



Full wwPDB X-ray Structure Validation Report

Dec 7, 2023 – 09:52 am GMT

PDB ID : 1GRW
Title : C. elegans major sperm protein
Authors : Baker, A.M.E.; Roberts, T.M.; Stewart, M.
Deposited on : 2001-12-18
Resolution : 2.60 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 : 2.36
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) : 2.36

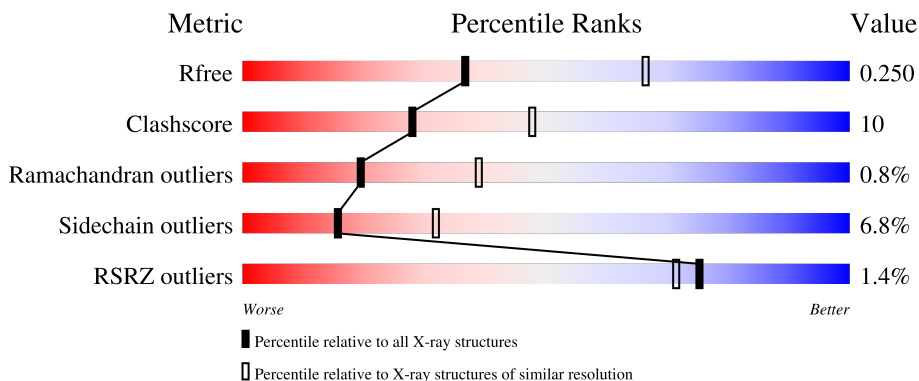
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	126	 60% 29% 9% 2% 2%
1	B	126	 2% 63% 29% 6% 2%
1	C	126	 1% 65% 30% 4% 2%
1	D	126	 2% 67% 26% 5% 2%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3986 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 MAJOR SPERM PROTEIN 31/40/142.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	124	978	614	175	185	4	36	0	0
1	B	124	978	614	175	185	4	56	0	0
1	C	124	978	614	175	185	4	41	0	0
1	D	124	978	614	175	185	4	34	0	0

- Molecule 2 is water.

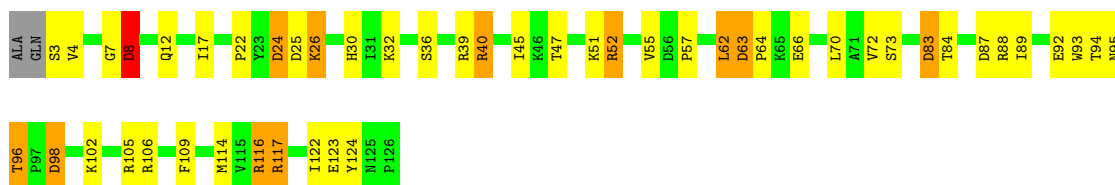
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	27	Total	O	1	1
			27	27		
2	B	14	Total	O	0	0
			14	14		
2	C	15	Total	O	0	0
			15	15		
2	D	18	Total	O	0	0
			18	18		

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.

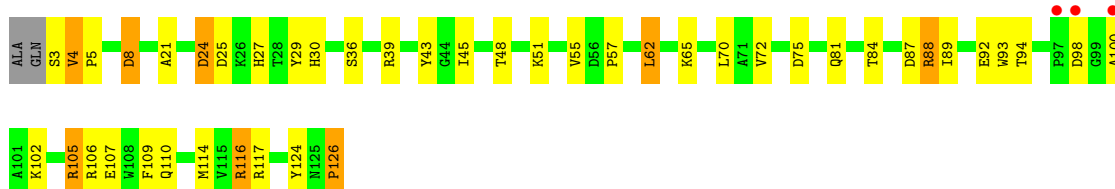
- Molecule 1: MAJOR SPERM PROTEIN 31/40/142

Chain A: 



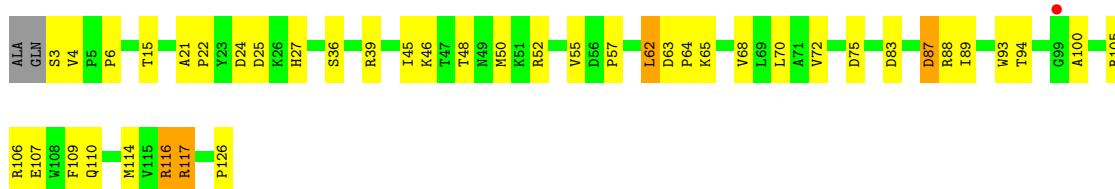
- Molecule 1: MAJOR SPERM PROTEIN 31/40/142

Chain B: 



- Molecule 1: MAJOR SPERM PROTEIN 31/40/142

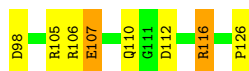
Chain C: 



- Molecule 1: MAJOR SPERM PROTEIN 31/40/142

Chain D: 





4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	53.48Å 53.48Å 457.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.60 37.69 – 2.43	Depositor EDS
% Data completeness (in resolution range)	99.6 (20.00-2.60) 97.5 (37.69-2.43)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.50 (at 2.42Å)	Xtrriage
Refinement program	REFMAC 4.0.0	Depositor
R, R_{free}	0.228 , 0.256 0.224 , 0.250	Depositor DCC
R_{free} test set	1293 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	46.8	Xtrriage
Anisotropy	0.480	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3986	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.12% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality 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	1.12	6/1002 (0.6%)	2.35	27/1360 (2.0%)
1	B	1.19	9/1002 (0.9%)	1.68	14/1360 (1.0%)
1	C	1.09	7/1002 (0.7%)	1.65	16/1360 (1.2%)
1	D	1.26	6/1002 (0.6%)	1.63	18/1360 (1.3%)
All	All	1.17	28/4008 (0.7%)	1.85	75/5440 (1.4%)

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	82	GLU	CB-CG	23.96	1.97	1.52
1	B	92	GLU	CG-CD	-18.60	1.24	1.51
1	C	117	ARG	CB-CG	-16.68	1.07	1.52
1	A	3	SER	C-N	15.36	1.69	1.34
1	D	105	ARG	NE-CZ	-14.53	1.14	1.33
1	B	39	ARG	CD-NE	-14.43	1.22	1.46
1	C	39	ARG	CG-CD	-13.85	1.17	1.51
1	A	39	ARG	CZ-NH1	13.02	1.50	1.33
1	D	106	ARG	CB-CG	-9.44	1.27	1.52
1	C	83	ASP	CB-CG	-9.07	1.32	1.51
1	A	39	ARG	CZ-NH2	-8.95	1.21	1.33
1	A	88	ARG	NE-CZ	-8.66	1.21	1.33
1	B	107	GLU	CB-CG	-8.37	1.36	1.52
1	B	51	LYS	CB-CG	8.35	1.75	1.52
1	D	51	LYS	CB-CG	8.35	1.75	1.52
1	A	3	SER	CA-C	8.29	1.74	1.52
1	C	46	LYS	CB-CG	7.30	1.72	1.52
1	B	126	PRO	N-CD	7.18	1.57	1.47
1	A	83	ASP	CB-CG	6.55	1.65	1.51
1	C	65	LYS	CD-CE	6.43	1.67	1.51
1	B	3	SER	C-N	-6.25	1.19	1.34
1	B	81	GLN	CB-CG	-6.22	1.35	1.52
1	B	75	ASP	CB-CG	-6.04	1.39	1.51
1	C	50	MET	CB-CG	-5.96	1.32	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	126	PRO	N-CD	5.89	1.56	1.47
1	B	110	GLN	CB-CG	5.61	1.67	1.52
1	D	126	PRO	N-CD	5.33	1.55	1.47
1	D	26	LYS	CD-CE	-5.17	1.38	1.51

All (75) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	88	ARG	NE-CZ-NH2	-25.57	107.52	120.30
1	A	3	SER	O-C-N	-25.42	82.03	122.70
1	A	117	ARG	NE-CZ-NH1	24.69	132.65	120.30
1	A	88	ARG	NE-CZ-NH1	22.09	131.35	120.30
1	A	3	SER	C-N-CA	20.25	172.34	121.70
1	A	39	ARG	NE-CZ-NH2	19.54	130.07	120.30
1	B	3	SER	O-C-N	-18.57	92.99	122.70
1	B	3	SER	C-N-CA	18.43	167.78	121.70
1	A	3	SER	CA-C-N	17.20	155.03	117.20
1	A	52	ARG	NE-CZ-NH2	17.09	128.85	120.30
1	C	117	ARG	CA-CB-CG	16.08	148.78	113.40
1	A	117	ARG	NE-CZ-NH2	-14.34	113.13	120.30
1	C	3	SER	C-N-CA	-13.96	86.80	121.70
1	A	39	ARG	NE-CZ-NH1	-13.47	113.56	120.30
1	B	3	SER	CA-C-N	12.55	144.81	117.20
1	D	39	ARG	NE-CZ-NH1	12.18	126.39	120.30
1	D	51	LYS	CA-CB-CG	-11.78	87.48	113.40
1	D	83	ASP	CB-CG-OD1	-11.77	107.71	118.30
1	A	117	ARG	CD-NE-CZ	10.60	138.44	123.60
1	D	105	ARG	NE-CZ-NH1	10.24	125.42	120.30
1	C	3	SER	O-C-N	10.12	138.90	122.70
1	B	116	ARG	NE-CZ-NH1	-10.12	115.24	120.30
1	A	52	ARG	NE-CZ-NH1	-9.79	115.41	120.30
1	B	39	ARG	CG-CD-NE	9.75	132.27	111.80
1	C	88	ARG	NE-CZ-NH1	-9.73	115.44	120.30
1	C	116	ARG	NE-CZ-NH2	9.41	125.00	120.30
1	A	116	ARG	NE-CZ-NH1	-9.24	115.68	120.30
1	D	8	ASP	CB-CG-OD2	9.19	126.57	118.30
1	C	3	SER	CA-C-N	-8.89	97.63	117.20
1	B	8	ASP	CB-CG-OD1	8.81	126.23	118.30
1	D	4	VAL	N-CA-CB	8.37	129.92	111.50
1	C	117	ARG	CB-CG-CD	8.05	132.52	111.60
1	B	105	ARG	NE-CZ-NH2	7.83	124.22	120.30
1	C	116	ARG	NE-CZ-NH1	-7.81	116.40	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	88	ARG	CD-NE-CZ	-7.75	112.75	123.60
1	C	52	ARG	NE-CZ-NH1	-7.63	116.48	120.30
1	D	83	ASP	CB-CG-OD2	7.60	125.14	118.30
1	A	105	ARG	NE-CZ-NH2	7.58	124.09	120.30
1	B	88	ARG	NE-CZ-NH2	-7.42	116.59	120.30
1	B	106	ARG	NE-CZ-NH2	7.41	124.01	120.30
1	A	63	ASP	CB-CG-OD2	-7.30	111.73	118.30
1	A	8	ASP	CB-CG-OD1	7.12	124.70	118.30
1	D	40	ARG	NE-CZ-NH1	-7.11	116.75	120.30
1	B	75	ASP	CB-CG-OD2	-7.10	111.91	118.30
1	A	106	ARG	NE-CZ-NH2	7.08	123.84	120.30
1	D	40	ARG	NE-CZ-NH2	-6.96	116.82	120.30
1	B	92	GLU	CA-CB-CG	6.94	128.68	113.40
1	D	8	ASP	CB-CG-OD1	-6.90	112.09	118.30
1	A	25	ASP	CB-CG-OD1	-6.89	112.10	118.30
1	C	88	ARG	NE-CZ-NH2	6.85	123.73	120.30
1	C	106	ARG	NE-CZ-NH2	6.63	123.62	120.30
1	A	40	ARG	NE-CZ-NH2	-6.52	117.04	120.30
1	D	52	ARG	NE-CZ-NH2	-6.51	117.05	120.30
1	C	39	ARG	NE-CZ-NH2	6.50	123.55	120.30
1	A	24	ASP	CB-CG-OD1	6.43	124.09	118.30
1	D	40	ARG	NH1-CZ-NH2	6.39	126.43	119.40
1	A	51	LYS	CA-CB-CG	-6.37	99.38	113.40
1	D	116	ARG	NE-CZ-NH1	-6.35	117.12	120.30
1	C	75	ASP	CB-CG-OD1	6.30	123.97	118.30
1	D	83	ASP	CA-CB-CG	6.24	127.14	113.40
1	A	7	GLY	CA-C-O	6.23	131.82	120.60
1	D	62	LEU	CA-CB-CG	6.20	129.55	115.30
1	A	123	GLU	OE1-CD-OE2	6.13	130.65	123.30
1	D	112	ASP	CB-CG-OD2	-5.83	113.05	118.30
1	D	4	VAL	CA-C-O	5.76	132.19	120.10
1	C	50	MET	CG-SD-CE	5.68	109.29	100.20
1	D	82	GLU	CA-CB-CG	-5.53	101.23	113.40
1	B	3	SER	N-CA-C	-5.51	96.13	111.00
1	A	116	ARG	NE-CZ-NH2	5.49	123.05	120.30
1	A	83	ASP	CA-CB-CG	-5.33	101.66	113.40
1	A	98	ASP	CA-CB-CG	-5.23	101.89	113.40
1	B	29	TYR	CB-CG-CD2	-5.11	117.93	121.00
1	C	87	ASP	CB-CG-OD2	-5.11	113.70	118.30
1	C	117	ARG	NE-CZ-NH1	5.08	122.84	120.30
1	B	24	ASP	CB-CG-OD2	-5.03	113.77	118.30

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	978	0	955	25	4
1	B	978	0	954	18	0
1	C	978	0	955	17	0
1	D	978	0	955	15	4
2	A	27	0	0	2	0
2	B	14	0	0	0	0
2	C	15	0	0	0	0
2	D	18	0	0	1	0
All	All	3986	0	3819	73	4

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 10.

All (73) 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:A:40:ARG:HD2	1:A:96:THR:HG21	1.55	0.88
1:C:105:ARG:HB3	1:C:107:GLU:OE1	1.75	0.87
1:A:26:LYS:O	2:A:2006:HOH:O	1.94	0.85
1:D:45:ILE:HG21	1:D:55:VAL:HG11	1.65	0.78
1:B:88:ARG:HH22	1:C:110:GLN:HB3	1.54	0.72
1:B:57:PRO:HD2	1:B:70:LEU:HD23	1.72	0.70
1:B:4:VAL:HG23	1:B:5:PRO:HD2	1.74	0.69
1:A:93:TRP:CZ2	1:A:116:ARG:HD3	2.28	0.68
1:A:72:VAL:HG11	1:A:89:ILE:HD11	1.76	0.67
1:A:12:GLN:OE1	1:A:32:LYS:HD3	1.95	0.66
1:C:21:ALA:HB1	1:C:22:PRO:HA	1.77	0.66
1:B:45:ILE:HG21	1:B:55:VAL:HG11	1.76	0.65
1:D:93:TRP:CZ2	1:D:116:ARG:HD3	2.33	0.64
1:A:52:ARG:NH1	1:A:87:ASP:OD2	2.31	0.63
1:B:72:VAL:HG11	1:B:89:ILE:HD11	1.80	0.63
1:A:57:PRO:HD2	1:A:70:LEU:HD23	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94:THR:HG22	1:A:109:PHE:HE1	1.65	0.61
1:B:94:THR:HG22	1:B:109:PHE:HE1	1.64	0.61
1:A:84:THR:HB	1:A:124:TYR:CD2	2.35	0.61
1:B:94:THR:HG22	1:B:109:PHE:CE1	2.37	0.60
1:D:3:SER:OG	1:D:4:VAL:N	2.38	0.56
1:A:40:ARG:CD	1:A:96:THR:HG21	2.32	0.56
1:C:94:THR:HG22	1:C:109:PHE:HE1	1.71	0.56
1:B:84:THR:HG23	1:B:124:TYR:CD2	2.41	0.55
1:A:45:ILE:HG21	1:A:55:VAL:HG11	1.88	0.54
1:D:48:THR:OG1	1:D:87:ASP:HA	2.08	0.53
1:D:8:ASP:OD2	1:D:8:ASP:N	2.42	0.53
1:A:22:PRO:HB2	1:A:24:ASP:OD1	2.09	0.52
1:A:63:ASP:HB3	1:A:64:PRO:HD2	1.90	0.52
1:C:57:PRO:HD2	1:C:70:LEU:HD23	1.90	0.52
1:D:10:GLN:HG2	1:D:34:ILE:HD12	1.90	0.52
1:D:21:ALA:HB1	1:D:22:PRO:HA	1.91	0.52
1:B:93:TRP:CZ2	1:B:116:ARG:HD3	2.45	0.51
1:A:89:ILE:CD1	1:A:122:ILE:HD11	2.39	0.51
1:C:72:VAL:HG11	1:C:89:ILE:HD11	1.92	0.51
1:D:78:ALA:HB3	1:D:81:GLN:HB2	1.92	0.51
1:D:107:GLU:HA	1:D:110:GLN:NE2	2.26	0.51
1:C:48:THR:OG1	1:C:87:ASP:HA	2.12	0.49
1:A:8:ASP:HB3	2:A:2002:HOH:O	2.13	0.49
1:C:15:THR:HB	2:D:2018:HOH:O	2.12	0.49
1:B:30:HIS:HA	1:B:70:LEU:O	2.12	0.48
1:C:94:THR:HG22	1:C:109:PHE:CE1	2.49	0.48
1:A:62:LEU:HD12	1:A:66:GLU:HG2	1.96	0.48
1:A:92:GLU:OE2	1:A:117:ARG:NH2	2.46	0.47
1:C:93:TRP:CZ2	1:C:116:ARG:HD3	2.49	0.47
1:C:62:LEU:HD11	1:C:68:VAL:HB	1.97	0.47
1:C:25:ASP:HB2	1:C:27:HIS:CE1	2.50	0.47
1:B:5:PRO:HB3	1:B:114:MET:SD	2.55	0.46
1:C:107:GLU:CD	1:C:107:GLU:H	2.18	0.46
1:D:6:PRO:HG2	1:D:116:ARG:HB2	1.99	0.46
1:D:12:GLN:OE1	1:D:32:LYS:HD3	2.16	0.45
1:D:15:THR:O	1:D:16:LYS:HB3	2.17	0.45
1:A:94:THR:HG22	1:A:109:PHE:CE1	2.47	0.45
1:A:17:ILE:CG2	1:A:122:ILE:HD13	2.47	0.44
1:C:63:ASP:HB3	1:C:64:PRO:HD2	1.99	0.44
1:D:43:TYR:CD2	1:D:62:LEU:HD22	2.52	0.43
1:A:8:ASP:OD1	1:A:8:ASP:N	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:6:PRO:HG2	1:C:116:ARG:HB2	2.01	0.43
1:B:43:TYR:CD2	1:B:62:LEU:HD22	2.55	0.42
1:B:48:THR:OG1	1:B:87:ASP:HA	2.19	0.42
1:B:25:ASP:O	1:B:27:HIS:CD2	2.73	0.42
1:A:52:ARG:HH11	1:A:52:ARG:HD2	1.62	0.42
1:A:4:VAL:HG23	1:A:95:ASN:HD22	1.85	0.41
1:A:30:HIS:HA	1:A:70:LEU:O	2.20	0.41
1:D:30:HIS:HA	1:D:70:LEU:O	2.21	0.41
1:B:21:ALA:HB2	1:B:126:PRO:HA	2.02	0.41
1:A:102:LYS:HG2	1:A:102:LYS:O	2.20	0.40
1:A:63:ASP:HB3	1:A:64:PRO:CD	2.52	0.40
1:B:48:THR:HG21	1:B:88:ARG:HG2	2.03	0.40
1:D:57:PRO:HD2	1:D:70:LEU:HD23	2.03	0.40
1:C:45:ILE:HG21	1:C:55:VAL:HG11	2.03	0.40
1:B:72:VAL:HG11	1:B:89:ILE:CD1	2.48	0.40
1:B:88:ARG:NH2	1:C:110:GLN:HB3	2.29	0.40

All (4) 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)
1:A:83:ASP:OD2	1:D:98:ASP:OD2[7_545]	1.53	0.67
1:A:83:ASP:OD1	1:D:98:ASP:OD1[7_545]	2.01	0.19
1:A:83:ASP:CG	1:D:98:ASP:OD2[7_545]	2.04	0.16
1:A:83:ASP:OD1	1:D:98:ASP:OD2[7_545]	2.12	0.08

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	122/126 (97%)	118 (97%)	4 (3%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	122/126 (97%)	114 (93%)	6 (5%)	2 (2%)	9	19
1	C	122/126 (97%)	118 (97%)	3 (2%)	1 (1%)	19	39
1	D	122/126 (97%)	117 (96%)	4 (3%)	1 (1%)	19	39
All	All	488/504 (97%)	467 (96%)	17 (4%)	4 (1%)	19	39

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	100	ALA
1	C	100	ALA
1	B	4	VAL
1	D	81	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	106/107 (99%)	97 (92%)	9 (8%)	10	21
1	B	106/107 (99%)	97 (92%)	9 (8%)	10	21
1	C	106/107 (99%)	100 (94%)	6 (6%)	20	41
1	D	106/107 (99%)	101 (95%)	5 (5%)	26	50
All	All	424/428 (99%)	395 (93%)	29 (7%)	16	32

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	ASP
1	A	26	LYS
1	A	36	SER
1	A	47	THR
1	A	62	LEU
1	A	73	SER
1	A	96	THR

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Mol	Chain	Res	Type
1	A	98	ASP
1	A	114	MET
1	B	8	ASP
1	B	24	ASP
1	B	36	SER
1	B	62	LEU
1	B	65	LYS
1	B	98	ASP
1	B	102	LYS
1	B	105	ARG
1	B	117	ARG
1	C	4	VAL
1	C	24	ASP
1	C	36	SER
1	C	62	LEU
1	C	114	MET
1	C	117	ARG
1	D	8	ASP
1	D	36	SER
1	D	62	LEU
1	D	83	ASP
1	D	107	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	27	HIS
1	C	27	HIS
1	C	110	GLN
1	D	110	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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	3:SER	C	4:VAL	N	1.69
1	B	3:SER	C	4:VAL	N	1.19

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	123/126 (97%)	-0.23	0 100 100	33, 47, 66, 88	8 (6%)
1	B	123/126 (97%)	-0.05	3 (2%) 59 53	34, 52, 78, 88	12 (9%)
1	C	123/126 (97%)	-0.27	1 (0%) 86 84	38, 50, 75, 90	9 (7%)
1	D	124/126 (98%)	-0.23	3 (2%) 59 53	33, 48, 78, 104	10 (8%)
All	All	493/504 (97%)	-0.20	7 (1%) 75 71	33, 50, 77, 104	39 (7%)

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	83	ASP	3.8
1	D	84	THR	3.3
1	B	100	ALA	3.2
1	B	98	ASP	2.7
1	B	97	PRO	2.5
1	C	99	GLY	2.5
1	D	81	GLN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no 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.