Full wwPDB X-ray Structure Validation Report (i)

Dec 16, 2023 – 11:54 pm GMT

PDB ID	:	4BZ5
Title	:	Crystal structure of Schistosoma mansoni HDAC8
Authors	:	Marek, M.; Romier, C.
Deposited on	:	2013-07-24
Resolution	:	1.78 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 (i)) were used in the production of this report:

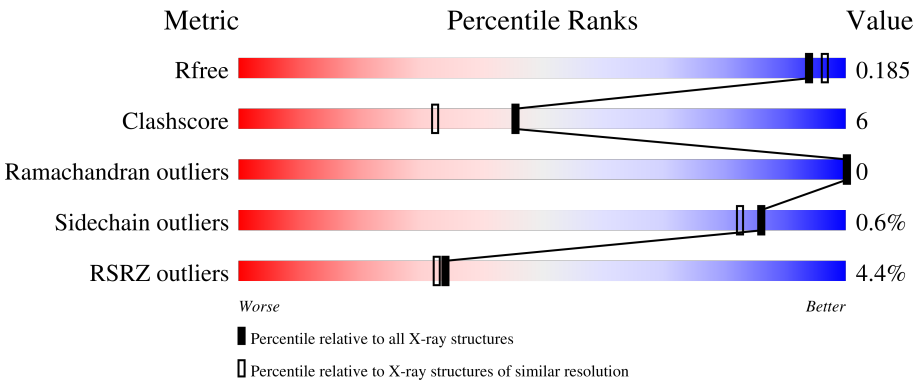
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 1.78 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	9185 (1.80-1.76)
Clashscore	141614	10184 (1.80-1.76)
Ramachandran outliers	138981	10051 (1.80-1.76)
Sidechain outliers	138945	10050 (1.80-1.76)
RSRZ outliers	127900	9032 (1.80-1.76)

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	446	4% 82% 9% 9%
1	B	446	3% 82% 12% 6%
1	C	446	3% 83% 10% 6%
1	D	446	6% 80% 10% 9%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 14813 atoms, of which 16 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 HISTONE DEACETYLASE 8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	404	3239	2088	539	596	16	0	1	0
1	B	419	3362	2166	560	619	17	0	4	0
1	C	418	3366	2169	561	619	17	0	6	0
1	D	404	3246	2096	538	596	16	0	3	0

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	441	GLY	-	expression tag	UNP A5H660
A	442	SER	-	expression tag	UNP A5H660
A	443	LEU	-	expression tag	UNP A5H660
A	444	VAL	-	expression tag	UNP A5H660
A	445	PRO	-	expression tag	UNP A5H660
A	446	ARG	-	expression tag	UNP A5H660
B	441	GLY	-	expression tag	UNP A5H660
B	442	SER	-	expression tag	UNP A5H660
B	443	LEU	-	expression tag	UNP A5H660
B	444	VAL	-	expression tag	UNP A5H660
B	445	PRO	-	expression tag	UNP A5H660
B	446	ARG	-	expression tag	UNP A5H660
C	441	GLY	-	expression tag	UNP A5H660
C	442	SER	-	expression tag	UNP A5H660
C	443	LEU	-	expression tag	UNP A5H660
C	444	VAL	-	expression tag	UNP A5H660
C	445	PRO	-	expression tag	UNP A5H660
C	446	ARG	-	expression tag	UNP A5H660
D	441	GLY	-	expression tag	UNP A5H660
D	442	SER	-	expression tag	UNP A5H660
D	443	LEU	-	expression tag	UNP A5H660

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Chain	Residue	Modelled	Actual	Comment	Reference
D	444	VAL	-	expression tag	UNP A5H660
D	445	PRO	-	expression tag	UNP A5H660
D	446	ARG	-	expression tag	UNP A5H660

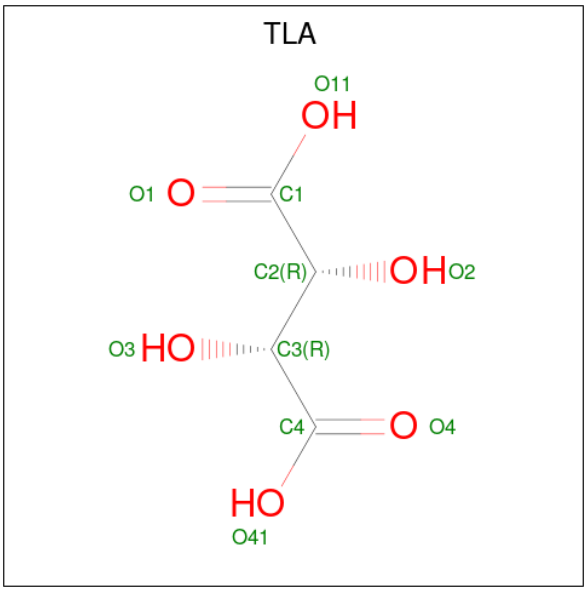
- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0
2	D	1	Total Zn 1 1	0	0

- Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total K 2 2	0	0
3	B	2	Total K 2 2	0	0
3	C	2	Total K 2 2	0	0
3	D	2	Total K 2 2	0	0

- Molecule 4 is L(+)-TARTARIC ACID (three-letter code: TLA) (formula: $C_4H_6O_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	H	O		
4	A	1	14	4	4	6	0	0
			Total	C	H	O		
4	B	1	14	4	4	6	0	0
			Total	C	H	O		
4	C	1	14	4	4	6	0	0
			Total	C	H	O		
4	D	1	14	4	4	6	0	0

- Molecule 5 is water.

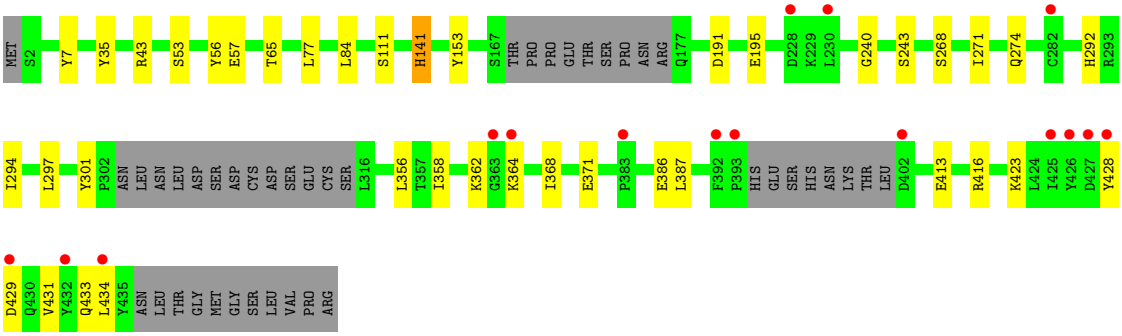
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	357	357	357	0	0
			Total	O		
5	B	416	416	416	0	0
			Total	O		
5	C	434	434	434	0	0
			Total	O		
5	D	325	325	325	0	0

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.

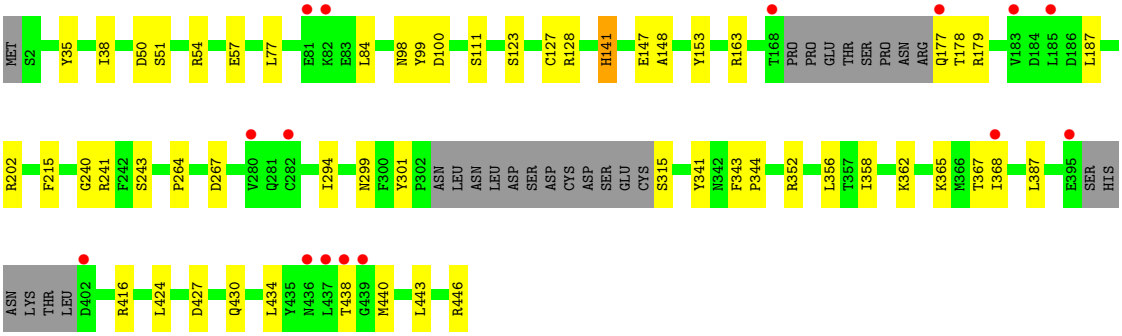
• Molecule 1: HISTONE DEACETYLASE 8

Chain A:



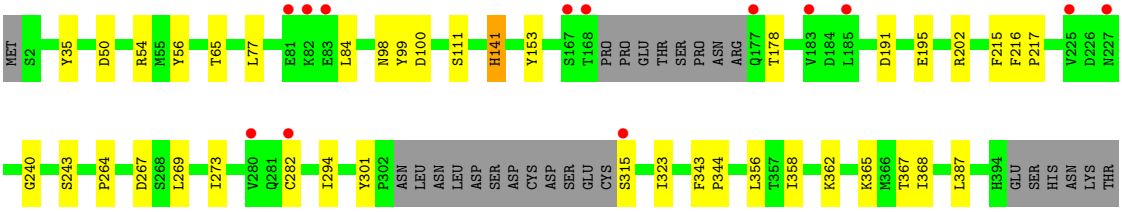
• Molecule 1: HISTONE DEACETYLASE 8

Chain B:



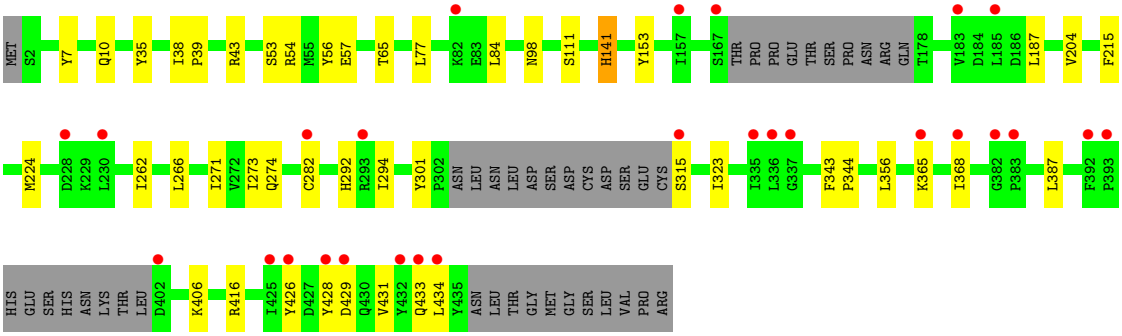
• Molecule 1: HISTONE DEACETYLASE 8

Chain C:



• Molecule 1: HISTONE DEACETYLASE 8

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	70.65Å 70.73Å 98.30Å 75.90° 78.32° 85.59°	Depositor
Resolution (Å)	31.37 – 1.78 31.37 – 1.79	Depositor EDS
% Data completeness (in resolution range)	96.8 (31.37-1.78) 96.8 (31.37-1.79)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.14 (at 1.78Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.154 , 0.188 0.151 , 0.185	Depositor DCC
R_{free} test set	8324 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å^2)	17.9	Xtriage
Anisotropy	0.321	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 59.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.064 for -k,-h,-l	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	14813	wwPDB-VP
Average B, all atoms (Å^2)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.56% 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)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, TLA, K

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	0.28	0/3333	0.45	0/4534
1	B	0.30	0/3467	0.47	0/4717
1	C	0.30	0/3477	0.47	0/4731
1	D	0.28	0/3346	0.45	0/4552
All	All	0.29	0/13623	0.46	0/18534

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	3239	0	3133	35	0
1	B	3362	0	3269	52	0
1	C	3366	0	3276	41	0
1	D	3246	0	3153	40	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
4	A	10	4	3	0	0
4	B	10	4	3	0	0
4	C	10	4	3	0	0
4	D	10	4	3	0	0
5	A	357	0	0	2	1
5	B	416	0	0	7	0
5	C	434	0	0	4	1
5	D	325	0	0	5	0
All	All	14797	16	12843	168	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 (168) 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:D:365:LYS:HD2	1:D:365:LYS:H	1.10	1.14
1:C:35:TYR:CE1	1:C:368[B]:ILE:HG23	2.03	0.93
1:B:35:TYR:CE1	1:B:368[A]:ILE:HG23	2.05	0.89
1:B:368[A]:ILE:HG21	1:B:387:LEU:HD22	1.55	0.89
1:A:35:TYR:CE1	1:A:368:ILE:HG23	2.08	0.88
1:D:429:ASP:O	1:D:433:GLN:HG2	1.76	0.86
1:B:35:TYR:CE1	1:B:368[B]:ILE:HG23	2.11	0.84
1:C:368[B]:ILE:HG21	1:C:387:LEU:HD22	1.59	0.82
1:A:35:TYR:CD1	1:A:368:ILE:HG23	2.15	0.81
1:C:267:ASP:HB3	1:C:434:LEU:HD11	1.62	0.80
1:C:35:TYR:CE1	1:C:368[A]:ILE:HG23	2.16	0.80
1:D:35:TYR:CE1	1:D:368[A]:ILE:HG23	2.15	0.80
1:D:365:LYS:H	1:D:365:LYS:CD	1.92	0.80
1:C:368[B]:ILE:HG21	1:C:387:LEU:CD2	2.11	0.80
1:C:35:TYR:CD1	1:C:368[A]:ILE:HG23	2.17	0.79
1:D:368[A]:ILE:HG21	1:D:387:LEU:CD2	2.12	0.78
1:A:429:ASP:O	1:A:433:GLN:HG2	1.82	0.78
1:C:267:ASP:CB	1:C:434:LEU:HD11	2.12	0.78
1:B:35:TYR:CD1	1:B:368[B]:ILE:HG23	2.19	0.77
1:D:35:TYR:CD1	1:D:368[B]:ILE:HG23	2.20	0.77
1:D:365:LYS:HD2	1:D:365:LYS:N	1.95	0.75
1:B:98:ASN:HB3	5:B:2159:HOH:O	1.86	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:368[A]:ILE:HG21	1:B:387:LEU:CD2	2.17	0.73
1:B:98:ASN:OD1	1:B:99:TYR:N	2.18	0.72
1:D:368[A]:ILE:HG21	1:D:387:LEU:HD22	1.72	0.71
1:A:362:LYS:HB3	1:A:364:LYS:HE3	1.74	0.70
1:C:98[B]:ASN:OD1	1:C:99:TYR:N	2.22	0.70
1:C:35:TYR:CE1	1:C:368[A]:ILE:CG2	2.75	0.70
1:D:315:SER:HB3	5:D:2252:HOH:O	1.92	0.69
1:C:178:THR:HB	1:C:202:ARG:HH21	1.57	0.69
1:A:35:TYR:CE1	1:A:368:ILE:CG2	2.75	0.68
1:D:35:TYR:CE1	1:D:368[B]:ILE:HG23	2.28	0.68
1:B:35:TYR:CE1	1:B:368[B]:ILE:CG2	2.75	0.68
1:A:368:ILE:HG21	1:A:387:LEU:CD2	2.25	0.67
1:B:368[A]:ILE:CG2	1:B:387:LEU:HD22	2.25	0.67
1:A:358:ILE:HG23	1:A:362:LYS:HD3	1.79	0.65
1:A:362:LYS:HE2	1:A:364:LYS:NZ	2.12	0.64
1:D:416:ARG:HD3	1:D:428:TYR:CE2	2.32	0.64
1:C:365:LYS:HE2	1:C:367:THR:OG1	1.96	0.64
1:B:267:ASP:HB3	1:B:434:LEU:HD11	1.80	0.64
1:D:54:ARG:NH1	5:D:2072:HOH:O	2.30	0.63
1:D:368[A]:ILE:CG2	1:D:387:LEU:HD22	2.28	0.63
1:C:368[B]:ILE:CG2	1:C:387:LEU:HD22	2.29	0.62
1:A:362:LYS:CB	1:A:364:LYS:HE3	2.31	0.60
1:B:440:MET:HE3	1:B:443:LEU:HD11	1.82	0.60
1:B:438:THR:O	1:B:438:THR:HG22	2.02	0.60
1:B:315:SER:HA	5:B:2330:HOH:O	2.02	0.59
1:A:271:ILE:HD12	1:A:434:LEU:HD11	1.85	0.59
1:A:423:LYS:HE2	5:A:2357:HOH:O	2.03	0.59
1:A:368:ILE:HG21	1:A:387:LEU:HD22	1.83	0.58
1:C:358:ILE:HG23	1:C:362:LYS:HD3	1.84	0.58
1:B:100:ASP:HB2	5:B:2164:HOH:O	2.03	0.58
1:B:356:LEU:C	1:B:356:LEU:HD23	2.24	0.58
1:D:35:TYR:CE1	1:D:368[B]:ILE:CG2	2.86	0.57
1:B:267:ASP:CB	1:B:434:LEU:HD11	2.35	0.57
1:A:371:GLU:HG3	1:A:386:GLU:OE2	2.05	0.57
1:D:271:ILE:HD12	1:D:434:LEU:HD11	1.85	0.57
1:D:426:TYR:CE2	1:D:431:VAL:HG21	2.40	0.56
1:B:77:LEU:HD13	1:B:84:LEU:HG	1.87	0.56
1:C:54:ARG:HD3	5:C:2113:HOH:O	2.04	0.56
1:C:178:THR:HB	1:C:202:ARG:NH2	2.20	0.56
1:C:315:SER:HB3	5:C:2325:HOH:O	2.03	0.56
1:A:368:ILE:CG2	1:A:387:LEU:HD22	2.36	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:368[B]:ILE:HG21	1:B:387:LEU:HD22	1.88	0.55
1:A:240:GLY:O	1:A:243:SER:OG	2.25	0.55
1:C:267:ASP:HB2	1:C:434:LEU:HD11	1.87	0.55
1:A:56:TYR:OH	1:A:65:THR:HB	2.06	0.55
1:B:358:ILE:HG23	1:B:362:LYS:HD3	1.89	0.55
1:C:50:ASP:HB2	1:C:54:ARG:HD2	1.89	0.55
1:B:111:SER:OG	1:B:153:TYR:HB2	2.06	0.54
1:A:7:TYR:CE2	1:A:43:ARG:HG3	2.43	0.54
1:A:416:ARG:HD3	1:A:428:TYR:CE2	2.42	0.53
1:C:240:GLY:O	1:C:243:SER:OG	2.26	0.53
1:D:56:TYR:OH	1:D:65:THR:HB	2.07	0.53
1:B:241:ARG:HB2	1:B:241:ARG:NH1	2.24	0.53
1:B:178:THR:HB	1:B:202:ARG:HH21	1.74	0.52
1:A:274:GLN:HG2	5:A:2288:HOH:O	2.09	0.52
1:B:54:ARG:HD3	5:B:2093:HOH:O	2.09	0.52
1:B:179:ARG:NH2	1:B:424:LEU:HD21	2.25	0.52
1:C:264:PRO:HA	1:C:434:LEU:HD21	1.91	0.52
1:C:100:ASP:HB2	5:C:2194:HOH:O	2.09	0.51
1:C:111:SER:OG	1:C:153:TYR:HB2	2.10	0.51
1:B:35:TYR:CD1	1:B:368[A]:ILE:HG23	2.46	0.51
1:A:53:SER:O	1:A:57:GLU:HG3	2.10	0.51
1:C:273[B]:ILE:HD12	1:C:418:TYR:CE2	2.46	0.51
1:D:426:TYR:CD2	1:D:431:VAL:HG21	2.46	0.51
1:A:292:HIS:HB3	1:A:294:ILE:HG12	1.92	0.51
1:D:53:SER:O	1:D:57:GLU:HG3	2.10	0.51
1:A:7:TYR:CE1	1:A:43:ARG:CD	2.94	0.51
1:A:356:LEU:C	1:A:356:LEU:HD23	2.32	0.50
1:B:57:GLU:HG3	5:B:2095:HOH:O	2.10	0.50
1:B:187:LEU:HD21	1:B:294[B]:ILE:HD12	1.93	0.50
1:D:416:ARG:HD3	1:D:428:TYR:HE2	1.75	0.50
1:A:7:TYR:CE1	1:A:43:ARG:HD2	2.48	0.49
1:B:427:ASP:OD2	1:B:430:GLN:NE2	2.25	0.49
1:D:215:PHE:CE1	1:D:294[B]:ILE:CD1	2.95	0.49
1:D:368[A]:ILE:CG2	1:D:387:LEU:CD2	2.86	0.49
1:C:267:ASP:HB3	1:C:434:LEU:CD1	2.39	0.49
1:C:191:ASP:O	1:C:195:GLU:HG3	2.13	0.49
1:C:440:MET:HE3	1:C:443:LEU:HD11	1.95	0.48
1:A:297:LEU:HD12	1:A:297:LEU:N	2.28	0.48
1:C:77:LEU:HD13	1:C:84:LEU:HG	1.95	0.48
1:B:240:GLY:O	1:B:243:SER:OG	2.32	0.48
1:B:416:ARG:HD2	1:B:446:ARG:O	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:365:LYS:HE3	1:B:367:THR:OG1	2.13	0.48
1:B:343:PHE:HB2	1:B:344:PRO:HD3	1.95	0.47
1:C:35:TYR:CD1	1:C:368[B]:ILE:HG23	2.47	0.47
1:B:264:PRO:HA	1:B:434:LEU:HD21	1.97	0.47
1:D:356:LEU:C	1:D:356:LEU:HD23	2.35	0.47
5:C:2148:HOH:O	1:D:10:GLN:HG2	2.14	0.47
1:D:7:TYR:CE1	1:D:43:ARG:HG3	2.50	0.46
1:B:147:GLU:HG2	1:B:148:ALA:O	2.15	0.46
1:B:368[B]:ILE:CG2	1:B:387:LEU:HD22	2.45	0.46
1:A:362:LYS:HE2	1:A:364:LYS:CE	2.45	0.46
1:C:141:HIS:CD2	1:C:141:HIS:H	2.33	0.46
1:C:282[B]:CYS:SG	1:C:323:ILE:HD11	2.55	0.46
1:B:51:SER:HB3	1:B:54:ARG:NH1	2.31	0.46
1:D:204:VAL:HG11	1:D:273:ILE:HD11	1.97	0.45
1:D:141:HIS:H	1:D:141:HIS:CD2	2.34	0.45
1:B:356:LEU:HD23	1:B:356:LEU:O	2.16	0.45
1:D:274:GLN:HG2	5:D:2261:HOH:O	2.16	0.45
1:C:368[A]:ILE:HG21	1:C:387:LEU:HD22	1.98	0.45
1:C:416:ARG:NH1	1:C:426:TYR:HB3	2.32	0.45
1:A:268:SER:HB2	1:A:431:VAL:HG13	1.99	0.45
1:C:343:PHE:HB2	1:C:344:PRO:HD3	1.99	0.45
1:B:177:GLN:HA	1:B:177:GLN:OE1	2.17	0.44
1:D:98:ASN:HB2	5:D:2133:HOH:O	2.17	0.44
1:A:362:LYS:HE2	1:A:364:LYS:HE3	1.99	0.44
1:B:440:MET:CE	1:B:443:LEU:CD1	2.96	0.44
1:D:406:LYS:HE2	5:D:2166:HOH:O	2.16	0.44
1:B:440:MET:HE3	1:B:443:LEU:CD1	2.47	0.44
1:B:299:ASN:O	1:B:352:ARG:HD2	2.18	0.44
1:A:368:ILE:CG2	1:A:387:LEU:CD2	2.93	0.44
1:C:215:PHE:CE1	1:C:294:ILE:CD1	3.00	0.43
1:C:356:LEU:HD23	1:C:356:LEU:C	2.38	0.43
1:D:38:ILE:N	1:D:39:PRO:CD	2.81	0.43
1:D:262:ILE:HG12	1:D:266:LEU:HG	1.99	0.43
1:B:38:ILE:HG23	5:B:2007:HOH:O	2.17	0.43
1:B:241:ARG:HB2	1:B:241:ARG:CZ	2.48	0.43
1:B:434:LEU:C	1:B:434:LEU:HD23	2.38	0.43
1:B:440:MET:HE2	1:B:443:LEU:HD12	2.00	0.43
1:A:141:HIS:H	1:A:141:HIS:CD2	2.35	0.43
1:D:292:HIS:HB3	1:D:294[A]:ILE:HG12	2.00	0.43
1:C:269:LEU:O	1:C:273[B]:ILE:HG12	2.18	0.43
1:C:56:TYR:OH	1:C:65:THR:HB	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:187:LEU:HD21	1:D:294[B]:ILE:HD12	2.01	0.43
1:A:111:SER:OG	1:A:153:TYR:HB2	2.18	0.42
1:D:38:ILE:HB	1:D:39:PRO:HD3	2.01	0.42
1:B:141:HIS:H	1:B:141:HIS:CD2	2.36	0.42
1:C:440:MET:HE2	1:C:443:LEU:HD12	2.01	0.42
1:B:123:SER:OG	1:B:163:ARG:NH1	2.52	0.42
1:D:224:MET:HA	1:D:224:MET:CE	2.50	0.42
1:B:50:ASP:HB2	1:B:54:ARG:HD2	2.02	0.42
1:B:215:PHE:CE1	1:B:294[B]:ILE:CD1	3.03	0.41
1:A:77:LEU:HD13	1:A:84:LEU:HG	2.02	0.41
1:D:343:PHE:HB2	1:D:344:PRO:HD3	2.02	0.41
1:B:127:CYS:O	1:B:128:ARG:HB2	2.20	0.41
1:B:341:TYR:OH	5:B:2037:HOH:O	2.21	0.41
1:C:216:PHE:CD1	1:C:217:PRO:HA	2.56	0.41
1:A:362:LYS:HE2	1:A:364:LYS:HZ1	1.82	0.41
1:A:191:ASP:O	1:A:195:GLU:HG3	2.20	0.41
1:C:368[B]:ILE:CG2	1:C:387:LEU:CD2	2.89	0.41
1:B:51:SER:CB	1:B:54:ARG:HH11	2.33	0.41
1:D:282[B]:CYS:SG	1:D:323:ILE:HD11	2.61	0.41
1:A:413:GLU:OE1	1:A:413:GLU:HA	2.22	0.40
1:D:77:LEU:HD13	1:D:84:LEU:HG	2.02	0.40
1:D:111:SER:OG	1:D:153:TYR:HB2	2.21	0.40
1:C:368[A]:ILE:HG21	1:C:387:LEU:CD2	2.52	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)
5:A:2153:HOH:O	5:C:2386:HOH:O[1_565]	2.18	0.02

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	397/446 (89%)	390 (98%)	7 (2%)	0	100	100
1	B	415/446 (93%)	408 (98%)	7 (2%)	0	100	100
1	C	416/446 (93%)	408 (98%)	8 (2%)	0	100	100
1	D	399/446 (90%)	395 (99%)	4 (1%)	0	100	100
All	All	1627/1784 (91%)	1601 (98%)	26 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	352/391 (90%)	350 (99%)	2 (1%)	86	82
1	B	368/391 (94%)	366 (100%)	2 (0%)	88	86
1	C	369/391 (94%)	367 (100%)	2 (0%)	88	86
1	D	354/391 (90%)	352 (99%)	2 (1%)	86	82
All	All	1443/1564 (92%)	1435 (99%)	8 (1%)	86	82

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	141	HIS
1	A	301	TYR
1	B	141	HIS
1	B	301	TYR
1	C	141	HIS
1	C	301	TYR
1	D	141	HIS
1	D	301	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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)

Of 16 ligands modelled in this entry, 12 are monoatomic - leaving 4 for Mogul analysis.

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	$\#\|Z\| > 2$
4	TLA	D	700	2	9,9,9	1.21	0	12,12,12	1.11	0
4	TLA	A	700	2	9,9,9	1.17	0	12,12,12	1.08	1 (8%)
4	TLA	B	700	2	9,9,9	1.15	0	12,12,12	1.22	2 (16%)
4	TLA	C	700	2	9,9,9	1.25	0	12,12,12	1.03	1 (8%)

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
4	TLA	D	700	2	-	4/12/12/12	-
4	TLA	A	700	2	-	4/12/12/12	-
4	TLA	B	700	2	-	2/12/12/12	-
4	TLA	C	700	2	-	2/12/12/12	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	B	700	TLA	O11-C1-C2	2.27	119.42	113.27
4	B	700	TLA	O41-C4-C3	2.20	119.22	113.27
4	A	700	TLA	O11-C1-C2	2.12	119.00	113.27
4	C	700	TLA	O11-C1-C2	2.01	118.69	113.27

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	700	TLA	O1-C1-C2-C3
4	D	700	TLA	O1-C1-C2-O2
4	D	700	TLA	O1-C1-C2-C3
4	A	700	TLA	O11-C1-C2-C3
4	A	700	TLA	O1-C1-C2-O2
4	D	700	TLA	O11-C1-C2-O2
4	B	700	TLA	C2-C3-C4-O4
4	D	700	TLA	O11-C1-C2-C3
4	A	700	TLA	O11-C1-C2-O2
4	B	700	TLA	C2-C3-C4-O41
4	C	700	TLA	O1-C1-C2-C3
4	C	700	TLA	O11-C1-C2-C3

There are no ring outliers.

No monomer is involved in short contacts.

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 (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	404/446 (90%)	0.08	16 (3%)	38	36	11, 20, 48, 80	0
1	B	419/446 (93%)	-0.04	15 (3%)	42	41	10, 17, 41, 75	0
1	C	418/446 (93%)	-0.06	14 (3%)	46	45	10, 17, 40, 68	0
1	D	404/446 (90%)	0.14	27 (6%)	17	17	12, 20, 53, 74	0
All	All	1645/1784 (92%)	0.03	72 (4%)	34	32	10, 19, 46, 80	0

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	392	PHE	7.1
1	A	425	ILE	7.1
1	D	426	TYR	6.3
1	D	432	TYR	5.7
1	A	428	TYR	5.5
1	B	168	THR	5.4
1	D	428	TYR	5.2
1	A	432	TYR	5.2
1	D	425	ILE	4.8
1	C	168	THR	4.5
1	A	426	TYR	4.5
1	A	392	PHE	4.4
1	C	81	GLU	4.3
1	C	177	GLN	4.1
1	A	230	LEU	3.8
1	D	282[A]	CYS	3.6
1	A	402	ASP	3.6
1	A	429	ASP	3.5
1	C	282[A]	CYS	3.4
1	D	402	ASP	3.4
1	B	282[A]	CYS	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	183	VAL	3.2
1	B	82	LYS	3.1
1	C	315	SER	3.1
1	D	429	ASP	3.1
1	B	402	ASP	3.0
1	B	437	LEU	3.0
1	C	227	ASN	3.0
1	B	81	GLU	3.0
1	B	395	GLU	3.0
1	C	82	LYS	3.0
1	D	383	PRO	2.9
1	D	434	LEU	2.9
1	D	82	LYS	2.8
1	D	365	LYS	2.8
1	D	337	GLY	2.7
1	D	368[A]	ILE	2.7
1	C	280	VAL	2.7
1	B	177	GLN	2.7
1	B	438	THR	2.6
1	B	439	GLY	2.6
1	A	434	LEU	2.6
1	D	433	GLN	2.6
1	A	282[A]	CYS	2.6
1	C	183	VAL	2.5
1	A	393	PRO	2.5
1	D	230	LEU	2.5
1	B	185	LEU	2.5
1	C	439	GLY	2.4
1	C	83	GLU	2.4
1	A	364	LYS	2.4
1	B	280	VAL	2.4
1	C	185	LEU	2.4
1	D	336	LEU	2.4
1	D	293	ARG	2.4
1	D	335	ILE	2.3
1	D	393	PRO	2.3
1	B	436	ASN	2.3
1	D	228	ASP	2.3
1	B	368[A]	ILE	2.3
1	A	363	GLY	2.3
1	D	185	LEU	2.2
1	D	167	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	383	PRO	2.2
1	A	427	ASP	2.2
1	C	225	VAL	2.2
1	D	315	SER	2.1
1	A	228	ASP	2.1
1	D	157	ILE	2.1
1	D	183	VAL	2.1
1	C	167	SER	2.1
1	D	382	GLY	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
4	TLA	D	700	10/10	0.91	0.11	20,30,38,38	0
4	TLA	B	700	10/10	0.93	0.12	16,28,45,46	0
4	TLA	C	700	10/10	0.95	0.10	17,27,38,41	0
4	TLA	A	700	10/10	0.95	0.09	20,28,35,40	0
3	K	A	600	1/1	0.99	0.04	17,17,17,17	0
3	K	C	601	1/1	0.99	0.07	15,15,15,15	0
3	K	B	600	1/1	1.00	0.06	17,17,17,17	0
3	K	B	601	1/1	1.00	0.09	10,10,10,10	0
3	K	C	600	1/1	1.00	0.08	11,11,11,11	0
2	ZN	B	500	1/1	1.00	0.03	18,18,18,18	0
3	K	D	600	1/1	1.00	0.04	18,18,18,18	0
3	K	D	601	1/1	1.00	0.07	12,12,12,12	0
2	ZN	C	500	1/1	1.00	0.03	18,18,18,18	0
2	ZN	D	500	1/1	1.00	0.02	19,19,19,19	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	ZN	A	500	1/1	1.00	0.02	19,19,19,19	0
3	K	A	601	1/1	1.00	0.08	11,11,11,11	0

6.5 Other polymers (i)

There are no such residues in this entry.