



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 26, 2014 – 07:05 PM GMT

PDB ID : 2C1J  
Title : Molecular basis for the recognition of phosphorylated and phosphoacetylated histone H3 by 14-3-3  
Authors : Welburn, J.P.I.; Macdonald, N.; Noble, M.E.M.; Nguyen, A.; Yaffe, M.B.; Clynes, D.; Moggs, J.G.; Orphanides, G.; Thomson, S.; Edmunds, J.W.; Clayton, A.L.; Endicott, J.A.; Mahadevan, L.C.  
Deposited on : 2005-09-15  
Resolution : 2.60 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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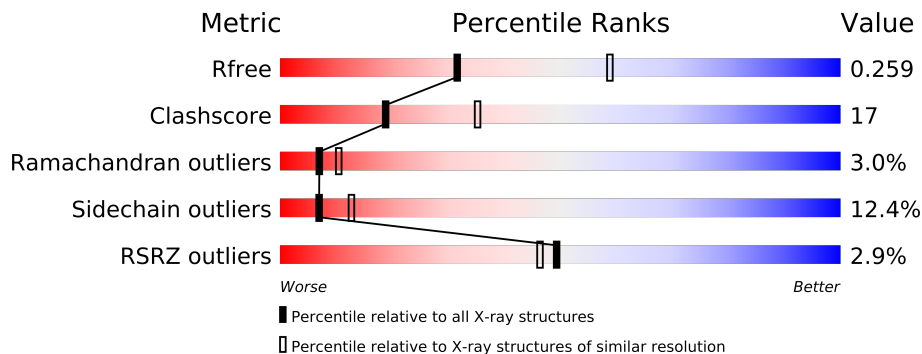
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.60 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))
$R_{free}$	66092	1718 (2.60-2.60)
Clashscore	79885	2154 (2.60-2.60)
Ramachandran outliers	78287	2113 (2.60-2.60)
Sidechain outliers	78261	2113 (2.60-2.60)
RSRZ outliers	66119	1718 (2.60-2.60)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	258	
1	B	258	
2	C	8	
2	D	8	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3892 atoms, of which 0 are hydrogen and 0 are deuterium.

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 14-3-3 PROTEIN ZETA/DELTA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	230	Total	C	N	O	S	0	0	0
			1841	1154	309	368	10			
1	B	230	Total	C	N	O	S	0	0	0
			1841	1154	309	368	10			

- Molecule 2 is a protein called HISTONE H3 ACETYLPHOSPHOPEPTIDE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	8	Total	C	N	O	P	0	0	0
			62	34	13	14	1			
2	D	8	Total	C	N	O	P	0	0	0
			62	34	13	14	1			

- Molecule 3 is water.

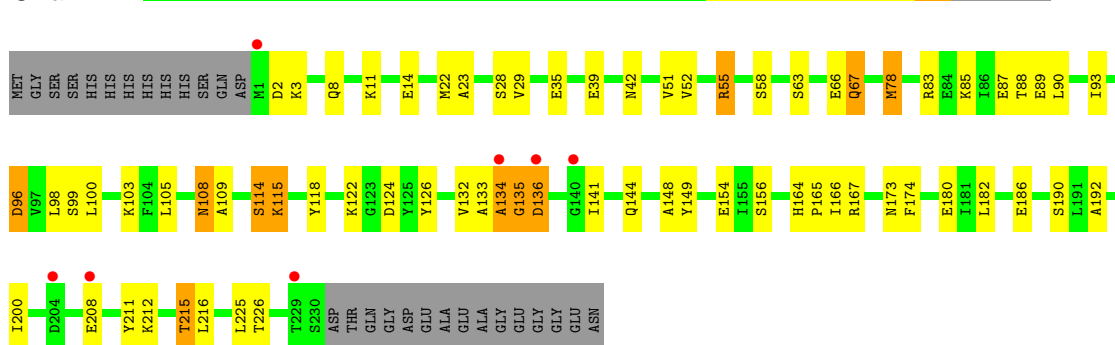
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	51	Total	O	0	0
			51	51		
3	B	27	Total	O	0	0
			27	27		
3	C	3	Total	O	0	0
			3	3		
3	D	5	Total	O	0	0
			5	5		

### 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.

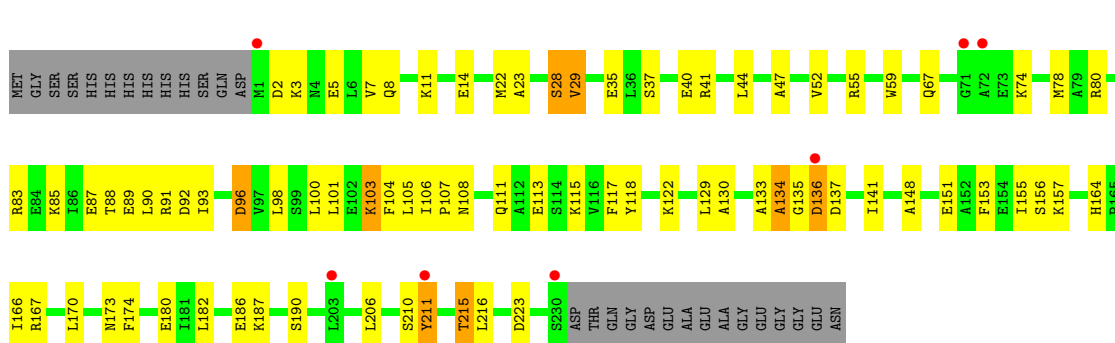
- Molecule 1: 14-3-3 PROTEIN ZETA/DELTA

Chain A:



- Molecule 1: 14-3-3 PROTEIN ZETA/DELTA

Chain B:



- Molecule 2: HISTONE H3 ACETYLPHOSPHOPEPTIDE

Chain C:



- Molecule 2: HISTONE H3 ACETYLPHOSPHOPEPTIDE

Chain D:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	71.19Å 72.41Å 71.16Å 90.00° 102.96° 90.00°	Depositor
Resolution (Å)	69.34 – 2.60 21.20 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.7 (69.34-2.60) 99.9 (21.20-2.60)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.19 (at 2.60Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.227 , 0.292 0.215 , 0.259	Depositor DCC
$R_{free}$ test set	1113 reflections (5.37%)	DCC
Wilson B-factor (Å <sup>2</sup> )	45.9	Xtriage
Anisotropy	0.330	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , -20.0	EDS
Estimated twinning fraction	0.459 for l,-k,h	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 21803 reflections	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	3892	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	10.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.63% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

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

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	0.65	0/1866	0.73	0/2509
1	B	0.64	0/1866	0.74	0/2509
2	C	0.96	0/38	1.34	0/46
2	D	0.83	0/38	1.52	1/46 (2.2%)
All	All	0.65	0/3808	0.76	1/5110 (0.0%)

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
2	D	0	3

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
2	D	14	LYS	N-CA-C	5.09	124.74	111.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	D	13	GLY	Peptide
2	D	8	ARG	Mainchain
2	D	9	ALY	Mainchain

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1841	0	1830	52	0
1	B	1841	0	1830	56	0
2	C	62	0	63	18	0
2	D	62	0	59	7	0
3	A	51	0	0	7	0
3	B	27	0	0	5	0
3	C	3	0	0	0	0
3	D	5	0	0	2	0
All	All	3892	0	3782	124	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 17.

All (124) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:C:9:ALY:C	2:C:10:SEP:H2	1.46	1.19
1:A:42:ASN:HB3	3:A:2010:HOH:O	1.45	1.12
1:A:133:ALA:HA	1:A:134:ALA:HB2	1.48	0.96
1:B:164:HIS:HD2	1:B:166:ILE:H	1.14	0.95
1:B:210:SER:HB3	3:B:2026:HOH:O	1.64	0.95
1:B:80:ARG:HD3	3:B:2012:HOH:O	1.65	0.95
1:A:83:ARG:HH11	1:A:83:ARG:HG2	1.31	0.94
2:C:8:ARG:C	2:C:9:ALY:H2	1.62	0.93
1:A:180:GLU:OE1	2:C:8:ARG:HD3	1.69	0.93
2:C:8:ARG:O	2:C:9:ALY:N	2.08	0.87
1:A:164:HIS:HD2	1:A:166:ILE:H	1.22	0.87
1:B:156:SER:HB2	1:B:167:ARG:HG3	1.58	0.85
1:B:133:ALA:HA	1:B:134:ALA:HB2	1.61	0.83
1:B:8:GLN:HB3	3:B:2002:HOH:O	1.79	0.83
1:A:156:SER:HB2	1:A:167:ARG:HG3	1.60	0.82
1:A:180:GLU:OE1	2:C:8:ARG:CD	2.28	0.81
1:B:83:ARG:HG2	1:B:83:ARG:HH11	1.44	0.81
1:A:164:HIS:CD2	1:A:165:PRO:HD2	2.16	0.80
1:A:83:ARG:HG2	1:A:83:ARG:NH1	1.97	0.78

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:C:9:ALY:CA	2:C:10:SEP:H2	1.96	0.78
2:C:9:ALY:CA	2:C:10:SEP:N	2.46	0.77
1:B:173:ASN:HD21	2:D:11:THR:HB	1.50	0.77
1:B:164:HIS:CD2	1:B:166:ILE:H	2.01	0.76
1:A:66:GLU:OE2	1:A:83:ARG:NE	2.20	0.74
1:A:85:LYS:HE2	3:A:2021:HOH:O	1.88	0.73
1:B:135:GLY:O	1:B:136:ASP:HB2	1.88	0.72
2:C:9:ALY:C	2:C:10:SEP:CA	2.68	0.71
1:A:133:ALA:CA	1:A:134:ALA:HB2	2.23	0.68
1:A:8:GLN:HG3	3:A:2003:HOH:O	1.93	0.67
1:A:173:ASN:HD21	2:C:11:THR:HB	1.61	0.66
1:B:103:LYS:HB2	1:B:104:PHE:CD1	2.31	0.65
1:B:180:GLU:OE1	2:D:8:ARG:HD3	1.98	0.64
2:C:9:ALY:O	2:C:10:SEP:N	2.29	0.63
1:B:153:PHE:HA	1:B:170:LEU:HD21	1.80	0.63
1:A:164:HIS:CD2	1:A:166:ILE:H	2.11	0.63
1:B:59:TRP:HE1	1:B:87:GLU:HG3	1.63	0.62
1:A:135:GLY:O	1:A:136:ASP:HB2	2.00	0.62
1:B:164:HIS:HD2	1:B:166:ILE:N	1.92	0.61
2:C:8:ARG:CA	2:C:9:ALY:N	2.64	0.60
1:A:11:LYS:HE2	1:A:39:GLU:OE2	2.02	0.60
1:B:7:VAL:O	1:B:11:LYS:HG3	2.00	0.60
1:B:3:LYS:HD3	1:B:40:GLU:OE2	2.02	0.60
1:B:133:ALA:CA	1:B:134:ALA:HB2	2.31	0.60
1:A:99:SER:HB2	3:A:2026:HOH:O	2.01	0.59
1:A:109:ALA:HB3	1:A:115:LYS:HD2	1.84	0.59
2:C:8:ARG:C	2:C:9:ALY:H	1.99	0.59
2:D:7:ALA:N	3:D:2002:HOH:O	2.35	0.59
1:A:78:MET:HA	1:A:78:MET:HE2	1.83	0.58
1:B:28:SER:O	1:B:29:VAL:C	2.42	0.58
1:B:157:LYS:HB2	3:B:2019:HOH:O	2.04	0.57
1:A:98:LEU:HD22	1:A:126:TYR:CE2	2.39	0.57
1:A:55:ARG:HD3	1:A:93:ILE:HD13	1.85	0.57
1:A:133:ALA:HB1	1:A:134:ALA:HB3	1.86	0.57
2:C:9:ALY:HCA	2:C:10:SEP:N	2.20	0.57
1:A:156:SER:CB	1:A:167:ARG:HG3	2.33	0.56
1:B:59:TRP:NE1	1:B:87:GLU:HG3	2.20	0.56
1:B:153:PHE:CZ	1:B:157:LYS:HE3	2.40	0.56
1:A:133:ALA:CA	1:A:134:ALA:CB	2.82	0.56
1:A:109:ALA:HB1	1:A:114:SER:HB3	1.87	0.55
1:A:103:LYS:HE2	3:A:2030:HOH:O	2.06	0.55
1:A:88:THR:O	1:A:89:GLU:C	2.45	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:37:SER:OG	1:B:40:GLU:HG3	2.08	0.53
1:B:122:LYS:HD3	1:B:148:ALA:HB2	1.91	0.52
1:A:108:ASN:N	1:A:108:ASN:HD22	2.09	0.51
1:A:124:ASP:HA	1:A:149:TYR:OH	2.10	0.51
1:B:55:ARG:HD2	1:B:93:ILE:HD13	1.92	0.51
1:B:105:LEU:HD22	1:B:118:TYR:CZ	2.46	0.51
1:B:211:TYR:O	1:B:215:THR:HB	2.11	0.51
1:A:14:GLU:HB2	1:A:22:MET:CE	2.41	0.51
1:B:83:ARG:HG2	1:B:83:ARG:NH1	2.19	0.50
1:A:180:GLU:OE1	2:C:8:ARG:HD2	2.11	0.50
1:B:106:ILE:HB	1:B:107:PRO:HD3	1.95	0.49
1:A:211:TYR:O	1:A:215:THR:HB	2.12	0.49
1:A:109:ALA:CB	1:A:114:SER:HB3	2.43	0.49
2:D:9:ALY:HCA	3:D:2004:HOH:O	2.13	0.49
1:A:122:LYS:HG2	1:A:148:ALA:HB2	1.95	0.48
1:B:41:ARG:NH1	1:B:113:GLU:OE2	2.42	0.48
1:A:135:GLY:O	1:A:136:ASP:CB	2.61	0.48
1:B:133:ALA:HA	1:B:134:ALA:CB	2.40	0.48
2:C:8:ARG:NH2	2:C:10:SEP:O1P	2.44	0.48
1:B:59:TRP:HE1	1:B:87:GLU:CG	2.27	0.47
2:C:9:ALY:HCA	2:C:10:SEP:H2	1.78	0.47
1:B:156:SER:CB	1:B:167:ARG:HG3	2.37	0.47
1:A:67:GLN:N	1:A:67:GLN:HE21	2.12	0.47
1:A:164:HIS:HD2	1:A:166:ILE:N	2.02	0.46
1:B:223:ASP:HB3	2:D:9:ALY:HH31	1.97	0.46
1:B:206:LEU:HD22	1:B:211:TYR:HA	1.97	0.46
1:B:151:GLU:O	1:B:155:ILE:HG13	2.15	0.46
1:B:22:MET:HG2	1:B:47:ALA:HB2	1.97	0.46
1:A:164:HIS:CD2	1:A:165:PRO:CD	2.93	0.46
1:B:91:ARG:HG2	1:B:129:LEU:CD2	2.45	0.45
2:C:11:THR:HG22	2:C:12:GLY:N	2.31	0.45
1:A:51:VAL:HG13	1:A:55:ARG:NH1	2.31	0.45
1:B:153:PHE:CA	1:B:170:LEU:HD21	2.46	0.45
1:B:23:ALA:HA	1:B:47:ALA:HB1	1.98	0.45
2:D:10:SEP:O	2:D:13:GLY:N	2.40	0.45
1:B:133:ALA:CA	1:B:134:ALA:CB	2.94	0.45
1:A:78:MET:CE	1:A:78:MET:HA	2.46	0.44
1:A:133:ALA:HB1	1:A:134:ALA:CB	2.47	0.44
1:B:173:ASN:ND2	2:D:11:THR:HB	2.25	0.43
1:B:135:GLY:O	1:B:136:ASP:CB	2.60	0.43
1:A:55:ARG:HD3	1:A:93:ILE:CD1	2.48	0.43
1:B:14:GLU:HB2	1:B:22:MET:SD	2.58	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:41:ARG:HD3	1:B:117:PHE:CD1	2.53	0.43
1:B:88:THR:O	1:B:89:GLU:C	2.57	0.43
1:A:154:GLU:CB	3:A:2036:HOH:O	2.67	0.43
1:A:78:MET:HE1	1:B:5:GLU:OE2	2.19	0.42
1:A:51:VAL:HG13	1:A:55:ARG:HH11	1.85	0.42
1:A:154:GLU:HB3	3:A:2036:HOH:O	2.20	0.41
1:A:28:SER:O	1:A:29:VAL:C	2.58	0.41
1:A:192:ALA:HB3	1:A:225:LEU:HD21	2.02	0.41
1:A:63:SER:O	1:A:67:GLN:NE2	2.54	0.41
1:B:157:LYS:HA	1:B:167:ARG:NH1	2.35	0.41
1:B:83:ARG:NH1	1:B:83:ARG:CG	2.84	0.41
1:B:186:GLU:HG3	1:B:187:LYS:H	1.84	0.41
2:C:14:LYS:H	2:C:14:LYS:HD3	1.85	0.41
1:B:35:GLU:HG3	1:B:108:ASN:HB3	2.03	0.41
1:B:98:LEU:HA	1:B:101:LEU:HD12	2.02	0.41
1:B:186:GLU:HG3	1:B:187:LYS:N	2.36	0.41
1:B:130:ALA:HA	1:B:141:ILE:HD11	2.03	0.40
1:B:59:TRP:CD1	1:B:87:GLU:HG3	2.56	0.40
1:B:85:LYS:HE3	3:B:2013:HOH:O	2.22	0.40
1:A:105:LEU:HD22	1:A:118:TYR:CZ	2.56	0.40
1:A:87:GLU:HG3	1:A:132:VAL:HG13	2.04	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

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	228/258 (88%)	208 (91%)	14 (6%)	6 (3%)	8	13
1	B	228/258 (88%)	207 (91%)	16 (7%)	5 (2%)	10	18
2	C	4/8 (50%)	2 (50%)	1 (25%)	1 (25%)	0	0
2	D	4/8 (50%)	2 (50%)	0	2 (50%)	0	0
All	All	464/532 (87%)	419 (90%)	31 (7%)	14 (3%)	7	10

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	134	ALA
1	B	134	ALA
1	B	29	VAL
1	B	96	ASP
1	A	96	ASP
2	D	8	ARG
2	C	11	THR
2	D	11	THR
1	A	23	ALA
1	A	58	SER
1	B	136	ASP
1	A	136	ASP
1	B	28	SER
1	A	135	GLY

### 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	199/221 (90%)	174 (87%)	25 (13%)	7	12
1	B	199/221 (90%)	179 (90%)	20 (10%)	11	20
2	C	3/3 (100%)	0	3 (100%)	0	0
2	D	3/3 (100%)	1 (33%)	2 (67%)	0	0
All	All	404/448 (90%)	354 (88%)	50 (12%)	7	13

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ASP
1	A	3	LYS
1	A	35	GLU
1	A	52	VAL
1	A	55	ARG
1	A	67	GLN
1	A	78	MET

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Mol	Chain	Res	Type
1	A	90	LEU
1	A	96	ASP
1	A	100	LEU
1	A	108	ASN
1	A	114	SER
1	A	115	LYS
1	A	141	ILE
1	A	144	GLN
1	A	174	PHE
1	A	182	LEU
1	A	186	GLU
1	A	190	SER
1	A	200	ILE
1	A	208	GLU
1	A	212	LYS
1	A	215	THR
1	A	216	LEU
1	A	226	THR
1	B	2	ASP
1	B	44	LEU
1	B	52	VAL
1	B	67	GLN
1	B	74	LYS
1	B	78	MET
1	B	90	LEU
1	B	92	ASP
1	B	96	ASP
1	B	100	LEU
1	B	103	LYS
1	B	111	GLN
1	B	115	LYS
1	B	137	ASP
1	B	174	PHE
1	B	182	LEU
1	B	190	SER
1	B	211	TYR
1	B	215	THR
1	B	216	LEU
2	C	8	ARG
2	C	11	THR
2	C	14	LYS
2	D	8	ARG

*Continued on next page...*

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Mol	Chain	Res	Type
2	D	11	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	67	GLN
1	A	108	ASN
1	A	164	HIS
1	B	76	GLN
1	B	108	ASN
1	B	164	HIS

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

4 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$
2	SEP	C	10	2	9,9,10	5.92	2 (22%)	10,12,14	2.95	2 (20%)
2	ALY	C	9	2	11,11,12	5.35	1 (9%)	10,12,14	4.03	2 (20%)
2	SEP	D	10	2	9,9,10	4.52	2 (22%)	10,12,14	2.14	3 (30%)
2	ALY	D	9	2	11,11,12	4.68	1 (9%)	10,12,14	2.04	3 (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
2	SEP	C	10	2	-	0/6/8/10	0/0/0/0
2	ALY	C	9	2	-	0/8/10/12	0/0/0/0
2	SEP	D	10	2	-	0/6/8/10	0/0/0/0
2	ALY	D	9	2	-	0/8/10/12	0/0/0/0

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	9	ALY	O-C	17.60	1.23	1.11
2	C	10	SEP	O-C	17.24	1.23	1.11
2	D	9	ALY	O-C	15.36	1.21	1.11
2	D	10	SEP	O-C	13.01	1.20	1.11
2	C	10	SEP	P-O1P	3.16	1.61	1.51
2	D	10	SEP	P-O1P	2.64	1.60	1.51

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	9	ALY	C-CA-N	-11.89	101.96	113.83
2	C	10	SEP	C-CA-N	-8.61	105.23	113.83
2	D	10	SEP	C-CA-N	-4.85	108.99	113.83
2	D	9	ALY	C-CA-N	-4.50	109.33	113.83
2	D	9	ALY	CE-NZ-CH	3.81	128.59	122.32
2	D	10	SEP	O3P-P-O1P	-2.93	100.87	110.44
2	C	9	ALY	CE-NZ-CH	2.75	126.84	122.32
2	D	10	SEP	P-OG-CB	-2.55	110.83	118.19
2	C	10	SEP	O2P-P-OG	2.33	113.07	106.65
2	D	9	ALY	CH3-CH-NZ	2.17	119.49	116.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.5 Carbohydrates

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

There are no ligands in this entry.

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	230/258 (89%)	0.41	7 (3%) 48 45	2, 9, 24, 35	0
1	B	230/258 (89%)	0.40	7 (3%) 48 45	2, 9, 24, 35	0
2	C	8/8 (100%)	0.47	0 100 100	8, 11, 15, 18	0
2	D	8/8 (100%)	0.20	0 100 100	8, 11, 15, 17	0
All	All	476/532 (89%)	0.41	14 (2%) 49 46	2, 9, 24, 35	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	4.2
1	B	71	GLY	4.1
1	A	136	ASP	3.8
1	A	204	ASP	3.4
1	B	230	SER	3.0
1	B	203	LEU	2.7
1	A	134	ALA	2.6
1	B	1	MET	2.5
1	A	229	THR	2.5
1	A	140	GLY	2.3
1	B	211	TYR	2.2
1	B	136	ASP	2.1
1	A	208	GLU	2.0
1	B	72	ALA	2.0

### 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, 95<sup>th</sup> 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	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	ALY	C	9	12/13	0.27	1.81	8,11,20,22	0
2	SEP	C	10	10/11	0.22	0.49	3,8,9,9	0
2	SEP	D	10	10/11	0.20	-0.19	2,7,9,9	0
2	ALY	D	9	12/13	0.17	-1.49	8,11,21,24	0

### 6.3 Carbohydrates

There are no carbohydrates in this entry.

### 6.4 Ligands

There are no ligands in this entry.

### 6.5 Other polymers

There are no such residues in this entry.