



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 26, 2024 – 06:08 PM JST

PDB ID : 8W8G
Title : Crystal structure of human TRF1 with PinX1
Authors : Lei, M.; Wu, J.
Deposited on : 2023-09-02
Resolution : 2.70 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
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 3.0
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.002 (Gargrove)
Density-Fitness : 1.0.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.38.2

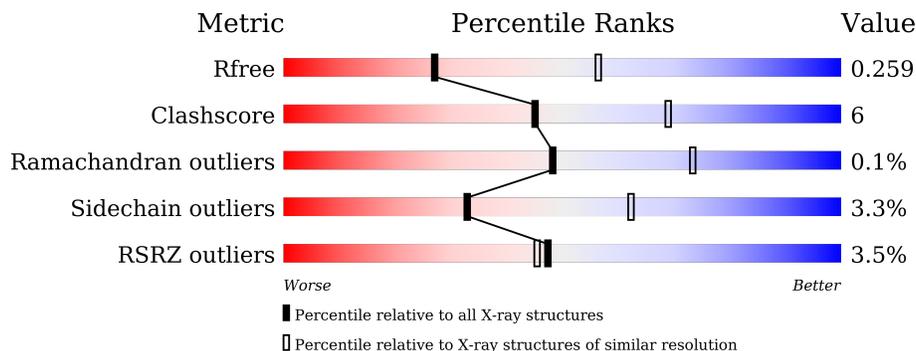
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	164625	3333 (2.70-2.70)
Clashscore	180529	3684 (2.70-2.70)
Ramachandran outliers	177936	3633 (2.70-2.70)
Sidechain outliers	177891	3633 (2.70-2.70)
RSRZ outliers	164620	3333 (2.70-2.70)

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	212	 4% (red), 80% (green), 16% (yellow), .. (grey)
1	B	212	 6% (red), 84% (green), 11% (yellow), .. (grey)
1	C	212	 % (red), 83% (green), 13% (yellow), .. (grey)
1	D	212	 2% (red), 79% (green), 13% (yellow), 8% (grey)
2	E	20	 60% (green), 5% (yellow), 35% (grey)
2	F	20	 40% (green), 15% (orange), 45% (grey)

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Mol	Chain	Length	Quality of chain
2	G	20	 55% 15% 5% 25%
2	H	20	 5% 50% 15% 35%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7064 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.

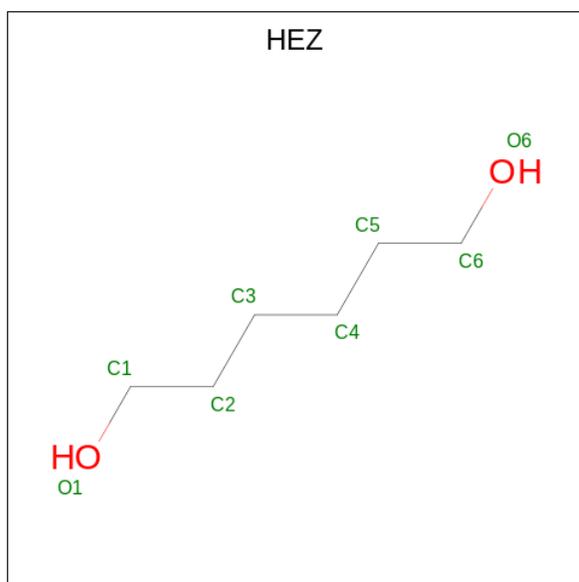
- Molecule 1 is a protein called Telomeric repeat-binding factor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	206	1648	1046	281	308	13	0	0	0
1	B	206	1648	1046	281	308	13	0	0	0
1	C	208	1666	1056	286	311	13	0	0	0
1	D	196	1575	1003	267	293	12	0	0	0

- Molecule 2 is a protein called PIN2/TERF1-interacting telomerase inhibitor 1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	E	13	111	69	25	17	0	0	0
2	F	11	101	64	23	14	0	0	0
2	G	15	124	77	28	19	0	0	0
2	H	13	111	69	25	17	0	0	0

- Molecule 3 is HEXANE-1,6-DIOL (three-letter code: HEZ) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 8 6 2	0	0
3	B	1	Total C O 8 6 2	0	0
3	C	1	Total C O 8 6 2	0	0
3	D	1	Total C O 8 6 2	0	0

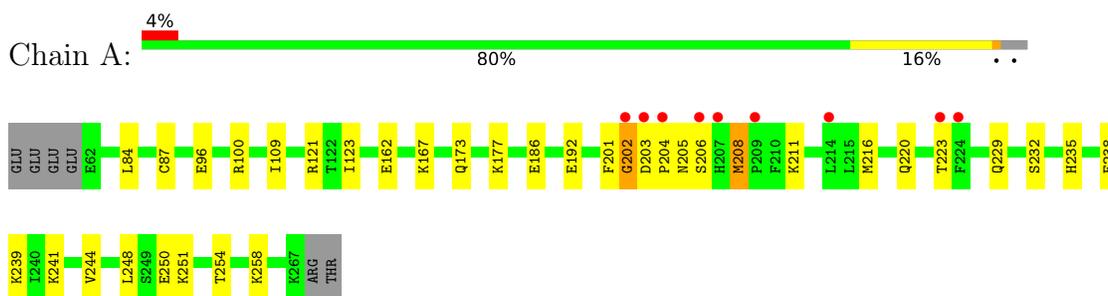
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	13	Total O 13 13	0	0
4	B	10	Total O 10 10	0	0
4	C	12	Total O 12 12	0	0
4	D	9	Total O 9 9	0	0
4	E	2	Total O 2 2	0	0
4	G	1	Total O 1 1	0	0
4	H	1	Total O 1 1	0	0

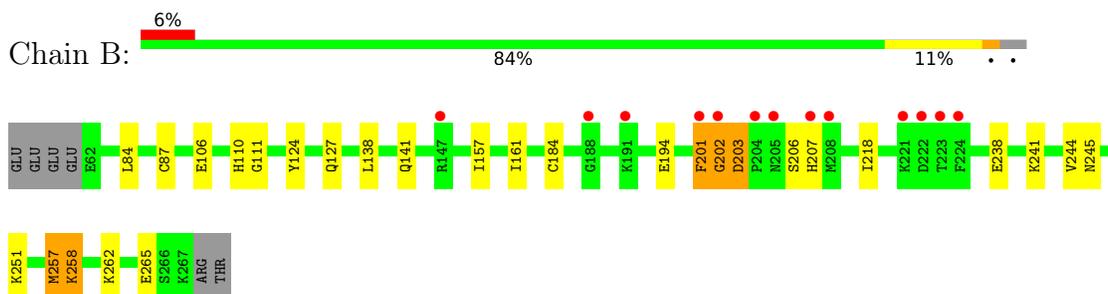
3 Residue-property plots

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.

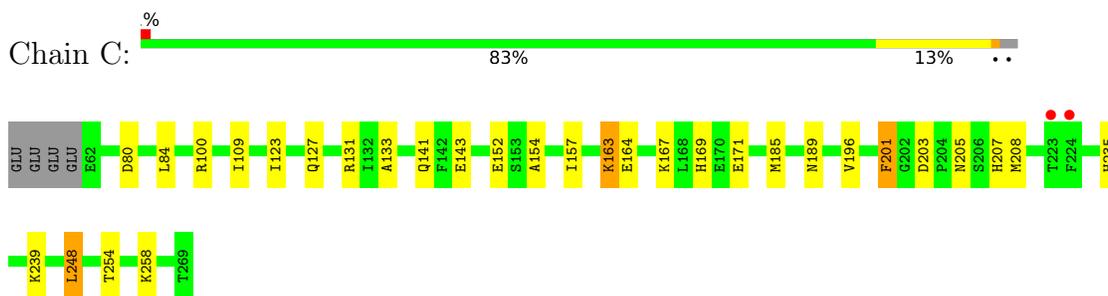
- Molecule 1: Telomeric repeat-binding factor 1



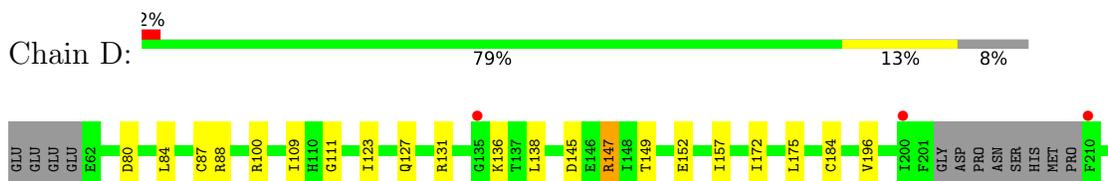
- Molecule 1: Telomeric repeat-binding factor 1



- Molecule 1: Telomeric repeat-binding factor 1



- Molecule 1: Telomeric repeat-binding factor 1





- Molecule 2: PIN2/TERF1-interacting telomerase inhibitor 1



- Molecule 2: PIN2/TERF1-interacting telomerase inhibitor 1



- Molecule 2: PIN2/TERF1-interacting telomerase inhibitor 1



- Molecule 2: PIN2/TERF1-interacting telomerase inhibitor 1



4 Data and refinement statistics i

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	87.40Å 89.19Å 113.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.95 – 2.70 47.95 – 2.70	Depositor EDS
% Data completeness (in resolution range)	96.7 (47.95-2.70) 96.7 (47.95-2.70)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.58 (at 2.69Å)	Xtrriage
Refinement program	PHENIX (1.19.2_4158: ???)	Depositor
R, R_{free}	0.227 , 0.263 0.229 , 0.259	Depositor DCC
R_{free} test set	1251 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	46.7	Xtrriage
Anisotropy	0.293	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 42.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.026 for k,h,-l	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7064	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 36.99 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.5957e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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](#)

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

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	0.26	0/1679	0.44	1/2255 (0.0%)
1	B	0.26	0/1679	0.69	6/2255 (0.3%)
1	C	0.23	0/1697	0.62	4/2279 (0.2%)
1	D	0.24	0/1602	0.42	0/2149
2	E	0.23	0/112	0.70	0/145
2	F	0.22	0/102	0.66	0/132
2	G	0.40	0/125	0.50	0/161
2	H	0.24	0/112	0.62	0/145
All	All	0.25	0/7108	0.56	11/9521 (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	B	0	1

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	201	PHE	CB-CA-C	-14.28	81.84	110.40
1	B	201	PHE	CB-CA-C	-13.46	83.48	110.40
1	B	201	PHE	N-CA-C	12.49	144.73	111.00
1	B	203	ASP	N-CA-C	11.28	141.47	111.00
1	C	203	ASP	N-CA-CB	-11.28	90.29	110.60
1	B	203	ASP	N-CA-CB	-9.90	92.77	110.60
1	C	203	ASP	N-CA-C	9.54	136.75	111.00
1	B	202	GLY	N-CA-C	-8.86	90.95	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	201	PHE	N-CA-C	6.63	128.90	111.00
1	B	203	ASP	CB-CA-C	-5.59	99.22	110.40
1	A	202	GLY	N-CA-C	-5.52	99.31	113.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	203	ASP	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	1648	0	1631	30	0
1	B	1648	0	1631	17	0
1	C	1666	0	1651	16	1
1	D	1575	0	1564	19	0
2	E	111	0	123	1	0
2	F	101	0	115	4	0
2	G	124	0	139	3	1
2	H	111	0	123	2	0
3	A	8	0	14	1	0
3	B	8	0	14	1	0
3	C	8	0	14	1	0
3	D	8	0	14	1	0
4	A	13	0	0	1	0
4	B	10	0	0	0	0
4	C	12	0	0	0	0
4	D	9	0	0	0	0
4	E	2	0	0	0	0
4	G	1	0	0	0	0
4	H	1	0	0	0	0
All	All	7064	0	7033	82	1

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 6.

All (82) 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:201:PHE:O	1:A:204:PRO:CD	1.97	1.13
1:A:201:PHE:O	1:A:204:PRO:HD3	1.54	1.07
1:B:201:PHE:C	1:B:202:GLY:O	2.11	0.80
1:B:194:GLU:HG3	1:B:218:ILE:HG21	1.71	0.72
1:A:202:GLY:O	1:A:211:LYS:HE3	1.91	0.71
1:A:202:GLY:O	1:A:211:LYS:CE	2.42	0.67
2:F:288:ASP:OD1	2:F:288:ASP:N	2.27	0.64
1:A:201:PHE:O	1:A:204:PRO:HD2	1.94	0.64
1:D:172:ILE:HG12	1:D:239:LYS:HB3	1.77	0.64
1:B:201:PHE:O	1:B:202:GLY:O	2.17	0.63
1:A:203:ASP:N	1:A:204:PRO:HD3	2.15	0.62
1:B:138:LEU:HD11	1:B:184:CYS:HA	1.81	0.61
1:C:141:GLN:HB2	2:G:297:ARG:HG3	1.82	0.60
1:A:201:PHE:O	1:A:204:PRO:CG	2.50	0.59
1:C:127:GLN:HB3	1:C:157:ILE:HD13	1.85	0.58
1:D:88:ARG:HG3	1:D:248:LEU:HD11	1.85	0.58
1:D:138:LEU:HD11	1:D:184:CYS:HA	1.86	0.57
1:B:124:TYR:HB3	1:B:161:ILE:HG12	1.86	0.56
1:A:186:GLU:HA	1:A:223:THR:HG23	1.87	0.56
1:C:152:GLU:OE2	1:C:196:VAL:HG22	2.05	0.56
1:A:202:GLY:C	1:A:204:PRO:HD3	2.26	0.55
1:B:241:LYS:HD2	3:B:301:HEZ:H62	1.89	0.55
1:C:205:ASN:O	1:C:207:HIS:ND1	2.28	0.55
1:C:84:LEU:HD12	1:C:248:LEU:HD12	1.89	0.54
1:D:80:ASP:HB3	1:D:254:THR:HG21	1.90	0.54
1:B:141:GLN:HB2	2:F:297:ARG:HA	1.91	0.53
1:D:84:LEU:HD12	1:D:248:LEU:HD12	1.89	0.53
1:B:87:CYS:HB3	1:B:244:VAL:HG13	1.89	0.53
1:B:110:HIS:HA	2:F:287:ARG:N	2.24	0.53
1:B:84:LEU:HD13	1:B:251:LYS:HG3	1.91	0.52
1:A:87:CYS:HB3	1:A:244:VAL:HG13	1.92	0.52
1:D:138:LEU:O	2:H:297:ARG:NH2	2.37	0.52
1:A:192:GLU:OE2	2:E:297:ARG:NH2	2.38	0.51
1:D:87:CYS:HB2	1:D:248:LEU:HD13	1.92	0.51
1:D:175:LEU:HD13	1:D:236:MET:HA	1.92	0.51
1:C:100:ARG:HD3	1:D:111:GLY:HA2	1.93	0.51
1:A:216:MET:O	1:A:220:GLN:HG2	2.11	0.51
1:B:84:LEU:HD23	1:B:257:MET:HG3	1.92	0.51
1:C:109:ILE:HD13	1:C:123:ILE:HG23	1.93	0.50
1:B:241:LYS:O	1:B:245:ASN:ND2	2.45	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:131:ARG:HB3	1:C:154:ALA:HB2	1.94	0.49
1:A:173:GLN:HG2	1:A:177:LYS:HE3	1.94	0.48
1:C:201:PHE:HE1	1:C:208:MET:HG3	1.77	0.48
1:D:239:LYS:HD2	3:D:301:HEZ:H42	1.95	0.48
1:C:167:LYS:HE3	1:C:171:GLU:OE2	2.14	0.47
1:A:109:ILE:HD13	1:A:123:ILE:HG23	1.97	0.47
1:A:121:ARG:NH2	1:A:250:GLU:OE1	2.42	0.47
1:C:189:ASN:OD1	1:C:189:ASN:N	2.46	0.47
1:A:84:LEU:HD11	1:A:251:LYS:HB3	1.96	0.47
2:F:287:ARG:HB3	2:F:288:ASP:H	1.46	0.47
1:B:127:GLN:HB3	1:B:157:ILE:HD13	1.97	0.47
1:A:121:ARG:HD2	1:A:162:GLU:HG3	1.97	0.47
1:D:149:THR:HA	2:H:297:ARG:HH22	1.79	0.47
1:C:254:THR:O	1:C:258:LYS:HB2	2.15	0.46
1:D:152:GLU:OE2	1:D:196:VAL:HG22	2.15	0.46
1:C:80:ASP:HB3	1:C:254:THR:HG21	1.97	0.46
1:A:258:LYS:NZ	4:A:401:HOH:O	2.48	0.46
1:A:96:GLU:HB3	1:A:100:ARG:NH2	2.31	0.45
1:D:109:ILE:HD13	1:D:123:ILE:HG23	1.97	0.45
1:D:127:GLN:HB3	1:D:157:ILE:HD13	1.99	0.45
1:D:100:ARG:HE	2:G:287:ARG:HE	1.64	0.45
1:B:157:ILE:O	1:B:161:ILE:HG13	2.17	0.45
1:D:138:LEU:HD23	1:D:138:LEU:HA	1.86	0.45
1:A:203:ASP:HA	1:A:211:LYS:HZ1	1.82	0.44
1:D:127:GLN:O	1:D:131:ARG:HG2	2.16	0.44
1:A:203:ASP:N	1:A:204:PRO:CD	2.81	0.44
1:B:207:HIS:O	1:B:207:HIS:CG	2.71	0.44
1:B:258:LYS:HD2	1:B:262:LYS:HE2	2.00	0.44
1:A:100:ARG:HD2	1:B:111:GLY:HA2	2.00	0.43
1:A:254:THR:O	1:A:258:LYS:HB2	2.18	0.43
1:A:248:LEU:HD12	1:A:248:LEU:HA	1.87	0.43
1:C:235:HIS:NE2	1:C:239:LYS:HE2	2.33	0.43
1:A:208:MET:HA	1:A:208:MET:HE2	2.01	0.43
1:D:100:ARG:HE	2:G:287:ARG:NE	2.17	0.42
1:D:145:ASP:OD1	1:D:147:ARG:HD2	2.19	0.42
1:C:163:LYS:HE3	1:C:169:HIS:ND1	2.34	0.41
1:A:206:SER:HA	1:A:211:LYS:NZ	2.35	0.41
1:A:229:GLN:O	1:A:232:SER:HB3	2.21	0.41
1:A:241:LYS:HZ2	3:A:301:HEZ:H41	1.86	0.41
1:C:133:ALA:HB1	3:C:301:HEZ:H11	2.01	0.41
1:A:235:HIS:CE1	1:A:239:LYS:HE2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:206:SER:HA	1:A:211:LYS:HZ3	1.85	0.40

All (1) 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:C:164:GLU:OE2	2:G:299:LYS:O[4_454]	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	204/212 (96%)	192 (94%)	12 (6%)	0	100	100
1	B	204/212 (96%)	196 (96%)	8 (4%)	0	100	100
1	C	206/212 (97%)	197 (96%)	9 (4%)	0	100	100
1	D	192/212 (91%)	186 (97%)	5 (3%)	1 (0%)	25	49
2	E	11/20 (55%)	9 (82%)	2 (18%)	0	100	100
2	F	9/20 (45%)	8 (89%)	1 (11%)	0	100	100
2	G	13/20 (65%)	11 (85%)	2 (15%)	0	100	100
2	H	11/20 (55%)	10 (91%)	1 (9%)	0	100	100
All	All	850/928 (92%)	809 (95%)	40 (5%)	1 (0%)	48	73

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	136	LYS

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	178/184 (97%)	174 (98%)	4 (2%)	47	76
1	B	178/184 (97%)	172 (97%)	6 (3%)	32	61
1	C	180/184 (98%)	176 (98%)	4 (2%)	47	76
1	D	169/184 (92%)	165 (98%)	4 (2%)	44	73
2	E	12/18 (67%)	12 (100%)	0	100	100
2	F	11/18 (61%)	8 (73%)	3 (27%)	0	1
2	G	13/18 (72%)	11 (85%)	2 (15%)	2	6
2	H	12/18 (67%)	10 (83%)	2 (17%)	2	5
All	All	753/808 (93%)	728 (97%)	25 (3%)	33	62

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	167	LYS
1	A	205	ASN
1	A	208	MET
1	A	238	GLU
1	B	106	GLU
1	B	206	SER
1	B	238	GLU
1	B	257	MET
1	B	258	LYS
1	B	265	GLU
1	C	143	GLU
1	C	163	LYS
1	C	185	MET
1	C	248	LEU
1	D	147	ARG
1	D	214	LEU
1	D	216	MET
1	D	238	GLU
2	F	287	ARG

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Mol	Chain	Res	Type
2	F	288	ASP
2	F	297	ARG
2	G	290	THR
2	G	297	ARG
2	H	287	ARG
2	H	290	THR

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

Mol	Chain	Res	Type
1	B	173	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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	$\# Z > 2$
3	HEZ	A	301	-	7,7,7	0.10	0	6,6,6	0.15	0
3	HEZ	D	301	-	7,7,7	0.11	0	6,6,6	0.11	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	HEZ	C	301	-	7,7,7	0.11	0	6,6,6	0.15	0
3	HEZ	B	301	-	7,7,7	0.09	0	6,6,6	0.19	0

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
3	HEZ	A	301	-	-	2/5/5/5	-
3	HEZ	D	301	-	-	3/5/5/5	-
3	HEZ	C	301	-	-	2/5/5/5	-
3	HEZ	B	301	-	-	2/5/5/5	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	301	HEZ	C2-C3-C4-C5
3	D	301	HEZ	C1-C2-C3-C4
3	B	301	HEZ	O1-C1-C2-C3
3	D	301	HEZ	C4-C5-C6-O6
3	B	301	HEZ	C3-C4-C5-C6
3	A	301	HEZ	C1-C2-C3-C4
3	D	301	HEZ	C3-C4-C5-C6
3	C	301	HEZ	C2-C3-C4-C5
3	C	301	HEZ	C1-C2-C3-C4

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	301	HEZ	1	0
3	D	301	HEZ	1	0
3	C	301	HEZ	1	0
3	B	301	HEZ	1	0

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, 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	206/212 (97%)	0.13	9 (4%) 39 38	32, 49, 89, 114	0
1	B	206/212 (97%)	0.34	13 (6%) 27 25	34, 51, 93, 137	0
1	C	208/212 (98%)	0.16	2 (0%) 79 79	33, 49, 102, 149	0
1	D	196/212 (92%)	0.20	5 (2%) 57 56	32, 53, 80, 112	0
2	E	13/20 (65%)	0.58	0 100 100	40, 53, 82, 101	0
2	F	11/20 (55%)	0.67	0 100 100	45, 53, 96, 104	0
2	G	15/20 (75%)	0.75	0 100 100	40, 57, 99, 104	0
2	H	13/20 (65%)	0.38	1 (7%) 21 19	38, 40, 85, 90	0
All	All	868/928 (93%)	0.23	30 (3%) 47 45	32, 51, 93, 149	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	204	PRO	3.5
1	D	210	PHE	3.3
1	A	204	PRO	3.2
1	A	202	GLY	3.2
1	A	206	SER	3.2
1	B	224	PHE	3.2
1	B	188	GLY	3.0
1	A	209	PRO	3.0
1	C	224	PHE	2.8
1	B	201	PHE	2.7
1	D	224	PHE	2.7
1	D	135	GLY	2.6
1	A	203	ASP	2.6
1	A	224	PHE	2.4
1	A	223	THR	2.4
1	B	222	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	208	MET	2.3
1	B	205	ASN	2.3
1	C	223	THR	2.3
1	B	147	ARG	2.2
1	A	207	HIS	2.2
1	D	200	ILE	2.2
1	B	207	HIS	2.2
1	B	202	GLY	2.2
1	B	223	THR	2.1
1	B	191	LYS	2.1
1	B	221	LYS	2.1
2	H	295	LYS	2.0
1	A	214	LEU	2.0
1	D	215	LEU	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](#)

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. The B-factors column lists the minimum, median, 95th 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	B-factors(Å ²)	Q<0.9
3	HEZ	C	301	8/8	0.80	0.14	39,45,46,56	0
3	HEZ	D	301	8/8	0.80	0.15	47,52,62,62	0
3	HEZ	A	301	8/8	0.89	0.15	34,38,45,45	0
3	HEZ	B	301	8/8	0.89	0.11	38,44,52,55	0

6.5 Other polymers [i](#)

There are no such residues in this entry.