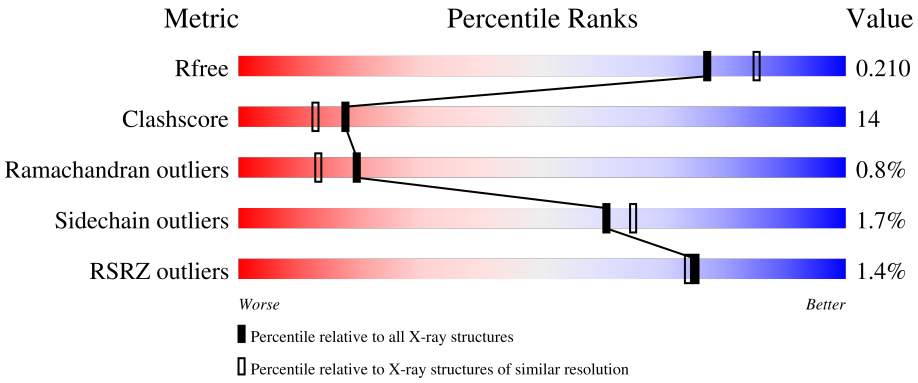
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	73	62% 26% • 11%
1	D	73	71% 18% 11%
2	B	70	74% 20% • •
2	E	70	% 69% 24% • 6%
3	C	197	% 69% 30% ••

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Mol	Chain	Length	Quality of chain
3	F	197	

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 5769 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 Surface presentation of antigens protein SpaO.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	65	Total	C	N	O	S	0	0	0
			494	316	90	86	2			
1	D	65	Total	C	N	O	S	0	1	0
			514	329	94	89	2			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP P40699
A	2	PRO	-	expression tag	UNP P40699
A	3	VAL	-	expression tag	UNP P40699
A	4	ASP	-	expression tag	UNP P40699
D	1	GLY	-	expression tag	UNP P40699
D	2	PRO	-	expression tag	UNP P40699
D	3	VAL	-	expression tag	UNP P40699
D	4	ASP	-	expression tag	UNP P40699

- Molecule 2 is a protein called Surface presentation of antigens protein SpaO.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	67	Total	C	N	O	S	0	0	0
			522	332	85	102	3			
2	E	66	Total	C	N	O	S	0	0	0
			515	327	84	101	3			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	GLY	-	expression tag	UNP P40699
B	2	PRO	-	expression tag	UNP P40699
B	3	VAL	-	expression tag	UNP P40699
B	4	ASP	-	expression tag	UNP P40699

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Chain	Residue	Modelled	Actual	Comment	Reference
E	1	GLY	-	expression tag	UNP P40699
E	2	PRO	-	expression tag	UNP P40699
E	3	VAL	-	expression tag	UNP P40699
E	4	ASP	-	expression tag	UNP P40699

- Molecule 3 is a protein called Oxygen-regulated invasion protein OrgB,Endolysin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	196	Total	C	N	O	S	0	0	0
			1543	981	279	278	5			
3	F	195	Total	C	N	O	S	0	0	0
			1524	967	274	278	5			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	1	GLY	-	expression tag	UNP P0CL45
C	2	PRO	-	expression tag	UNP P0CL45
C	3	VAL	-	expression tag	UNP P0CL45
C	4	ASP	-	expression tag	UNP P0CL45
C	45	GLY	ARG	conflict	UNP P00720
C	53	ASN	ASP	engineered mutation	UNP P00720
C	87	THR	CYS	engineered mutation	UNP P00720
C	130	ALA	CYS	engineered mutation	UNP P00720
C	170	ARG	ILE	conflict	UNP P00720
C	195	ALA	LYS	engineered mutation	UNP P00720
C	196	ALA	ASN	engineered mutation	UNP P00720
C	197	ALA	LEU	engineered mutation	UNP P00720
F	1	GLY	-	expression tag	UNP P0CL45
F	2	PRO	-	expression tag	UNP P0CL45
F	3	VAL	-	expression tag	UNP P0CL45
F	4	ASP	-	expression tag	UNP P0CL45
F	45	GLY	ARG	conflict	UNP P00720
F	53	ASN	ASP	engineered mutation	UNP P00720
F	87	THR	CYS	engineered mutation	UNP P00720
F	130	ALA	CYS	engineered mutation	UNP P00720
F	170	ARG	ILE	conflict	UNP P00720
F	195	ALA	LYS	engineered mutation	UNP P00720
F	196	ALA	ASN	engineered mutation	UNP P00720
F	197	ALA	LEU	engineered mutation	UNP P00720

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	55	Total O 55 55	0	0
4	B	65	Total O 65 65	0	0
4	C	218	Total O 218 218	0	0
4	D	67	Total O 67 67	0	0
4	E	76	Total O 76 76	0	0
4	F	176	Total O 176 176	0	0

3 Residue-property plots (i)

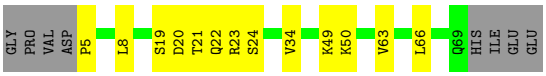
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Surface presentation of antigens protein SpaO

Chain A:

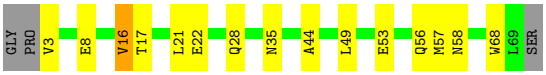
• Molecule 1: Surface presentation of antigens protein SpaO

Chain D:



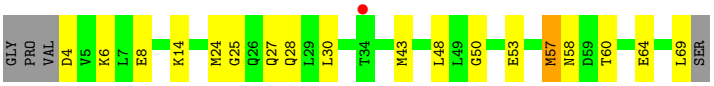
• Molecule 2: Surface presentation of antigens protein SpaO

Chain B:



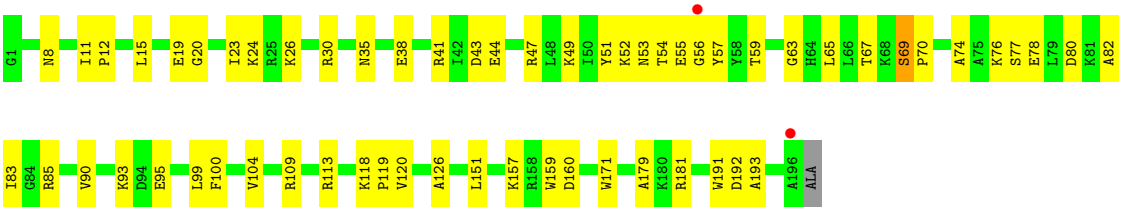
• Molecule 2: Surface presentation of antigens protein SpaO

Chain E:



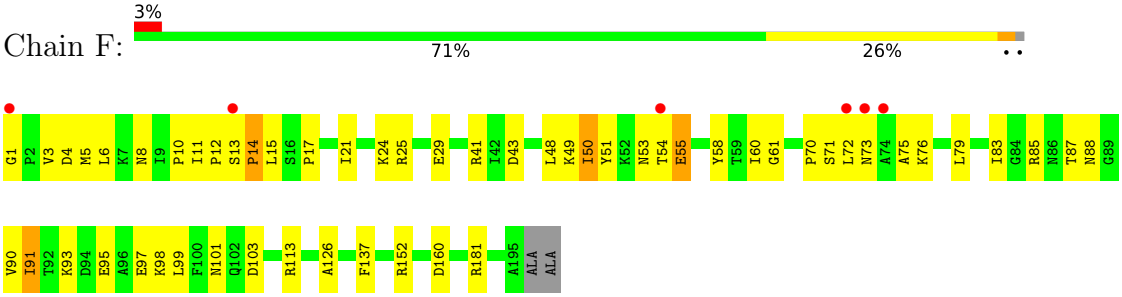
• Molecule 3: Oxygen-regulated invasion protein OrgB,Endolysin

Chain C:



• Molecule 3: Oxygen-regulated invasion protein OrgB,Endolysin

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	62.09Å 89.07Å 62.09Å 90.00° 114.94° 90.00°	Depositor
Resolution (Å)	47.59 – 2.00 47.59 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.4 (47.59-2.00) 99.4 (47.59-2.00)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.98 (at 2.00Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.157 , 0.210 0.159 , 0.210	Depositor DCC
R_{free} test set	1424 reflections (3.46%)	wwPDB-VP
Wilson B-factor (Å^2)	24.3	Xtriage
Anisotropy	0.157	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 43.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.43$, $< L^2 > = 0.25$	Xtriage
Estimated twinning fraction	0.226 for l,-k,h	Xtriage
Reported twinning fraction	0.230 for l,-k,h	Depositor
Outliers	0 of 41181 reflections	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	5769	wwPDB-VP
Average B, all atoms (Å^2)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.95% 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

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	0.37	0/501	0.63	1/676 (0.1%)
1	D	0.35	0/522	0.55	0/703
2	B	0.45	0/527	0.61	0/715
2	E	0.40	0/520	0.54	0/705
3	C	0.38	0/1570	0.57	0/2121
3	F	0.38	0/1551	0.59	0/2099
All	All	0.39	0/5191	0.58	1/7019 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	22	GLN	N-CA-C	-5.84	95.23	111.00

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	494	0	498	20	0
1	D	514	0	529	12	0
2	B	522	0	529	12	0
2	E	515	0	520	17	0
3	C	1543	0	1590	47	0
3	F	1524	0	1543	41	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	55	0	0	4	1
4	B	65	0	0	6	0
4	C	218	0	0	14	8
4	D	67	0	0	4	4
4	E	76	0	0	10	4
4	F	176	0	0	13	7
All	All	5769	0	5209	141	12

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 14.

All (141) 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)
3:F:113:ARG:NH2	4:F:201:HOH:O	1.96	0.97
3:C:192:ASP:OD2	4:C:201:HOH:O	1.92	0.85
2:E:4:ASP:OD1	4:E:101:HOH:O	1.96	0.83
3:C:43:ASP:OD1	3:C:181:ARG:NH1	2.11	0.83
3:F:41:ARG:NH1	4:F:202:HOH:O	1.99	0.82
2:B:35:ASN:OD1	4:B:101:HOH:O	2.00	0.78
1:A:43:GLU:OE2	4:A:101:HOH:O	2.01	0.77
3:C:74:ALA:HA	3:C:77:SER:HB2	1.66	0.76
3:F:98:LYS:NZ	4:F:209:HOH:O	2.20	0.74
1:A:18:SER:OG	1:A:41:ARG:NH1	2.20	0.73
3:F:3:VAL:O	4:F:204:HOH:O	2.05	0.73
2:E:53:GLU:OE2	4:E:102:HOH:O	2.06	0.73
2:B:53:GLU:OE2	4:B:102:HOH:O	2.08	0.72
3:F:88:ASN:OD1	4:F:205:HOH:O	2.07	0.72
2:B:53:GLU:OE1	4:B:103:HOH:O	2.08	0.72
4:D:157:HOH:O	2:E:60:THR:HG21	1.91	0.70
2:E:58:ASN:OD1	4:E:103:HOH:O	2.10	0.69
3:F:160:ASP:OD2	4:F:206:HOH:O	2.09	0.69
3:C:113:ARG:NH2	4:C:210:HOH:O	2.26	0.69
1:D:5:PRO:HA	1:D:8:LEU:HD13	1.75	0.68
3:C:160:ASP:OD2	4:C:203:HOH:O	2.12	0.68
3:C:53:ASN:OD1	3:C:57:TYR:N	2.26	0.68
3:F:160:ASP:OD2	4:F:207:HOH:O	2.11	0.67
2:E:28:GLN:OE1	4:E:104:HOH:O	2.13	0.67
1:A:27:GLY:HA3	1:D:23:ARG:HD2	1.77	0.67
3:F:103:ASP:OD1	4:F:208:HOH:O	2.14	0.66
3:C:30:ARG:NH2	4:C:205:HOH:O	2.17	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:64:GLU:OE1	4:E:105:HOH:O	2.14	0.65
1:D:50:LYS:NZ	4:D:103:HOH:O	2.29	0.65
1:A:53:HIS:ND1	4:A:102:HOH:O	2.30	0.65
3:F:1:GLY:N	4:F:214:HOH:O	2.29	0.64
1:D:22:GLN:NE2	1:D:24:SER:OG	2.26	0.63
3:C:41:ARG:NH1	4:C:214:HOH:O	2.31	0.63
1:A:52:GLY:HA3	1:A:65:THR:HB	1.80	0.63
3:C:54:THR:HG23	4:C:354:HOH:O	1.99	0.62
3:F:85:ARG:HE	3:F:87:THR:HA	1.65	0.62
1:A:64:GLU:HG2	4:A:110:HOH:O	1.99	0.62
2:E:6:LYS:NZ	4:E:106:HOH:O	2.16	0.62
3:F:72:LEU:O	3:F:76:LYS:N	2.29	0.61
3:F:11:ILE:HG22	3:F:14:PRO:HB2	1.83	0.60
2:B:56:GLN:NE2	2:B:57:MET:O	2.32	0.59
1:A:38:ARG:NH1	3:C:20:GLY:O	2.35	0.58
3:F:43:ASP:OD1	3:F:181:ARG:NH1	2.37	0.58
2:E:50:GLY:HA2	2:E:69:LEU:HD13	1.87	0.57
3:F:83:ILE:HG13	3:F:91:ILE:HD11	1.87	0.56
3:C:49:LYS:HG2	3:C:90:VAL:HG22	1.87	0.56
3:F:25:ARG:NH1	3:F:29:GLU:OE2	2.38	0.56
3:C:8:ASN:HD21	3:C:126:ALA:HB3	1.71	0.56
1:A:23:ARG:HE	2:B:3:VAL:HG23	1.70	0.55
2:E:60:THR:HG23	4:E:141:HOH:O	2.07	0.55
1:D:19:SER:HB3	3:F:21:ILE:HD12	1.87	0.55
2:B:28:GLN:NE2	4:B:107:HOH:O	2.31	0.55
1:D:20[A]:ASP:OD2	4:D:102:HOH:O	2.18	0.55
3:F:72:LEU:HA	3:F:75:ALA:HB3	1.88	0.54
3:F:48:LEU:HD23	3:F:90:VAL:HG13	1.88	0.54
3:C:109:ARG:O	3:C:113:ARG:HG3	2.08	0.54
3:C:74:ALA:O	3:C:78:GLU:HG3	2.08	0.54
3:C:83:ILE:HD12	3:C:95:GLU:HB3	1.90	0.54
2:E:8:GLU:OE2	4:E:107:HOH:O	2.18	0.54
3:F:93:LYS:NZ	3:F:97:GLU:OE2	2.41	0.53
1:A:56:ARG:NH2	4:A:105:HOH:O	2.41	0.53
3:C:38:GLU:HG2	4:C:357:HOH:O	2.10	0.51
3:F:71:SER:OG	3:F:73:ASN:O	2.25	0.51
3:C:55:GLU:N	3:C:56:GLY:HA2	2.26	0.51
3:C:181:ARG:HG2	3:C:193:ALA:HB1	1.94	0.50
3:F:13:SER:N	3:F:14:PRO:CD	2.75	0.50
3:F:50:ILE:CB	3:F:51:TYR:HA	2.42	0.50
2:E:25:GLY:O	2:E:27:GLN:NE2	2.45	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:28:ARG:HG2	1:A:28:ARG:HH11	1.77	0.49
3:F:152:ARG:NH1	4:F:223:HOH:O	2.45	0.49
3:C:26:LYS:HE3	4:C:331:HOH:O	2.11	0.49
2:B:16:VAL:HG11	2:B:21:LEU:HD13	1.95	0.49
3:C:85:ARG:NE	3:C:95:GLU:OE2	2.24	0.49
3:F:4:ASP:HB2	3:F:5:MET:HE2	1.94	0.49
1:D:63:VAL:HG11	1:D:66:LEU:HD12	1.95	0.49
1:A:26:LEU:HD22	2:B:68:TRP:HZ2	1.77	0.49
3:C:59:THR:HG22	3:C:65:LEU:HA	1.93	0.48
2:B:8:GLU:OE2	4:B:104:HOH:O	2.20	0.48
1:A:29:ILE:HD13	1:A:35:LEU:HD21	1.95	0.48
3:C:67:THR:HB	3:C:69:SER:H	1.78	0.48
3:F:54:THR:HG23	3:F:55:GLU:HG2	1.95	0.48
1:D:22:GLN:HB2	4:E:124:HOH:O	2.14	0.47
2:B:44:ALA:HB2	2:B:49:LEU:HD11	1.95	0.47
3:C:93:LYS:NZ	4:C:224:HOH:O	2.46	0.47
3:C:8:ASN:HB3	3:C:191:TRP:HE1	1.79	0.47
3:F:101:ASN:ND2	4:F:224:HOH:O	2.46	0.47
3:F:79:LEU:O	3:F:83:ILE:HG12	2.15	0.47
1:A:25:LEU:HD11	3:C:19:GLU:HG3	1.97	0.46
3:C:8:ASN:HB2	4:C:233:HOH:O	2.15	0.46
2:B:16:VAL:HG13	2:B:17:THR:O	2.15	0.46
3:C:69:SER:HA	3:C:70:PRO:HD3	1.85	0.46
3:F:8:ASN:O	3:F:10:PRO:HD3	2.16	0.46
3:F:8:ASN:HD21	3:F:126:ALA:HB3	1.81	0.45
1:A:26:LEU:CD2	1:A:29:ILE:HD12	2.47	0.45
3:C:47:ARG:HG3	3:C:51:TYR:CE1	2.52	0.45
3:F:11:ILE:CG2	3:F:14:PRO:HB2	2.46	0.45
1:D:34:VAL:HG11	2:E:57:MET:HE1	1.99	0.45
1:A:21:THR:HG23	1:A:25:LEU:HD22	1.99	0.44
3:F:14:PRO:CB	3:F:15:LEU:HB2	2.47	0.44
3:C:44:GLU:OE1	3:C:63:GLY:HA3	2.17	0.44
1:A:28:ARG:O	3:C:24:LYS:NZ	2.50	0.44
2:E:14:LYS:HD2	2:E:30:LEU:HD11	2.00	0.44
1:A:26:LEU:HA	1:A:27:GLY:HA2	1.60	0.44
1:D:21:THR:HA	4:D:110:HOH:O	2.17	0.44
2:E:57:MET:HE2	2:E:57:MET:HB2	1.66	0.44
3:F:15:LEU:O	3:F:17:PRO:HD3	2.17	0.43
3:C:126:ALA:HB2	4:C:380:HOH:O	2.18	0.43
3:C:8:ASN:ND2	3:C:126:ALA:HB3	2.33	0.43
1:D:22:GLN:HE21	1:D:24:SER:HG	1.58	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:8:ASN:ND2	3:F:126:ALA:HB3	2.34	0.43
2:E:24:MET:HE3	2:E:30:LEU:HD22	2.00	0.43
2:E:43:MET:HG2	2:E:48:LEU:HA	2.01	0.43
3:C:82:ALA:HB1	3:C:99:LEU:HD11	2.01	0.43
3:C:109:ARG:HD3	4:C:278:HOH:O	2.18	0.43
3:F:58:TYR:HD2	4:F:244:HOH:O	2.01	0.43
3:C:120:VAL:HG21	3:C:151:LEU:HB3	2.01	0.42
3:C:157:LYS:HG2	3:C:159:TRP:CZ2	2.55	0.42
1:D:49:LYS:HD3	1:D:49:LYS:HA	1.70	0.42
3:F:60:ILE:HG13	3:F:61:GLY:H	1.84	0.42
3:C:171:TRP:CZ2	3:C:179:ALA:HA	2.54	0.42
3:F:91:ILE:HG13	3:F:95:GLU:HG3	2.02	0.42
3:C:11:ILE:HG22	3:C:12:PRO:O	2.19	0.42
3:F:60:ILE:HG13	3:F:61:GLY:N	2.35	0.42
3:F:76:LYS:HA	3:F:79:LEU:HD12	2.01	0.42
3:C:52:LYS:NZ	4:C:225:HOH:O	2.52	0.42
3:C:76:LYS:O	3:C:80:ASP:HB2	2.20	0.42
3:C:118:LYS:HB3	3:C:119:PRO:HD3	2.02	0.41
2:B:22:GLU:HG3	4:B:137:HOH:O	2.21	0.41
3:C:15:LEU:HD23	3:C:23:ILE:HD11	2.03	0.41
3:C:53:ASN:OD1	3:C:56:GLY:HA2	2.20	0.41
1:A:38:ARG:HA	1:A:38:ARG:HD2	1.88	0.41
1:A:66:LEU:HD23	1:A:66:LEU:HA	1.85	0.41
3:C:109:ARG:NH1	4:C:231:HOH:O	2.54	0.41
3:C:24:LYS:HA	3:C:24:LYS:HD2	1.95	0.41
3:F:3:VAL:HA	3:F:6:LEU:HG	2.03	0.41
1:A:21:THR:HG22	1:A:22:GLN:O	2.21	0.40
3:C:100:PHE:O	3:C:104:VAL:HG23	2.22	0.40
3:F:99:LEU:HD23	3:F:99:LEU:HA	1.78	0.40
3:C:35:ASN:OD1	3:C:38:GLU:HG3	2.21	0.40
2:E:28:GLN:HB2	4:E:104:HOH:O	2.21	0.40
3:F:24:LYS:NZ	4:F:203:HOH:O	2.03	0.40

All (12) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:C:221:HOH:O	4:D:122:HOH:O[2_547]	1.98	0.22
4:C:305:HOH:O	4:F:316:HOH:O[2_546]	2.01	0.19
4:C:386:HOH:O	4:D:141:HOH:O[2_547]	2.02	0.18

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:C:395:HOH:O	4:F:343:HOH:O[2_546]	2.05	0.15
4:D:161:HOH:O	4:F:373:HOH:O[2_646]	2.06	0.14
4:C:232:HOH:O	4:E:109:HOH:O[2_547]	2.10	0.10
4:A:106:HOH:O	4:F:271:HOH:O[2_646]	2.14	0.06
4:C:332:HOH:O	4:E:151:HOH:O[2_647]	2.14	0.06
4:E:150:HOH:O	4:F:333:HOH:O[2_646]	2.14	0.06
4:C:396:HOH:O	4:D:157:HOH:O[2_547]	2.15	0.05
4:E:123:HOH:O	4:F:282:HOH:O[2_646]	2.15	0.05
4:C:405:HOH:O	4:F:321:HOH:O[2_546]	2.17	0.03

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	63/73 (86%)	56 (89%)	7 (11%)	0	100 100
1	D	64/73 (88%)	58 (91%)	6 (9%)	0	100 100
2	B	65/70 (93%)	61 (94%)	4 (6%)	0	100 100
2	E	64/70 (91%)	61 (95%)	3 (5%)	0	100 100
3	C	194/197 (98%)	187 (96%)	7 (4%)	0	100 100
3	F	193/197 (98%)	181 (94%)	7 (4%)	5 (3%)	5 2
All	All	643/680 (95%)	604 (94%)	34 (5%)	5 (1%)	19 13

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	F	12	PRO
3	F	14	PRO
3	F	50	ILE
3	F	53	ASN
3	F	70	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	50/63 (79%)	49 (98%)	1 (2%)	55	58
1	D	54/63 (86%)	54 (100%)	0	100	100
2	B	58/60 (97%)	56 (97%)	2 (3%)	37	36
2	E	57/60 (95%)	56 (98%)	1 (2%)	59	63
3	C	162/165 (98%)	161 (99%)	1 (1%)	86	90
3	F	158/165 (96%)	154 (98%)	4 (2%)	47	49
All	All	539/576 (94%)	530 (98%)	9 (2%)	60	65

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	SER
2	B	16	VAL
2	B	58	ASN
3	C	69	SER
2	E	57	MET
3	F	49	LYS
3	F	55	GLU
3	F	91	ILE
3	F	137	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
3	C	8	ASN
1	D	22	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

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, 95^{th} 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	65/73 (89%)	-0.38	0	100	100	16, 34, 61, 81	0
1	D	65/73 (89%)	-0.48	0	100	100	16, 29, 61, 74	0
2	B	67/70 (95%)	-0.51	0	100	100	13, 26, 61, 63	0
2	E	66/70 (94%)	-0.59	1 (1%)	73	72	16, 28, 50, 58	0
3	C	196/197 (99%)	-0.58	2 (1%)	82	81	12, 26, 59, 71	0
3	F	195/197 (98%)	-0.30	6 (3%)	49	48	14, 32, 73, 95	0
All	All	654/680 (96%)	-0.46	9 (1%)	75	74	12, 29, 63, 95	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	F	74	ALA	3.7
3	F	72	LEU	3.4
3	F	73	ASN	3.3
3	F	13	SER	2.9
3	C	56	GLY	2.8
3	C	196	ALA	2.7
3	F	1	GLY	2.4
3	F	54	THR	2.1
2	E	34	THR	2.0

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

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

6.3 Carbohydrates

There are no monosaccharides in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.