



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 7, 2023 – 01:57 PM JST

PDB ID : 5HZS
Title : Crystal structure of Dronpa-Co2+
Authors : Hwang, K.Y.; Nam, K.H.
Deposited on : 2016-02-03
Resolution : 2.17 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)
Xtrriage (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

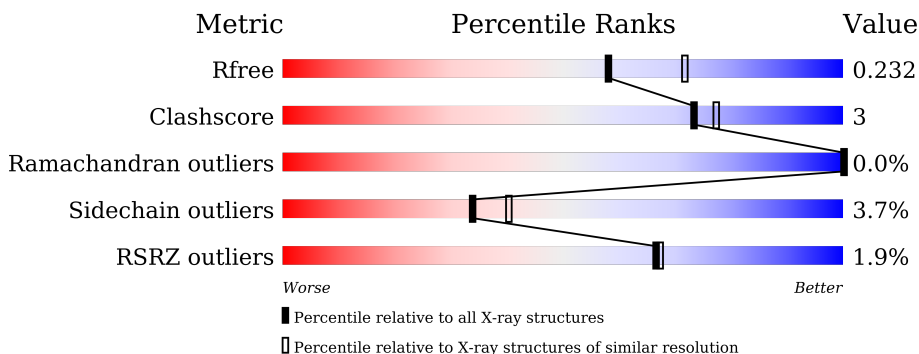
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.17 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	6864 (2.20-2.16)
Clashscore	141614	7689 (2.20-2.16)
Ramachandran outliers	138981	7564 (2.20-2.16)
Sidechain outliers	138945	7564 (2.20-2.16)
RSRZ outliers	127900	6738 (2.20-2.16)

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	215	 2% 87% 11%
1	B	215	 92% 7%
1	C	215	 91% 7%
1	D	215	 89% 11%
1	E	215	 2% 88% 10%
1	F	215	 2% 88% 10%

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Mol	Chain	Length	Quality of chain
1	G	215	 <p>% 93% 7%</p>
1	H	215	 <p>89% 9%</p>
1	I	215	 <p>% 86% 13%</p>
1	J	215	 <p>6% 86% 12%</p>
1	K	215	 <p>4% 89% 9%</p>
1	L	215	 <p>2% 87% 13%</p>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 22118 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 Fluorescent protein Dronpa.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	215	1736	1108	292	327	9	0	0	0
1	B	213	1721	1100	290	322	9	0	0	0
1	C	215	1736	1108	292	327	9	0	0	0
1	D	215	1736	1108	292	327	9	0	0	0
1	E	215	1736	1108	292	327	9	0	0	0
1	F	214	1730	1105	291	325	9	0	0	0
1	G	214	1727	1103	291	324	9	0	0	0
1	H	214	1727	1103	291	324	9	0	0	0
1	I	213	1721	1100	290	322	9	0	0	0
1	J	214	1727	1103	291	324	9	0	0	0
1	K	213	1721	1100	290	322	9	0	0	0
1	L	214	1730	1105	291	325	9	0	0	0

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	CYS	chromophore	UNP Q5TLG6
A	?	-	TYR	chromophore	UNP Q5TLG6
A	62	GYS	GLY	chromophore	UNP Q5TLG6
B	?	-	CYS	chromophore	UNP Q5TLG6
B	?	-	TYR	chromophore	UNP Q5TLG6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	62	GYS	GLY	chromophore	UNP Q5TLG6
C	?	-	CYS	chromophore	UNP Q5TLG6
C	?	-	TYR	chromophore	UNP Q5TLG6
C	62	GYS	GLY	chromophore	UNP Q5TLG6
D	?	-	CYS	chromophore	UNP Q5TLG6
D	?	-	TYR	chromophore	UNP Q5TLG6
D	62	GYS	GLY	chromophore	UNP Q5TLG6
E	?	-	CYS	chromophore	UNP Q5TLG6
E	?	-	TYR	chromophore	UNP Q5TLG6
E	62	GYS	GLY	chromophore	UNP Q5TLG6
F	?	-	CYS	chromophore	UNP Q5TLG6
F	?	-	TYR	chromophore	UNP Q5TLG6
F	62	GYS	GLY	chromophore	UNP Q5TLG6
G	?	-	CYS	chromophore	UNP Q5TLG6
G	?	-	TYR	chromophore	UNP Q5TLG6
G	62	GYS	GLY	chromophore	UNP Q5TLG6
H	?	-	CYS	chromophore	UNP Q5TLG6
H	?	-	TYR	chromophore	UNP Q5TLG6
H	62	GYS	GLY	chromophore	UNP Q5TLG6
I	?	-	CYS	chromophore	UNP Q5TLG6
I	?	-	TYR	chromophore	UNP Q5TLG6
I	62	GYS	GLY	chromophore	UNP Q5TLG6
J	?	-	CYS	chromophore	UNP Q5TLG6
J	?	-	TYR	chromophore	UNP Q5TLG6
J	62	GYS	GLY	chromophore	UNP Q5TLG6
K	?	-	CYS	chromophore	UNP Q5TLG6
K	?	-	TYR	chromophore	UNP Q5TLG6
K	62	GYS	GLY	chromophore	UNP Q5TLG6
L	?	-	CYS	chromophore	UNP Q5TLG6
L	?	-	TYR	chromophore	UNP Q5TLG6
L	62	GYS	GLY	chromophore	UNP Q5TLG6

- Molecule 2 is COBALT (II) ION (three-letter code: CO) (formula: Co).

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

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	E	2	Total 2	Co 2	0	0
2	F	2	Total 2	Co 2	0	0
2	G	2	Total 2	Co 2	0	0
2	H	2	Total 2	Co 2	0	0
2	I	2	Total 2	Co 2	0	0
2	J	2	Total 2	Co 2	0	0
2	K	2	Total 2	Co 2	0	0
2	L	2	Total 2	Co 2	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	115	Total 115	O 115	0	0
3	B	122	Total 122	O 122	0	0
3	C	128	Total 128	O 128	0	0
3	D	118	Total 118	O 118	0	0
3	E	125	Total 125	O 125	0	0
3	F	122	Total 122	O 122	0	0
3	G	119	Total 119	O 119	0	0
3	H	149	Total 149	O 149	0	0
3	I	77	Total 77	O 77	0	0
3	J	77	Total 77	O 77	0	0
3	K	93	Total 93	O 93	0	0

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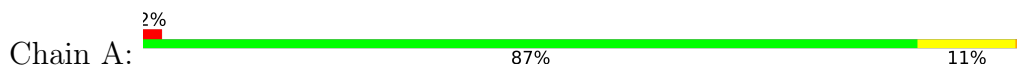
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	L	101	Total 101	O 101	0	0

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.

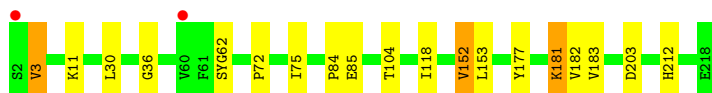
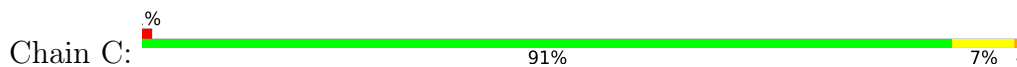
- Molecule 1: Fluorescent protein Dronpa



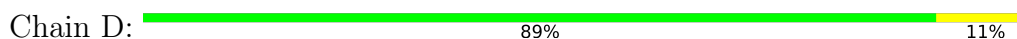
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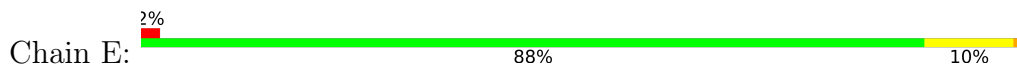
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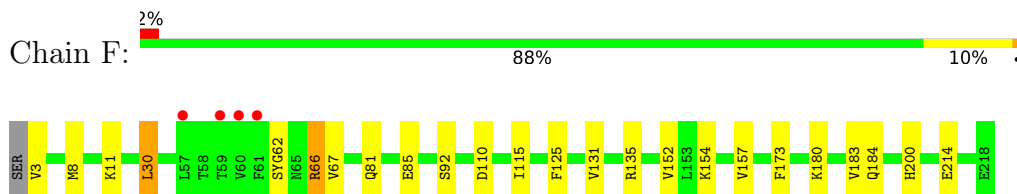
- Molecule 1: Fluorescent protein Dronpa



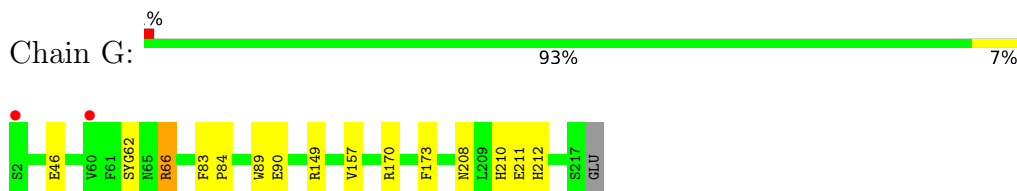
- Molecule 1: Fluorescent protein Dronpa



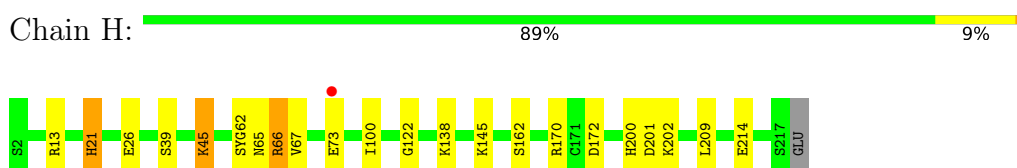
- Molecule 1: Fluorescent protein Dronpa



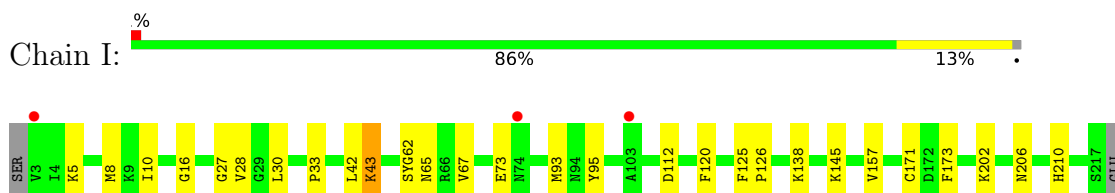
- Molecule 1: Fluorescent protein Dronpa



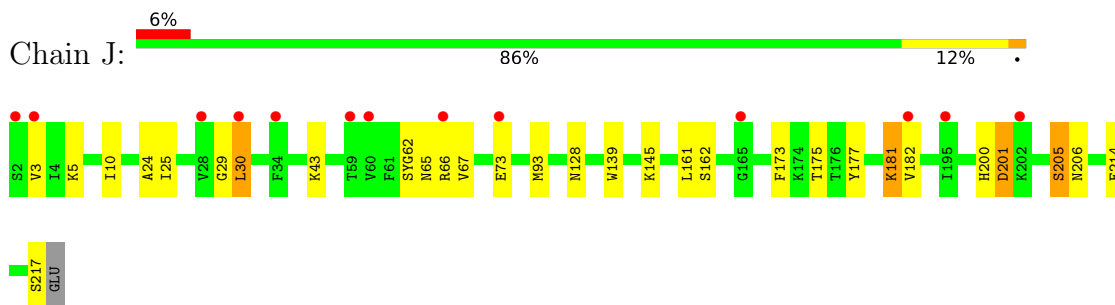
- Molecule 1: Fluorescent protein Dronpa



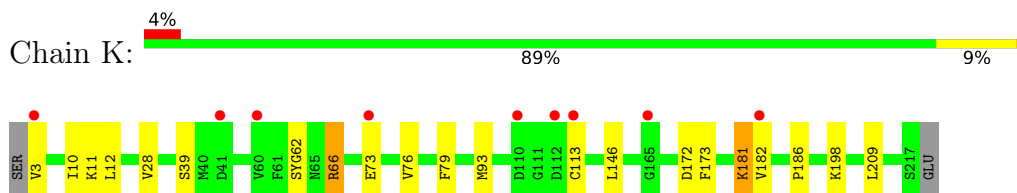
- Molecule 1: Fluorescent protein Dronpa



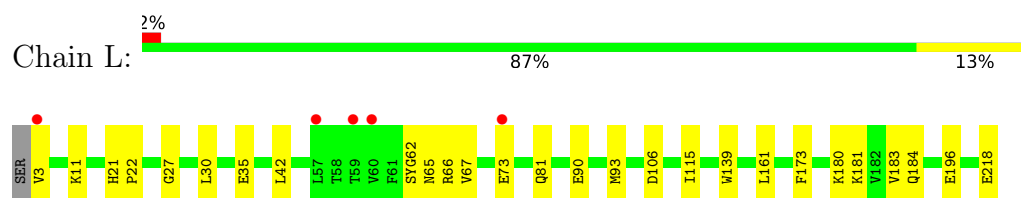
- Molecule 1: Fluorescent protein Dronpa



- Molecule 1: Fluorescent protein Dronpa



- Molecule 1: Fluorescent protein Dronpa



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	72.33Å 105.48Å 109.78Å 115.79° 109.51° 93.95°	Depositor
Resolution (Å)	29.57 – 2.17 29.11 – 2.17	Depositor EDS
% Data completeness (in resolution range)	93.1 (29.57-2.17) 93.2 (29.11-2.17)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 2.18Å)	Xtrriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.166 , 0.225 0.177 , 0.232	Depositor DCC
R_{free} test set	6546 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	28.0	Xtrriage
Anisotropy	0.100	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 40.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	0.002 for -h,-k,h+k+l	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	22118	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.69% 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

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

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	0.91	0/1760	0.93	1/2375 (0.0%)
1	B	0.88	0/1745	0.89	1/2355 (0.0%)
1	C	0.91	0/1760	0.92	0/2375
1	D	0.86	0/1760	0.92	3/2375 (0.1%)
1	E	0.92	1/1760 (0.1%)	0.93	2/2375 (0.1%)
1	F	0.93	0/1754	0.94	1/2367 (0.0%)
1	G	0.87	1/1751 (0.1%)	0.91	3/2363 (0.1%)
1	H	0.93	0/1751	0.97	2/2363 (0.1%)
1	I	0.83	0/1745	0.89	0/2355
1	J	0.76	0/1751	0.87	0/2363
1	K	0.74	0/1745	0.85	1/2355 (0.0%)
1	L	0.80	0/1754	0.88	2/2367 (0.1%)
All	All	0.86	2/21036 (0.0%)	0.91	16/28388 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	89	TRP	CB-CG	-6.42	1.38	1.50
1	G	89	TRP	CB-CG	6.13	1.61	1.50

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	112	ASP	CB-CG-OD1	6.84	124.45	118.30
1	D	170	ARG	NE-CZ-NH1	6.81	123.70	120.30
1	E	41	ASP	CB-CG-OD1	5.79	123.51	118.30
1	G	66	ARG	NE-CZ-NH2	-5.48	117.56	120.30
1	L	106	ASP	CB-CG-OD1	5.40	123.16	118.30
1	A	187	ASP	CB-CG-OD1	5.39	123.16	118.30
1	E	119	ARG	NE-CZ-NH2	-5.39	117.60	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	66	ARG	NE-CZ-NH2	-5.38	117.61	120.30
1	H	170	ARG	NE-CZ-NH1	5.36	122.98	120.30
1	L	106	ASP	CB-CG-OD2	-5.23	113.59	118.30
1	G	149	ARG	NE-CZ-NH2	5.23	122.91	120.30
1	H	172	ASP	CB-CG-OD2	-5.22	113.60	118.30
1	G	170	ARG	NE-CZ-NH1	5.22	122.91	120.30
1	K	172	ASP	CB-CG-OD2	-5.20	113.62	118.30
1	B	170	ARG	NE-CZ-NH1	5.04	122.82	120.30
1	D	162	SER	N-CA-CB	-5.02	102.98	110.50

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	1736	0	1658	14	0
1	B	1721	0	1647	5	0
1	C	1736	0	1658	13	0
1	D	1736	0	1658	12	0
1	E	1736	0	1657	14	0
1	F	1730	0	1653	11	0
1	G	1727	0	1652	5	0
1	H	1727	0	1651	13	0
1	I	1721	0	1646	17	0
1	J	1727	0	1652	14	0
1	K	1721	0	1646	10	0
1	L	1730	0	1652	11	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	2	0	0	0	0
2	G	2	0	0	0	0
2	H	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	I	2	0	0	0	0
2	J	2	0	0	0	0
2	K	2	0	0	0	0
2	L	2	0	0	1	0
3	A	115	0	0	1	0
3	B	122	0	0	1	0
3	C	128	0	0	1	0
3	D	118	0	0	0	0
3	E	125	0	0	3	0
3	F	122	0	0	3	0
3	G	119	0	0	1	0
3	H	149	0	0	6	0
3	I	77	0	0	0	0
3	J	77	0	0	1	0
3	K	93	0	0	0	0
3	L	101	0	0	2	0
All	All	22118	0	19830	135	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 3.

All (135) 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:H:21:HIS:NE2	3:H:401:HOH:O	1.95	0.99
1:H:21:HIS:CD2	3:H:401:HOH:O	2.31	0.79
1:A:65:ASN:OD1	1:A:67:VAL:HG22	1.86	0.76
1:A:135:ARG:NE	1:E:48:GLY:O	2.20	0.75
1:B:21:HIS:NE2	3:B:401:HOH:O	2.22	0.72
1:I:43:LYS:HD2	1:I:206:ASN:HD21	1.54	0.72
1:F:110:ASP:HB3	1:F:115:ILE:HD11	1.73	0.71
1:E:21:HIS:CD2	3:E:403:HOH:O	2.42	0.70
1:J:214:GLU:HG3	3:J:445:HOH:O	1.90	0.70
2:L:302:CO:CO	3:L:485:HOH:O	1.38	0.69
1:E:217:SER:O	1:E:218:GLU:HB2	1.91	0.68
1:E:2:SER:O	1:E:5:LYS:NZ	2.28	0.66
1:L:65:ASN:OD1	1:L:67:VAL:HG22	1.96	0.66
1:E:21:HIS:NE2	3:E:403:HOH:O	2.29	0.65
1:C:181:LYS:HE3	1:C:183:VAL:HG12	1.77	0.65
1:C:72:PRO:HG2	1:C:75:ILE:HD12	1.80	0.62
1:F:125:PHE:CE1	1:F:131:VAL:HG21	2.34	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:30:LEU:N	1:L:30:LEU:HD23	2.17	0.60
1:A:217:SER:O	1:A:218:GLU:HB2	2.02	0.59
1:K:146:LEU:N	1:K:146:LEU:HD12	2.17	0.59
1:G:157:VAL:HG13	1:G:173:PHE:HB2	1.84	0.59
1:D:143:THR:H	1:E:145:LYS:NZ	2.02	0.57
1:I:65:ASN:OD1	1:I:67:VAL:HG22	2.03	0.57
1:F:85:GLU:OE1	1:F:85:GLU:N	2.33	0.57
1:C:181:LYS:CE	1:C:183:VAL:HG12	2.33	0.57
1:K:93:MET:HG2	1:K:173:PHE:CD1	2.38	0.57
1:J:181:LYS:HG2	1:J:182:VAL:N	2.19	0.57
1:J:200:HIS:ND1	1:J:201:ASP:O	2.37	0.56
1:G:46:GLU:OE2	3:G:401:HOH:O	2.18	0.56
1:C:85:GLU:H	1:C:85:GLU:CD	2.10	0.55
1:C:72:PRO:CG	1:C:75:ILE:HD12	2.36	0.55
1:E:218:GLU:HA	1:E:218:GLU:OE1	2.05	0.55
1:G:83:PHE:HB3	1:G:84:PRO:HA	1.89	0.55
1:D:143:THR:H	1:E:145:LYS:HZ3	1.56	0.54
1:J:139:TRP:CZ3	1:J:161:LEU:HG	2.43	0.54
1:K:93:MET:HG2	1:K:173:PHE:CE1	2.44	0.53
1:I:8:MET:HG2	1:I:112:ASP:O	2.09	0.53
1:D:148:VAL:HG11	1:D:185:LEU:HD13	1.89	0.53
1:J:93:MET:HG2	1:J:173:PHE:CD1	2.44	0.52
1:A:93:MET:HG2	1:A:173:PHE:CD1	2.45	0.51
1:C:152:VAL:HG22	1:C:177:TYR:O	2.10	0.51
1:H:21:HIS:CE1	3:H:403:HOH:O	2.64	0.51
1:A:2:SER:O	1:A:3:VAL:CG1	2.58	0.51
1:J:128:ASN:OD1	1:J:128:ASN:C	2.48	0.51
1:D:36:GLY:O	1:D:212:HIS:HA	2.10	0.51
1:J:30:LEU:HD23	1:J:30:LEU:N	2.26	0.51
1:A:93:MET:HG2	1:A:173:PHE:CE1	2.46	0.51
1:I:8:MET:HE3	1:I:33:PRO:HG2	1.93	0.50
1:I:65:ASN:CG	1:I:67:VAL:HG22	2.32	0.50
1:J:175:THR:HG21	1:J:177:TYR:CE1	2.47	0.50
1:J:24:ALA:C	1:J:25:ILE:HG13	2.34	0.48
1:L:3:VAL:N	3:L:407:HOH:O	2.46	0.48
1:C:152:VAL:HG22	1:C:153:LEU:H	1.78	0.48
1:I:93:MET:HG2	1:I:173:PHE:CD1	2.49	0.48
1:J:10:ILE:N	1:J:29:GLY:O	2.46	0.48
1:F:214:GLU:OE2	3:F:401:HOH:O	2.20	0.47
1:J:175:THR:HG21	1:J:177:TYR:CZ	2.49	0.47
1:B:94:ASN:HA	1:B:100:ILE:HD13	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:8:MET:CE	1:I:33:PRO:HG2	2.45	0.47
1:H:21:HIS:ND1	3:H:403:HOH:O	2.35	0.47
1:H:45:LYS:HE3	3:H:529:HOH:O	2.15	0.47
1:K:66:ARG:HG2	1:K:79:PHE:CE2	2.50	0.47
1:B:138:LYS:C	1:B:161:LEU:HD23	2.35	0.47
1:L:65:ASN:CG	1:L:67:VAL:HG22	2.34	0.46
1:L:139:TRP:CZ3	1:L:161:LEU:HG	2.50	0.46
1:K:11:LYS:HG2	1:K:113:CYS:SG	2.56	0.46
1:H:100:ILE:O	1:H:122:GLY:HA2	2.16	0.46
1:G:208:ASN:OD1	1:G:210:HIS:HE1	1.99	0.45
1:A:143:THR:H	1:H:145:LYS:NZ	2.14	0.45
1:B:72:PRO:HD2	1:B:75:ILE:HD12	1.98	0.45
1:F:154:LYS:NZ	3:F:403:HOH:O	2.30	0.45
1:J:65:ASN:OD1	1:J:67:VAL:HG22	2.16	0.45
1:C:3:VAL:HG21	1:C:84:PRO:HB3	1.99	0.45
1:A:76:VAL:HG22	1:A:186:PRO:HB3	1.97	0.45
1:I:27:GLY:HA3	1:I:42:LEU:HD23	1.99	0.45
1:D:85:GLU:OE1	1:D:85:GLU:N	2.47	0.44
1:I:157:VAL:HG12	1:I:173:PHE:HB2	1.98	0.44
1:K:12:LEU:C	1:K:12:LEU:HD12	2.37	0.44
1:A:2:SER:O	1:A:3:VAL:HG13	2.18	0.44
1:K:10:ILE:O	1:K:28:VAL:HA	2.18	0.44
1:I:157:VAL:CG1	1:I:173:PHE:HB2	2.48	0.43
1:A:65:ASN:CG	1:A:67:VAL:HG22	2.38	0.43
1:D:144:GLU:HG2	1:D:146:LEU:CD1	2.48	0.43
1:H:200:HIS:ND1	1:H:201:ASP:O	2.47	0.43
1:I:145:LYS:N	1:I:145:LYS:HD2	2.32	0.43
1:L:93:MET:HG2	1:L:173:PHE:CD1	2.53	0.43
1:H:39:SER:HA	1:H:209:LEU:O	2.18	0.43
1:A:157:VAL:HG13	1:A:173:PHE:HB2	2.00	0.43
1:K:181:LYS:HD2	1:K:182:VAL:O	2.19	0.43
1:L:81:GLN:NE2	1:L:184:GLN:H	2.16	0.43
1:I:5:LYS:O	1:I:8:MET:HE2	2.18	0.43
1:I:43:LYS:HD2	1:I:206:ASN:ND2	2.29	0.43
1:C:85:GLU:CD	1:C:85:GLU:N	2.72	0.43
1:C:104:THR:O	1:C:118:ILE:HA	2.18	0.43
1:E:17:ALA:HA	1:E:21:HIS:O	2.19	0.42
1:A:36:GLY:O	1:A:212:HIS:HA	2.19	0.42
1:E:125:PHE:CE1	1:E:131:VAL:HG21	2.54	0.42
1:E:157:VAL:CG1	1:E:173:PHE:HB2	2.49	0.42
1:L:81:GLN:HG3	1:L:183:VAL:HB	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:39:SER:HA	1:E:209:LEU:O	2.19	0.42
1:I:16:GLY:HA2	1:I:120:PHE:O	2.19	0.42
1:D:95:TYR:CD1	1:D:171:CYS:HB2	2.54	0.42
1:J:201:ASP:OD2	1:J:206:ASN:N	2.52	0.42
1:H:66:ARG:CZ	1:H:66:ARG:HB2	2.49	0.42
1:H:138:LYS:HE3	3:H:444:HOH:O	2.20	0.42
1:F:157:VAL:HG13	1:F:173:PHE:HB2	2.01	0.42
1:F:200:HIS:CE1	3:F:415:HOH:O	2.72	0.42
3:E:500:HOH:O	1:F:135:ARG:HG2	2.20	0.41
1:H:65:ASN:CG	1:H:67:VAL:HG12	2.40	0.41
1:F:81:GLN:HB3	1:F:183:VAL:HG13	2.01	0.41
1:C:182:VAL:HA	3:C:401:HOH:O	2.20	0.41
1:I:125:PHE:HA	1:I:126:PRO:HD3	1.93	0.41
1:L:21:HIS:HA	1:L:22:PRO:HD3	1.91	0.41
1:C:203:ASP:OD1	1:C:203:ASP:N	2.54	0.41
1:D:30:LEU:N	1:D:30:LEU:HD23	2.36	0.41
1:G:211:GLU:HG2	1:G:212:HIS:N	2.36	0.41
1:C:36:GLY:O	1:C:212:HIS:HA	2.21	0.41
1:D:84:PRO:HD2	1:D:85:GLU:OE1	2.21	0.41
1:E:128:ASN:O	1:E:135:ARG:NH2	2.54	0.41
1:E:147:TYR:HB3	1:E:188:TYR:CD1	2.56	0.41
1:K:39:SER:HA	1:K:209:LEU:O	2.21	0.41
1:L:27:GLY:HA3	1:L:42:LEU:HD23	2.01	0.41
1:B:139:TRP:CE3	1:B:161:LEU:HG	2.56	0.40
1:D:59:THR:O	1:D:62:GYS:C2	2.70	0.40
1:D:93:MET:HG2	1:D:173:PHE:CD1	2.56	0.40
1:I:10:ILE:O	1:I:28:VAL:HA	2.20	0.40
1:K:76:VAL:HB	1:K:186:PRO:HB3	2.02	0.40
1:D:198:LYS:HD3	1:D:198:LYS:HA	1.98	0.40
1:H:13:ARG:NH2	1:H:26:GLU:OE1	2.54	0.40
1:J:43:LYS:HA	1:J:205:SER:O	2.22	0.40
1:F:183:VAL:HG12	1:F:184:GLN:N	2.36	0.40
1:A:123:VAL:HB	1:L:90:GLU:HB3	2.03	0.40
1:I:95:TYR:CD1	1:I:171:CYS:HB2	2.56	0.40
1:A:210:HIS:HE1	3:A:401:HOH:O	2.05	0.40
1:F:30:LEU:N	1:F:30:LEU:CD2	2.84	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	210/215 (98%)	205 (98%)	5 (2%)	0	100	100
1	B	208/215 (97%)	204 (98%)	4 (2%)	0	100	100
1	C	210/215 (98%)	208 (99%)	2 (1%)	0	100	100
1	D	210/215 (98%)	207 (99%)	3 (1%)	0	100	100
1	E	210/215 (98%)	207 (99%)	3 (1%)	0	100	100
1	F	209/215 (97%)	206 (99%)	3 (1%)	0	100	100
1	G	209/215 (97%)	202 (97%)	7 (3%)	0	100	100
1	H	209/215 (97%)	206 (99%)	3 (1%)	0	100	100
1	I	208/215 (97%)	202 (97%)	6 (3%)	0	100	100
1	J	209/215 (97%)	207 (99%)	1 (0%)	1 (0%)	29	28
1	K	208/215 (97%)	205 (99%)	3 (1%)	0	100	100
1	L	209/215 (97%)	203 (97%)	6 (3%)	0	100	100
All	All	2509/2580 (97%)	2462 (98%)	46 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	J	201	ASP

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	185/185 (100%)	174 (94%)	11 (6%)	19	21
1	B	183/185 (99%)	178 (97%)	5 (3%)	44	54
1	C	185/185 (100%)	180 (97%)	5 (3%)	44	54
1	D	185/185 (100%)	181 (98%)	4 (2%)	52	62
1	E	185/185 (100%)	177 (96%)	8 (4%)	29	34
1	F	184/185 (100%)	175 (95%)	9 (5%)	25	28
1	G	184/185 (100%)	182 (99%)	2 (1%)	73	83
1	H	184/185 (100%)	177 (96%)	7 (4%)	33	39
1	I	183/185 (99%)	177 (97%)	6 (3%)	38	46
1	J	184/185 (100%)	174 (95%)	10 (5%)	22	24
1	K	183/185 (99%)	178 (97%)	5 (3%)	44	54
1	L	184/185 (100%)	175 (95%)	9 (5%)	25	28
All	All	2209/2220 (100%)	2128 (96%)	81 (4%)	34	40

All (81) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	VAL
1	A	8	MET
1	A	11	LYS
1	A	30	LEU
1	A	66	ARG
1	A	67	VAL
1	A	70	LYS
1	A	73	GLU
1	A	138	LYS
1	A	202	LYS
1	A	218	GLU
1	B	5	LYS
1	B	30	LEU
1	B	66	ARG
1	B	67	VAL
1	B	92	SER
1	C	3	VAL
1	C	11	LYS
1	C	30	LEU
1	C	152	VAL
1	C	181	LYS
1	D	11	LYS

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Mol	Chain	Res	Type
1	D	66	ARG
1	D	90	GLU
1	D	164	GLU
1	E	2	SER
1	E	21	HIS
1	E	66	ARG
1	E	67	VAL
1	E	92	SER
1	E	134	LYS
1	E	145	LYS
1	E	218	GLU
1	F	3	VAL
1	F	8	MET
1	F	11	LYS
1	F	30	LEU
1	F	66	ARG
1	F	67	VAL
1	F	92	SER
1	F	152	VAL
1	F	180	LYS
1	G	66	ARG
1	G	90	GLU
1	H	21	HIS
1	H	45	LYS
1	H	66	ARG
1	H	73	GLU
1	H	162	SER
1	H	202	LYS
1	H	214	GLU
1	I	30	LEU
1	I	43	LYS
1	I	73	GLU
1	I	138	LYS
1	I	202	LYS
1	I	210	HIS
1	J	3	VAL
1	J	5	LYS
1	J	30	LEU
1	J	66	ARG
1	J	73	GLU
1	J	145	LYS
1	J	162	SER

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Mol	Chain	Res	Type
1	J	181	LYS
1	J	205	SER
1	J	217	SER
1	K	3	VAL
1	K	66	ARG
1	K	73	GLU
1	K	181	LYS
1	K	198	LYS
1	L	11	LYS
1	L	35	GLU
1	L	66	ARG
1	L	73	GLU
1	L	115	ILE
1	L	180	LYS
1	L	181	LYS
1	L	196	GLU
1	L	218	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	210	HIS
1	B	38	GLN
1	D	208	ASN
1	D	210	HIS
1	F	158	ASN
1	G	38	GLN
1	G	210	HIS
1	H	210	HIS
1	J	38	GLN
1	L	81	GLN
1	L	194	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

12 non-standard protein/DNA/RNA residues 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
1	GYS	J	62	1	22,22,23	3.40	6 (27%)	27,30,32	3.86	8 (29%)
1	GYS	B	62	1	22,22,23	3.22	6 (27%)	27,30,32	3.90	10 (37%)
1	GYS	H	62	1	22,22,23	4.38	6 (27%)	27,30,32	5.50	7 (25%)
1	GYS	G	62	1	22,22,23	3.24	6 (27%)	27,30,32	4.01	11 (40%)
1	GYS	L	62	1	22,22,23	2.59	6 (27%)	27,30,32	3.60	13 (48%)
1	GYS	C	62	1	22,22,23	3.18	5 (22%)	27,30,32	3.43	11 (40%)
1	GYS	I	62	1	22,22,23	3.22	5 (22%)	27,30,32	2.70	6 (22%)
1	GYS	F	62	1	22,22,23	3.04	6 (27%)	27,30,32	3.13	7 (25%)
1	GYS	D	62	1	22,22,23	3.57	6 (27%)	27,30,32	3.30	5 (18%)
1	GYS	K	62	1	22,22,23	3.60	6 (27%)	27,30,32	3.94	5 (18%)
1	GYS	E	62	1	22,22,23	4.03	5 (22%)	27,30,32	4.01	8 (29%)
1	GYS	A	62	1	22,22,23	3.88	6 (27%)	27,30,32	3.69	10 (37%)

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
1	GYS	J	62	1	-	2/9/29/30	0/2/2/2
1	GYS	B	62	1	-	1/9/29/30	0/2/2/2
1	GYS	H	62	1	-	1/9/29/30	0/2/2/2
1	GYS	G	62	1	-	3/9/29/30	0/2/2/2
1	GYS	L	62	1	-	3/9/29/30	0/2/2/2
1	GYS	C	62	1	-	2/9/29/30	0/2/2/2
1	GYS	I	62	1	-	3/9/29/30	0/2/2/2
1	GYS	F	62	1	-	3/9/29/30	0/2/2/2
1	GYS	D	62	1	-	3/9/29/30	0/2/2/2
1	GYS	K	62	1	-	1/9/29/30	0/2/2/2
1	GYS	E	62	1	-	2/9/29/30	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	GYS	A	62	1	-	3/9/29/30	0/2/2/2

All (69) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	62	GYS	CB2-CA2	18.37	1.50	1.35
1	E	62	GYS	CB2-CA2	17.32	1.49	1.35
1	A	62	GYS	CB2-CA2	16.32	1.48	1.35
1	K	62	GYS	CB2-CA2	15.08	1.47	1.35
1	D	62	GYS	CB2-CA2	15.00	1.47	1.35
1	J	62	GYS	CB2-CA2	14.20	1.47	1.35
1	B	62	GYS	CB2-CA2	12.96	1.46	1.35
1	I	62	GYS	CB2-CA2	12.59	1.45	1.35
1	G	62	GYS	CB2-CA2	12.54	1.45	1.35
1	C	62	GYS	CB2-CA2	12.50	1.45	1.35
1	F	62	GYS	CB2-CA2	11.53	1.44	1.35
1	L	62	GYS	CB2-CA2	8.53	1.42	1.35
1	L	62	GYS	C1-N2	5.75	1.40	1.32
1	H	62	GYS	C1-N2	5.64	1.40	1.32
1	C	62	GYS	CA2-C2	-4.82	1.43	1.48
1	H	62	GYS	CA2-C2	-4.76	1.43	1.48
1	E	62	GYS	C1-N2	4.55	1.38	1.32
1	F	62	GYS	C2-N3	-4.47	1.29	1.39
1	G	62	GYS	CA2-C2	-4.41	1.44	1.48
1	A	62	GYS	C1-N2	4.37	1.38	1.32
1	I	62	GYS	C1-N2	4.36	1.38	1.32
1	J	62	GYS	C1-N2	4.34	1.38	1.32
1	G	62	GYS	O2-C2	4.30	1.32	1.23
1	K	62	GYS	C1-N2	4.28	1.38	1.32
1	I	62	GYS	CA2-C2	-4.07	1.44	1.48
1	F	62	GYS	C1-N2	3.92	1.38	1.32
1	A	62	GYS	CA2-C2	-3.81	1.44	1.48
1	D	62	GYS	C1-N2	3.66	1.37	1.32
1	C	62	GYS	O2-C2	3.60	1.30	1.23
1	I	62	GYS	O2-C2	3.58	1.30	1.23
1	A	62	GYS	C2-N3	-3.47	1.31	1.39
1	K	62	GYS	C2-N3	-3.44	1.31	1.39
1	B	62	GYS	C1-N2	3.44	1.37	1.32
1	E	62	GYS	C2-N3	-3.38	1.31	1.39
1	L	62	GYS	O2-C2	3.37	1.30	1.23
1	B	62	GYS	O2-C2	3.37	1.30	1.23
1	J	62	GYS	C2-N3	-3.36	1.32	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	62	GYS	O2-C2	3.30	1.30	1.23
1	D	62	GYS	C2-N3	-3.22	1.32	1.39
1	I	62	GYS	C2-N3	-3.12	1.32	1.39
1	E	62	GYS	O2-C2	3.10	1.29	1.23
1	L	62	GYS	CA2-C2	-3.09	1.45	1.48
1	C	62	GYS	C1-N2	3.09	1.36	1.32
1	D	62	GYS	CA2-C2	-3.05	1.45	1.48
1	A	62	GYS	O2-C2	2.98	1.29	1.23
1	H	62	GYS	C2-N3	-2.97	1.32	1.39
1	G	62	GYS	C1-N2	2.97	1.36	1.32
1	K	62	GYS	CA2-C2	-2.93	1.45	1.48
1	B	62	GYS	CA2-C2	-2.93	1.45	1.48
1	G	62	GYS	C2-N3	-2.86	1.33	1.39
1	K	62	GYS	O2-C2	2.84	1.29	1.23
1	H	62	GYS	CA1-C1	-2.84	1.45	1.51
1	F	62	GYS	O2-C2	2.76	1.28	1.23
1	C	62	GYS	C2-N3	-2.72	1.33	1.39
1	B	62	GYS	CE1-CD1	2.66	1.43	1.38
1	F	62	GYS	CA2-C2	-2.60	1.46	1.48
1	J	62	GYS	CA2-C2	-2.59	1.46	1.48
1	J	62	GYS	O2-C2	2.52	1.28	1.23
1	B	62	GYS	C2-N3	-2.50	1.34	1.39
1	F	62	GYS	CA2-N2	-2.42	1.33	1.38
1	K	62	GYS	CA2-N2	-2.41	1.33	1.38
1	G	62	GYS	CA2-N2	-2.37	1.33	1.38
1	A	62	GYS	CA2-N2	-2.26	1.33	1.38
1	E	62	GYS	CA2-N2	-2.26	1.33	1.38
1	D	62	GYS	CA2-N2	-2.25	1.33	1.38
1	J	62	GYS	CA2-N2	-2.24	1.33	1.38
1	H	62	GYS	O2-C2	2.24	1.27	1.23
1	L	62	GYS	CA2-N2	-2.13	1.34	1.38
1	L	62	GYS	C2-N3	-2.08	1.35	1.39

All (101) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	62	GYS	O2-C2-CA2	-18.90	120.35	130.96
1	H	62	GYS	CA2-C2-N3	17.89	111.83	103.37
1	G	62	GYS	O2-C2-CA2	-15.89	122.04	130.96
1	E	62	GYS	O2-C2-CA2	-14.33	122.91	130.96
1	C	62	GYS	O2-C2-CA2	-12.90	123.71	130.96
1	K	62	GYS	CA2-C2-N3	12.88	109.46	103.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	J	62	GYS	CA2-C2-N3	12.86	109.45	103.37
1	K	62	GYS	O2-C2-CA2	-12.74	123.81	130.96
1	J	62	GYS	O2-C2-CA2	-12.44	123.97	130.96
1	A	62	GYS	O2-C2-CA2	-12.24	124.09	130.96
1	B	62	GYS	O2-C2-CA2	-11.87	124.29	130.96
1	E	62	GYS	CA2-C2-N3	11.57	108.84	103.37
1	A	62	GYS	CA2-C2-N3	11.36	108.74	103.37
1	D	62	GYS	O2-C2-CA2	-11.24	124.65	130.96
1	B	62	GYS	CA2-C2-N3	10.63	108.40	103.37
1	F	62	GYS	O2-C2-CA2	-10.57	125.03	130.96
1	L	62	GYS	O2-C2-CA2	-10.56	125.03	130.96
1	D	62	GYS	CA2-C2-N3	10.46	108.32	103.37
1	L	62	GYS	CA2-C2-N3	10.12	108.16	103.37
1	G	62	GYS	CA2-C2-N3	9.42	107.83	103.37
1	F	62	GYS	CA2-C2-N3	9.24	107.74	103.37
1	I	62	GYS	O2-C2-CA2	-8.60	126.13	130.96
1	H	62	GYS	C2-N3-C1	-7.82	104.01	107.97
1	I	62	GYS	CA2-C2-N3	7.51	106.92	103.37
1	C	62	GYS	CA2-C2-N3	7.35	106.85	103.37
1	B	62	GYS	C2-N3-C1	-6.68	104.58	107.97
1	L	62	GYS	C2-N3-C1	-6.33	104.76	107.97
1	K	62	GYS	C2-N3-C1	-5.89	104.98	107.97
1	E	62	GYS	C2-N3-C1	-5.79	105.04	107.97
1	B	62	GYS	O3-C3-CA3	-5.47	109.89	126.39
1	J	62	GYS	C2-N3-C1	-5.11	105.38	107.97
1	I	62	GYS	O3-C3-CA3	-4.89	111.61	126.39
1	G	62	GYS	O3-C3-CA3	-4.72	112.14	126.39
1	A	62	GYS	C2-N3-C1	-4.64	105.62	107.97
1	K	62	GYS	O3-C3-CA3	-4.63	112.41	126.39
1	B	62	GYS	CA1-C1-N3	-4.50	118.99	124.85
1	L	62	GYS	O3-C3-CA3	-4.43	113.03	126.39
1	C	62	GYS	O3-C3-CA3	-4.30	113.40	126.39
1	D	62	GYS	O3-C3-CA3	-4.22	113.64	126.39
1	E	62	GYS	O3-C3-CA3	-4.12	113.94	126.39
1	B	62	GYS	N3-C1-N2	4.04	114.25	111.45
1	H	62	GYS	C2-CA2-N2	-4.01	106.12	108.93
1	H	62	GYS	O3-C3-CA3	-3.75	115.07	126.39
1	A	62	GYS	O3-C3-CA3	-3.68	115.27	126.39
1	H	62	GYS	CA1-C1-N3	-3.66	120.08	124.85
1	F	62	GYS	O3-C3-CA3	-3.55	115.68	126.39
1	F	62	GYS	CD1-CE1-CZ	3.37	123.58	119.88
1	C	62	GYS	CD1-CE1-CZ	3.28	123.47	119.88

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	62	GYS	C2-N3-C1	-3.25	106.32	107.97
1	C	62	GYS	CB1-CA1-C1	3.20	117.75	110.19
1	J	62	GYS	O3-C3-CA3	-3.08	117.08	126.39
1	G	62	GYS	C2-N3-C1	-2.98	106.46	107.97
1	G	62	GYS	CD2-CE2-CZ	2.97	123.14	119.88
1	F	62	GYS	CE1-CD1-CG2	-2.94	117.42	121.25
1	G	62	GYS	O2-C2-N3	2.91	130.14	124.35
1	C	62	GYS	N3-C1-N2	-2.89	109.45	111.45
1	L	62	GYS	CD2-CG2-CB2	-2.88	111.43	121.22
1	A	62	GYS	CE1-CD1-CG2	-2.84	117.55	121.25
1	J	62	GYS	CA1-C1-N3	-2.83	121.16	124.85
1	L	62	GYS	CE1-CD1-CG2	-2.83	117.56	121.25
1	E	62	GYS	CA1-C1-N3	-2.80	121.20	124.85
1	G	62	GYS	CB1-CA1-C1	2.79	116.78	110.19
1	L	62	GYS	CA2-N2-C1	-2.73	103.76	105.77
1	G	62	GYS	CD2-CG2-CB2	-2.72	111.94	121.22
1	L	62	GYS	CA3-N3-C1	2.71	130.41	127.16
1	C	62	GYS	CE1-CD1-CG2	-2.64	117.80	121.25
1	A	62	GYS	CB1-CA1-C1	2.60	116.33	110.19
1	I	62	GYS	CG2-CB2-CA2	-2.59	126.77	129.94
1	J	62	GYS	C2-CA2-N2	-2.58	107.13	108.93
1	C	62	GYS	O2-C2-N3	2.56	129.43	124.35
1	E	62	GYS	CB1-CA1-C1	2.55	116.23	110.19
1	F	62	GYS	CD2-CG2-CB2	-2.49	112.75	121.22
1	L	62	GYS	CA1-C1-N3	-2.42	121.70	124.85
1	L	62	GYS	OG1-CB1-CA1	-2.41	104.48	110.85
1	A	62	GYS	CD2-CG2-CD1	2.38	121.16	117.64
1	H	62	GYS	CG2-CB2-CA2	-2.37	127.04	129.94
1	L	62	GYS	CD2-CG2-CD1	2.33	121.09	117.64
1	L	62	GYS	CB1-CA1-C1	2.33	115.70	110.19
1	E	62	GYS	N3-C1-N2	2.29	113.04	111.45
1	J	62	GYS	CB1-CA1-C1	2.26	115.53	110.19
1	B	62	GYS	CD2-CE2-CZ	2.25	122.35	119.88
1	D	62	GYS	CD2-CG2-CB2	-2.25	113.57	121.22
1	C	62	GYS	CD2-CG2-CB2	-2.25	113.57	121.22
1	G	62	GYS	CD1-CE1-CZ	2.22	122.31	119.88
1	I	62	GYS	CD2-CG2-CB2	-2.21	113.68	121.22
1	A	62	GYS	CD2-CG2-CB2	-2.20	113.74	121.22
1	E	62	GYS	CA3-N3-C2	2.17	128.78	123.80
1	L	62	GYS	N3-C1-N2	2.16	112.95	111.45
1	G	62	GYS	CE2-CZ-CE1	-2.16	116.14	119.77
1	C	62	GYS	OG1-CB1-CA1	2.15	116.55	110.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	62	GYS	CB1-CA1-C1	2.15	115.28	110.19
1	G	62	GYS	CD1-CG2-CB2	2.15	128.55	121.22
1	A	62	GYS	CA3-N3-C1	2.15	129.75	127.16
1	B	62	GYS	CD1-CG2-CB2	2.14	128.50	121.22
1	J	62	GYS	CD2-CG2-CB2	-2.13	113.97	121.22
1	F	62	GYS	C2-N3-C1	-2.11	106.90	107.97
1	B	62	GYS	CD2-CG2-CB2	-2.11	114.05	121.22
1	C	62	GYS	CA3-N3-C1	-2.10	124.64	127.16
1	B	62	GYS	CA2-N2-C1	-2.10	104.22	105.77
1	K	62	GYS	CA1-C1-N3	-2.09	122.13	124.85
1	A	62	GYS	CA1-C1-N3	-2.03	122.21	124.85

There are no chirality outliers.

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	62	GYS	C3-CA3-N3-C2
1	F	62	GYS	C3-CA3-N3-C2
1	G	62	GYS	C3-CA3-N3-C2
1	H	62	GYS	C3-CA3-N3-C2
1	I	62	GYS	N1-CA1-CB1-OG1
1	I	62	GYS	C3-CA3-N3-C2
1	J	62	GYS	C3-CA3-N3-C2
1	K	62	GYS	C3-CA3-N3-C2
1	L	62	GYS	N1-CA1-CB1-OG1
1	A	62	GYS	N1-CA1-CB1-OG1
1	A	62	GYS	C3-CA3-N3-C2
1	C	62	GYS	C3-CA3-N3-C2
1	D	62	GYS	N1-CA1-CB1-OG1
1	D	62	GYS	C3-CA3-N3-C2
1	E	62	GYS	C3-CA3-N3-C2
1	A	62	GYS	C1-CA1-CB1-OG1
1	L	62	GYS	C3-CA3-N3-C2
1	L	62	GYS	C1-CA1-CB1-OG1
1	D	62	GYS	C1-CA1-CB1-OG1
1	F	62	GYS	N1-CA1-CB1-OG1
1	G	62	GYS	N1-CA1-CB1-OG1
1	J	62	GYS	C3-CA3-N3-C1
1	F	62	GYS	C3-CA3-N3-C1
1	G	62	GYS	C3-CA3-N3-C1
1	I	62	GYS	C3-CA3-N3-C1
1	C	62	GYS	C3-CA3-N3-C1

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Mol	Chain	Res	Type	Atoms
1	E	62	GYS	C3-CA3-N3-C1

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	62	GYS	1	0

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 24 ligands modelled in this entry, 24 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

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

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	214/215 (99%)	-0.14	4 (1%) 66 67	19, 29, 50, 69	0
1	B	212/215 (98%)	-0.24	1 (0%) 91 91	18, 27, 45, 61	0
1	C	214/215 (99%)	-0.21	2 (0%) 84 84	16, 25, 43, 75	0
1	D	214/215 (99%)	-0.27	1 (0%) 91 91	17, 29, 46, 60	0
1	E	214/215 (99%)	-0.21	4 (1%) 66 67	17, 26, 41, 59	0
1	F	213/215 (99%)	-0.22	4 (1%) 66 67	16, 25, 44, 69	0
1	G	213/215 (99%)	-0.23	2 (0%) 84 84	18, 27, 45, 64	0
1	H	213/215 (99%)	-0.34	1 (0%) 91 91	17, 25, 38, 60	0
1	I	212/215 (98%)	-0.12	3 (1%) 75 75	21, 34, 54, 78	0
1	J	213/215 (99%)	0.11	13 (6%) 21 22	21, 36, 58, 76	0
1	K	212/215 (98%)	0.01	9 (4%) 36 37	21, 35, 55, 74	0
1	L	213/215 (99%)	-0.09	5 (2%) 60 61	20, 32, 51, 77	0
All	All	2557/2580 (99%)	-0.16	49 (1%) 66 67	16, 29, 50, 78	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	J	182	VAL	3.6
1	J	73	GLU	3.4
1	L	3	VAL	3.4
1	B	3	VAL	3.2
1	C	60	VAL	3.2
1	D	2	SER	3.2
1	F	60	VAL	3.0
1	K	73	GLU	2.9
1	E	60	VAL	2.9
1	C	2	SER	2.9
1	L	60	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	K	60	VAL	2.8
1	I	3	VAL	2.8
1	K	112	ASP	2.7
1	I	74	ASN	2.6
1	A	60	VAL	2.6
1	L	73	GLU	2.6
1	J	165	GLY	2.5
1	L	57	LEU	2.5
1	F	61	PHE	2.5
1	K	3	VAL	2.5
1	K	165	GLY	2.5
1	G	2	SER	2.5
1	F	59	THR	2.4
1	A	73	GLU	2.4
1	H	73	GLU	2.4
1	J	59	THR	2.4
1	J	202	LYS	2.4
1	F	57	LEU	2.4
1	J	2	SER	2.3
1	E	59	THR	2.3
1	J	195	ILE	2.3
1	J	3	VAL	2.2
1	K	113	CYS	2.2
1	A	202	LYS	2.2
1	J	66	ARG	2.2
1	J	30	LEU	2.2
1	G	60	VAL	2.2
1	E	61	PHE	2.2
1	K	182	VAL	2.2
1	K	110	ASP	2.1
1	J	34	PHE	2.1
1	I	103	ALA	2.1
1	A	59	THR	2.1
1	L	59	THR	2.1
1	E	73	GLU	2.1
1	J	60	VAL	2.0
1	J	28	VAL	2.0
1	K	41	ASP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	GYS	F	62	21/22	0.94	0.23	17,21,23,29	0
1	GYS	L	62	21/22	0.94	0.16	20,23,27,28	0
1	GYS	J	62	21/22	0.96	0.27	29,36,39,47	0
1	GYS	I	62	21/22	0.96	0.12	22,25,28,33	0
1	GYS	C	62	21/22	0.96	0.18	16,21,25,26	0
1	GYS	K	62	21/22	0.97	0.20	24,28,36,41	0
1	GYS	B	62	21/22	0.97	0.16	17,19,23,31	0
1	GYS	A	62	21/22	0.97	0.19	20,23,27,31	0
1	GYS	H	62	21/22	0.97	0.15	16,20,24,26	0
1	GYS	G	62	21/22	0.98	0.15	18,21,23,32	0
1	GYS	D	62	21/22	0.98	0.13	22,27,31,33	0
1	GYS	E	62	21/22	0.98	0.23	18,21,23,24	0

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
2	CO	J	302	1/1	0.90	0.05	69,69,69,69	0
2	CO	K	301	1/1	0.93	0.05	48,48,48,48	0
2	CO	L	301	1/1	0.94	0.03	49,49,49,49	0
2	CO	L	302	1/1	0.94	0.03	66,66,66,66	0
2	CO	K	302	1/1	0.95	0.03	60,60,60,60	0
2	CO	A	302	1/1	0.96	0.03	52,52,52,52	0
2	CO	J	301	1/1	0.97	0.04	54,54,54,54	0
2	CO	F	302	1/1	0.97	0.04	53,53,53,53	0
2	CO	B	302	1/1	0.98	0.03	47,47,47,47	0
2	CO	I	301	1/1	0.98	0.03	56,56,56,56	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	CO	D	301	1/1	0.98	0.04	41,41,41,41	0
2	CO	D	302	1/1	0.98	0.04	48,48,48,48	0
2	CO	H	302	1/1	0.99	0.02	47,47,47,47	0
2	CO	A	301	1/1	0.99	0.03	39,39,39,39	0
2	CO	I	302	1/1	0.99	0.04	61,61,61,61	0
2	CO	C	301	1/1	0.99	0.03	34,34,34,34	0
2	CO	E	301	1/1	0.99	0.02	32,32,32,32	0
2	CO	E	302	1/1	0.99	0.02	45,45,45,45	0
2	CO	F	301	1/1	0.99	0.04	41,41,41,41	0
2	CO	C	302	1/1	0.99	0.02	49,49,49,49	0
2	CO	G	302	1/1	0.99	0.02	45,45,45,45	0
2	CO	G	301	1/1	1.00	0.02	37,37,37,37	0
2	CO	B	301	1/1	1.00	0.02	37,37,37,37	0
2	CO	H	301	1/1	1.00	0.03	34,34,34,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.