



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:37 PM GMT

PDB ID : 4CRY
Title : Direct visualisation of strain-induced protein post-translational modification
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Deposited on : 2014-03-02
Resolution : 1.61 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

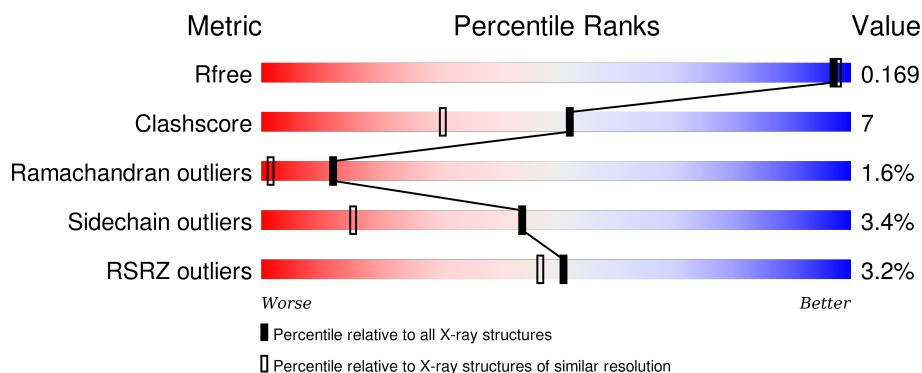
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.61 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}	91344	3202 (1.64-1.60)
Clashscore	102246	3500 (1.64-1.60)
Ramachandran outliers	100387	3411 (1.64-1.60)
Sidechain outliers	100360	3410 (1.64-1.60)
RSRZ outliers	91569	3207 (1.64-1.60)

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	41	
2	B	137	
3	G	102	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 2271 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 ASPARTATE 1-DECARBOXYLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	25	Total	C	N	O	S	0	1	0
			209	129	41	37	2			

There are 17 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-16	MET	-	EXPRESSION TAG	UNP P0A790
A	-15	ARG	-	EXPRESSION TAG	UNP P0A790
A	-14	GLY	-	EXPRESSION TAG	UNP P0A790
A	-13	SER	-	EXPRESSION TAG	UNP P0A790
A	-12	HIS	-	EXPRESSION TAG	UNP P0A790
A	-11	HIS	-	EXPRESSION TAG	UNP P0A790
A	-10	HIS	-	EXPRESSION TAG	UNP P0A790
A	-9	HIS	-	EXPRESSION TAG	UNP P0A790
A	-8	HIS	-	EXPRESSION TAG	UNP P0A790
A	-7	HIS	-	EXPRESSION TAG	UNP P0A790
A	-6	GLY	-	EXPRESSION TAG	UNP P0A790
A	-5	LEU	-	EXPRESSION TAG	UNP P0A790
A	-4	VAL	-	EXPRESSION TAG	UNP P0A790
A	-3	PRO	-	EXPRESSION TAG	UNP P0A790
A	-2	ARG	-	EXPRESSION TAG	UNP P0A790
A	-1	GLY	-	EXPRESSION TAG	UNP P0A790
A	0	SER	-	EXPRESSION TAG	UNP P0A790

- Molecule 2 is a protein called PANZ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	127	Total	C	N	O	S	0	9	0
			1079	678	204	190	7			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	128	SER	-	EXPRESSION TAG	UNP P37613
B	129	GLY	-	EXPRESSION TAG	UNP P37613
B	130	LEU	-	EXPRESSION TAG	UNP P37613
B	131	GLU	-	EXPRESSION TAG	UNP P37613
B	132	HIS	-	EXPRESSION TAG	UNP P37613
B	133	HIS	-	EXPRESSION TAG	UNP P37613
B	134	HIS	-	EXPRESSION TAG	UNP P37613
B	135	HIS	-	EXPRESSION TAG	UNP P37613
B	136	HIS	-	EXPRESSION TAG	UNP P37613
B	137	HIS	-	EXPRESSION TAG	UNP P37613

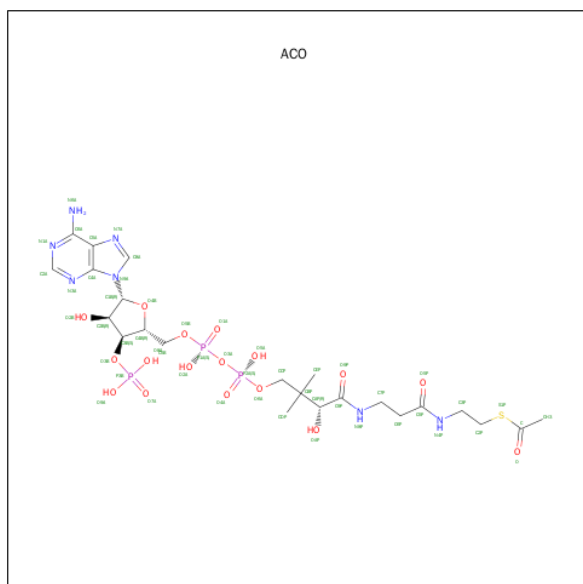
- Molecule 3 is a protein called ASPARTATE 1-DECARBOXYLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	G	102	Total	C	N	O	S	0	5	0
			806	504	142	156	4			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	57	VAL	THR	ENGINEERED MUTATION	UNP P0A790

- Molecule 4 is ACETYL COENZYME *A (three-letter code: ACO) (formula: C₂₃H₃₈N₇O₁₇P₃S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	B	1	Total	C	N	O	P	S	0	0
			51	23	7	17	3	1		

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total 1	Mg 1	0	0

- Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total 1	Cl 1	0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	19	Total 19	O 19	0	0
7	B	56	Total 56	O 56	0	0
7	G	49	Total 49	O 49	0	0

- Molecule 1: ASPARTATE 1-DECARBOXYLASE



S25
D29
D34
E40
D45
G52
F55
S56
V57
R67
V71
S89
D95
R102
E113
K119
A126

4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α , β , γ	86.40Å 86.40Å 81.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.19 – 1.61 43.20 – 1.61	Depositor EDS
% Data completeness (in resolution range)	97.3 (43.19-1.61) 98.8 (43.20-1.61)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.40 (at 1.61Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.129 , 0.150 0.151 , 0.169	Depositor DCC
R_{free} test set	1942 reflections (5.38%)	DCC
Wilson B-factor (Å ²)	20.4	Xtriage
Anisotropy	0.041	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 50.2	EDS
Estimated twinning fraction	0.841 for H, K, L 0.159 for K, H, -L 0.014 for -k,-h,-l	Xtriage
Reported twinning fraction	0.841 for H, K, L 0.159 for K, H, -L	Depositor
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 38084 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	2271	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.78% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CSO, ACO, MG, CL

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	1.83	6/211 (2.8%)	1.80	6/276 (2.2%)
2	B	1.30	2/1124 (0.2%)	1.47	15/1512 (1.0%)
3	G	1.41	1/826 (0.1%)	1.53	10/1118 (0.9%)
All	All	1.40	9/2161 (0.4%)	1.53	31/2906 (1.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	24[A]	GLY	N-CA	9.64	1.60	1.46
1	A	24[B]	GLY	N-CA	9.64	1.60	1.46
2	B	57	SER	CB-OG	-7.25	1.32	1.42
1	A	9	LYS	CB-CG	-6.17	1.35	1.52
1	A	3	ARG	NE-CZ	-5.96	1.25	1.33
1	A	23	GLU	CD-OE1	5.79	1.32	1.25
3	G	52	GLY	N-CA	-5.28	1.38	1.46
1	A	23	GLU	C-O	5.22	1.33	1.23
2	B	91	GLY	N-CA	5.06	1.53	1.46

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	38	ARG	NE-CZ-NH1	16.47	128.54	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	34	ASP	CB-CG-OD2	-13.98	105.72	118.30
2	B	38	ARG	NE-CZ-NH2	-11.60	114.50	120.30
1	A	12	ARG	NE-CZ-NH2	-11.59	114.50	120.30
2	B	47	ARG	NE-CZ-NH2	-10.59	115.01	120.30
3	G	29	ASP	CB-CG-OD2	-9.33	109.91	118.30
2	B	69[A]	ARG	NE-CZ-NH2	-8.51	116.04	120.30
2	B	69[B]	ARG	NE-CZ-NH2	-8.51	116.04	120.30
3	G	25[A]	SER	N-CA-CB	7.51	121.76	110.50
3	G	25[B]	SER	N-CA-CB	7.51	121.76	110.50
1	A	3	ARG	NE-CZ-NH2	-7.47	116.57	120.30
3	G	102[A]	ARG	NE-CZ-NH1	-6.96	116.82	120.30
3	G	102[B]	ARG	NE-CZ-NH1	-6.96	116.82	120.30
1	A	12	ARG	NE-CZ-NH1	6.94	123.77	120.30
2	B	97[A]	MET	CG-SD-CE	-6.82	89.29	100.20
2	B	97[B]	MET	CG-SD-CE	-6.82	89.29	100.20
3	G	55	PHE	CB-CG-CD1	6.72	125.51	120.80
1	A	24[A]	GLY	N-CA-C	6.60	129.59	113.10
1	A	24[B]	GLY	N-CA-C	6.60	129.59	113.10
2	B	34	ASP	CB-CG-OD1	6.16	123.84	118.30
2	B	47	ARG	NE-CZ-NH1	6.15	123.37	120.30
2	B	67	ARG	NE-CZ-NH2	-5.93	117.33	120.30
2	B	73	ARG	NE-CZ-NH1	-5.87	117.36	120.30
2	B	75[A]	ARG	NE-CZ-NH1	-5.79	117.41	120.30
2	B	75[B]	ARG	NE-CZ-NH1	-5.79	117.41	120.30
3	G	34	ASP	CB-CG-OD1	5.67	123.40	118.30
2	B	80	TYR	CB-CG-CD1	5.57	124.34	121.00
1	A	1	MET	CG-SD-CE	-5.32	91.68	100.20
2	B	43	ARG	NE-CZ-NH2	-5.29	117.65	120.30
3	G	45	ASP	CB-CG-OD2	5.19	122.97	118.30
3	G	95	ASP	CB-CG-OD1	5.10	122.89	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	23	GLU	Peptide

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	209	0	215	2	0
2	B	1079	0	1108	24	1
3	G	806	0	796	5	1
4	B	51	0	34	0	0
5	B	1	0	0	0	0
6	B	1	0	0	0	0
7	A	19	0	0	0	0
7	B	56	0	0	2	0
7	G	49	0	0	0	0
All	All	2271	0	2153	29	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (29) 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)
3:G:67[A]:ARG:HH22	3:G:113[A]:GLU:HG2	1.19	1.00
2:B:28:PRO:HA	2:B:29:SER:C	1.88	0.94
3:G:67[A]:ARG:NH2	3:G:113[A]:GLU:HG2	1.83	0.93
2:B:28:PRO:HG2	2:B:31:LEU:HD22	1.65	0.79
2:B:27:SER:N	2:B:28:PRO:HD2	2.01	0.76
2:B:12:SER:O	2:B:16[A]:ARG:HG3	1.92	0.70
2:B:28:PRO:HG3	2:B:31:LEU:HD13	1.75	0.68
2:B:27:SER:N	2:B:28:PRO:CD	2.57	0.66
2:B:28:PRO:HA	2:B:29:SER:O	1.97	0.64
2:B:71:VAL:HG13	7:B:2033:HOH:O	1.98	0.63
2:B:99:ASP:OD2	2:B:105[A]:ARG:NH1	2.31	0.63
2:B:28:PRO:HB3	2:B:31:LEU:HB2	1.88	0.56
1:A:24[A]:GLY:HA2	3:G:25[A]:SER:HA	1.88	0.55
2:B:28:PRO:CG	2:B:31:LEU:HD22	2.35	0.53
2:B:75[B]:ARG:HH11	2:B:75[B]:ARG:HG3	1.75	0.51
2:B:22:ILE:O	2:B:67:ARG:HD2	2.11	0.50
2:B:31:LEU:HG	2:B:39:ILE:HD11	1.95	0.48
2:B:33:VAL:HG13	7:B:2011:HOH:O	2.13	0.48
2:B:28:PRO:CG	2:B:31:LEU:HD13	2.44	0.48
2:B:84:GLU:OE2	2:B:87[A]:ARG:NH2	2.48	0.46
2:B:26:TYR:C	2:B:28:PRO:HD2	2.36	0.46
2:B:69[B]:ARG:HH11	2:B:69[B]:ARG:HD3	1.63	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:63:LEU:O	2:B:97[B]:MET:HG3	2.17	0.44
3:G:102[A]:ARG:HB3	3:G:119:LYS:O	2.18	0.43
2:B:6:ILE:HD13	2:B:6:ILE:HG21	1.81	0.42
2:B:89:ASN:N	2:B:90:PRO:CD	2.82	0.42
2:B:29:SER:O	2:B:30:SER:CB	2.68	0.41
1:A:18:ALA:HA	3:G:71:VAL:O	2.21	0.41

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:75[A]:ARG:NH1	3:G:40:GLU:OE2[4_555]	1.85	0.35

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	23/41 (56%)	22 (96%)	1 (4%)	0	100	100
2	B	134/137 (98%)	128 (96%)	3 (2%)	3 (2%)	8	1
3	G	103/102 (101%)	102 (99%)	0	1 (1%)	19	4
All	All	260/280 (93%)	252 (97%)	4 (2%)	4 (2%)	12	2

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	28	PRO
2	B	30	SER
2	B	27	SER
3	G	57	VAL

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	22/35 (63%)	21 (96%)	1 (4%)	34 9
2	B	116/116 (100%)	111 (96%)	5 (4%)	35 10
3	G	83/78 (106%)	82 (99%)	1 (1%)	78 59
All	All	221/229 (96%)	214 (97%)	7 (3%)	44 17

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
2	B	25	GLU
2	B	28	PRO
2	B	31	LEU
2	B	32	GLN
2	B	35	ASP
3	G	89	SER

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

Mol	Chain	Res	Type
2	B	120	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](#)

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link

column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	CSO	G	78	3	3,6,7	1.42	1 (33%)	1,6,8	2.28	1 (100%)

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	CSO	G	78	3	-	0/1/5/7	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	G	78	CSO	CB-CA	-2.34	1.47	1.53

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	78	CSO	O-C-CA	-2.28	119.56	125.49

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	ACO	B	1128	5	43,53,53	1.14	2 (4%)	55,79,79	2.29	17 (30%)

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	ACO	B	1128	5	-	0/47/67/67	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1128	ACO	C4A-N3A	2.30	1.39	1.35
4	B	1128	ACO	P3B-O3B	2.38	1.67	1.60

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1128	ACO	N3A-C2A-N1A	-9.89	121.32	128.89
4	B	1128	ACO	O6A-CCP-CBP	-3.59	104.78	110.55
4	B	1128	ACO	O8A-P3B-O7A	-3.45	99.46	110.58
4	B	1128	ACO	O9P-C9P-N8P	-3.30	116.47	123.08
4	B	1128	ACO	C1B-N9A-C4A	-3.16	122.18	126.94
4	B	1128	ACO	C4A-C5A-N7A	-3.10	106.63	109.48
4	B	1128	ACO	O5P-C5P-N4P	-2.85	117.28	122.94
4	B	1128	ACO	C2P-C3P-N4P	-2.83	106.70	112.36
4	B	1128	ACO	OAP-CAP-C9P	-2.75	104.08	110.38
4	B	1128	ACO	O2B-C2B-C3B	-2.20	104.79	111.16
4	B	1128	ACO	O2A-P1A-O3A	2.00	114.17	105.09
4	B	1128	ACO	C2A-N1A-C6A	2.03	122.40	118.77
4	B	1128	ACO	O3B-C3B-C2B	2.14	119.83	111.51
4	B	1128	ACO	O5P-C5P-C6P	2.91	127.00	121.98
4	B	1128	ACO	C4B-O4B-C1B	3.74	113.83	109.72
4	B	1128	ACO	O9A-P3B-O8A	4.00	122.59	107.38
4	B	1128	ACO	O9A-P3B-O7A	4.45	124.91	110.58

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

There are no such residues in this entry.

5.8 Polymer linkage issues

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	25/41 (60%)	-0.04	0 100 100	14, 17, 32, 52	0
2	B	127/137 (92%)	0.04	8 (6%) 23 20	16, 29, 51, 100	0
3	G	101/102 (99%)	-0.18	0 100 100	13, 17, 33, 36	0
All	All	253/280 (90%)	-0.05	8 (3%) 51 47	13, 22, 45, 100	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	29	SER	5.8
2	B	28	PRO	5.4
2	B	27	SER	5.0
2	B	32	GLN	4.9
2	B	31	LEU	3.1
2	B	30	SER	2.8
2	B	26	TYR	2.4
2	B	10	LYS	2.2

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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
3	CSO	G	78	7/8	0.99	0.08	-	15,17,24,33	0

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(\AA^2)	Q<0.9
5	MG	B	1129	1/1	1.00	0.07	0.60	26,26,26,26	0
4	ACO	B	1128	51/51	0.98	0.06	-0.79	18,28,36,37	0
6	CL	B	1130	1/1	0.92	0.48	-	30,30,30,30	1

6.5 Other polymers [i](#)

There are no such residues in this entry.