Full wwPDB X-ray Structure Validation Report (i)

Jun 16, 2024 – 03:20 AM EDT

PDB ID	:	2AQ1
Title	:	Crystal structure of T-cell receptor V beta domain variant complexed with superantigen SEC3 mutant
Authors	:	Cho, S.; Swaminathan, C.P.; Yang, J.; Kerzic, M.C.; Guan, R.; Kieke, M.C.; Kranz, D.M.; Mariuzza, R.A.; Sundberg, E.J.
Deposited on	:	2005-08-17
Resolution	:	2.10 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 (i) 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 (1)) were used in the production of this report:

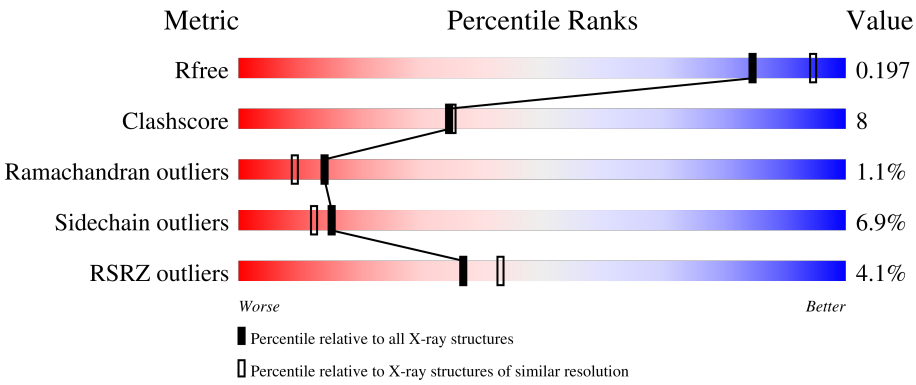
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	112	88% 6% • •
1	C	112	% 76% 18% • •
1	E	112	3% 79% 12% 5% •
1	G	112	2% 71% 24% • •
2	B	237	3% 80% 15% • •

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
2	D	237	8% 76% 19%
2	F	237	5% 79% 19%
2	H	237	6% 78% 17%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 11673 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 T-cell receptor beta chain V.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	109	832	517	147	165	3	0	0	0
1	C	109	832	517	147	165	3	0	0	0
1	E	109	832	517	147	165	3	0	0	0
1	G	109	832	517	147	165	3	0	0	0

- Molecule 2 is a protein called Enterotoxin type C-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	234	1926	1225	316	375	10	0	0	0
2	D	236	1939	1231	319	379	10	0	0	0
2	F	237	1948	1236	320	382	10	0	0	0
2	H	237	1948	1236	320	382	10	0	0	0

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	GLY	SEE REMARK 999	UNP P0A0L5
B	?	-	LYS	SEE REMARK 999	UNP P0A0L5
B	102	TRP	VAL	SEE REMARK 999	UNP P0A0L5
B	103	TRP	THR	SEE REMARK 999	UNP P0A0L5
B	104	HIS	GLY	SEE REMARK 999	UNP P0A0L5
D	?	-	GLY	SEE REMARK 999	UNP P0A0L5
D	?	-	LYS	SEE REMARK 999	UNP P0A0L5
D	102	TRP	VAL	SEE REMARK 999	UNP P0A0L5

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	103	TRP	THR	SEE REMARK 999	UNP P0A0L5
D	104	HIS	GLY	SEE REMARK 999	UNP P0A0L5
F	?	-	GLY	SEE REMARK 999	UNP P0A0L5
F	?	-	LYS	SEE REMARK 999	UNP P0A0L5
F	102	TRP	VAL	SEE REMARK 999	UNP P0A0L5
F	103	TRP	THR	SEE REMARK 999	UNP P0A0L5
F	104	HIS	GLY	SEE REMARK 999	UNP P0A0L5
H	?	-	GLY	SEE REMARK 999	UNP P0A0L5
H	?	-	LYS	SEE REMARK 999	UNP P0A0L5
H	102	TRP	VAL	SEE REMARK 999	UNP P0A0L5
H	103	TRP	THR	SEE REMARK 999	UNP P0A0L5
H	104	HIS	GLY	SEE REMARK 999	UNP P0A0L5

- Molecule 3 is water.

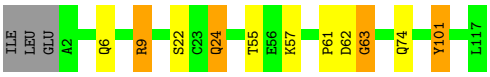
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	64	Total O 64 64	0	0
3	B	118	Total O 118 118	0	0
3	C	56	Total O 56 56	0	0
3	D	85	Total O 85 85	0	0
3	E	50	Total O 50 50	0	0
3	F	91	Total O 91 91	0	0
3	G	45	Total O 45 45	0	0
3	H	75	Total O 75 75	0	0

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.

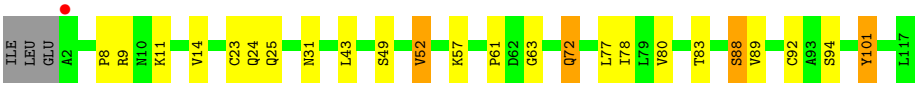
• Molecule 1: T-cell receptor beta chain V

Chain A:



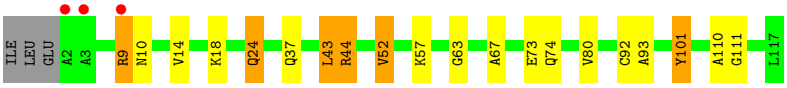
• Molecule 1: T-cell receptor beta chain V

Chain C:



• Molecule 1: T-cell receptor beta chain V

Chain E:

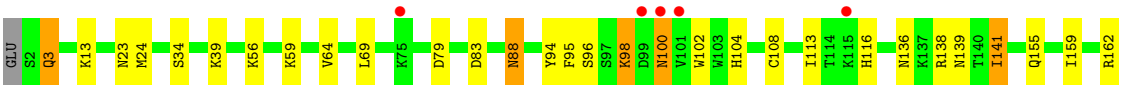


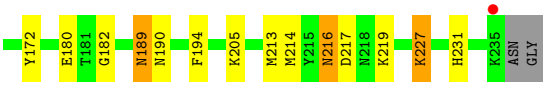
• Molecule 1: T-cell receptor beta chain V

Chain G:

• Molecule 2: Enterotoxin type C-3

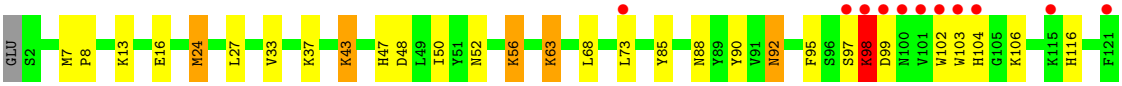
Chain B:

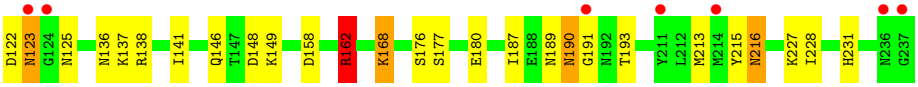




• Molecule 2: Enterotoxin type C-3

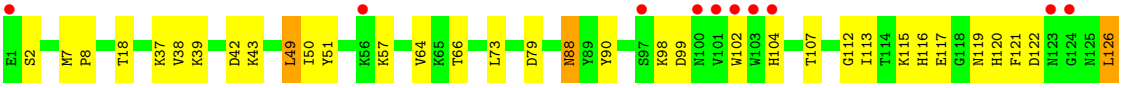
Chain D:

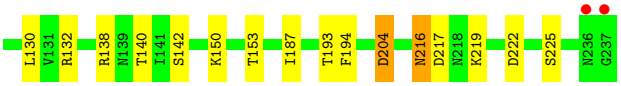




• Molecule 2: Enterotoxin type C-3

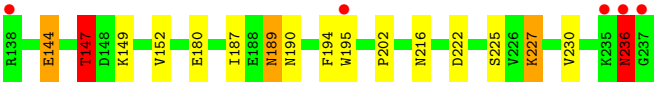
Chain F:





• Molecule 2: Enterotoxin type C-3

Chain H:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	63.20Å 70.19Å 98.40Å 74.79° 75.05° 88.54°	Depositor
Resolution (Å)	35.00 – 2.10 35.21 – 2.10	Depositor EDS
% Data completeness (in resolution range)	96.0 (35.00-2.10) 88.6 (35.21-2.10)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.95 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.185 , 0.243 0.196 , 0.197	Depositor DCC
R_{free} test set	4085 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	39.4	Xtriage
Anisotropy	0.711	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 53.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	11673	wwPDB-VP
Average B, all atoms (Å^2)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.91% 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 (i)

5.1 Standard geometry (i)

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	1.32	2/850 (0.2%)	1.07	1/1153 (0.1%)
1	C	1.26	4/850 (0.5%)	1.09	5/1153 (0.4%)
1	E	1.25	4/850 (0.5%)	1.00	1/1153 (0.1%)
1	G	1.33	5/850 (0.6%)	0.99	1/1153 (0.1%)
2	B	1.17	2/1972 (0.1%)	0.94	2/2660 (0.1%)
2	D	1.07	1/1985 (0.1%)	0.95	2/2676 (0.1%)
2	F	1.10	1/1994 (0.1%)	0.94	3/2688 (0.1%)
2	H	1.04	0/1994	0.91	1/2688 (0.0%)
All	All	1.16	19/11345 (0.2%)	0.97	16/15324 (0.1%)

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	2
1	C	0	1
1	G	0	1
2	D	0	1
2	F	0	1
2	H	0	1
All	All	0	7

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	63	GLY	C-N	12.58	1.62	1.34
1	A	63	GLY	C-N	11.99	1.61	1.34
1	C	63	GLY	C-N	11.50	1.60	1.34
1	G	101	TYR	C-N	11.38	1.60	1.34
1	C	101	TYR	C-N	9.16	1.55	1.34
1	E	63	GLY	C-N	9.12	1.55	1.34

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	92	CYS	CB-SG	-7.99	1.68	1.82
1	E	67	ALA	CA-CB	7.60	1.68	1.52
1	E	101	TYR	C-N	7.23	1.50	1.34
1	A	101	TYR	C-N	7.08	1.50	1.34
2	D	215	TYR	CE1-CZ	6.24	1.46	1.38
1	C	23	CYS	CB-SG	6.01	1.92	1.82
2	B	172	TYR	CE1-CZ	6.00	1.46	1.38
2	F	51	TYR	CD1-CE1	-5.81	1.30	1.39
1	E	92	CYS	CB-SG	-5.67	1.72	1.81
1	G	73	GLU	CG-CD	5.36	1.59	1.51
1	G	93	ALA	CA-CB	5.33	1.63	1.52
1	G	35	TYR	CD2-CE2	5.17	1.47	1.39
2	B	172	TYR	CD2-CE2	5.09	1.47	1.39

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	162	ARG	NE-CZ-NH1	14.78	127.69	120.30
1	A	63	GLY	O-C-N	-12.35	102.95	122.70
2	D	162	ARG	NE-CZ-NH2	-10.68	114.96	120.30
1	C	101	TYR	O-C-N	-10.03	106.66	122.70
2	F	132	ARG	NE-CZ-NH2	-6.85	116.88	120.30
1	G	36	ARG	NE-CZ-NH2	-6.73	116.93	120.30
1	C	63	GLY	C-N-CA	-6.40	105.69	121.70
1	E	63	GLY	O-C-N	-6.35	112.54	122.70
2	F	49	LEU	CA-CB-CG	6.31	129.82	115.30
2	B	162	ARG	NE-CZ-NH1	-6.13	117.24	120.30
2	H	147	THR	CB-CA-C	-5.79	95.96	111.60
1	C	63	GLY	O-C-N	5.55	131.57	122.70
1	C	101	TYR	CA-C-N	5.28	128.82	117.20
1	C	23	CYS	CA-CB-SG	5.14	123.26	114.00
2	F	126	LEU	CB-CG-CD1	5.07	119.62	111.00
2	B	24	MET	CG-SD-CE	-5.07	92.09	100.20

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	101	TYR	Mainchain
1	A	63	GLY	Mainchain
1	C	101	TYR	Mainchain
2	D	98	LYS	Peptide

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Group
2	F	121	PHE	Peptide
1	G	63	GLY	Mainchain
2	H	97	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	832	0	793	8	0
1	C	832	0	793	12	0
1	E	832	0	793	16	0
1	G	832	0	793	19	0
2	B	1926	0	1854	29	0
2	D	1939	0	1863	47	0
2	F	1948	0	1872	31	0
2	H	1948	0	1872	28	0
3	A	64	0	0	1	0
3	B	118	0	0	3	0
3	C	56	0	0	0	0
3	D	85	0	0	9	0
3	E	50	0	0	2	0
3	F	91	0	0	6	0
3	G	45	0	0	4	0
3	H	75	0	0	2	0
All	All	11673	0	10633	181	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 8.

All (181) 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)
2:D:24:MET:HA	2:D:24:MET:HE3	1.25	1.13
2:D:158:ASP:OD1	2:D:162:ARG:HD3	1.62	0.98
2:F:120:HIS:O	2:F:150:LYS:HE3	1.72	0.88
1:E:80:VAL:HG12	3:E:131:HOH:O	1.74	0.87

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:THR:H	2:B:23:ASN:HD21	1.21	0.86
2:D:24:MET:HA	2:D:24:MET:CE	2.06	0.83
2:D:158:ASP:O	2:D:162:ARG:HG2	1.79	0.81
2:B:100:ASN:HB3	2:B:102:TRP:HD1	1.44	0.81
2:F:88:ASN:HD22	2:F:88:ASN:H	1.27	0.80
2:D:92:ASN:HB2	3:D:296:HOH:O	1.79	0.80
2:H:88:ASN:H	2:H:88:ASN:HD22	1.30	0.79
2:D:168:LYS:HE3	2:D:168:LYS:HA	1.63	0.79
2:H:227:LYS:HA	2:H:227:LYS:HE2	1.66	0.78
2:D:24:MET:HE1	2:D:27:LEU:HD12	1.63	0.78
1:E:44:ARG:HG3	1:E:44:ARG:HH21	1.48	0.78
2:D:168:LYS:HE3	2:D:168:LYS:CA	2.12	0.78
1:G:42:GLY:HA2	3:G:148:HOH:O	1.82	0.77
1:A:9:ARG:HB3	1:A:9:ARG:NH1	1.99	0.76
1:A:9:ARG:HB3	1:A:9:ARG:HH11	1.52	0.75
2:B:189:ASN:HD22	2:B:190:ASN:N	1.83	0.75
2:B:214:MET:HE3	3:B:283:HOH:O	1.87	0.73
2:D:138:ARG:HD3	3:D:273:HOH:O	1.88	0.73
2:B:100:ASN:HB3	2:B:102:TRP:CD1	2.25	0.72
2:H:102:TRP:HB2	2:H:104:HIS:HD2	1.55	0.72
2:D:136:ASN:O	2:D:137:LYS:HB2	1.90	0.71
2:B:205:LYS:HD3	3:B:289:HOH:O	1.91	0.71
2:H:13:LYS:HB3	2:H:180:GLU:OE1	1.91	0.71
2:D:56:LYS:CD	2:D:56:LYS:H	2.04	0.70
3:F:321:HOH:O	1:G:113:ARG:HB3	1.90	0.70
2:D:168:LYS:HA	2:D:168:LYS:CE	2.21	0.70
2:H:128:ASN:HB3	2:H:144:GLU:HG2	1.73	0.70
1:E:80:VAL:HG11	1:G:83:THR:CG2	2.23	0.69
1:G:28:ASN:ND2	1:G:72:GLN:HG2	2.07	0.68
1:E:57:LYS:HD2	3:F:294:HOH:O	1.92	0.67
1:G:21:LEU:HD22	1:G:112:THR:HG21	1.76	0.67
2:D:13:LYS:HB2	2:D:16:GLU:HG3	1.76	0.66
2:D:24:MET:HE3	2:D:24:MET:CA	2.11	0.66
2:F:88:ASN:H	2:F:88:ASN:ND2	1.92	0.66
2:D:24:MET:CE	2:D:27:LEU:HD12	2.26	0.65
2:H:202:PRO:HG3	3:H:312:HOH:O	1.97	0.65
2:H:101:VAL:O	2:H:102:TRP:HB2	1.98	0.64
2:H:147:THR:HG21	2:H:152:VAL:CG2	2.28	0.64
1:C:9:ARG:HB3	1:C:9:ARG:NH1	2.12	0.63
2:H:189:ASN:HD22	2:H:190:ASN:N	1.96	0.63
2:H:189:ASN:HD22	2:H:190:ASN:H	1.46	0.62

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:13:LYS:HB3	2:D:180:GLU:OE1	2.00	0.62
2:H:115:LYS:HG3	3:H:281:HOH:O	1.98	0.62
2:H:123:ASN:HB3	2:H:125:ASN:H	1.65	0.61
2:H:147:THR:HG21	2:H:152:VAL:HG21	1.81	0.61
2:F:66:THR:HG21	2:F:113:ILE:HD11	1.82	0.61
2:H:236:ASN:N	2:H:236:ASN:OD1	2.34	0.60
2:B:83:ASP:OD1	2:B:116:HIS:HD2	1.84	0.60
1:E:80:VAL:CG1	1:G:83:THR:HG21	2.31	0.60
2:D:97:SER:C	2:D:99:ASP:H	2.03	0.60
2:D:176:SER:OG	2:D:177:SER:N	2.31	0.59
1:E:80:VAL:HG11	1:G:83:THR:HG21	1.83	0.59
2:F:39:LYS:HD2	2:F:79:ASP:O	2.02	0.59
2:H:227:LYS:HA	2:H:227:LYS:CE	2.29	0.59
2:B:88:ASN:HD22	2:B:88:ASN:H	1.51	0.59
1:C:77:LEU:C	1:C:78:ILE:HD13	2.23	0.59
2:D:123:ASN:HA	3:D:275:HOH:O	2.01	0.58
2:H:55:ASP:OD1	2:H:57:LYS:HG2	2.02	0.58
2:H:68:LEU:HD21	2:H:113:ILE:HD12	1.85	0.58
1:E:52:VAL:HG22	2:F:90:TYR:CD1	2.39	0.57
2:F:112:GLY:HA2	2:F:153:THR:HG21	1.85	0.57
2:B:189:ASN:HD22	2:B:190:ASN:H	1.51	0.56
2:F:88:ASN:HD22	2:F:88:ASN:N	2.01	0.56
1:C:52:VAL:HG22	2:D:90:TYR:CD1	2.40	0.56
2:D:189:ASN:C	2:D:191:GLY:H	2.08	0.56
2:D:47:HIS:H	2:D:47:HIS:CD2	2.25	0.55
2:F:104:HIS:HB3	3:F:328:HOH:O	2.06	0.55
2:B:94:TYR:O	2:B:95:PHE:HB3	2.06	0.54
1:E:80:VAL:CG1	1:G:83:THR:CG2	2.86	0.54
1:E:9:ARG:HH22	1:E:110:ALA:HB3	1.73	0.54
1:A:62:ASP:HB2	3:A:124:HOH:O	2.08	0.53
2:D:43:LYS:HA	2:D:50:ILE:HG12	1.89	0.53
2:D:92:ASN:CB	3:D:296:HOH:O	2.45	0.53
2:F:39:LYS:CD	2:F:79:ASP:O	2.57	0.53
2:F:138:ARG:HD2	2:F:140:THR:HG22	1.91	0.53
2:H:88:ASN:H	2:H:88:ASN:ND2	2.04	0.53
1:G:31:ASN:ND2	1:G:50:TYR:HD1	2.07	0.52
2:F:187:ILE:HG23	2:F:193:THR:HG22	1.90	0.52
2:F:115:LYS:HE3	2:F:117:GLU:N	2.24	0.52
2:F:18:THR:OG1	2:F:204:ASP:OD2	2.22	0.51
2:F:119:ASN:O	2:F:150:LYS:HG3	2.11	0.51
2:F:138:ARG:HD3	3:F:327:HOH:O	2.09	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:THR:H	2:B:23:ASN:ND2	2.00	0.51
2:D:168:LYS:CA	2:D:168:LYS:CE	2.86	0.51
2:B:213:MET:HE2	3:B:279:HOH:O	2.11	0.50
2:H:147:THR:HB	2:H:149:LYS:H	1.76	0.50
2:F:115:LYS:HE3	2:F:117:GLU:HB2	1.94	0.50
1:G:22:SER:HB3	3:G:162:HOH:O	2.10	0.50
2:D:231:HIS:ND1	3:D:286:HOH:O	2.35	0.50
2:H:187:ILE:HB	2:H:227:LYS:HB3	1.94	0.50
2:F:115:LYS:CE	2:F:117:GLU:HB2	2.41	0.49
2:B:182:GLY:HA2	2:B:231:HIS:O	2.13	0.49
1:E:44:ARG:HH21	1:E:44:ARG:CG	2.21	0.49
3:E:149:HOH:O	1:G:84:PRO:HG2	2.12	0.49
2:B:194:PHE:CE2	2:B:219:LYS:HE2	2.48	0.49
2:B:136:ASN:O	2:B:138:ARG:NH1	2.47	0.48
2:D:213:MET:O	2:D:216:ASN:HB2	2.13	0.48
2:F:117:GLU:CG	3:F:305:HOH:O	2.60	0.48
2:F:130:LEU:HD11	2:F:142:SER:HB3	1.94	0.48
2:F:222:ASP:OD2	2:F:225:SER:OG	2.17	0.48
2:D:37:LYS:HG3	2:D:116:HIS:CD2	2.49	0.48
2:D:227:LYS:HZ2	2:D:228:ILE:H	1.61	0.48
1:A:24:GLN:OE1	1:A:74:GLN:NE2	2.44	0.48
1:C:57:LYS:HB3	1:C:61:PRO:HG3	1.96	0.48
2:F:194:PHE:CE2	2:F:219:LYS:HE3	2.49	0.47
1:C:31:ASN:HA	1:C:49:SER:O	2.13	0.47
2:B:227:LYS:HA	2:B:227:LYS:HD2	1.68	0.47
2:B:213:MET:O	2:B:216:ASN:HB2	2.14	0.47
2:F:115:LYS:HE3	2:F:117:GLU:H	1.80	0.47
2:B:3:GLN:HE21	2:B:194:PHE:HB2	1.80	0.47
2:D:138:ARG:CD	3:D:273:HOH:O	2.54	0.47
2:D:97:SER:C	2:D:99:ASP:N	2.68	0.47
2:B:96:SER:HB3	2:B:98:LYS:HG2	1.96	0.47
2:H:37:LYS:HG3	2:H:116:HIS:CD2	2.50	0.47
2:H:222:ASP:OD2	2:H:225:SER:OG	2.29	0.47
2:B:88:ASN:H	2:B:88:ASN:ND2	2.12	0.46
2:F:64:VAL:HG22	2:F:107:THR:HG22	1.97	0.46
2:D:148:ASP:OD2	2:D:149:LYS:HG3	2.15	0.46
2:F:37:LYS:HG3	2:F:116:HIS:CD2	2.50	0.46
2:D:95:PHE:CZ	2:D:106:LYS:HB2	2.51	0.46
2:F:120:HIS:O	2:F:150:LYS:CE	2.55	0.46
2:F:187:ILE:HA	2:F:193:THR:HG22	1.97	0.46
1:A:6:GLN:HB2	1:A:9:ARG:HH21	1.81	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:88:SER:OG	1:C:89:VAL:N	2.49	0.46
2:D:24:MET:CE	2:D:24:MET:CA	2.80	0.46
1:C:9:ARG:HB3	1:C:9:ARG:HH11	1.78	0.45
2:H:57:LYS:N	2:H:57:LYS:HD3	2.31	0.45
2:F:117:GLU:HG2	3:F:305:HOH:O	2.17	0.45
2:B:56:LYS:HD3	2:B:56:LYS:HA	1.82	0.45
2:D:187:ILE:HG23	2:D:193:THR:HG22	1.98	0.45
1:G:66:GLU:HB2	1:G:78:ILE:HB	2.00	0.44
2:H:77:TYR:O	2:H:78:LYS:C	2.56	0.44
2:H:147:THR:CG2	2:H:152:VAL:HG21	2.48	0.44
1:C:8:PRO:HG2	1:C:11:LYS:HB2	1.99	0.44
1:C:25:GLN:O	1:C:25:GLN:HG3	2.18	0.44
2:F:42:ASP:HB3	2:F:50:ILE:HB	2.00	0.44
1:E:80:VAL:HG11	1:G:83:THR:HG22	1.99	0.44
1:G:42:GLY:CA	3:G:148:HOH:O	2.52	0.44
2:B:13:LYS:HB3	2:B:180:GLU:OE1	2.18	0.43
2:B:39:LYS:HD2	2:B:79:ASP:O	2.18	0.43
1:G:69:ARG:HD2	1:G:74:GLN:O	2.18	0.43
2:D:189:ASN:OD1	2:D:190:ASN:N	2.51	0.43
2:D:43:LYS:NZ	2:D:48:ASP:O	2.38	0.43
2:D:227:LYS:NZ	3:D:319:HOH:O	2.52	0.43
2:F:7:MET:O	2:F:8:PRO:C	2.56	0.43
2:D:52:ASN:OD1	2:D:63:LYS:CD	2.66	0.43
2:D:123:ASN:HB2	2:D:125:ASN:OD1	2.19	0.43
2:F:216:ASN:HD22	2:F:217:ASP:N	2.15	0.43
2:D:98:LYS:NZ	3:D:283:HOH:O	2.47	0.43
1:C:72:GLN:HE21	1:C:72:GLN:HB3	1.21	0.42
1:C:78:ILE:HD13	1:C:78:ILE:N	2.33	0.42
2:D:146:GLN:HE21	2:D:146:GLN:HB2	1.61	0.42
1:E:93:ALA:HA	1:E:101:TYR:O	2.19	0.42
2:D:7:MET:HG2	3:D:303:HOH:O	2.18	0.42
1:G:9:ARG:HG2	3:G:160:HOH:O	2.19	0.42
2:B:88:ASN:HA	2:B:108:CYS:O	2.20	0.42
1:C:31:ASN:OD1	1:C:31:ASN:N	2.51	0.42
2:H:3:GLN:HE21	2:H:194:PHE:HA	1.85	0.42
2:D:227:LYS:HD2	2:D:227:LYS:HA	1.73	0.42
1:E:9:ARG:NH2	1:E:111:GLY:O	2.53	0.42
2:B:69:LEU:HG	2:B:217:ASP:HA	2.02	0.41
2:B:141:ILE:H	2:B:141:ILE:HD13	1.84	0.41
2:D:162:ARG:HG2	2:D:162:ARG:H	1.78	0.41
2:H:6:PRO:HD3	2:H:195:TRP:CE2	2.54	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:113:ILE:HD13	2:B:113:ILE:HG21	1.78	0.41
1:G:87:SER:O	1:G:88:SER:HB2	2.21	0.41
1:E:24:GLN:HG2	1:E:74:GLN:NE2	2.36	0.41
1:G:11:LYS:HE3	1:G:19:VAL:HG13	2.02	0.41
1:G:2:ALA:O	1:G:3:ALA:HB3	2.21	0.41
1:E:24:GLN:HA	1:E:73:GLU:O	2.21	0.41
1:A:57:LYS:HB3	1:A:61:PRO:HG3	2.02	0.40
1:E:37:GLN:HB2	1:E:43:LEU:CD1	2.51	0.40
2:B:155:GLN:O	2:B:159:ILE:HG13	2.21	0.40
2:D:33:VAL:O	2:D:85:TYR:HA	2.21	0.40
2:B:216:ASN:HD22	2:B:216:ASN:HA	1.45	0.40
2:D:56:LYS:H	2:D:56:LYS:HD3	1.83	0.40
2:H:38:VAL:O	2:H:82:VAL:HG22	2.22	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	107/112 (96%)	106 (99%)	1 (1%)	0	100 100
1	C	107/112 (96%)	104 (97%)	2 (2%)	1 (1%)	17 12
1	E	107/112 (96%)	104 (97%)	3 (3%)	0	100 100
1	G	107/112 (96%)	102 (95%)	4 (4%)	1 (1%)	17 12
2	B	232/237 (98%)	217 (94%)	13 (6%)	2 (1%)	17 12
2	D	234/237 (99%)	208 (89%)	21 (9%)	5 (2%)	7 3
2	F	235/237 (99%)	215 (92%)	18 (8%)	2 (1%)	17 12
2	H	235/237 (99%)	209 (89%)	22 (9%)	4 (2%)	9 4
All	All	1364/1396 (98%)	1265 (93%)	84 (6%)	15 (1%)	14 9

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	98	LYS
2	D	102	TRP
2	D	122	ASP
2	D	190	ASN
2	D	98	LYS
1	G	3	ALA
2	H	102	TRP
2	B	100	ASN
2	F	57	LYS
2	F	102	TRP
2	H	98	LYS
2	D	8	PRO
2	H	236	ASN
1	C	88	SER
2	H	3	GLN

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	90/93 (97%)	87 (97%)	3 (3%)	38	40
1	C	90/93 (97%)	82 (91%)	8 (9%)	9	6
1	E	90/93 (97%)	82 (91%)	8 (9%)	9	6
1	G	90/93 (97%)	87 (97%)	3 (3%)	38	40
2	B	217/219 (99%)	206 (95%)	11 (5%)	24	22
2	D	218/219 (100%)	202 (93%)	16 (7%)	14	11
2	F	219/219 (100%)	207 (94%)	12 (6%)	21	19
2	H	219/219 (100%)	195 (89%)	24 (11%)	6	3
All	All	1233/1248 (99%)	1148 (93%)	85 (7%)	15	12

All (85) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ARG
1	A	22	SER
1	A	24	GLN
2	B	3	GLN
2	B	34	SER
2	B	59	LYS
2	B	64	VAL
2	B	88	ASN
2	B	104	HIS
2	B	139	ASN
2	B	141	ILE
2	B	189	ASN
2	B	216	ASN
2	B	227	LYS
1	C	14	VAL
1	C	24	GLN
1	C	43	LEU
1	C	52	VAL
1	C	72	GLN
1	C	80	VAL
1	C	83	THR
1	C	94	SER
2	D	24	MET
2	D	43	LYS
2	D	56	LYS
2	D	63	LYS
2	D	68	LEU
2	D	73	LEU
2	D	88	ASN
2	D	92	ASN
2	D	98	LYS
2	D	103	TRP
2	D	104	HIS
2	D	123	ASN
2	D	141	ILE
2	D	162	ARG
2	D	168	LYS
2	D	216	ASN
1	E	9	ARG
1	E	10	ASN
1	E	14	VAL
1	E	18	LYS
1	E	24	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	E	43	LEU
1	E	44	ARG
1	E	52	VAL
2	F	2	SER
2	F	38	VAL
2	F	43	LYS
2	F	49	LEU
2	F	73	LEU
2	F	88	ASN
2	F	98	LYS
2	F	99	ASP
2	F	122	ASP
2	F	126	LEU
2	F	204	ASP
2	F	216	ASN
1	G	9	ARG
1	G	14	VAL
1	G	43	LEU
2	H	1	GLU
2	H	2	SER
2	H	3	GLN
2	H	9	ASP
2	H	43	LYS
2	H	45	LEU
2	H	47	HIS
2	H	56	LYS
2	H	57	LYS
2	H	71	GLU
2	H	75	LYS
2	H	88	ASN
2	H	96	SER
2	H	97	SER
2	H	100	ASN
2	H	104	HIS
2	H	117	GLU
2	H	144	GLU
2	H	147	THR
2	H	189	ASN
2	H	216	ASN
2	H	227	LYS
2	H	230	VAL
2	H	236	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (44) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	28	ASN
1	A	31	ASN
1	A	74	GLN
2	B	3	GLN
2	B	23	ASN
2	B	88	ASN
2	B	116	HIS
2	B	123	ASN
2	B	139	ASN
2	B	189	ASN
2	B	216	ASN
1	C	28	ASN
1	C	30	ASN
1	C	37	GLN
1	C	72	GLN
2	D	47	HIS
2	D	88	ASN
2	D	92	ASN
2	D	104	HIS
2	D	127	GLN
2	D	128	ASN
2	D	139	ASN
2	D	146	GLN
2	D	167	ASN
2	D	170	ASN
2	D	216	ASN
1	E	24	GLN
1	E	74	GLN
2	F	88	ASN
2	F	92	ASN
2	F	100	ASN
2	F	127	GLN
2	F	139	ASN
2	F	216	ASN
1	G	28	ASN
1	G	31	ASN
2	H	3	GLN
2	H	88	ASN
2	H	104	HIS
2	H	146	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	H	170	ASN
2	H	189	ASN
2	H	216	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no 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)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	G	1
1	A	1
1	C	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	G	63:GLY	C	65:TYR	N	1.63
1	A	63:GLY	C	65:TYR	N	1.61

Continued on next page...

Continued from previous page...

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	63:GLY	C	65:TYR	N	1.60

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	109/112 (97%)	-0.26	0	100	100	24, 35, 46, 56	0
1	C	109/112 (97%)	-0.27	1 (0%)	84	86	29, 38, 51, 59	0
1	E	109/112 (97%)	-0.10	3 (2%)	53	59	29, 43, 61, 72	0
1	G	109/112 (97%)	-0.12	2 (1%)	68	72	26, 37, 54, 66	0
2	B	234/237 (98%)	0.04	6 (2%)	56	61	25, 39, 64, 87	0
2	D	236/237 (99%)	0.49	18 (7%)	13	18	26, 48, 79, 92	0
2	F	237/237 (100%)	0.39	12 (5%)	28	33	27, 43, 81, 104	0
2	H	237/237 (100%)	0.44	15 (6%)	20	24	29, 49, 83, 101	0
All	All	1380/1396 (98%)	0.17	57 (4%)	37	43	24, 42, 73, 104	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	102	TRP	12.5
2	H	102	TRP	10.0
2	F	237	GLY	9.4
2	F	101	VAL	8.4
2	H	237	GLY	8.3
2	H	236	ASN	6.9
2	F	104	HIS	6.5
2	H	101	VAL	6.4
2	H	103	TRP	6.3
2	D	237	GLY	6.2
2	F	103	TRP	6.1
2	D	236	ASN	5.4
2	F	100	ASN	5.3
2	H	100	ASN	5.0
2	F	1	GLU	5.0
2	H	104	HIS	4.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	H	1	GLU	4.8
1	G	2	ALA	4.8
2	D	123	ASN	4.2
2	D	99	ASP	4.2
2	D	102	TRP	3.9
2	B	101	VAL	3.8
2	D	103	TRP	3.8
2	D	101	VAL	3.7
2	B	235	LYS	3.4
2	F	236	ASN	3.3
2	B	100	ASN	3.3
2	D	121	PHE	3.1
2	F	97	SER	3.0
2	D	100	ASN	3.0
2	F	56	LYS	3.0
2	D	98	LYS	2.9
2	D	97	SER	2.8
2	D	115	LYS	2.7
2	F	123	ASN	2.6
2	D	211	TYR	2.6
2	H	8	PRO	2.6
2	D	73	LEU	2.6
2	H	7	MET	2.5
1	E	2	ALA	2.5
2	H	138	ARG	2.5
2	H	235	LYS	2.4
2	H	99	ASP	2.4
2	D	191	GLY	2.4
2	D	124	GLY	2.4
1	E	9	ARG	2.3
2	F	124	GLY	2.3
1	C	2	ALA	2.3
2	B	75	LYS	2.2
2	B	115	LYS	2.2
2	D	214	MET	2.2
2	H	195	TRP	2.2
2	D	104	HIS	2.2
2	H	122	ASP	2.2
1	E	3	ALA	2.2
2	B	99	ASP	2.2
1	G	3	ALA	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 monosaccharides 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.