Full wwPDB X-ray Structure Validation Report (i)

Mar 13, 2024 – 02:32 PM JST

PDB ID	:	4XM1
Title	:	N,N'-diacetylchitobiose deacetylase from Pyrococcus furiosus in the presence of cadmium
Authors	:	Nakamura, T.; Niiyama, M.; Hashimoto, W.; Ida, K.; Uegaki, K.
Deposited on	:	2015-01-14
Resolution	:	1.80 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:

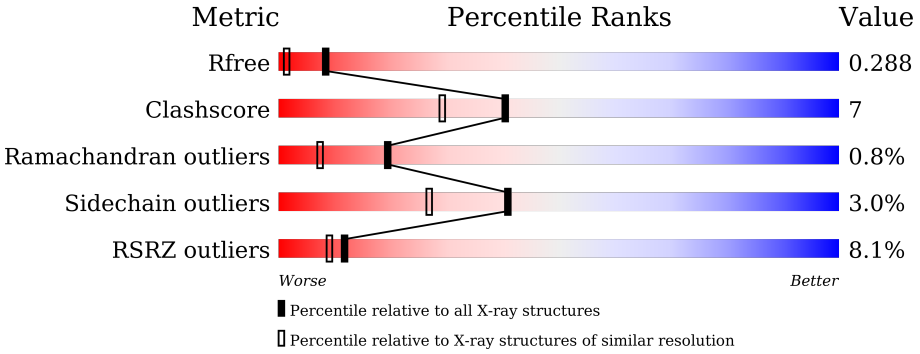
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, 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 1.80 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	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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	267	6% 81% 16%
1	B	267	11% 81% 16%
1	C	267	4% 79% 19%
1	D	267	5% 79% 16%
1	E	267	13% 76% 19%
1	F	267	8% 80% 19%

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 13350 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.

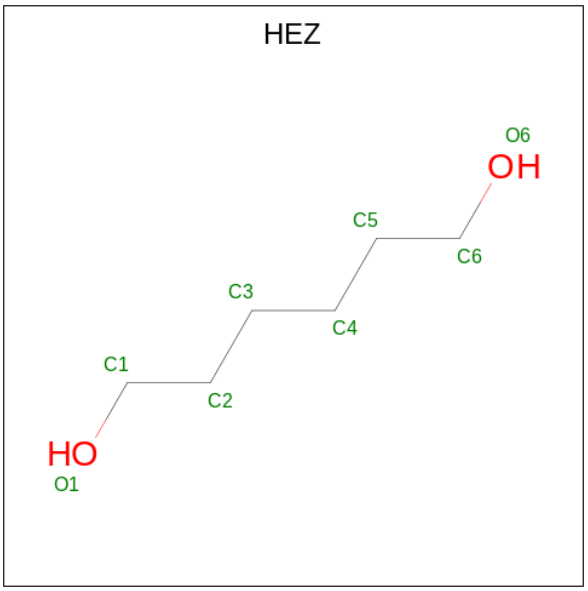
- Molecule 1 is a protein called Uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	263	2154	1393	355	397	9	0	0	0
			Total	C	N	O	S			
1	B	261	2141	1385	353	394	9	0	0	0
			Total	C	N	O	S			
1	C	267	2190	1416	360	404	10	0	0	0
			Total	C	N	O	S			
1	D	260	2134	1381	352	392	9	0	0	0
			Total	C	N	O	S			
1	E	257	2105	1362	349	385	9	0	0	0
			Total	C	N	O	S			
1	F	267	2190	1416	360	404	10	0	0	0

- Molecule 2 is CADMIUM ION (three-letter code: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Cd		
2	A	2	2	2	0	0
			Total	Cd		
2	B	1	1	1	0	0
			Total	Cd		
2	C	1	1	1	0	0
			Total	Cd		
2	D	2	2	2	0	0
			Total	Cd		
2	E	1	1	1	0	0
			Total	Cd		
2	F	1	1	1	0	0

- Molecule 3 is HEXANE-1,6-DIOL (three-letter code: HEZ) (formula: $C_6H_{14}O_2$).

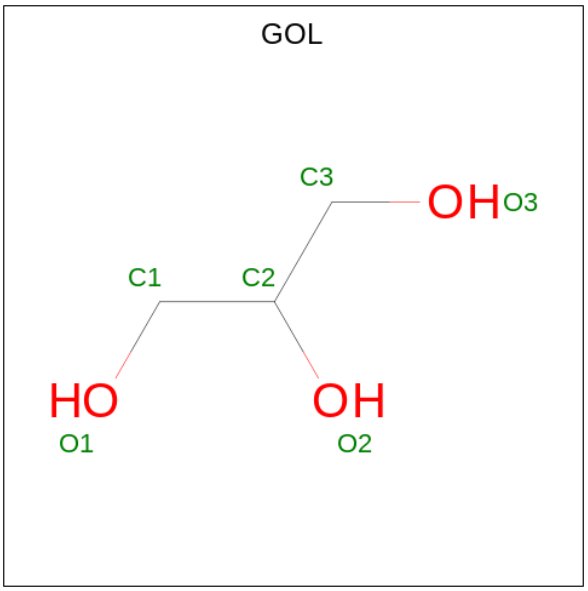


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	A	1	8	6	2	0	0
3	B	1	8	6	2	0	0
3	C	1	8	6	2	0	0
3	D	1	8	6	2	0	0
3	E	1	8	6	2	0	0
3	F	1	8	6	2	0	0

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Cl		
4	A	1	1	1	0	0
4	C	2	2	2	0	0
4	D	1	1	1	0	0
4	F	2	2	2	0	0

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
5	A	1	6	3	3	0	0
			Total	C	O		
5	D	1	6	3	3	0	0

- Molecule 6 is water.

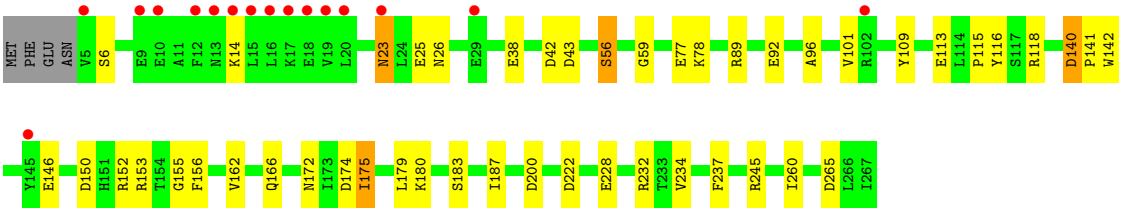
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	A	83	83	83	0	0
			Total	O		
6	B	56	56	56	0	0
			Total	O		
6	C	67	67	67	0	0
			Total	O		
6	D	65	65	65	0	0
			Total	O		
6	E	43	43	43	0	0
			Total	O		
6	F	48	48	48	0	0

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.

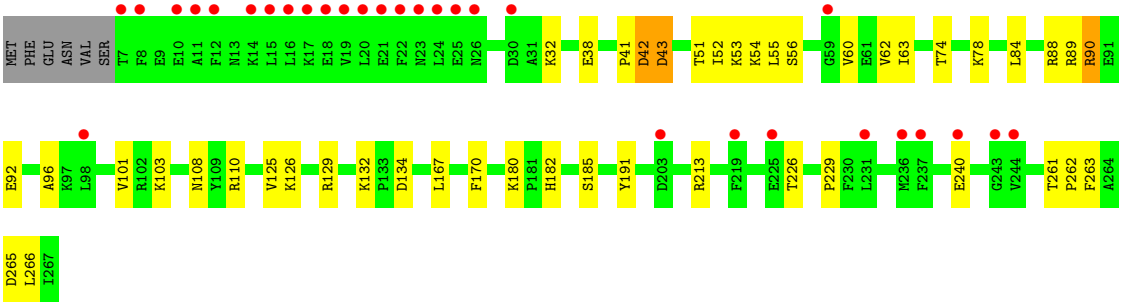
• Molecule 1: Uncharacterized protein

Chain A:



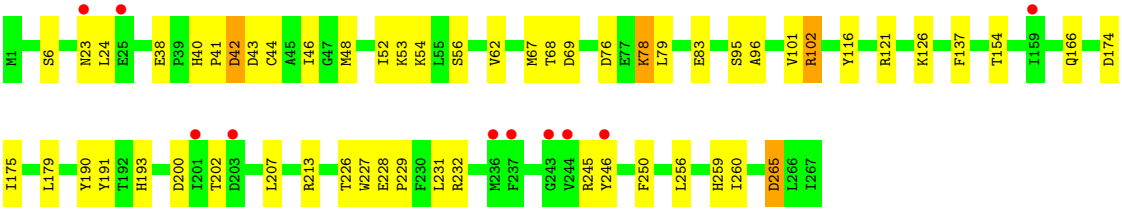
• Molecule 1: Uncharacterized protein

Chain B:



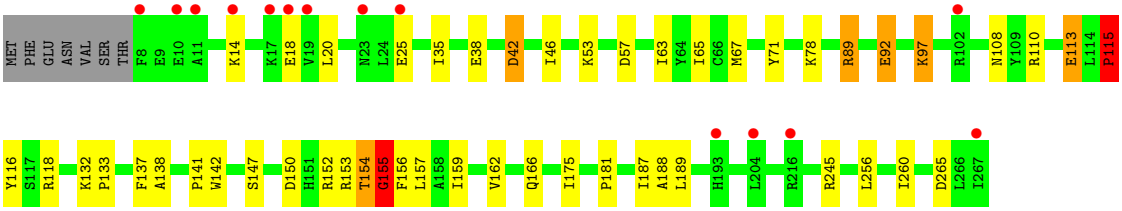
• Molecule 1: Uncharacterized protein

Chain C:



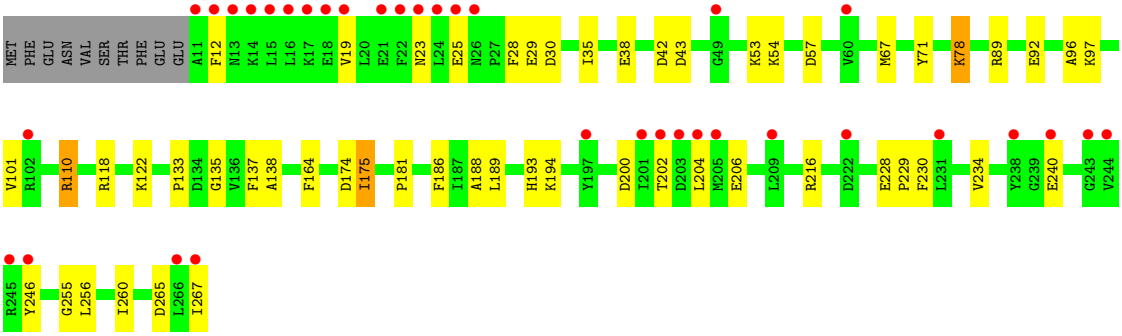
• Molecule 1: Uncharacterized protein

Chain D:



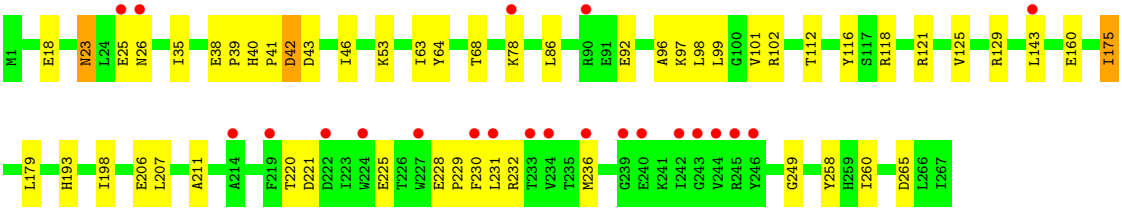
• Molecule 1: Uncharacterized protein

Chain E:



• Molecule 1: Uncharacterized protein

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	74.35Å 77.85Å 93.51Å 66.39° 84.70° 61.95°	Depositor
Resolution (Å)	41.13 – 1.80 41.13 – 1.80	Depositor EDS
% Data completeness (in resolution range)	97.6 (41.13-1.80) 97.6 (41.13-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.44 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.238 , 0.290 0.237 , 0.288	Depositor DCC
R_{free} test set	7656 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å^2)	26.0	Xtriage
Anisotropy	0.229	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 48.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	13350	wwPDB-VP
Average B, all atoms (Å^2)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.88% 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: CL, GOL, CD, 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	1.15	4/2211 (0.2%)	1.08	11/2992 (0.4%)
1	B	0.92	0/2198	0.93	2/2974 (0.1%)
1	C	1.06	2/2248 (0.1%)	0.99	5/3041 (0.2%)
1	D	1.16	9/2191 (0.4%)	1.05	13/2964 (0.4%)
1	E	0.87	0/2161	0.86	2/2924 (0.1%)
1	F	0.94	1/2248 (0.0%)	0.89	2/3041 (0.1%)
All	All	1.02	16/13257 (0.1%)	0.97	35/17936 (0.2%)

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	155	GLY	N-CA	11.86	1.63	1.46
1	A	155	GLY	N-CA	7.05	1.56	1.46
1	D	71	TYR	CD2-CE2	6.77	1.49	1.39
1	C	116	TYR	CE2-CZ	6.43	1.47	1.38
1	D	154	THR	CA-CB	-6.37	1.36	1.53
1	A	153	ARG	CA-CB	6.24	1.67	1.53
1	D	155	GLY	CA-C	-6.12	1.42	1.51
1	D	154	THR	CB-OG1	5.56	1.54	1.43
1	A	146	GLU	CD-OE2	5.43	1.31	1.25
1	D	115	PRO	CB-CG	5.27	1.76	1.50
1	D	116	TYR	CD1-CE1	5.15	1.47	1.39
1	D	113	GLU	CB-CG	-5.13	1.42	1.52
1	A	109	TYR	CE1-CZ	5.11	1.45	1.38
1	F	160	GLU	CD-OE1	-5.09	1.20	1.25
1	D	92	GLU	CB-CG	5.06	1.61	1.52
1	C	250	PHE	CE1-CZ	5.05	1.47	1.37

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	43	ASP	CB-CG-OD1	14.80	131.62	118.30
1	A	153	ARG	NE-CZ-NH1	-13.25	113.67	120.30
1	D	118	ARG	NE-CZ-NH2	-9.52	115.54	120.30
1	D	154	THR	O-C-N	8.96	138.44	123.20
1	E	43	ASP	CB-CG-OD1	8.57	126.01	118.30
1	C	69	ASP	CB-CG-OD1	8.42	125.88	118.30
1	F	118	ARG	NE-CZ-NH1	7.78	124.19	120.30
1	D	155	GLY	N-CA-C	-7.65	93.98	113.10
1	A	140	ASP	CB-CG-OD2	-7.12	111.89	118.30
1	A	150	ASP	CB-CG-OD2	6.57	124.22	118.30
1	F	118	ARG	NE-CZ-NH2	-6.41	117.09	120.30
1	E	110	ARG	NE-CZ-NH2	6.34	123.47	120.30
1	B	110	ARG	NE-CZ-NH2	6.32	123.46	120.30
1	D	157	LEU	CB-CG-CD2	6.15	121.46	111.00
1	D	154	THR	CA-C-N	-5.96	104.28	116.20
1	A	174	ASP	CB-CG-OD2	5.84	123.56	118.30
1	A	245	ARG	NE-CZ-NH2	5.71	123.15	120.30
1	D	245	ARG	NE-CZ-NH2	5.69	123.15	120.30
1	C	256	LEU	CA-CB-CG	5.62	128.23	115.30
1	C	121	ARG	NE-CZ-NH2	-5.61	117.50	120.30
1	D	118	ARG	NE-CZ-NH1	5.60	123.10	120.30
1	A	118	ARG	NE-CZ-NH2	-5.57	117.51	120.30
1	A	140	ASP	CB-CG-OD1	5.55	123.30	118.30
1	A	200	ASP	CB-CG-OD1	5.50	123.25	118.30
1	A	156	PHE	CB-CG-CD2	5.45	124.62	120.80
1	D	166	GLN	CB-CA-C	-5.45	99.51	110.40
1	D	89	ARG	NE-CZ-NH2	-5.42	117.59	120.30
1	C	265	ASP	CB-CG-OD1	5.37	123.13	118.30
1	D	110	ARG	NE-CZ-NH1	-5.27	117.67	120.30
1	C	69	ASP	CB-CG-OD2	-5.24	113.59	118.30
1	D	150	ASP	CB-CG-OD2	5.24	123.01	118.30
1	A	152	ARG	O-C-N	-5.22	114.34	122.70
1	D	67	MET	CG-SD-CE	5.08	108.34	100.20
1	A	116	TYR	CA-CB-CG	-5.08	103.75	113.40
1	D	153	ARG	NE-CZ-NH2	-5.02	117.79	120.30

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	2154	0	2128	21	0
1	B	2141	0	2114	37	0
1	C	2190	0	2161	39	0
1	D	2134	0	2107	30	0
1	E	2105	0	2086	26	0
1	F	2190	0	2161	35	0
2	A	2	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	2	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	8	0	14	2	0
3	B	8	0	14	1	0
3	C	8	0	14	1	0
3	D	8	0	14	1	0
3	E	8	0	14	3	0
3	F	8	0	14	1	0
4	A	1	0	0	0	0
4	C	2	0	0	0	0
4	D	1	0	0	0	0
4	F	2	0	0	0	0
5	A	6	0	7	2	0
5	D	6	0	8	1	0
6	A	83	0	0	3	0
6	B	56	0	0	8	0
6	C	67	0	0	4	0
6	D	65	0	0	8	0
6	E	43	0	0	4	0
6	F	48	0	0	6	0
All	All	13350	0	12856	186	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 7.

All (186) 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:115:PRO:CB	1:D:115:PRO:CG	1.76	1.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:159:ILE:HD12	6:D:437:HOH:O	1.53	1.08
1:A:23:ASN:HD21	1:A:25:GLU:HB2	1.24	1.02
1:D:155:GLY:C	6:D:437:HOH:O	2.01	0.98
1:D:155:GLY:O	6:D:437:HOH:O	1.88	0.92
1:B:90:ARG:HG3	1:B:90:ARG:HH11	1.35	0.91
5:D:305:GOL:H12	6:F:446:HOH:O	1.73	0.89
1:B:90:ARG:HH11	1:B:90:ARG:CG	1.88	0.85
1:D:89:ARG:HH21	1:D:108:ASN:HD21	1.25	0.82
1:C:102:ARG:HB2	6:C:440:HOH:O	1.83	0.78
1:A:175:ILE:HD13	6:A:444:HOH:O	1.85	0.76
1:D:159:ILE:CD1	6:D:437:HOH:O	2.23	0.74
1:B:129:ARG:NH1	1:B:167:LEU:O	2.19	0.74
1:A:23:ASN:ND2	1:A:25:GLU:HB2	2.00	0.73
1:B:126:LYS:NZ	6:B:448:HOH:O	2.22	0.72
1:E:256:LEU:HD12	6:E:416:HOH:O	1.90	0.70
3:E:302:HEZ:H42	3:F:302:HEZ:H31	1.75	0.69
1:C:265:ASP:OD1	6:C:455:HOH:O	2.12	0.67
1:D:154:THR:O	1:D:155:GLY:O	2.12	0.67
3:A:303:HEZ:H22	3:C:302:HEZ:H21	1.76	0.67
1:C:226:THR:O	1:C:229:PRO:HD2	1.94	0.67
1:D:152:ARG:HD2	3:D:303:HEZ:H42	1.77	0.66
1:E:35:ILE:HD12	1:E:133:PRO:HG3	1.77	0.66
1:D:265:ASP:OD1	6:D:408:HOH:O	2.13	0.66
1:A:175:ILE:HD12	1:A:179:LEU:O	1.94	0.65
1:B:51:THR:O	1:B:55:LEU:HG	1.97	0.65
1:E:228:GLU:HB3	1:E:229:PRO:HD3	1.77	0.65
1:B:38:GLU:OE2	1:B:92:GLU:OE1	2.15	0.64
1:F:193:HIS:HB3	6:F:428:HOH:O	1.98	0.63
1:D:89:ARG:NH2	1:D:108:ASN:HD21	1.95	0.62
1:F:23:ASN:HD21	1:F:25:GLU:HB2	1.65	0.62
1:F:42:ASP:O	1:F:46:ILE:HG13	2.00	0.61
1:A:23:ASN:ND2	1:A:25:GLU:H	1.97	0.61
1:B:132:LYS:HE3	6:B:438:HOH:O	2.00	0.61
1:B:226:THR:O	1:B:229:PRO:HD2	2.00	0.60
1:E:53:LYS:NZ	1:E:57:ASP:OD1	2.34	0.60
1:C:38:GLU:HB2	1:C:43:ASP:HB2	1.83	0.60
1:C:96:ALA:HB1	1:C:101:VAL:HB	1.84	0.60
1:B:180:LYS:O	6:B:424:HOH:O	2.17	0.59
1:F:23:ASN:ND2	1:F:25:GLU:H	1.99	0.59
1:A:140:ASP:OD2	1:A:140:ASP:C	2.41	0.59
1:F:96:ALA:HB1	1:F:101:VAL:HB	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:202:THR:HA	1:E:246:TYR:HB2	1.84	0.58
5:A:305:GOL:H12	1:C:259:HIS:NE2	2.19	0.58
1:D:14:LYS:NZ	1:D:18:GLU:OE1	2.32	0.58
1:C:42:ASP:O	1:C:46:ILE:HG13	2.03	0.58
1:E:96:ALA:HB1	1:E:101:VAL:HB	1.85	0.57
1:D:141:PRO:HG2	1:D:142:TRP:CE3	2.39	0.57
1:C:23:ASN:HB3	6:C:439:HOH:O	2.04	0.57
1:B:52:ILE:HG23	1:B:62:VAL:HG11	1.87	0.57
1:F:64:TYR:CE1	1:F:101:VAL:HG21	2.39	0.57
1:E:23:ASN:HD21	1:E:25:GLU:HB3	1.70	0.57
1:B:90:ARG:HG3	1:B:90:ARG:NH1	2.10	0.57
3:A:303:HEZ:H11	3:B:302:HEZ:H42	1.86	0.56
1:C:227:TRP:O	1:C:231:LEU:HG	2.06	0.56
1:C:23:ASN:CG	1:C:24:LEU:H	2.10	0.55
1:F:41:PRO:O	1:F:42:ASP:CG	2.45	0.55
1:F:53:LYS:HE3	1:F:207:LEU:HD21	1.89	0.55
1:D:42:ASP:O	1:D:46:ILE:HG13	2.08	0.54
1:E:135:GLY:HA3	1:E:186:PHE:CZ	2.43	0.54
1:A:234:VAL:O	1:A:237:PHE:HB3	2.08	0.54
1:B:84:LEU:O	1:B:88:ARG:HG3	2.07	0.54
1:B:191:TYR:HA	6:B:422:HOH:O	2.07	0.53
1:D:53:LYS:NZ	1:D:57:ASP:OD1	2.36	0.53
1:E:265:ASP:OD1	6:E:434:HOH:O	2.18	0.53
1:F:265:ASP:OD1	6:F:402:HOH:O	2.18	0.53
1:B:263:PHE:HB2	1:B:266:LEU:HD12	1.89	0.53
1:B:262:PRO:O	1:C:193:HIS:HD2	1.92	0.53
1:E:230:PHE:O	1:E:234:VAL:HG23	2.08	0.52
1:D:132:LYS:HD3	6:D:425:HOH:O	2.09	0.52
1:C:41:PRO:O	1:C:42:ASP:CG	2.48	0.52
1:B:96:ALA:HB1	1:B:101:VAL:HB	1.92	0.52
1:B:53:LYS:O	1:B:54:LYS:C	2.46	0.52
1:C:260:ILE:O	1:C:260:ILE:HG13	2.09	0.52
1:A:96:ALA:HB1	1:A:101:VAL:HB	1.92	0.51
1:F:260:ILE:HG13	1:F:260:ILE:O	2.09	0.51
1:D:138:ALA:O	1:D:189:LEU:HA	2.10	0.51
1:D:159:ILE:CG1	6:D:437:HOH:O	2.56	0.51
1:C:53:LYS:HG2	1:C:207:LEU:HD23	1.92	0.51
1:A:38:GLU:OE2	1:A:92:GLU:OE1	2.28	0.51
1:D:137:PHE:CD2	1:D:188:ALA:HB3	2.46	0.51
1:F:125:VAL:O	1:F:129:ARG:HG3	2.11	0.50
1:B:38:GLU:HB2	1:B:43:ASP:HB2	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:262:PRO:O	1:C:193:HIS:CD2	2.65	0.50
1:F:46:ILE:HG23	1:F:231:LEU:HD21	1.94	0.49
1:B:129:ARG:NH1	1:B:170:PHE:O	2.45	0.49
1:D:20:LEU:HD21	1:D:256:LEU:HD22	1.94	0.49
1:D:14:LYS:HG3	1:D:18:GLU:OE1	2.13	0.49
1:A:265:ASP:OD1	6:A:407:HOH:O	2.20	0.49
1:B:55:LEU:O	1:B:60:VAL:HG23	2.12	0.49
1:B:265:ASP:OD1	6:B:402:HOH:O	2.20	0.48
1:B:132:LYS:HB3	1:B:182:HIS:CE1	2.48	0.48
1:D:113:GLU:O	1:D:115:PRO:HD3	2.13	0.48
1:B:90:ARG:CG	1:B:90:ARG:NH1	2.57	0.48
1:A:162:VAL:HG11	1:A:187:ILE:HD11	1.95	0.48
1:B:62:VAL:O	1:B:101:VAL:HG13	2.14	0.48
3:E:302:HEZ:H11	6:E:442:HOH:O	2.13	0.48
1:F:38:GLU:OE2	1:F:92:GLU:OE1	2.32	0.48
1:F:175:ILE:HD12	1:F:179:LEU:O	2.14	0.48
1:F:265:ASP:CG	6:F:402:HOH:O	2.51	0.48
1:A:166:GLN:HE21	1:A:183:SER:HA	1.78	0.48
1:E:118:ARG:HG2	1:E:122:LYS:HE3	1.96	0.48
1:C:102:ARG:CG	1:C:102:ARG:HH11	2.27	0.47
1:E:38:GLU:O	1:E:67:MET:HG2	2.15	0.47
1:E:28:PHE:O	1:E:30:ASP:N	2.48	0.47
1:B:41:PRO:O	1:B:42:ASP:CB	2.63	0.47
1:C:54:LYS:HE3	1:C:200:ASP:O	2.15	0.47
1:B:265:ASP:OD2	6:B:402:HOH:O	2.21	0.47
1:C:126:LYS:NZ	6:C:465:HOH:O	2.46	0.46
1:D:35:ILE:HG21	1:D:65:ILE:HD12	1.97	0.46
5:A:305:GOL:C1	1:C:259:HIS:NE2	2.78	0.46
1:E:256:LEU:HD21	1:F:230:PHE:CD1	2.50	0.46
1:F:41:PRO:O	1:F:42:ASP:CB	2.64	0.46
1:D:162:VAL:HG11	1:D:187:ILE:HD11	1.96	0.46
1:C:226:THR:C	1:C:229:PRO:HD2	2.36	0.46
1:E:54:LYS:NZ	1:E:200:ASP:O	2.49	0.46
1:F:53:LYS:HD3	1:F:99:LEU:HA	1.96	0.46
1:A:141:PRO:HG2	1:A:142:TRP:CE3	2.51	0.46
1:E:164:PHE:CZ	1:F:112:THR:HG22	2.51	0.46
1:B:42:ASP:OD1	1:B:42:ASP:C	2.55	0.45
1:C:48:MET:CE	1:C:137:PHE:CD2	2.99	0.45
1:E:12:PHE:HZ	1:E:267:ILE:HG23	1.80	0.45
1:E:38:GLU:OE2	1:E:92:GLU:OE1	2.34	0.45
1:A:23:ASN:HD21	1:A:25:GLU:CB	2.11	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:53:LYS:O	1:B:56:SER:N	2.49	0.45
1:E:260:ILE:HA	1:F:143:LEU:HD21	1.99	0.45
1:C:41:PRO:O	1:C:42:ASP:CB	2.64	0.45
1:D:147:SER:HB2	1:F:258:TYR:O	2.16	0.45
1:C:40:HIS:HD2	1:C:68:THR:OG1	2.00	0.45
1:C:79:LEU:HD11	1:C:83:GLU:HG2	1.99	0.45
1:F:38:GLU:HA	1:F:39:PRO:HD3	1.83	0.44
1:B:261:THR:HA	1:B:262:PRO:HD2	1.77	0.44
1:C:83:GLU:HB2	1:F:86:LEU:HD13	1.98	0.44
1:A:172:ASN:HA	6:A:444:HOH:O	2.18	0.44
3:E:302:HEZ:C1	6:E:442:HOH:O	2.65	0.44
1:F:116:TYR:CZ	1:F:121:ARG:NH2	2.85	0.44
1:B:89:ARG:HH21	1:B:108:ASN:HD21	1.65	0.44
1:D:38:GLU:OE2	1:D:92:GLU:OE1	2.35	0.44
1:A:38:GLU:HB2	1:A:43:ASP:HB2	2.00	0.43
1:D:97:LYS:HE2	1:D:97:LYS:HB2	1.93	0.43
1:F:18:GLU:HB3	6:F:401:HOH:O	2.19	0.43
1:C:54:LYS:CE	1:C:200:ASP:O	2.66	0.43
1:D:35:ILE:HD12	1:D:133:PRO:HG3	2.00	0.43
1:A:180:LYS:HE3	1:A:180:LYS:HB3	1.68	0.43
1:C:67:MET:HG3	1:C:154:THR:HG23	2.01	0.43
1:D:175:ILE:CG1	1:D:181:PRO:HD3	2.48	0.43
1:B:63:ILE:HG23	1:B:103:LYS:O	2.17	0.43
1:C:126:LYS:NZ	1:E:174:ASP:OD2	2.46	0.43
1:F:228:GLU:HB3	1:F:229:PRO:HD3	2.01	0.43
1:F:265:ASP:OD2	6:F:402:HOH:O	2.21	0.43
1:B:265:ASP:CG	6:B:402:HOH:O	2.54	0.42
1:E:137:PHE:CD2	1:E:188:ALA:HB3	2.53	0.42
1:E:138:ALA:O	1:E:189:LEU:HA	2.19	0.42
1:C:76:ASP:OD2	1:C:78:LYS:HG2	2.18	0.42
1:C:78:LYS:HG2	1:C:78:LYS:H	1.57	0.42
1:B:74:THR:HB	1:B:84:LEU:HD22	1.99	0.42
1:E:175:ILE:HD11	1:E:181:PRO:HD3	2.02	0.42
1:F:40:HIS:HD2	1:F:68:THR:OG1	2.02	0.42
1:C:228:GLU:O	1:C:232:ARG:HG3	2.19	0.42
1:A:113:GLU:O	1:A:115:PRO:HD3	2.19	0.42
1:E:71:TYR:CD2	1:E:110:ARG:HD3	2.54	0.42
1:F:38:GLU:HB2	1:F:43:ASP:HB2	2.02	0.42
1:F:102:ARG:HH11	1:F:102:ARG:CG	2.33	0.42
1:F:232:ARG:O	1:F:236:MET:HG3	2.20	0.42
1:C:23:ASN:CG	1:C:24:LEU:N	2.72	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:125:VAL:HG21	1:B:170:PHE:CE2	2.55	0.42
1:F:35:ILE:HA	1:F:63:ILE:O	2.20	0.42
1:F:221:ASP:O	1:F:225:GLU:HG3	2.19	0.42
1:C:52:ILE:HG23	1:C:62:VAL:HG11	2.02	0.41
1:B:74:THR:OG1	6:B:419:HOH:O	2.19	0.41
1:D:35:ILE:HA	1:D:63:ILE:O	2.21	0.41
1:F:98:LEU:HD12	1:F:211:ALA:HA	2.02	0.41
1:C:102:ARG:CG	1:C:102:ARG:NH1	2.83	0.41
1:A:228:GLU:OE1	1:A:232:ARG:NH2	2.48	0.41
1:E:78:LYS:HE3	1:E:78:LYS:HB2	1.77	0.41
1:A:162:VAL:CG1	1:A:187:ILE:HD11	2.51	0.41
1:C:44:CYS:HB3	1:C:52:ILE:HD11	2.01	0.41
1:C:174:ASP:O	1:C:179:LEU:HB2	2.20	0.41
1:D:156:PHE:N	6:D:437:HOH:O	2.45	0.41
1:F:198:ILE:HD13	1:F:249:GLY:HA2	2.02	0.41
1:B:134:ASP:O	1:B:185:SER:HB2	2.21	0.40
1:C:102:ARG:HH11	1:C:102:ARG:HG2	1.86	0.40
1:C:202:THR:HA	1:C:246:TYR:HB2	2.03	0.40
1:E:193:HIS:CE1	1:E:194:LYS:HE2	2.56	0.40
1:A:56:SER:O	1:A:59:GLY:N	2.50	0.40
1:C:190:TYR:O	1:C:191:TYR:C	2.58	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	261/267 (98%)	249 (95%)	10 (4%)	2 (1%)	19 7
1	B	259/267 (97%)	246 (95%)	12 (5%)	1 (0%)	34 21
1	C	265/267 (99%)	258 (97%)	6 (2%)	1 (0%)	34 21
1	D	258/267 (97%)	241 (93%)	14 (5%)	3 (1%)	13 3

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	E	255/267 (96%)	235 (92%)	16 (6%)	4 (2%)	9 2
1	F	265/267 (99%)	255 (96%)	8 (3%)	2 (1%)	19 7
All	All	1563/1602 (98%)	1484 (95%)	66 (4%)	13 (1%)	19 7

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	29	GLU
1	A	42	ASP
1	B	42	ASP
1	D	42	ASP
1	D	155	GLY
1	E	42	ASP
1	E	240	GLU
1	F	220	THR
1	C	42	ASP
1	F	42	ASP
1	A	260	ILE
1	E	255	GLY
1	D	260	ILE

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	235/239 (98%)	225 (96%)	10 (4%)	29 14
1	B	233/239 (98%)	228 (98%)	5 (2%)	53 42
1	C	239/239 (100%)	230 (96%)	9 (4%)	33 18
1	D	232/239 (97%)	228 (98%)	4 (2%)	60 51
1	E	229/239 (96%)	221 (96%)	8 (4%)	36 21
1	F	239/239 (100%)	233 (98%)	6 (2%)	47 34
All	All	1407/1434 (98%)	1365 (97%)	42 (3%)	41 27

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	SER
1	A	14	LYS
1	A	23	ASN
1	A	26	ASN
1	A	56	SER
1	A	77	GLU
1	A	78	LYS
1	A	89	ARG
1	A	175	ILE
1	A	222	ASP
1	B	32	LYS
1	B	78	LYS
1	B	90	ARG
1	B	213	ARG
1	B	240	GLU
1	C	6	SER
1	C	56	SER
1	C	78	LYS
1	C	95	SER
1	C	102	ARG
1	C	166	GLN
1	C	175	ILE
1	C	213	ARG
1	C	245	ARG
1	D	25	GLU
1	D	78	LYS
1	D	97	LYS
1	D	115	PRO
1	E	19	VAL
1	E	78	LYS
1	E	89	ARG
1	E	97	LYS
1	E	175	ILE
1	E	204	LEU
1	E	206	GLU
1	E	216	ARG
1	F	23	ASN
1	F	26	ASN
1	F	78	LYS
1	F	97	LYS
1	F	175	ILE
1	F	206	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	ASN
1	A	108	ASN
1	A	166	GLN
1	B	23	ASN
1	B	108	ASN
1	C	108	ASN
1	C	166	GLN
1	D	23	ASN
1	D	26	ASN
1	D	108	ASN
1	D	166	GLN
1	E	26	ASN
1	E	108	ASN
1	E	166	GLN
1	F	23	ASN
1	F	108	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)

Of 22 ligands modelled in this entry, 14 are monoatomic - leaving 8 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	$\#\|Z\| > 2$
3	HEZ	D	303	-	7,7,7	0.38	0	6,6,6	0.51	0
3	HEZ	F	302	-	7,7,7	0.43	0	6,6,6	0.40	0
5	GOL	D	305	2	5,5,5	0.40	0	5,5,5	0.46	0
3	HEZ	C	302	-	7,7,7	0.77	0	6,6,6	0.87	0
3	HEZ	A	303	-	7,7,7	0.54	0	6,6,6	0.82	0
5	GOL	A	305	2	5,5,5	0.58	0	5,5,5	1.00	0
3	HEZ	B	302	-	7,7,7	0.48	0	6,6,6	0.38	0
3	HEZ	E	302	-	7,7,7	0.28	0	6,6,6	0.89	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	D	303	-	-	4/5/5/5	-
3	HEZ	F	302	-	-	4/5/5/5	-
5	GOL	D	305	2	-	3/4/4/4	-
3	HEZ	C	302	-	-	1/5/5/5	-
3	HEZ	A	303	-	-	2/5/5/5	-
5	GOL	A	305	2	-	2/4/4/4	-
3	HEZ	B	302	-	-	3/5/5/5	-
3	HEZ	E	302	-	-	4/5/5/5	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	D	305	GOL	C1-C2-C3-O3
3	F	302	HEZ	C2-C3-C4-C5
5	A	305	GOL	O1-C1-C2-C3
5	D	305	GOL	O1-C1-C2-C3
3	F	302	HEZ	C1-C2-C3-C4
5	D	305	GOL	O2-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
3	E	302	HEZ	C2-C3-C4-C5
3	D	303	HEZ	C3-C4-C5-C6
3	B	302	HEZ	C2-C3-C4-C5
3	D	303	HEZ	O1-C1-C2-C3
5	A	305	GOL	O1-C1-C2-O2
3	E	302	HEZ	O1-C1-C2-C3
3	F	302	HEZ	O1-C1-C2-C3
3	B	302	HEZ	O1-C1-C2-C3
3	D	303	HEZ	C4-C5-C6-O6
3	E	302	HEZ	C3-C4-C5-C6
3	F	302	HEZ	C4-C5-C6-O6
3	D	303	HEZ	C2-C3-C4-C5
3	E	302	HEZ	C4-C5-C6-O6
3	A	303	HEZ	C3-C4-C5-C6
3	A	303	HEZ	C4-C5-C6-O6
3	C	302	HEZ	C1-C2-C3-C4
3	B	302	HEZ	C1-C2-C3-C4

There are no ring outliers.

8 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	303	HEZ	1	0
3	F	302	HEZ	1	0
5	D	305	GOL	1	0
3	C	302	HEZ	1	0
3	A	303	HEZ	2	0
5	A	305	GOL	2	0
3	B	302	HEZ	1	0
3	E	302	HEZ	3	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 (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	263/267 (98%)	0.25	16 (6%)	21	16	13, 31, 61, 71	0
1	B	261/267 (97%)	0.64	30 (11%)	4	3	19, 44, 65, 76	0
1	C	267/267 (100%)	0.37	10 (3%)	41	36	17, 35, 52, 58	0
1	D	260/267 (97%)	0.30	14 (5%)	25	20	9, 33, 62, 70	0
1	E	257/267 (96%)	0.85	35 (13%)	3	2	21, 46, 81, 90	0
1	F	267/267 (100%)	0.52	22 (8%)	11	9	18, 39, 60, 76	0
All	All	1575/1602 (98%)	0.49	127 (8%)	12	9	9, 37, 64, 90	0

All (127) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	16	LEU	5.8
1	E	15	LEU	5.1
1	F	245	ARG	5.1
1	A	5	VAL	5.0
1	B	244	VAL	5.0
1	D	11	ALA	5.0
1	E	17	LYS	4.6
1	A	13	ASN	4.6
1	B	14	LYS	4.5
1	B	30	ASP	4.4
1	A	17	LYS	4.4
1	E	202	THR	4.1
1	D	23	ASN	4.1
1	A	10	GLU	4.1
1	E	244	VAL	4.0
1	E	25	GLU	3.9
1	E	209	LEU	3.8
1	E	11	ALA	3.8
1	E	19	VAL	3.7

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Mol	Chain	Res	Type	RSRZ
1	C	243	GLY	3.7
1	E	23	ASN	3.7
1	D	17	LYS	3.7
1	F	25	GLU	3.7
1	B	20	LEU	3.6
1	E	231	LEU	3.6
1	F	242	ILE	3.6
1	F	246	TYR	3.6
1	B	22	PHE	3.5
1	D	8	PHE	3.5
1	D	267	ILE	3.5
1	B	21	GLU	3.5
1	B	243	GLY	3.5
1	C	236	MET	3.5
1	B	8	PHE	3.4
1	E	24	LEU	3.4
1	D	10	GLU	3.4
1	A	18	GLU	3.4
1	E	21	GLU	3.3
1	E	267	ILE	3.3
1	E	102	ARG	3.3
1	B	11	ALA	3.2
1	B	17	LYS	3.2
1	E	13	ASN	3.2
1	E	240	GLU	3.2
1	D	14	LYS	3.2
1	B	59	GLY	3.2
1	E	26	ASN	3.2
1	E	246	TYR	3.1
1	C	201	ILE	3.1
1	E	201	ILE	3.1
1	B	25	GLU	3.1
1	F	227	TRP	3.0
1	E	243	GLY	3.0
1	B	15	LEU	3.0
1	E	266	LEU	2.9
1	C	25	GLU	2.9
1	F	236	MET	2.9
1	B	219	PHE	2.9
1	F	240	GLU	2.9
1	E	203	ASP	2.8
1	D	19	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	246	TYR	2.8
1	F	239	GLY	2.8
1	E	204	LEU	2.7
1	A	102	ARG	2.7
1	E	22	PHE	2.7
1	F	244	VAL	2.7
1	F	230	PHE	2.7
1	F	90	ARG	2.7
1	B	26	ASN	2.7
1	D	216	ARG	2.6
1	E	197	TYR	2.6
1	E	12	PHE	2.5
1	E	49	GLY	2.5
1	A	20	LEU	2.5
1	B	19	VAL	2.4
1	B	7	THR	2.4
1	F	243	GLY	2.4
1	B	236	MET	2.4
1	A	19	VAL	2.4
1	C	23	ASN	2.4
1	F	78	LYS	2.3
1	E	222	ASP	2.3
1	B	10	GLU	2.3
1	E	245	ARG	2.3
1	E	18	GLU	2.3
1	F	233	THR	2.3
1	F	234	VAL	2.3
1	A	145	TYR	2.2
1	F	219	PHE	2.2
1	A	14	LYS	2.2
1	E	14	LYS	2.2
1	A	9	GLU	2.2
1	B	203	ASP	2.2
1	C	203	ASP	2.2
1	B	225	GLU	2.2
1	D	18	GLU	2.2
1	F	224	TRP	2.2
1	A	15	LEU	2.2
1	A	12	PHE	2.2
1	C	237	PHE	2.2
1	B	240	GLU	2.2
1	B	237	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
1	F	222	ASP	2.2
1	E	205	MET	2.1
1	B	18	GLU	2.1
1	E	60	VAL	2.1
1	B	16	LEU	2.1
1	A	29	GLU	2.1
1	A	23	ASN	2.1
1	D	102	ARG	2.1
1	D	193	HIS	2.1
1	E	238	TYR	2.1
1	C	159	ILE	2.1
1	B	231	LEU	2.1
1	D	25	GLU	2.1
1	D	204	LEU	2.1
1	F	143	LEU	2.1
1	F	26	ASN	2.0
1	C	244	VAL	2.0
1	B	12	PHE	2.0
1	F	214	ALA	2.0
1	A	16	LEU	2.0
1	B	23	ASN	2.0
1	B	24	LEU	2.0
1	B	98	LEU	2.0
1	F	231	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	F	302	8/8	0.81	0.36	47,47,50,50	0
3	HEZ	A	303	8/8	0.83	0.33	31,36,43,44	0
3	HEZ	C	302	8/8	0.85	0.26	28,33,40,45	0
3	HEZ	D	303	8/8	0.87	0.33	36,40,42,43	0
3	HEZ	E	302	8/8	0.88	0.32	29,35,36,39	0
3	HEZ	B	302	8/8	0.90	0.23	34,36,41,47	0
5	GOL	A	305	6/6	0.91	0.16	26,28,32,37	0
5	GOL	D	305	6/6	0.94	0.09	26,32,33,37	0
4	CL	F	303	1/1	0.96	0.08	40,40,40,40	0
4	CL	F	304	1/1	0.98	0.04	38,38,38,38	0
4	CL	D	304	1/1	0.98	0.05	36,36,36,36	0
4	CL	C	303	1/1	0.98	0.05	34,34,34,34	0
2	CD	E	301	1/1	0.99	0.05	36,36,36,36	0
2	CD	A	301	1/1	0.99	0.09	23,23,23,23	0
2	CD	B	301	1/1	0.99	0.04	30,30,30,30	0
2	CD	D	302	1/1	0.99	0.04	31,31,31,31	0
4	CL	C	304	1/1	0.99	0.09	43,43,43,43	0
2	CD	D	301	1/1	1.00	0.10	27,27,27,27	0
2	CD	A	302	1/1	1.00	0.07	37,37,37,37	0
4	CL	A	304	1/1	1.00	0.04	35,35,35,35	0
2	CD	C	301	1/1	1.00	0.06	31,31,31,31	0
2	CD	F	301	1/1	1.00	0.07	40,40,40,40	0

6.5 Other polymers (i)

There are no such residues in this entry.