



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 08:50 PM GMT

PDB ID : 4YYI
Title : Crystal structure of BRD9 Bromodomain bound to an acetylated peptide
Authors : Tang, Y.; Bellon, S.; Cochran, A.G.; Poy, F.
Deposited on : 2015-03-23
Resolution : 1.50 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
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) : rb-20026982

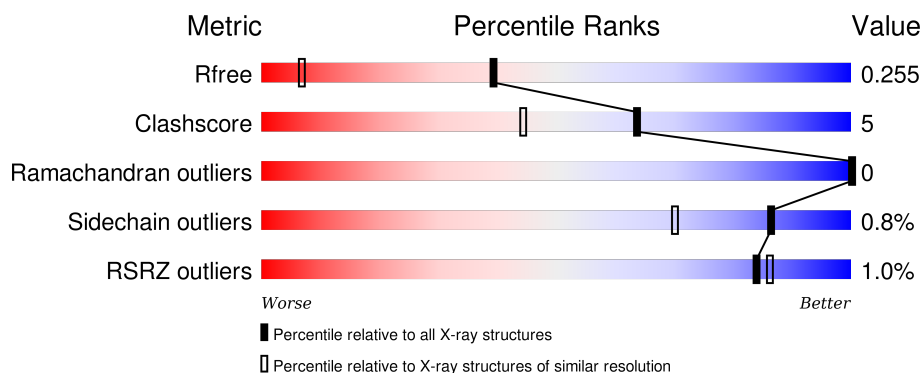
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


The reported resolution of this entry is 1.50 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.



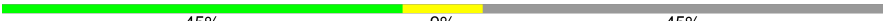
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	2072 (1.50-1.50)
Clashscore	102246	2274 (1.50-1.50)
Ramachandran outliers	100387	2218 (1.50-1.50)
Sidechain outliers	100360	2216 (1.50-1.50)
RSRZ outliers	91569	2075 (1.50-1.50)

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	108	 74% 18% • 7%
1	B	108	 70% 22% 7%
1	D	108	 73% 19% • 6%
1	E	108	 74% 18% • 6%
2	C	11	 36% 18% 45%

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Mol	Chain	Length	Quality of chain
2	F	11	 A horizontal bar chart showing the quality of chain F. The bar is divided into three segments: a green segment on the left labeled '45%', a yellow segment in the middle labeled '9%', and a grey segment on the right labeled '45%'. The segments are separated by thin white lines.

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3639 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 Bromodomain-containing protein 9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	100	Total	C	N	O	S	0	0	0
			815	531	135	141	8			
1	B	100	Total	C	N	O	S	0	0	0
			815	531	135	141	8			
1	D	101	Total	C	N	O	S	0	0	0
			821	534	136	143	8			
1	E	101	Total	C	N	O	S	0	0	0
			821	534	136	143	8			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	16	GLY	-	expression tag	UNP Q9H8M2
B	16	GLY	-	expression tag	UNP Q9H8M2
D	16	GLY	-	expression tag	UNP Q9H8M2
E	16	GLY	-	expression tag	UNP Q9H8M2

- Molecule 2 is a protein called Histone H4.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	6	Total	C	N	O	0	0	0
			40	24	8	8			
2	F	6	Total	C	N	O	0	0	0
			40	24	8	8			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	76	Total	O	0	0
			76	76		
3	B	61	Total	O	0	0
			61	61		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	5	Total 5	O 5	0	0
3	D	66	Total 66	O 66	0	0
3	E	76	Total 76	O 76	0	0
3	F	3	Total 3	O 3	0	0

3 Residue-property plots [i](#)

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.

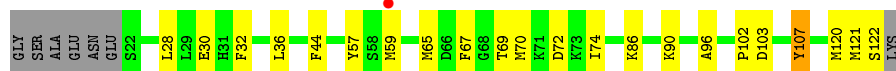
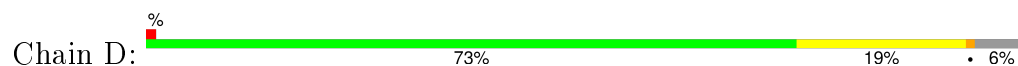
- Molecule 1: Bromodomain-containing protein 9



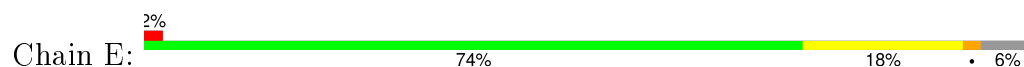
- Molecule 1: Bromodomain-containing protein 9



- Molecule 1: Bromodomain-containing protein 9



- Molecule 1: Bromodomain-containing protein 9



- Molecule 2: Histone H4



- Molecule 2: Histone H4



SER	GLY	ARG	G4	K5	G6	G9	LEU	GLY
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	24.75Å 34.60Å 129.36Å 88.79° 90.00° 68.92°	Depositor
Resolution (Å)	32.28 – 1.50 32.28 – 1.50	Depositor EDS
% Data completeness (in resolution range)	94.5 (32.28-1.50) 98.7 (32.28-1.50)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.38 (at 1.50Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.206 , 0.242 0.225 , 0.255	Depositor DCC
R_{free} test set	3185 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	22.7	Xtriage
Anisotropy	0.255	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 27.0	EDS
Estimated twinning fraction	0.556 for H, K, L 0.444 for H, H-K, -L 0.279 for h,h-k,-l 0.180 for -h,-k,l 0.177 for -h,-h+k,-l	Xtriage
Reported twinning fraction	0.556 for H, K, L 0.444 for H, H-K, -L	Depositor
L-test for twinning ²	$\langle L \rangle = 0.54$, $\langle L^2 \rangle = 0.38$	Xtriage
Outliers	0 of 63617 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3639	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 8.55% 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, 0.375 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: ALY

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	1.40	3/837 (0.4%)	1.40	7/1125 (0.6%)
1	B	1.34	4/837 (0.5%)	1.44	11/1125 (1.0%)
1	D	1.34	6/843 (0.7%)	1.45	8/1133 (0.7%)
1	E	1.42	1/843 (0.1%)	1.43	8/1133 (0.7%)
2	C	2.47	1/13 (7.7%)	2.18	0/11
2	F	2.66	1/13 (7.7%)	2.88	2/11 (18.2%)
All	All	1.39	16/3386 (0.5%)	1.44	36/4538 (0.8%)

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	67	PHE	CG-CD2	8.84	1.52	1.38
2	F	6	GLY	C-O	7.40	1.35	1.23
1	D	67	PHE	CG-CD2	6.50	1.48	1.38
1	B	67	PHE	C-O	6.45	1.35	1.23
1	D	96	ALA	C-O	6.20	1.35	1.23
2	C	6	GLY	N-CA	6.07	1.55	1.46
1	D	36	LEU	C-O	5.89	1.34	1.23
1	D	107	TYR	CE1-CZ	5.82	1.46	1.38
1	E	79	TYR	CB-CG	5.81	1.60	1.51
1	A	47	PHE	CG-CD2	5.69	1.47	1.38
1	B	89	PHE	CG-CD1	5.51	1.47	1.38
1	B	57	TYR	CB-CG	5.46	1.59	1.51
1	A	98	THR	C-O	5.44	1.33	1.23
1	D	70	MET	C-O	5.36	1.33	1.23
1	B	106	TYR	CE2-CZ	-5.13	1.31	1.38
1	D	32	PHE	CG-CD2	5.09	1.46	1.38

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	103	ASP	CB-CG-OD1	10.36	127.62	118.30
1	D	103	ASP	CB-CG-OD1	8.64	126.08	118.30
1	A	103	ASP	CB-CG-OD2	-8.59	110.57	118.30
1	E	67	PHE	CB-CG-CD1	8.56	126.80	120.80
1	B	70	MET	CG-SD-CE	8.45	113.71	100.20
1	D	72	ASP	CB-CG-OD1	8.30	125.77	118.30
1	D	103	ASP	CB-CG-OD2	-8.16	110.95	118.30
1	E	29	LEU	CB-CG-CD2	7.85	124.35	111.00
1	B	44	PHE	CB-CG-CD2	-7.72	115.40	120.80
1	A	44	PHE	CB-CG-CD2	-7.56	115.51	120.80
1	D	28	LEU	CB-CG-CD2	7.42	123.61	111.00
1	B	57	TYR	CD1-CE1-CZ	-7.33	113.21	119.80
1	D	57	TYR	CG-CD2-CE2	-7.05	115.66	121.30
1	D	44	PHE	CB-CG-CD2	-6.82	116.03	120.80
1	E	33	LEU	CB-CG-CD2	-6.70	99.62	111.00
2	F	6	GLY	CA-C-N	-6.66	102.89	116.20
1	E	80	LYS	CD-CE-NZ	6.56	126.78	111.70
1	E	94	ASP	CB-CG-OD2	-6.47	112.48	118.30
1	E	57	TYR	CB-CG-CD2	-6.46	117.13	121.00
1	B	66	ASP	CB-CG-OD2	-6.43	112.51	118.30
1	B	57	TYR	CB-CG-CD1	-6.25	117.25	121.00
1	B	28	LEU	CB-CG-CD2	6.14	121.44	111.00
1	B	57	TYR	CG-CD2-CE2	-5.96	116.54	121.30
1	B	66	ASP	CB-CG-OD1	5.90	123.61	118.30
1	D	67	PHE	CB-CG-CD2	5.84	124.89	120.80
1	B	65	MET	CG-SD-CE	5.81	109.49	100.20
1	A	97	MET	CG-SD-CE	5.59	109.15	100.20
1	B	106	TYR	CB-CG-CD2	-5.57	117.66	121.00
1	A	105	VAL	CA-CB-CG1	-5.53	102.60	110.90
2	F	6	GLY	O-C-N	5.34	132.27	123.20
1	B	67	PHE	CB-CA-C	-5.17	100.07	110.40
1	E	40	ASP	CB-CG-OD2	-5.15	113.67	118.30
1	E	106	TYR	CB-CG-CD2	-5.14	117.91	121.00
1	A	29	LEU	CB-CG-CD2	5.04	119.57	111.00
1	D	65	MET	CG-SD-CE	5.04	108.27	100.20
1	A	89	PHE	CD1-CE1-CZ	-5.04	114.05	120.10

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	815	0	820	7	0
1	B	815	0	820	10	0
1	D	821	0	825	9	0
1	E	821	0	825	9	0
2	C	40	0	38	1	0
2	F	40	0	38	0	0
3	A	76	0	0	0	0
3	B	61	0	0	2	0
3	C	5	0	0	0	0
3	D	66	0	0	3	0
3	E	76	0	0	4	0
3	F	3	0	0	0	0
All	All	3639	0	3366	35	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 5.

All (35) 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:B:59:MET:HB2	3:B:214:HOH:O	1.72	0.88
1:E:115:HIS:HD2	3:E:254:HOH:O	1.63	0.80
1:E:101:ARG:NH1	1:E:103:ASP:OD2	2.20	0.74
1:A:37:GLN:HG2	1:A:67:PHE:CD2	2.29	0.67
1:E:25:ILE:HG22	3:E:225:HOH:O	1.94	0.67
1:B:103:ASP:O	1:B:108:LYS:HE3	2.00	0.61
1:E:66:ASP:OD1	1:E:69:THR:OG1	2.18	0.59
1:E:115:HIS:HE1	3:E:256:HOH:O	1.87	0.57
1:D:59:MET:HB2	3:D:211:HOH:O	2.04	0.57
1:B:104:THR:O	1:B:108:LYS:HG3	2.05	0.57
1:A:99:TYR:OH	2:C:7:GLY:HA2	2.05	0.57
1:D:86:LYS:HE2	1:D:121:MET:CB	2.39	0.52
1:D:90:LYS:HE2	3:D:222:HOH:O	2.09	0.51
1:A:34:ARG:O	1:A:38:ARG:HG3	2.11	0.50
1:B:30:GLU:HG2	1:B:74:ILE:HG21	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:102:PRO:HA	1:D:107:TYR:CG	2.48	0.49
1:A:90:LYS:NZ	1:A:94:ASP:OD2	2.45	0.47
1:E:37:GLN:HG2	1:E:67:PHE:CD2	2.50	0.47
1:B:102:PRO:HA	1:B:107:TYR:CG	2.50	0.47
1:A:35:GLN:OE1	1:A:38:ARG:NH2	2.48	0.46
1:D:30:GLU:HG2	1:D:74:ILE:HG21	1.99	0.45
1:E:102:PRO:HA	1:E:107:TYR:CG	2.52	0.44
1:E:115:HIS:CD2	3:E:254:HOH:O	2.51	0.43
1:E:93:CYS:O	1:E:97:MET:HG3	2.19	0.43
1:D:86:LYS:HE2	1:D:121:MET:HB2	2.01	0.43
1:B:33:LEU:HD21	1:B:71:LYS:HB2	2.00	0.43
1:B:56:GLY:O	1:B:60:ILE:HG13	2.19	0.42
1:D:69:THR:HG23	3:D:242:HOH:O	2.18	0.42
1:D:86:LYS:HE2	1:D:121:MET:HB3	2.00	0.42
1:A:102:PRO:HA	1:A:107:TYR:CG	2.55	0.42
1:A:24:PRO:O	1:A:25:ILE:C	2.55	0.41
1:B:90:LYS:HE2	3:B:206:HOH:O	2.19	0.41
1:D:120:MET:C	1:D:122:SER:H	2.24	0.41
1:B:109:LEU:HA	1:B:109:LEU:HD12	1.83	0.40
1:B:33:LEU:CD2	1:B:71:LYS:HB2	2.52	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	98/108 (91%)	97 (99%)	1 (1%)	0	100	100
1	B	98/108 (91%)	98 (100%)	0	0	100	100
1	D	99/108 (92%)	98 (99%)	1 (1%)	0	100	100
1	E	99/108 (92%)	99 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	C	2/11 (18%)	2 (100%)	0	0	100	100
2	F	2/11 (18%)	2 (100%)	0	0	100	100
All	All	398/454 (88%)	396 (100%)	2 (0%)	0	100	100

There are no Ramachandran outliers to report.

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	88/94 (94%)	88 (100%)	0	100	100
1	B	88/94 (94%)	87 (99%)	1 (1%)	80	58
1	D	89/94 (95%)	89 (100%)	0	100	100
1	E	89/94 (95%)	87 (98%)	2 (2%)	60	25
All	All	354/376 (94%)	351 (99%)	3 (1%)	86	70

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	105	VAL
1	E	34	ARG
1	E	103	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	35	GLN
1	E	115	HIS

5.3.3 RNA ⓘ

There are no RNA molecules 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	ALY	C	5	2	9,11,12	2.06	5 (55%)	10,12,14	3.17	5 (50%)
2	ALY	C	8	2	9,11,12	2.58	2 (22%)	10,12,14	3.80	3 (30%)
2	ALY	F	5	2	9,11,12	1.48	1 (11%)	10,12,14	4.01	4 (40%)
2	ALY	F	8	2	9,11,12	1.84	2 (22%)	10,12,14	3.43	4 (40%)

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	ALY	C	5	2	-	0/8/10/12	0/0/0/0
2	ALY	C	8	2	-	0/8/10/12	0/0/0/0
2	ALY	F	5	2	-	0/8/10/12	0/0/0/0
2	ALY	F	8	2	-	0/8/10/12	0/0/0/0

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	5	ALY	OH-CH	-2.80	1.16	1.23
2	C	5	ALY	CH-NZ	-2.31	1.28	1.33
2	C	5	ALY	CB-CA	2.13	1.56	1.53
2	C	5	ALY	CH3-CH	2.50	1.55	1.50
2	C	5	ALY	CE-NZ	3.10	1.53	1.46
2	F	5	ALY	CH3-CH	3.20	1.57	1.50
2	F	8	ALY	CE-NZ	3.42	1.54	1.46
2	F	8	ALY	CH3-CH	3.52	1.57	1.50
2	C	8	ALY	CE-NZ	3.69	1.54	1.46
2	C	8	ALY	CH3-CH	6.33	1.63	1.50

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	8	ALY	CH3-CH-NZ	-8.91	103.44	116.28
2	F	5	ALY	OH-CH-CH3	-8.05	107.25	122.07
2	F	8	ALY	CH3-CH-NZ	-7.65	105.25	116.28
2	C	5	ALY	OH-CH-CH3	-7.00	109.19	122.07
2	F	5	ALY	CE-NZ-CH	-4.72	114.28	122.38
2	C	5	ALY	CE-NZ-CH	-3.54	116.30	122.38
2	F	5	ALY	O-C-CA	-2.78	118.27	125.72
2	F	8	ALY	CE-NZ-CH	-2.66	117.81	122.38
2	C	5	ALY	O-C-CA	-2.38	119.33	125.72
2	F	8	ALY	O-C-CA	-2.37	119.35	125.72
2	C	8	ALY	O-C-CA	-2.27	119.63	125.72
2	C	5	ALY	CH3-CH-NZ	3.30	121.04	116.28
2	C	5	ALY	OH-CH-NZ	3.87	128.80	121.71
2	F	8	ALY	OH-CH-NZ	6.23	133.11	121.71
2	C	8	ALY	OH-CH-NZ	7.26	134.99	121.71
2	F	5	ALY	OH-CH-NZ	7.98	136.30	121.71

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](#)

There are no ligands in this entry.

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	100/108 (92%)	0.01	1 (1%) 84 86	17, 24, 35, 44	0
1	B	100/108 (92%)	0.03	0 100 100	19, 27, 39, 45	0
1	D	101/108 (93%)	-0.03	1 (0%) 84 86	19, 25, 37, 53	0
1	E	101/108 (93%)	0.02	2 (1%) 68 72	17, 25, 36, 56	0
2	C	4/11 (36%)	0.52	0 100 100	24, 25, 37, 38	0
2	F	4/11 (36%)	0.39	0 100 100	24, 24, 40, 45	0
All	All	410/454 (90%)	0.02	4 (0%) 84 86	17, 26, 38, 56	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	22	SER	3.8
1	A	59	MET	2.7
1	D	59	MET	2.3
1	E	59	MET	2.2

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. 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	ALY	C	5	12/13	0.94	0.08	-	18,24,31,37	0
2	ALY	F	8	12/13	0.90	0.11	-	15,24,35,44	0
2	ALY	F	5	12/13	0.91	0.09	-	18,24,36,38	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	ALY	C	8	12/13	0.90	0.10	-	16,25,32,43	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.