Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 09:06 PM GMT

PDB ID	:	4UW5
Title	:	Human galectin-7 in complex with a galactose based dendron D2-2.
Authors	:	Ramaswamy, S.; Sleiman, M.H.; Masuyer, G.; Arbez-Gindre, C.; Micha-Screttas, M.; Calogeropoulou, T.; Steele, B.R.; Acharya, K.R.
Deposited on	:	2014-08-08
Resolution	:	2.04 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

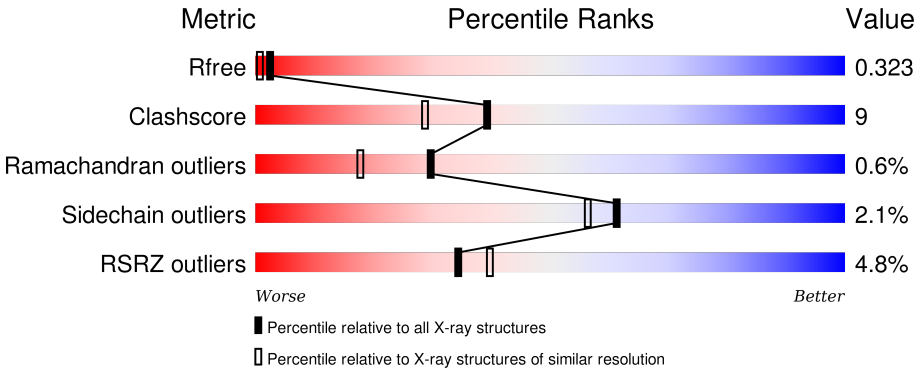
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.04 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	1192 (2.04-2.04)
Clashscore	102246	1269 (2.04-2.04)
Ramachandran outliers	100387	1258 (2.04-2.04)
Sidechain outliers	100360	1258 (2.04-2.04)
RSRZ outliers	91569	1194 (2.04-2.04)

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 <=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	84% 11%
1	B	136	% 87% 11%
1	C	136	82% 14%
1	D	136	% 76% 21%
1	E	136	10% 71% 27%

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Mol	Chain	Length	Quality of chain
1	F	136	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	4S0	A	1136	X	-	-	-
2	4S0	F	1136	X	-	-	-

2 Entry composition

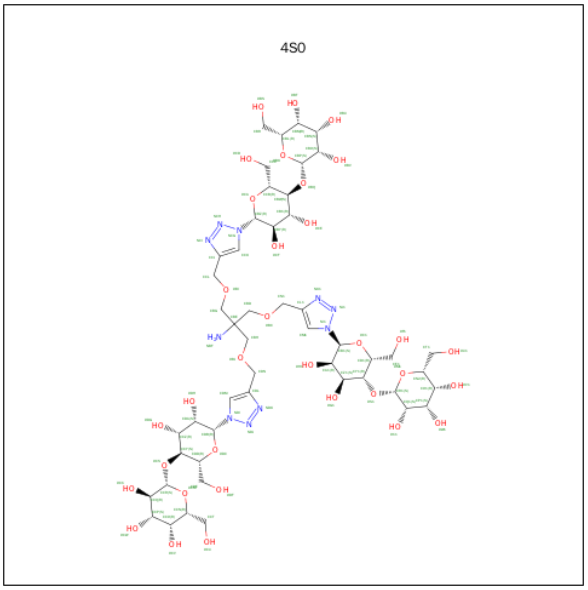
There are 3 unique types of molecules in this entry. The entry contains 6846 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 HUMAN GALECTIN-7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	132	Total	C	N	O	S	0	0	0
			1037	652	197	187	1			
1	B	133	Total	C	N	O	S	0	0	0
			1044	657	198	188	1			
1	C	132	Total	C	N	O	S	0	0	0
			1036	652	197	186	1			
1	D	132	Total	C	N	O	S	0	0	0
			1037	652	197	187	1			
1	E	133	Total	C	N	O	S	0	0	0
			1043	657	198	187	1			
1	F	132	Total	C	N	O	S	0	1	0
			1043	655	198	189	1			

- Molecule 2 is DENDRON D2-1 (three-letter code: 4S0) (formula: $C_{49}H_{80}N_{10}O_{33}$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			92	49	10	33		
2	F	1	Total	C	N	O	0	0
			92	49	10	33		

- Molecule 3 is water.

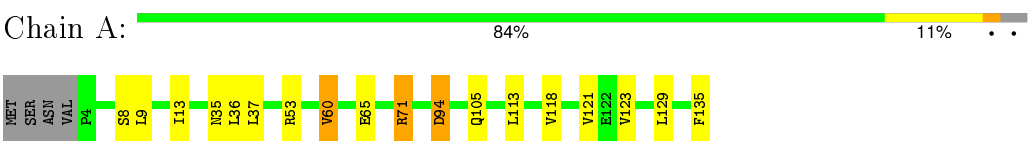
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	117	Total	O	0	0
			117	117		
3	B	53	Total	O	0	0
			53	53		
3	C	72	Total	O	0	0
			72	72		
3	D	51	Total	O	0	0
			51	51		
3	E	69	Total	O	0	0
			69	69		
3	F	60	Total	O	0	0
			60	60		

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.

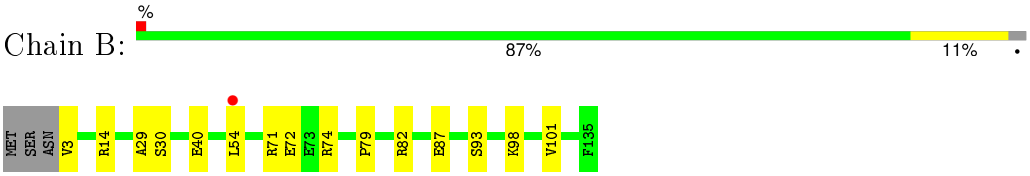
- Molecule 1: HUMAN GALECTIN-7

Chain A:



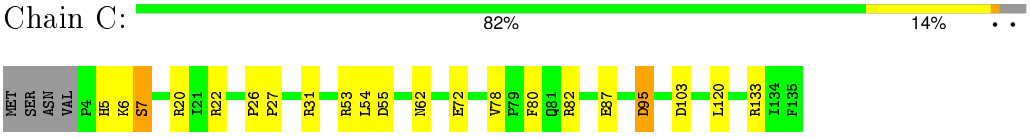
- Molecule 1: HUMAN GALECTIN-7

Chain B:



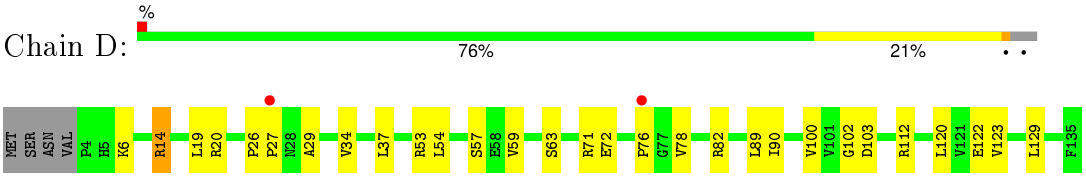
- Molecule 1: HUMAN GALECTIN-7

Chain C:



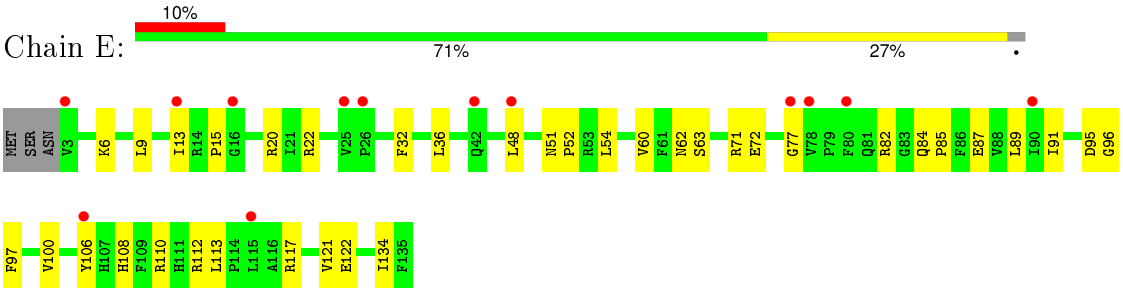
- Molecule 1: HUMAN GALECTIN-7

Chain D:



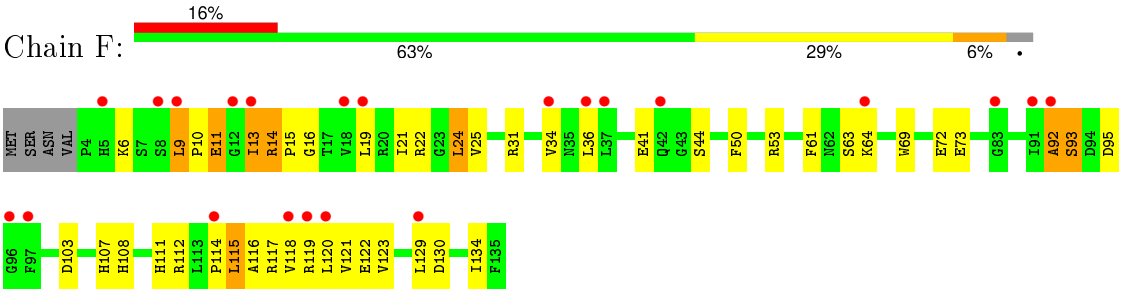
- Molecule 1: HUMAN GALECTIN-7

Chain E:



- Molecule 1: HUMAN GALECTIN-7

Chain F:


16%
63%
29%
6%
MET SER ASN VAL P4 H5 K6 S7 S8 L9 P10 E11 G12 I13 R14 P15 G16 T17 V18 L19 R20 I21 R22 G23 L24 V25 R31 V34 N35 L36 L37 E41 Q42 G43 S44 F50 R53 F61 W62 S63 K64 W69 E72 E73 G83 I91 A92 S93 D94 D95
G96 F97 D103 H107 H108 H111 R112 L113 P114 L115 A116 R117 V118 R119 L120 V121 E122 V123 L129 D130 I134 F135

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	63.44Å 115.65Å 116.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	55.69 – 2.04 55.69 – 2.04	Depositor EDS
% Data completeness (in resolution range)	98.7 (55.69-2.04) 98.7 (55.69-2.04)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.17 (at 2.05Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.238 , 0.319 0.245 , 0.323	Depositor DCC
R_{free} test set	1384 reflections (2.61%)	DCC
Wilson B-factor (Å^2)	32.5	Xtriage
Anisotropy	0.728	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 42.8	EDS
Estimated twinning fraction	0.010 for -h,l,k	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Outliers	2 of 54509 reflections (0.004%)	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	6846	wwPDB-VP
Average B, all atoms (Å^2)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.18% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 4S0

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	0.47	0/1063	0.63	0/1437
1	B	0.42	0/1070	0.56	0/1448
1	C	0.39	0/1062	0.56	0/1437
1	D	0.39	0/1063	0.55	0/1437
1	E	0.42	0/1069	0.59	0/1448
1	F	0.41	0/1069	0.65	1/1445 (0.1%)
All	All	0.42	0/6396	0.59	1/8652 (0.0%)

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	F	0	1

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	F	115	LEU	CA-CB-CG	5.09	127.00	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	F	93	SER	Peptide

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	1037	0	1019	13	0
1	B	1044	0	1027	9	1
1	C	1036	0	1019	15	1
1	D	1037	0	1019	19	0
1	E	1043	0	1027	26	0
1	F	1043	0	1023	40	0
2	A	92	0	80	0	1
2	F	92	0	80	5	0
3	A	117	0	0	3	4
3	B	53	0	0	2	2
3	C	72	0	0	2	1
3	D	51	0	0	1	0
3	E	69	0	0	6	0
3	F	60	0	0	3	0
All	All	6846	0	6294	113	5

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 (113) 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:F:117:ARG:O	1:F:119:ARG:NH1	2.05	0.88
1:E:9:LEU:HD13	1:E:13:ILE:HD11	1.59	0.83
1:A:105:GLN:OE1	3:A:2084:HOH:O	1.96	0.82
1:C:95:ASP:OD2	1:D:14:ARG:NH2	2.12	0.81
1:E:95:ASP:OD2	1:F:14:ARG:NH2	2.13	0.80
1:A:94:ASP:OD2	1:B:14:ARG:NH2	2.18	0.76
1:C:31:ARG:NH2	1:C:55:ASP:OD2	2.18	0.75
1:F:93:SER:HB3	1:F:111:HIS:HE1	1.51	0.74
1:F:13:ILE:HD12	1:F:116:ALA:HB2	1.69	0.72
1:C:78:VAL:HG12	1:C:80:PHE:H	1.55	0.70
1:C:87:GLU:OE1	3:C:2055:HOH:O	2.09	0.69
1:F:6:LYS:NZ	1:F:122:GLU:OE2	2.25	0.67
1:C:20:ARG:NH2	1:D:102:GLY:O	2.28	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:51:ASN:HB3	1:E:60:VAL:HG22	1.79	0.65
1:D:89:LEU:HB2	1:D:100:VAL:HB	1.80	0.62
1:F:36:LEU:HD23	1:F:118:VAL:HG11	1.82	0.62
1:F:13:ILE:HA	1:F:115:LEU:HD23	1.81	0.61
1:C:133:ARG:NH1	1:D:103:ASP:O	2.34	0.61
1:E:54:LEU:HD22	1:E:82:ARG:HG2	1.82	0.61
1:F:93:SER:CB	1:F:111:HIS:HE1	2.13	0.60
1:A:35:ASN:HB3	1:A:37:LEU:HD21	1.83	0.60
1:E:9:LEU:HD11	1:E:134:ILE:HD13	1.84	0.60
1:C:22:ARG:NE	3:C:2012:HOH:O	2.35	0.60
1:F:11:GLU:HG3	1:F:13:ILE:HG22	1.83	0.59
1:F:61:PHE:HB2	1:F:73:GLU:HB2	1.85	0.59
1:E:48:LEU:HD12	3:E:2030:HOH:O	2.02	0.59
1:D:71:ARG:NH2	3:D:2030:HOH:O	2.25	0.58
1:E:62:ASN:O	3:E:2030:HOH:O	2.16	0.57
1:F:13:ILE:HB	1:F:116:ALA:HA	1.85	0.57
1:E:77:GLY:O	1:E:106:TYR:OH	2.21	0.56
1:D:37:LEU:HD12	1:D:120:LEU:HD23	1.87	0.56
1:F:9:LEU:HD21	1:F:134:ILE:HD13	1.88	0.56
1:F:53:ARG:HH12	2:F:1136:4S0:HDG	1.50	0.56
1:E:72:GLU:O	3:E:2041:HOH:O	2.18	0.54
1:E:96:GLY:HA3	1:E:108:HIS:HE1	1.73	0.54
1:B:54:LEU:HD22	1:B:82:ARG:HG3	1.90	0.53
1:A:135:PHE:OXT	1:B:98:LYS:HE3	2.08	0.53
1:E:22:ARG:HG2	1:E:87:GLU:HG3	1.90	0.53
1:F:9:LEU:HD11	1:F:134:ILE:HD12	1.90	0.53
1:B:72:GLU:OE1	1:B:74:ARG:NH2	2.42	0.53
1:C:6:LYS:HG2	1:C:120:LEU:HD11	1.90	0.53
1:F:6:LYS:HB3	1:F:120:LEU:HD11	1.91	0.53
1:D:6:LYS:HE2	1:D:122:GLU:OE1	2.09	0.52
1:D:63:SER:OG	1:D:112:ARG:HD2	2.10	0.52
1:A:8:SER:OG	3:A:2008:HOH:O	2.19	0.52
1:E:96:GLY:HA3	1:E:110:ARG:HA	1.92	0.51
1:A:71:ARG:HG2	1:A:71:ARG:HH11	1.75	0.51
1:D:54:LEU:HD22	1:D:82:ARG:HD2	1.91	0.51
1:E:89:LEU:HB2	1:E:100:VAL:HB	1.91	0.51
1:F:53:ARG:NH1	2:F:1136:4S0:ODG	2.27	0.51
1:F:117:ARG:O	1:F:119:ARG:HG3	2.11	0.51
1:E:108:HIS:NE2	1:E:110:ARG:HG2	2.27	0.50
1:E:112:ARG:NH1	3:E:2034:HOH:O	2.25	0.50
1:E:84:GLN:HG3	1:E:85:PRO:HD2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:63:SER:HB3	1:E:112:ARG:HD2	1.94	0.49
1:A:53:ARG:CZ	1:A:60:VAL:HG11	2.43	0.49
1:E:113:LEU:HD11	1:E:117:ARG:NH1	2.27	0.49
1:E:71:ARG:O	1:E:112:ARG:NH1	2.36	0.48
1:F:41:GLU:HG2	1:F:44:SER:HB2	1.96	0.48
1:D:76:PRO:HB3	1:F:108:HIS:ND1	2.29	0.48
1:E:20:ARG:NH2	1:F:103:ASP:OD1	2.47	0.48
1:C:31:ARG:HB3	1:C:53:ARG:HA	1.96	0.47
1:F:41:GLU:HB2	1:F:44:SER:N	2.29	0.47
1:D:72:GLU:OE2	2:F:1136:4S0:HYZ	2.14	0.47
1:A:36:LEU:HD12	1:A:118:VAL:HG11	1.97	0.47
1:F:63[A]:SER:OG	1:F:112:ARG:HD3	2.14	0.47
1:A:36:LEU:HD13	1:A:121:VAL:HG22	1.95	0.47
1:B:79:PRO:HB2	1:B:101:VAL:HG11	1.96	0.47
1:F:13:ILE:HA	1:F:115:LEU:CD2	2.44	0.46
1:F:72:GLU:OE2	2:F:1136:4S0:HCZ	2.16	0.46
1:E:95:ASP:OD1	3:E:2060:HOH:O	2.21	0.46
1:F:123:VAL:HG11	1:F:129:LEU:HD21	1.97	0.46
1:B:30:SER:HB2	3:B:2052:HOH:O	2.16	0.46
1:D:34:VAL:HG22	1:D:123:VAL:HG22	1.98	0.46
1:C:103:ASP:OD1	1:D:20:ARG:NH2	2.50	0.45
1:E:36:LEU:CD2	1:E:121:VAL:HG22	2.47	0.45
1:C:26:PRO:HA	1:C:27:PRO:HD3	1.83	0.45
1:F:114:PRO:HG2	1:F:117:ARG:HE	1.81	0.45
1:F:19:LEU:HD21	1:F:21:ILE:CD1	2.47	0.45
1:D:26:PRO:HA	1:D:27:PRO:HD3	1.87	0.44
1:F:19:LEU:HD11	1:F:121:VAL:HG11	1.99	0.44
1:E:15:PRO:HG3	3:E:2058:HOH:O	2.17	0.44
1:D:19:LEU:HB3	1:D:90:ILE:HB	1.98	0.44
1:C:62:ASN:ND2	1:C:72:GLU:HG2	2.33	0.43
1:F:64:LYS:HB2	1:F:69:TRP:CZ3	2.53	0.43
1:F:22:ARG:O	1:F:130:ASP:HB2	2.18	0.43
1:F:41:GLU:H	1:F:44:SER:HB2	1.83	0.43
1:D:129:LEU:HA	1:D:129:LEU:HD12	1.88	0.43
1:B:29:ALA:O	1:B:82:ARG:NE	2.47	0.43
1:B:3:VAL:N	3:B:2003:HOH:O	2.50	0.43
1:F:13:ILE:HD12	1:F:116:ALA:CB	2.43	0.42
1:A:9:LEU:HD13	1:A:13:ILE:HG23	2.01	0.42
1:C:31:ARG:CZ	1:C:53:ARG:HD3	2.49	0.42
1:F:31:ARG:NH2	3:F:2013:HOH:O	2.42	0.42
1:E:6:LYS:HE2	1:E:122:GLU:OE2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:91:ILE:O	1:E:97:PHE:HA	2.20	0.42
1:F:9:LEU:HD21	1:F:134:ILE:CD1	2.49	0.41
1:F:61:PHE:CZ	1:F:107:HIS:HB3	2.55	0.41
1:D:26:PRO:HG2	1:D:29:ALA:HB2	2.02	0.41
1:D:53:ARG:HH12	2:F:1136:4S0:HZG	1.59	0.41
1:F:24:LEU:HG	1:F:25:VAL:N	2.36	0.41
1:E:32:PHE:CZ	1:E:52:PRO:HG2	2.56	0.41
1:A:123:VAL:HG11	1:A:129:LEU:HD21	2.02	0.41
1:F:112:ARG:HG3	3:F:2050:HOH:O	2.21	0.41
1:A:65:GLU:HG2	1:A:113:LEU:HD21	2.02	0.41
1:F:92:ALA:HB3	3:F:2006:HOH:O	2.20	0.41
1:C:27:PRO:O	1:C:82:ARG:HD3	2.21	0.40
3:A:2020:HOH:O	1:B:87:GLU:OE1	2.20	0.40
1:D:59:VAL:HG12	1:D:78:VAL:HG13	2.03	0.40
1:F:34:VAL:HG22	1:F:50:PHE:HB3	2.02	0.40
1:F:10:PRO:HA	1:F:11:GLU:HA	1.87	0.40
1:A:9:LEU:HD11	1:A:121:VAL:HG23	2.03	0.40
1:C:5:HIS:HE2	1:C:7:SER:HG	1.69	0.40

All (5) 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)
3:A:2108:HOH:O	3:B:2023:HOH:O[4_577]	2.03	0.17
3:A:2106:HOH:O	3:B:2033:HOH:O[4_577]	2.09	0.11
1:B:40:GLU:OE1	3:A:2008:HOH:O[4_577]	2.14	0.06
1:C:53:ARG:NH1	2:A:1136:4S0:OCD[2_784]	2.14	0.06
3:A:2113:HOH:O	3:C:2035:HOH:O[2_785]	2.19	0.01

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	130/136 (96%)	128 (98%)	2 (2%)	0	100 100
1	B	131/136 (96%)	127 (97%)	4 (3%)	0	100 100
1	C	130/136 (96%)	127 (98%)	3 (2%)	0	100 100
1	D	130/136 (96%)	126 (97%)	3 (2%)	1 (1%)	24 12
1	E	131/136 (96%)	127 (97%)	4 (3%)	0	100 100
1	F	131/136 (96%)	111 (85%)	16 (12%)	4 (3%)	5 0
All	All	783/816 (96%)	746 (95%)	32 (4%)	5 (1%)	30 18

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	13	ILE
1	D	57	SER
1	F	92	ALA
1	F	16	GLY
1	F	15	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	111/115 (96%)	108 (97%)	3 (3%)	52 45
1	B	112/115 (97%)	110 (98%)	2 (2%)	66 62
1	C	111/115 (96%)	108 (97%)	3 (3%)	52 45
1	D	111/115 (96%)	110 (99%)	1 (1%)	84 84
1	E	112/115 (97%)	112 (100%)	0	100 100
1	F	112/115 (97%)	107 (96%)	5 (4%)	34 25
All	All	669/690 (97%)	655 (98%)	14 (2%)	61 56

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	VAL
1	A	71	ARG
1	A	94	ASP
1	B	71	ARG
1	B	93	SER
1	C	7	SER
1	C	54	LEU
1	C	95	ASP
1	D	14	ARG
1	F	9	LEU
1	F	11	GLU
1	F	14	ARG
1	F	24	LEU
1	F	95	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	F	111	HIS

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)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	4S0	A	1136	-	94,100,100	2.24	10 (10%)	120,147,147	1.46	20 (16%)
2	4S0	F	1136	-	94,100,100	2.01	12 (12%)	120,147,147	1.37	15 (12%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	4S0	A	1136	-	17/17/31/31	0/39/177/177	0/9/9/9
2	4S0	F	1136	-	17/17/31/31	0/39/177/177	0/9/9/9

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	1136	4S0	NK1-NJ1	-14.53	1.13	1.34
2	A	1136	4S0	NCI-NCH	-12.45	1.16	1.34
2	A	1136	4S0	NK1-NJ1	-10.70	1.18	1.34
2	A	1136	4S0	NCH-NCG	-7.36	1.21	1.34
2	A	1136	4S0	NDK-NDJ	-5.62	1.26	1.34
2	F	1136	4S0	NDJ-NDI	-5.30	1.25	1.34
2	F	1136	4S0	NCI-NCH	-4.74	1.27	1.34
2	A	1136	4S0	NDJ-NDI	-4.39	1.26	1.34
2	A	1136	4S0	NJ1-NI1	-4.16	1.27	1.34
2	F	1136	4S0	NJ1-NI1	-4.13	1.27	1.34
2	A	1136	4S0	CDM-CDL	-2.85	1.32	1.36
2	F	1136	4S0	CCK-CCJ	-2.73	1.32	1.36
2	A	1136	4S0	CDM-NDI	-2.58	1.33	1.35
2	F	1136	4S0	CDM-CDL	-2.40	1.32	1.36
2	F	1136	4S0	NCH-NCG	-2.29	1.30	1.34
2	F	1136	4S0	CM1-CL1	-2.24	1.32	1.36
2	A	1136	4S0	CCK-CCJ	-2.20	1.33	1.36
2	A	1136	4S0	CM1-CL1	-2.04	1.33	1.36
2	F	1136	4S0	CM1-NI1	-2.01	1.33	1.35
2	F	1136	4S0	OBQ-CBP	2.13	1.47	1.41
2	F	1136	4S0	NDK-NDJ	2.28	1.37	1.34
2	F	1136	4S0	OCA-CBZ	2.35	1.45	1.42

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	F	1136	4S0	CL1-CM1-NI1	-3.95	100.62	107.14
2	A	1136	4S0	CCJ-CCK-NCG	-3.94	100.64	107.14
2	A	1136	4S0	CL1-CM1-NI1	-3.02	102.16	107.14
2	F	1136	4S0	CDA-CDB-NDI	-2.54	109.04	111.98
2	A	1136	4S0	OCS-CCR-CCQ	-2.49	102.05	108.10
2	A	1136	4S0	OBJ-CDN-CDL	-2.42	103.73	110.87
2	A	1136	4S0	CP1-CO1-CN2	-2.41	106.00	110.20
2	A	1136	4S0	CCK-CCJ-NCI	-2.36	107.64	111.42
2	A	1136	4S0	CDA-CCZ-CCY	-2.33	104.48	109.60
2	A	1136	4S0	CBP-OBQ-CBW	-2.32	111.94	118.01
2	A	1136	4S0	CBY-CBZ-NCG	-2.32	109.29	111.98
2	F	1136	4S0	CDA-CCZ-CCY	-2.25	104.65	109.60
2	A	1136	4S0	CCQ-CCP-CCO	-2.23	106.62	110.79
2	F	1136	4S0	CBN-CBM-CBL	-2.23	106.31	110.20
2	A	1136	4S0	OF1-CE1-CD1	-2.22	104.00	111.33
2	A	1136	4S0	CZ1-CA1-CB1	-2.21	105.61	109.21
2	F	1136	4S0	CZ1-CA1-CB1	-2.21	105.61	109.21
2	F	1136	4S0	CR1-OM1-CN2	-2.13	109.62	113.75
2	A	1136	4S0	OS1-CR1-CQ1	-2.10	102.99	108.10
2	A	1136	4S0	CCZ-CCY-CDD	-2.08	106.14	110.84
2	F	1136	4S0	CDM-CDL-NDK	-2.05	108.13	111.42
2	F	1136	4S0	CDD-ODC-CDB	2.17	112.55	108.71
2	A	1136	4S0	OCS-CCY-CCZ	2.24	112.94	107.17
2	F	1136	4S0	OBQ-CBW-CBX	2.27	113.02	107.17
2	A	1136	4S0	CDA-CDB-NDI	2.49	114.86	111.98
2	F	1136	4S0	OS1-CY1-CD1	2.75	116.56	109.32
2	A	1136	4S0	OCS-CCY-CDD	2.99	117.18	109.32
2	F	1136	4S0	CA1-CB1-NI1	3.19	115.66	111.98
2	F	1136	4S0	OCS-CCY-CDD	3.57	118.71	109.32
2	A	1136	4S0	OS1-CY1-CD1	3.60	118.78	109.32
2	A	1136	4S0	CCR-OCS-CCY	3.70	127.68	118.01
2	F	1136	4S0	CBP-OBQ-CBW	3.71	127.71	118.01
2	F	1136	4S0	CR1-OS1-CY1	4.00	128.46	118.01
2	F	1136	4S0	CCR-OCS-CCY	4.42	129.55	118.01
2	A	1136	4S0	CR1-OS1-CY1	4.86	130.71	118.01

All (34) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	F	1136	4S0	CCY
2	F	1136	4S0	CB1
2	F	1136	4S0	CDD

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Mol	Chain	Res	Type	Atom
2	F	1136	4S0	CCZ
2	F	1136	4S0	CDB
2	F	1136	4S0	CBY
2	F	1136	4S0	CBX
2	F	1136	4S0	CN2
2	F	1136	4S0	CP1
2	F	1136	4S0	CR1
2	F	1136	4S0	CBZ
2	F	1136	4S0	CBO
2	F	1136	4S0	CY1
2	F	1136	4S0	CZ1
2	F	1136	4S0	CO1
2	F	1136	4S0	CBW
2	F	1136	4S0	CCB
2	A	1136	4S0	CCY
2	A	1136	4S0	CB1
2	A	1136	4S0	CDD
2	A	1136	4S0	CCZ
2	A	1136	4S0	CDB
2	A	1136	4S0	CBY
2	A	1136	4S0	CBX
2	A	1136	4S0	CN2
2	A	1136	4S0	CP1
2	A	1136	4S0	CR1
2	A	1136	4S0	CBZ
2	A	1136	4S0	CBO
2	A	1136	4S0	CY1
2	A	1136	4S0	CZ1
2	A	1136	4S0	CO1
2	A	1136	4S0	CBW
2	A	1136	4S0	CCB

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1136	4S0	0	1
2	F	1136	4S0	5	0

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

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	132/136 (97%)	0.01	0 100 100	17, 28, 39, 47	0
1	B	133/136 (97%)	0.07	1 (0%) 87 90	19, 31, 47, 59	0
1	C	132/136 (97%)	0.05	0 100 100	22, 32, 44, 57	0
1	D	132/136 (97%)	0.14	2 (1%) 76 81	23, 34, 52, 63	0
1	E	133/136 (97%)	0.83	13 (9%) 10 10	24, 40, 51, 59	0
1	F	132/136 (97%)	0.92	22 (16%) 2 2	21, 42, 69, 81	0
All	All	794/816 (97%)	0.34	38 (4%) 34 40	17, 34, 56, 81	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	13	ILE	10.5
1	F	19	LEU	4.8
1	E	13	ILE	4.8
1	F	92	ALA	4.3
1	E	80	PHE	3.9
1	F	37	LEU	3.8
1	E	25	VAL	3.7
1	E	106	TYR	3.2
1	F	64	LYS	3.0
1	F	18	VAL	2.9
1	B	54	LEU	2.8
1	F	36	LEU	2.8
1	F	119	ARG	2.7
1	F	97	PHE	2.7
1	F	8	SER	2.7
1	E	78	VAL	2.6
1	E	42	GLN	2.6
1	E	16	GLY	2.6
1	E	26	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
1	F	83	GLY	2.5
1	F	12	GLY	2.5
1	F	114	PRO	2.4
1	F	120	LEU	2.4
1	E	90	ILE	2.4
1	D	27	PRO	2.4
1	F	129	LEU	2.4
1	F	96	GLY	2.3
1	F	5	HIS	2.3
1	F	118	VAL	2.2
1	F	91	ILE	2.2
1	E	115	LEU	2.2
1	F	9	LEU	2.1
1	F	34	VAL	2.1
1	D	76	PRO	2.1
1	E	3	VAL	2.1
1	E	48	LEU	2.1
1	F	42	GLN	2.1
1	E	77	GLY	2.1

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)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
2	4S0	A	1136	92/92	0.92	0.12	0.03	17,29,40,49	0
2	4S0	F	1136	92/92	0.89	0.13	-0.21	26,38,48,54	0

6.5 Other polymers (i)

There are no such residues in this entry.