



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:46 PM GMT

PDB ID : 1YMC
Title : THREE-DIMENSIONAL STRUCTURE OF CYANOMET-SULFMYOGLOBIN C
Authors : Evans, S.V.; Brayer, G.D.
Deposited on : 1993-09-27
Resolution : 2.00 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

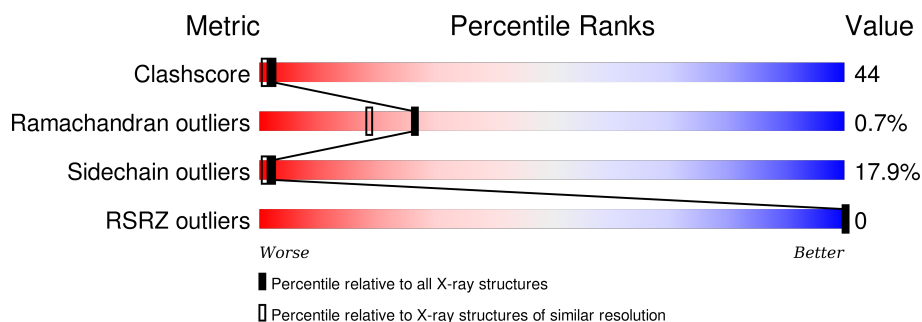
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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))
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	153	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	A	155	-	-	X	X
4	CLN	A	154	X	-	-	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 1423 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 CYANOMET-SULFMYOGLOBIN.

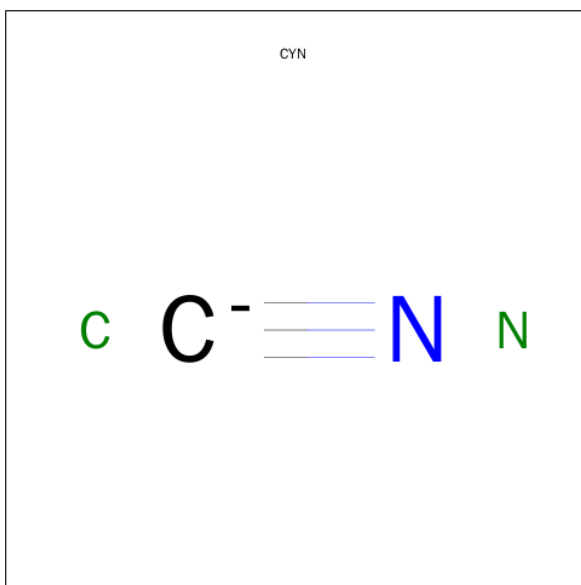
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	153	Total	C	N	O	S	0	0	0
			1199	769	210	218	2			

- Molecule 2 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



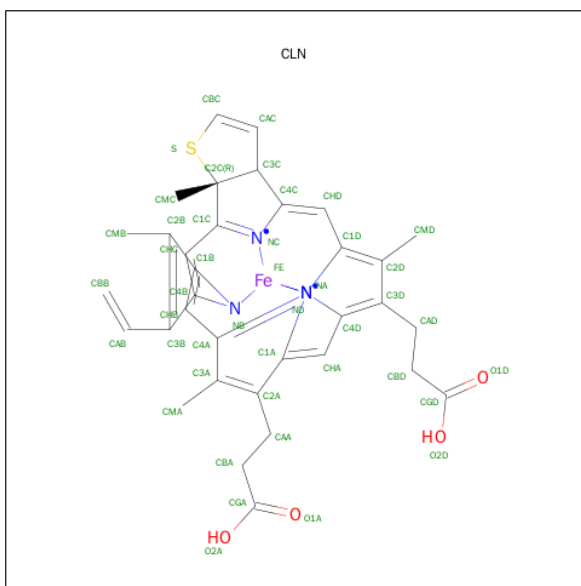
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is CYANIDE ION (three-letter code: CYN) (formula: CN).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			2	1	1		

- Molecule 4 is SULFUR SUBSTITUTED PROTOPORPHYRIN IX (three-letter code: CLN) (formula: $C_{34}H_{32}FeN_4O_4S$).

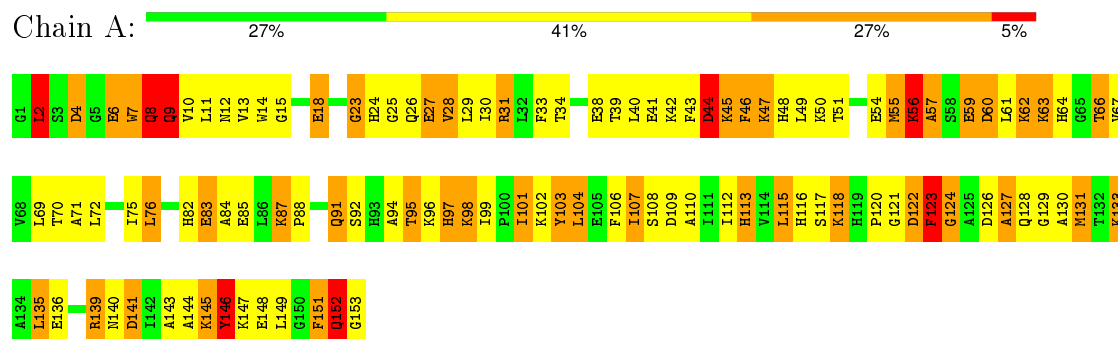


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	173	Total 173	O 173	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: CYANOMET-SULFMYOGLOBIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	64.50 Å 28.91 Å 35.66 Å 90.00° 107.00° 90.00°	Depositor
Resolution (Å)	6.00 – 2.00 6.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (6.00-2.00) 63.9 (6.00-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	PROLSQ	Depositor
R, R_{free}	0.129 , (Not available) 0.129 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	9.8	Xtriage
Anisotropy	0.146	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.18 , 173.7	EDS
Estimated twinning fraction	0.096 for -h,-l,-k,l	Xtriage
L-test for twinning ¹	$\langle L \rangle = 0.37$, $\langle L^2 \rangle = 0.20$	Xtriage
Outliers	0 of 5369 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	1423	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.03% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: CLN, CYN, SO4

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	1.80	15/1226 (1.2%)	2.67	90/1645 (5.5%)

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	144	ALA	C-O	8.16	1.38	1.23
1	A	18	GLU	CD-OE2	-7.58	1.17	1.25
1	A	121	GLY	C-O	6.76	1.34	1.23
1	A	124	GLY	N-CA	-6.75	1.35	1.46
1	A	23	GLY	N-CA	6.42	1.55	1.46
1	A	133	LYS	C-O	6.11	1.34	1.23
1	A	31	ARG	CD-NE	-5.83	1.36	1.46
1	A	27	GLU	CD-OE1	5.77	1.31	1.25
1	A	11	LEU	C-O	5.56	1.33	1.23
1	A	62	LYS	CA-CB	-5.43	1.42	1.53
1	A	15	GLY	N-CA	-5.34	1.38	1.46
1	A	60	ASP	CB-CG	-5.31	1.40	1.51
1	A	91	GLN	N-CA	5.30	1.56	1.46
1	A	83	GLU	CD-OE1	-5.21	1.20	1.25
1	A	10	VAL	CB-CG1	5.17	1.63	1.52

All (90) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	62	LYS	CA-CB-CG	18.00	153.00	113.40
1	A	139	ARG	CD-NE-CZ	16.98	147.38	123.60
1	A	139	ARG	NE-CZ-NH1	15.46	128.03	120.30
1	A	139	ARG	NH1-CZ-NH2	-12.22	105.96	119.40
1	A	31	ARG	CD-NE-CZ	12.00	140.41	123.60
1	A	60	ASP	CA-CB-CG	11.83	139.42	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	139	ARG	NE-CZ-NH2	11.42	126.01	120.30
1	A	126	ASP	CB-CG-OD1	11.31	128.48	118.30
1	A	106	PHE	CB-CG-CD1	11.12	128.58	120.80
1	A	2	LEU	CA-CB-CG	11.00	140.59	115.30
1	A	85	GLU	CG-CD-OE1	10.72	139.74	118.30
1	A	55	MET	CG-SD-CE	10.61	117.17	100.20
1	A	66	THR	CA-CB-CG2	10.13	126.59	112.40
1	A	103	TYR	CB-CG-CD2	-9.87	115.08	121.00
1	A	8	GLN	CA-CB-CG	9.84	135.04	113.40
1	A	103	TYR	CA-CB-CG	9.33	131.13	113.40
1	A	76	LEU	CA-CB-CG	9.04	136.10	115.30
1	A	31	ARG	CG-CD-NE	8.71	130.09	111.80
1	A	141	ASP	CB-CG-OD1	8.43	125.89	118.30
1	A	33	PHE	CB-CG-CD2	-8.30	114.99	120.80
1	A	135	LEU	CA-CB-CG	8.26	134.29	115.30
1	A	148	GLU	O-C-N	7.92	135.38	122.70
1	A	144	ALA	CB-CA-C	7.87	121.91	110.10
1	A	85	GLU	CG-CD-OE2	-7.79	102.73	118.30
1	A	27	GLU	OE1-CD-OE2	-7.75	114.00	123.30
1	A	9	GLN	O-C-N	-7.59	110.56	122.70
1	A	83	GLU	OE1-CD-OE2	-7.53	114.27	123.30
1	A	104	LEU	CA-CB-CG	7.44	132.40	115.30
1	A	127	ALA	N-CA-CB	7.41	120.48	110.10
1	A	152	GLN	CB-CG-CD	7.39	130.81	111.60
1	A	27	GLU	CG-CD-OE2	7.18	132.66	118.30
1	A	33	PHE	CB-CG-CD1	7.08	125.75	120.80
1	A	67	VAL	CB-CA-C	7.05	124.79	111.40
1	A	146	TYR	CZ-CE2-CD2	-6.96	113.54	119.80
1	A	135	LEU	CB-CA-C	6.81	123.14	110.20
1	A	146	TYR	CG-CD2-CE2	6.81	126.75	121.30
1	A	57	ALA	CB-CA-C	6.72	120.18	110.10
1	A	123	PHE	CB-CA-C	6.61	123.62	110.40
1	A	126	ASP	CA-CB-CG	6.58	127.88	113.40
1	A	38	GLU	O-C-N	6.56	133.19	122.70
1	A	106	PHE	CG-CD1-CE1	6.44	127.88	120.80
1	A	31	ARG	NE-CZ-NH1	6.39	123.50	120.30
1	A	4	ASP	N-CA-CB	6.34	122.01	110.60
1	A	7	TRP	CA-C-O	6.30	133.34	120.10
1	A	148	GLU	CA-C-O	-6.26	106.96	120.10
1	A	64	HIS	CA-CB-CG	6.24	124.20	113.60
1	A	113	HIS	CA-C-O	6.15	133.02	120.10
1	A	34	THR	C-N-CA	6.14	135.20	122.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	62	LYS	CB-CA-C	6.06	122.52	110.40
1	A	91	GLN	CB-CA-C	6.04	122.47	110.40
1	A	102	LYS	CA-CB-CG	6.02	126.65	113.40
1	A	148	GLU	N-CA-CB	6.02	121.43	110.60
1	A	43	PHE	CB-CG-CD2	5.97	124.98	120.80
1	A	113	HIS	N-CA-CB	-5.97	99.86	110.60
1	A	113	HIS	CB-CA-C	5.97	122.33	110.40
1	A	84	ALA	N-CA-CB	-5.93	101.80	110.10
1	A	64	HIS	N-CA-CB	5.81	121.06	110.60
1	A	106	PHE	CB-CG-CD2	-5.79	116.74	120.80
1	A	44	ASP	CB-CG-OD1	-5.78	113.09	118.30
1	A	122	ASP	N-CA-CB	5.74	120.93	110.60
1	A	4	ASP	CB-CG-OD1	5.73	123.46	118.30
1	A	54	GLU	N-CA-CB	5.67	120.81	110.60
1	A	71	ALA	N-CA-CB	5.61	117.95	110.10
1	A	143	ALA	N-CA-CB	5.57	117.90	110.10
1	A	63	LYS	CA-CB-CG	5.56	125.63	113.40
1	A	85	GLU	O-C-N	-5.48	113.93	122.70
1	A	118	LYS	CD-CE-NZ	-5.42	99.22	111.70
1	A	126	ASP	CB-CG-OD2	-5.40	113.44	118.30
1	A	91	GLN	CA-CB-CG	5.38	125.24	113.40
1	A	11	LEU	CB-CA-C	5.38	120.42	110.20
1	A	147	LYS	CB-CG-CD	5.35	125.50	111.60
1	A	131	MET	CB-CA-C	5.31	121.02	110.40
1	A	103	TYR	CB-CA-C	5.30	121.00	110.40
1	A	97	HIS	N-CA-CB	5.28	120.11	110.60
1	A	39	THR	CA-CB-CG2	5.26	119.77	112.40
1	A	70	THR	N-CA-CB	5.24	120.26	110.30
1	A	133	LYS	CA-CB-CG	-5.21	101.94	113.40
1	A	151	PHE	C-N-CA	5.18	134.66	121.70
1	A	41	GLU	CG-CD-OE1	-5.16	107.98	118.30
1	A	139	ARG	CB-CG-CD	5.16	125.02	111.60
1	A	6	GLU	CA-CB-CG	5.16	124.75	113.40
1	A	28	VAL	CG1-CB-CG2	-5.15	102.66	110.90
1	A	57	ALA	N-CA-CB	-5.13	102.91	110.10
1	A	98	LYS	CG-CD-CE	5.13	127.28	111.90
1	A	69	LEU	CB-CG-CD2	5.11	119.69	111.00
1	A	67	VAL	N-CA-CB	-5.09	100.30	111.50
1	A	56	LYS	N-CA-CB	5.08	119.74	110.60
1	A	139	ARG	CG-CD-NE	5.06	122.43	111.80
1	A	140	ASN	N-CA-CB	-5.02	101.56	110.60
1	A	46	PHE	CB-CG-CD2	-5.00	117.30	120.80

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	1199	0	1211	109	5
2	A	5	0	0	2	0
3	A	2	0	0	0	0
4	A	44	0	29	5	0
5	A	173	0	0	52	5
All	All	1423	0	1240	110	8

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 44.

All (110) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:HIS:HA	1:A:118:LYS:HZ1	1.24	1.00
1:A:62:LYS:HB3	5:A:276:HOH:O	1.66	0.95
1:A:104:LEU:HD23	4:A:154:CLN:HMC3	1.51	0.89
1:A:152:GLN:HG2	5:A:228:HOH:O	1.79	0.83
1:A:104:LEU:HD23	4:A:154:CLN:CMC	2.10	0.81
1:A:115:LEU:HD13	5:A:321:HOH:O	1.79	0.81
1:A:87:LYS:HA	5:A:273:HOH:O	1.82	0.80
1:A:4:ASP:O	1:A:8:GLN:HB2	1.82	0.80
1:A:131:MET:HB3	5:A:274:HOH:O	1.81	0.79
1:A:51:THR:O	1:A:55:MET:HG3	1.82	0.79
1:A:30:ILE:CG1	1:A:55:MET:HE3	2.14	0.77
1:A:24:HIS:HA	1:A:118:LYS:NZ	2.00	0.77
1:A:9:GLN:HG2	5:A:182:HOH:O	1.83	0.77
1:A:94:ALA:O	1:A:98:LYS:HD3	1.86	0.76
1:A:30:ILE:HG12	1:A:55:MET:HE3	1.68	0.75
1:A:122:ASP:HB2	5:A:249:HOH:O	1.86	0.74
1:A:124:GLY:HA3	5:A:252:HOH:O	1.88	0.73
1:A:60:ASP:HB2	2:A:155:SO4:O4	1.90	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:GLU:OE2	1:A:118:LYS:HD2	1.90	0.71
1:A:6:GLU:HA	5:A:245:HOH:O	1.92	0.69
1:A:110:ALA:HA	5:A:189:HOH:O	1.93	0.68
1:A:146:TYR:HA	5:A:271:HOH:O	1.95	0.66
1:A:123:PHE:HA	5:A:299:HOH:O	1.95	0.65
1:A:145:LYS:HD2	5:A:222:HOH:O	1.97	0.65
1:A:112:ILE:HD13	5:A:327:HOH:O	1.97	0.64
1:A:45:LYS:HD2	5:A:202:HOH:O	1.98	0.64
1:A:130:ALA:HB2	5:A:245:HOH:O	1.98	0.63
1:A:59:GLU:HA	5:A:276:HOH:O	1.97	0.63
1:A:27:GLU:CD	1:A:118:LYS:HD2	2.20	0.61
1:A:24:HIS:O	1:A:28:VAL:HG23	2.01	0.60
1:A:48:HIS:HD2	5:A:310:HOH:O	1.84	0.60
1:A:44:ASP:HA	1:A:47:LYS:HD2	1.82	0.59
1:A:40:LEU:HD11	1:A:46:PHE:O	2.03	0.59
1:A:112:ILE:HD11	1:A:135:LEU:HD12	1.85	0.58
1:A:30:ILE:CD1	1:A:56:LYS:HG3	2.33	0.58
1:A:95:THR:O	1:A:98:LYS:HE2	2.04	0.58
1:A:128:GLN:HA	5:A:274:HOH:O	2.04	0.57
1:A:87:LYS:N	1:A:88:PRO:HD2	2.20	0.57
1:A:24:HIS:CD2	1:A:118:LYS:NZ	2.73	0.56
1:A:127:ALA:HB3	5:A:188:HOH:O	2.04	0.56
1:A:99:ILE:HB	5:A:248:HOH:O	2.05	0.56
1:A:99:ILE:HG21	4:A:154:CLN:HMC2	1.87	0.56
1:A:91:GLN:HB3	5:A:305:HOH:O	2.05	0.56
1:A:127:ALA:N	5:A:188:HOH:O	2.39	0.55
1:A:113:HIS:HB3	5:A:189:HOH:O	2.06	0.55
1:A:151:PHE:C	1:A:152:GLN:HG3	2.27	0.54
1:A:82:HIS:HE1	1:A:141:ASP:OD2	1.90	0.54
1:A:59:GLU:CA	5:A:276:HOH:O	2.55	0.54
1:A:108:SER:O	1:A:112:ILE:HD12	2.07	0.54
1:A:49:LEU:HD21	5:A:254:HOH:O	2.09	0.53
1:A:112:ILE:CD1	1:A:135:LEU:HD12	2.38	0.53
1:A:101:ILE:O	1:A:104:LEU:HB2	2.10	0.52
1:A:87:LYS:HD3	5:A:305:HOH:O	2.08	0.52
1:A:29:LEU:HB3	1:A:61:LEU:HD11	1.91	0.52
1:A:47:LYS:HD3	5:A:231:HOH:O	2.10	0.51
1:A:83:GLU:HG3	5:A:172:HOH:O	2.09	0.51
1:A:75:ILE:HG23	1:A:82:HIS:CB	2.41	0.51
1:A:110:ALA:HB2	5:A:313:HOH:O	2.11	0.50
1:A:13:VAL:HG21	1:A:127:ALA:HB1	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:97:HIS:HB2	5:A:248:HOH:O	2.10	0.50
1:A:25:GLY:O	1:A:29:LEU:HB2	2.12	0.49
1:A:57:ALA:HB2	5:A:277:HOH:O	2.11	0.49
1:A:63:LYS:HG2	5:A:261:HOH:O	2.11	0.49
1:A:2:LEU:HD11	1:A:133:LYS:HB3	1.93	0.49
1:A:62:LYS:HE2	5:A:307:HOH:O	2.12	0.49
1:A:26:GLN:HG2	1:A:30:ILE:HD12	1.96	0.48
1:A:30:ILE:HD11	1:A:56:LYS:HG3	1.93	0.48
1:A:117:SER:HB2	5:A:234:HOH:O	2.13	0.48
1:A:60:ASP:N	2:A:155:SO4:O4	2.47	0.48
1:A:24:HIS:CA	1:A:118:LYS:HZ1	2.11	0.48
1:A:59:GLU:C	5:A:276:HOH:O	2.52	0.48
1:A:88:PRO:HG3	5:A:318:HOH:O	2.12	0.48
1:A:42:LYS:HG3	1:A:99:ILE:HD11	1.95	0.48
1:A:14:TRP:O	1:A:18:GLU:HG3	2.14	0.48
1:A:124:GLY:C	5:A:188:HOH:O	2.52	0.47
1:A:30:ILE:HD13	1:A:56:LYS:HG3	1.95	0.47
1:A:30:ILE:HG13	1:A:55:MET:HE3	1.95	0.47
1:A:44:ASP:HA	1:A:47:LYS:CD	2.45	0.47
1:A:96:LYS:HD3	5:A:165:HOH:O	2.15	0.47
1:A:116:HIS:NE2	1:A:128:GLN:NE2	2.63	0.46
1:A:141:ASP:O	1:A:145:LYS:HE2	2.15	0.46
1:A:42:LYS:HG3	1:A:99:ILE:CD1	2.46	0.46
1:A:149:LEU:HD12	5:A:271:HOH:O	2.15	0.46
1:A:24:HIS:CD2	1:A:118:LYS:HZ1	2.34	0.46
1:A:99:ILE:HD12	1:A:103:TYR:HD2	1.81	0.46
1:A:60:ASP:HB3	5:A:283:HOH:O	2.16	0.46
1:A:24:HIS:CG	1:A:118:LYS:HZ1	2.35	0.45
1:A:145:LYS:HB3	5:A:281:HOH:O	2.17	0.44
1:A:23:GLY:HA2	5:A:307:HOH:O	2.17	0.44
1:A:109:ASP:HB3	5:A:298:HOH:O	2.18	0.44
1:A:87:LYS:HG2	5:A:273:HOH:O	2.17	0.44
1:A:83:GLU:CD	1:A:83:GLU:H	2.20	0.44
1:A:31:ARG:CG	5:A:189:HOH:O	2.65	0.43
1:A:107:ILE:O	1:A:110:ALA:HB3	2.18	0.43
1:A:127:ALA:HB2	5:A:182:HOH:O	2.18	0.43
1:A:98:LYS:HD2	1:A:151:PHE:HE1	1.84	0.43
1:A:99:ILE:CD1	4:A:154:CLN:HAC	2.49	0.43
1:A:82:HIS:CE1	1:A:141:ASP:OD2	2.71	0.43
1:A:98:LYS:HA	1:A:151:PHE:CZ	2.54	0.42
1:A:87:LYS:HD3	1:A:91:GLN:OE1	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:SER:O	1:A:96:LYS:HB3	2.19	0.42
4:A:154:CLN:HAA1	5:A:170:HOH:O	2.19	0.42
1:A:2:LEU:CD1	1:A:133:LYS:HB3	2.49	0.42
1:A:12:ASN:ND2	5:A:324:HOH:O	2.52	0.42
1:A:4:ASP:HA	1:A:7:TRP:HB2	2.02	0.41
1:A:141:ASP:O	1:A:145:LYS:CE	2.68	0.41
1:A:61:LEU:O	1:A:61:LEU:HD12	2.21	0.41
1:A:12:ASN:HA	5:A:241:HOH:O	2.21	0.40
1:A:97:HIS:CD2	5:A:223:HOH:O	2.74	0.40
1:A:152:GLN:N	5:A:312:HOH:O	2.53	0.40

All (8) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:GLY:CA	5:A:318:HOH:O[1_545]	1.42	0.78
5:A:165:HOH:O	5:A:289:HOH:O[1_565]	1.54	0.66
1:A:91:GLN:NE2	1:A:153:GLY:C[2_555]	2.00	0.20
1:A:91:GLN:NE2	1:A:153:GLY:O[2_555]	2.05	0.15
1:A:129:GLY:N	5:A:318:HOH:O[1_545]	2.09	0.11
5:A:192:HOH:O	5:A:272:HOH:O[1_546]	2.12	0.08
5:A:255:HOH:O	5:A:326:HOH:O[2_555]	2.14	0.06
1:A:136:GLU:OE1	1:A:153:GLY:OXT[2_545]	2.19	0.01

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	151/153 (99%)	142 (94%)	8 (5%)	1 (1%)	26 19

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	152	GLN

5.3.2 Protein sidechains ⓘ

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	123/123 (100%)	101 (82%)	22 (18%)	2 1

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LEU
1	A	8	GLN
1	A	9	GLN
1	A	44	ASP
1	A	45	LYS
1	A	47	LYS
1	A	50	LYS
1	A	56	LYS
1	A	59	GLU
1	A	66	THR
1	A	72	LEU
1	A	76	LEU
1	A	87	LYS
1	A	95	THR
1	A	101	ILE
1	A	107	ILE
1	A	115	LEU
1	A	120	PRO
1	A	123	PHE
1	A	139	ARG
1	A	145	LYS
1	A	146	TYR

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

Mol	Chain	Res	Type
1	A	24	HIS
1	A	36	HIS
1	A	119	HIS
1	A	128	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

3 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	CLN	A	154	1,3	25,52,52	3.84	11 (44%)	23,87,87	3.73	12 (52%)
2	SO4	A	155	-	4,4,4	1.73	1 (25%)	6,6,6	1.17	0
3	CYN	A	156	4	0,1,1	0.00	-	0,0,0	0.00	-

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	CLN	A	154	1,3	1/1/8/15	0/6/83/83	0/0/9/9
2	SO4	A	155	-	-	0/0/0/0	0/0/0/0
3	CYN	A	156	4	-	0/0/0/0	0/0/0/0

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	154	CLN	C3C-CAC	-7.63	1.34	1.52
4	A	154	CLN	C2C-S	-6.65	1.77	1.85
4	A	154	CLN	C3C-C4C	-5.82	1.46	1.51
4	A	154	CLN	C3B-C2B	-5.22	1.33	1.40
4	A	154	CLN	CBC-CAC	2.46	1.45	1.33
4	A	154	CLN	CAD-C3D	2.49	1.56	1.52
4	A	154	CLN	FE-NC	2.77	2.10	1.97
2	A	155	SO4	O4-S	2.88	1.57	1.47
4	A	154	CLN	C1D-ND	3.04	1.40	1.36
4	A	154	CLN	CMB-C2B	3.89	1.59	1.51
4	A	154	CLN	C3B-CAB	4.29	1.56	1.47
4	A	154	CLN	CBC-S	11.36	1.89	1.74

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	154	CLN	C3B-CAB-CBB	-7.86	110.24	126.32
4	A	154	CLN	CAA-C2A-C1A	-5.68	120.84	127.01
4	A	154	CLN	CMC-C2C-S	-4.64	102.28	108.94
4	A	154	CLN	CMB-C2B-C1B	-3.43	122.68	128.36
4	A	154	CLN	C3C-C2C-C1C	-3.05	97.43	101.03
4	A	154	CLN	CMB-C2B-C3B	2.14	129.28	125.09
4	A	154	CLN	CBA-CAA-C2A	2.96	117.84	112.53
4	A	154	CLN	C4C-NC-C1C	3.58	108.94	105.13
4	A	154	CLN	C2C-C1C-CHC	3.86	127.10	123.60
4	A	154	CLN	CMA-C3A-C2A	4.02	133.64	125.24
4	A	154	CLN	CAD-CBD-CGD	7.04	125.64	112.75
4	A	154	CLN	C2C-C3C-C4C	8.23	111.04	100.35

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	A	154	CLN	C3C

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	154	CLN	5	0
2	A	155	SO4	2	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	153/153 (100%)	-1.07	0 100 100	3, 12, 24, 31	0

There are no RSRZ outliers to report.

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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	SO4	A	155	5/5	0.94	0.18	6.36	53,53,54,54	0
4	CLN	A	154	44/44	0.98	0.07	1.11	13,18,20,22	0
3	CYN	A	156	2/2	0.98	0.10	-	14,14,14,15	0

6.5 Other polymers [i](#)

There are no such residues in this entry.