



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 12, 2018 – 07:06 PM EST

PDB ID : 6BM7
Title : Crystal structure of Trypanosoma brucei AdoMetDC/prozyme heterodimer in complex with pyrimidineamine inhibitor UTSAM568
Authors : Volkov, O.A.; Chen, Z.; Phillips, M.A.
Deposited on : 2017-11-13
Resolution : 2.98 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030736
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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-20030736

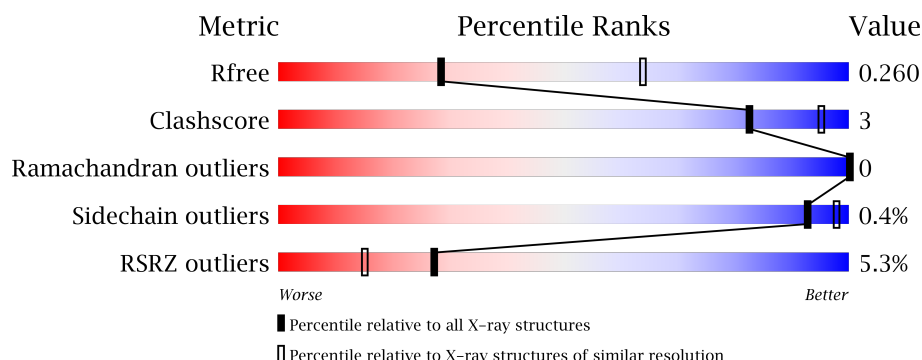
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.98 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}	100719	2168 (3.00-2.96)
Clashscore	112137	2535 (3.00-2.96)
Ramachandran outliers	110173	2451 (3.00-2.96)
Sidechain outliers	110143	2454 (3.00-2.96)
RSRZ outliers	101464	2192 (3.00-2.96)

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	85	<div> <div>6%</div> <div>87%</div> <div>9%</div> </div>
1	C	85	<div> <div>6%</div> <div>84%</div> <div>9%</div> <div>7%</div> </div>
2	B	285	<div> <div>6%</div> <div>89%</div> <div>6%</div> <div>5%</div> </div>
2	D	285	<div> <div>4%</div> <div>87%</div> <div>8%</div> <div>5%</div> </div>
3	E	325	<div> <div>6%</div> <div>81%</div> <div>9%</div> <div>10%</div> </div>

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Mol	Chain	Length	Quality of chain
3	F	325	

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
7	PGE	E	402	-	-	-	X
8	PG4	F	402	-	-	-	X

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 20224 atoms, of which 9907 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 S-adenosylmethionine decarboxylase beta chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	77	Total	C	H	N	O	S	0	0	0
			1217	391	600	105	117	4			
1	C	79	Total	C	H	N	O	S	0	0	0
			1263	403	626	111	119	4			

- Molecule 2 is a protein called S-adenosylmethionine decarboxylase alpha chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	271	Total	C	H	N	O	S	0	0	0
			4214	1386	2047	354	411	16			
2	D	271	Total	C	H	N	O	S	0	0	0
			4215	1386	2048	354	411	16			

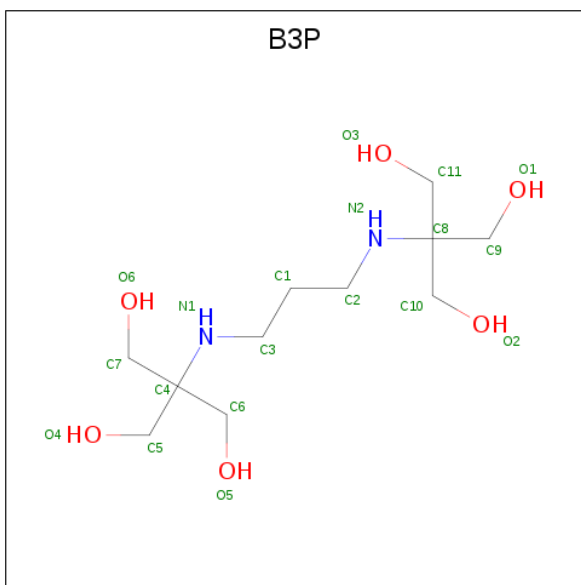
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	86	PYR	-	modified residue	UNP Q587A7
D	86	PYR	-	modified residue	UNP Q587A7

- Molecule 3 is a protein called Inactive S-adenosylmethionine decarboxylase prozyme.

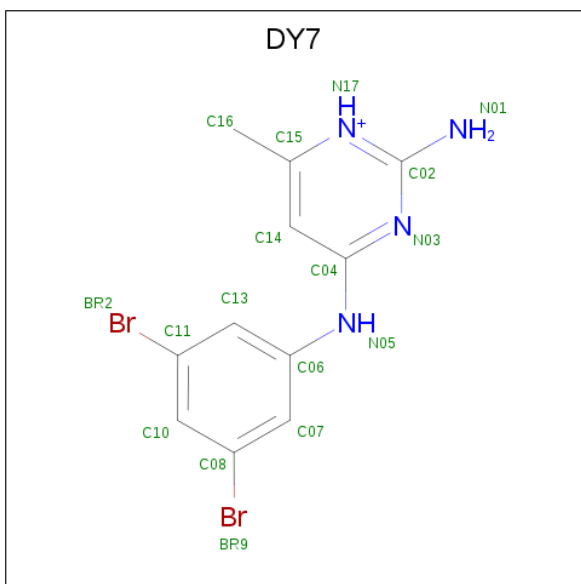
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	E	291	Total	C	H	N	O	S	0	0	0
			4536	1452	2227	405	438	14			
3	F	289	Total	C	H	N	O	S	0	0	0
			4511	1445	2215	403	434	14			

- Molecule 4 is 2-[3-(2-HYDROXY-1,1-DIHYDROXYMETHYL-ETHYLAMINO)-PROPYL AMINO]-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: B3P) (formula: C₁₁H₂₆N₂O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	H	N	O	0
			45	11	26	2	6	
4	D	1	Total	C	H	N	O	0
			45	11	26	2	6	

- Molecule 5 is 2-amino-4-[(3,5-dibromophenyl)amino]-6-methylpyrimidin-1-ium (three-letter code: DY7) (formula: C₁₁H₁₁Br₂N₄).



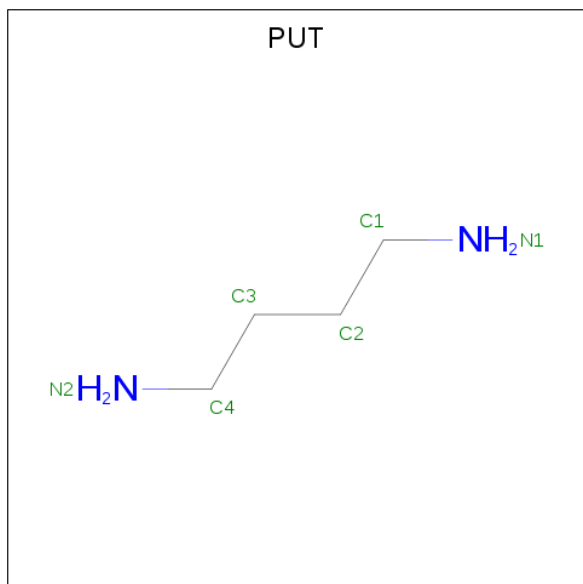
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	Br	C	H	N	0
			28	2	11	11	4	

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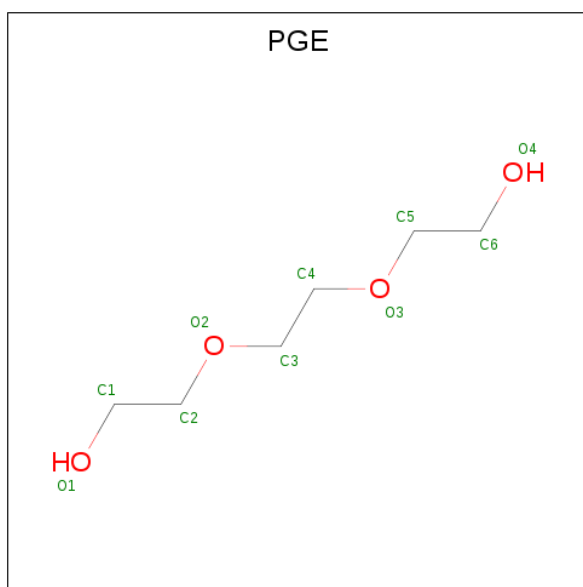
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	D	1	Total	Br	C	H	N	0	0
			28	2	11	11	4		

- Molecule 6 is 1,4-DIAMINOBTANE (three-letter code: PUT) (formula: $C_4H_{12}N_2$).



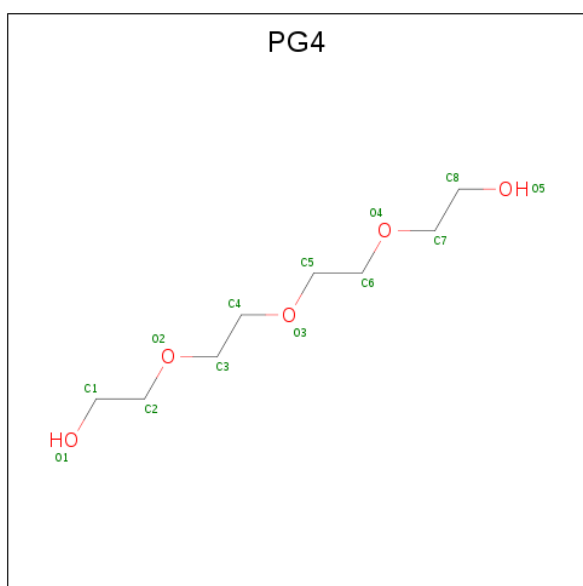
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	E	1	Total	C	H	N	0	0
			18	4	12	2		
6	F	1	Total	C	H	N	0	0
			18	4	12	2		

- Molecule 7 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	E	1	Total	C	H	O	0	0
			24	6	14	4		
7	E	1	Total	C	H	O	0	0
			24	6	14	4		

- Molecule 8 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	F	1	Total	C	H	O	0	0
			31	8	18	5		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	1	Total 1	O 1	0	0
9	B	1	Total 1	O 1	0	0
9	C	1	Total 1	O 1	0	0
9	D	2	Total 2	O 2	0	0
9	F	2	Total 2	O 2	0	0

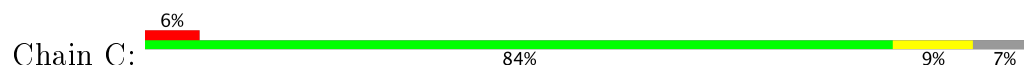
3 Residue-property plots [i](#)

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

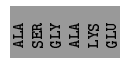
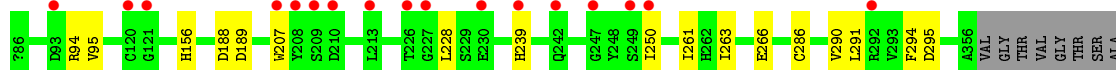
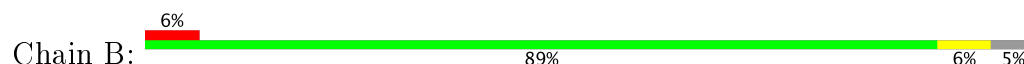
- Molecule 1: S-adenosylmethionine decarboxylase beta chain



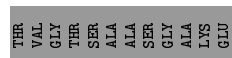
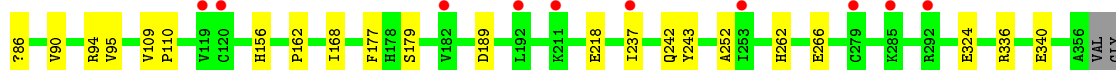
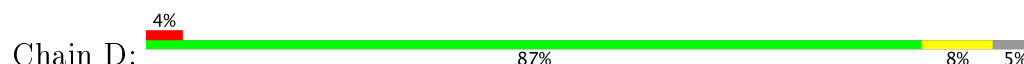
- Molecule 1: S-adenosylmethionine decarboxylase beta chain



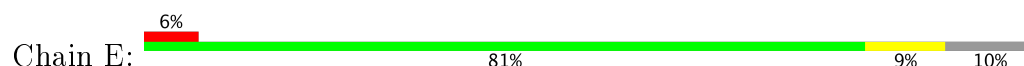
- Molecule 2: S-adenosylmethionine decarboxylase alpha chain

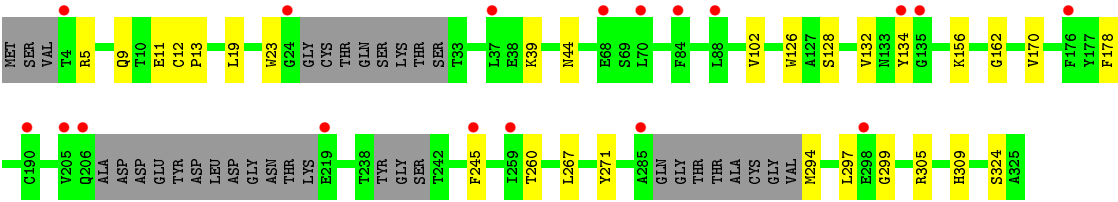


- Molecule 2: S-adenosylmethionine decarboxylase alpha chain

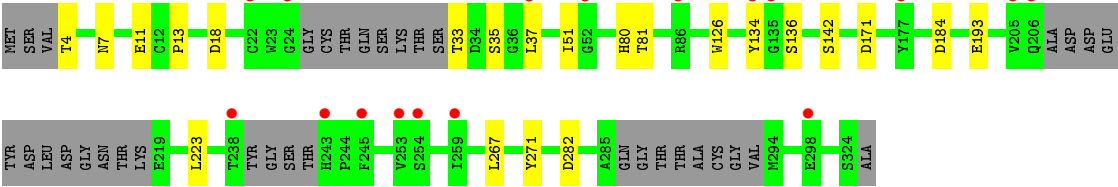
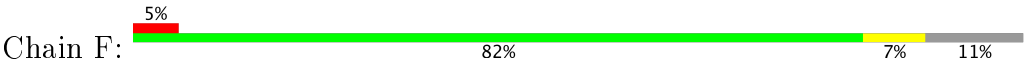


- Molecule 3: Inactive S-adenosylmethionine decarboxylase prozyme





● Molecule 3: Inactive S-adenosylmethionine decarboxylase prozyme



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	81.09Å 96.25Å 98.84Å 90.00° 102.43° 90.00°	Depositor
Resolution (Å)	48.26 – 2.98 48.26 – 2.98	Depositor EDS
% Data completeness (in resolution range)	84.6 (48.26-2.98) 98.0 (48.26-2.98)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.61 (at 2.96Å)	Xtriage
Refinement program	PHENIX 1.10.1 _2155	Depositor
R, R_{free}	0.200 , 0.255 0.206 , 0.260	Depositor DCC
R_{free} test set	1498 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	54.8	Xtriage
Anisotropy	0.149	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 38.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	20224	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.60% 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: PGE, PYR, DY7, B3P, PG4, PUT

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	0.27	0/629	0.42	0/847
1	C	0.25	0/649	0.43	0/872
2	B	0.27	0/2229	0.44	0/3038
2	D	0.28	0/2229	0.45	0/3038
3	E	0.26	0/2358	0.45	0/3191
3	F	0.26	0/2345	0.44	0/3174
All	All	0.27	0/10439	0.44	0/14160

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	617	600	600	5	0
1	C	637	626	626	6	0
2	B	2167	2047	2050	13	0
2	D	2167	2048	2050	15	0
3	E	2309	2227	2227	19	0
3	F	2296	2215	2215	15	0
4	B	19	26	25	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	19	26	26	0	0
5	B	17	11	0	0	0
5	D	17	11	0	1	0
6	E	6	12	12	0	0
6	F	6	12	12	3	0
7	E	20	28	28	0	0
8	F	13	18	18	1	0
9	A	1	0	0	0	0
9	B	1	0	0	0	0
9	C	1	0	0	0	0
9	D	2	0	0	0	0
9	F	2	0	0	0	0
All	All	10317	9907	9889	58	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 (58) 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)
2:D:86:PYR:H32	2:D:262:HIS:CE1	2.26	0.70
1:C:75:ASN:OD1	1:C:76:GLU:N	2.33	0.61
1:C:10:LEU:HD11	3:E:19:LEU:HG	1.82	0.61
2:B:94:ARG:NH2	4:B:401:B3P:O5	2.37	0.58
2:D:218:GLU:N	2:D:218:GLU:OE1	2.36	0.58
1:A:82:VAL:HG23	2:B:239:HIS:ND1	2.21	0.56
3:E:156:LYS:NZ	3:E:162:GLY:O	2.42	0.52
1:C:50:LEU:HD21	2:D:95:VAL:HG22	1.91	0.52
3:E:134:TYR:HH	3:E:245:PHE:HE1	1.59	0.51
3:F:184:ASP:HA	8:F:402:PG4:H22	1.93	0.51
3:E:13:PRO:HG2	3:E:19:LEU:HD13	1.94	0.50
1:A:82:VAL:HG23	2:B:239:HIS:CE1	2.47	0.49
3:F:33:THR:O	3:F:142:SER:OG	2.26	0.49
2:D:94:ARG:NH1	2:D:189:ASP:OD2	2.45	0.49
3:F:267:LEU:O	3:F:271:TYR:N	2.43	0.48
2:D:242:GLN:NE2	2:D:243:TYR:O	2.45	0.47
2:B:291:LEU:O	2:B:295:ASP:N	2.39	0.47
1:C:79:ARG:HB2	2:D:90:VAL:HB	1.97	0.47
3:F:37:LEU:HD11	3:F:134:TYR:HB3	1.97	0.47
3:E:267:LEU:HD11	3:E:297:LEU:HD13	1.97	0.46
2:B:94:ARG:NH1	2:B:189:ASP:OD2	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:VAL:HG23	2:B:239:HIS:HD1	1.81	0.44
3:E:11:GLU:HA	3:F:4:THR:O	2.18	0.44
3:E:128:SER:HA	3:E:178:PHE:O	2.17	0.44
2:B:263:ILE:HD13	2:B:294:PHE:HE1	1.81	0.44
3:E:44:ASN:HB2	3:E:126:TRP:HB3	2.00	0.44
2:D:266:GLU:OE2	5:D:402:DY7:N17	2.51	0.44
3:E:132:VAL:HG12	3:E:134:TYR:CE2	2.53	0.44
3:E:267:LEU:O	3:E:271:TYR:N	2.42	0.43
2:D:109:VAL:HB	2:D:110:PRO:HD3	2.01	0.43
3:F:81:THR:OG1	3:F:223:LEU:HD13	2.19	0.43
1:C:22:PRO:O	1:C:23:LYS:HB2	2.18	0.43
2:D:340:GLU:OE1	3:E:305:ARG:NH1	2.32	0.43
2:D:324:GLU:OE1	2:D:324:GLU:N	2.44	0.43
3:E:19:LEU:HD12	3:E:23:TRP:CG	2.54	0.43
2:D:237:ILE:O	2:D:252:ALA:HA	2.19	0.43
1:A:73:GLU:HG2	2:B:239:HIS:CD2	2.54	0.42
3:E:299:GLY:C	3:E:324:SER:HB2	2.39	0.42
2:B:250:ILE:CG2	2:B:261:ILE:HB	2.50	0.42
2:D:177:PHE:CZ	2:D:179:SER:HB3	2.55	0.42
3:F:193:GLU:HB3	6:F:401:PUT:HN22	1.83	0.42
2:D:162:PRO:HG2	3:E:170:VAL:HB	2.02	0.42
3:E:39:LYS:NZ	3:E:102:VAL:O	2.47	0.42
3:E:5:ARG:NH1	3:F:11:GLU:HG3	2.34	0.42
3:E:260:THR:HG22	3:E:294:MET:HG2	2.01	0.42
2:D:336:ARG:HB2	3:E:309:HIS:HB2	2.02	0.41
3:F:35:SER:HB2	3:F:134:TYR:CZ	2.55	0.41
3:F:126:TRP:CE2	6:F:401:PUT:H32	2.55	0.41
3:F:13:PRO:HB3	3:F:18:ASP:HB2	2.02	0.41
3:E:9:GLN:NE2	3:F:7:ASN:OD1	2.53	0.41
1:A:36:GLU:HA	2:B:95:VAL:O	2.21	0.41
3:F:51:ILE:HG22	3:F:80:HIS:CE1	2.56	0.41
2:B:286:CYS:O	2:B:290:VAL:HG23	2.21	0.41
2:B:207:TRP:CH2	2:B:228:LEU:HD11	2.56	0.41
2:B:266:GLU:N	2:B:266:GLU:OE1	2.50	0.41
3:F:136:SER:HB2	3:F:171:ASP:HA	2.02	0.40
1:C:8:LEU:HD22	2:D:168:ILE:HD13	2.03	0.40
3:F:282:ASP:OD1	6:F:401:PUT:N2	2.54	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	73/85 (86%)	71 (97%)	2 (3%)	0	100	100
1	C	75/85 (88%)	71 (95%)	4 (5%)	0	100	100
2	B	268/285 (94%)	256 (96%)	12 (4%)	0	100	100
2	D	268/285 (94%)	264 (98%)	4 (2%)	0	100	100
3	E	281/325 (86%)	276 (98%)	5 (2%)	0	100	100
3	F	279/325 (86%)	275 (99%)	4 (1%)	0	100	100
All	All	1244/1390 (90%)	1213 (98%)	31 (2%)	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	68/76 (90%)	68 (100%)	0	100	100
1	C	70/76 (92%)	70 (100%)	0	100	100
2	B	243/251 (97%)	241 (99%)	2 (1%)	85	95
2	D	243/251 (97%)	242 (100%)	1 (0%)	93	98
3	E	254/281 (90%)	253 (100%)	1 (0%)	93	98
3	F	253/281 (90%)	253 (100%)	0	100	100
All	All	1131/1216 (93%)	1127 (100%)	4 (0%)	93	98

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	156	HIS
2	B	188	ASP
2	D	156	HIS
3	E	12	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

9 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	B3P	B	401	-	18,18,18	2.42	4 (22%)	21,23,23	1.59	6 (28%)
5	DY7	B	402	-	18,18,18	2.65	8 (44%)	24,25,25	2.56	7 (29%)
4	B3P	D	401	-	18,18,18	2.97	4 (22%)	21,23,23	1.66	6 (28%)
5	DY7	D	402	-	18,18,18	2.56	8 (44%)	24,25,25	2.54	7 (29%)
6	PUT	E	401	-	5,5,5	0.28	0	4,4,4	0.58	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	PGE	E	402	-	9,9,9	0.52	0	8,8,8	0.31	0
7	PGE	E	403	-	9,9,9	0.48	0	8,8,8	0.40	0
6	PUT	F	401	-	5,5,5	0.16	0	4,4,4	0.50	0
8	PG4	F	402	-	12,12,12	0.51	0	11,11,11	0.38	0

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	B3P	B	401	-	-	0/28/28/28	0/0/0/0
5	DY7	B	402	-	-	0/4/4/4	0/2/2/2
4	B3P	D	401	-	-	0/28/28/28	0/0/0/0
5	DY7	D	402	-	-	0/4/4/4	0/2/2/2
6	PUT	E	401	-	-	0/3/3/3	0/0/0/0
7	PGE	E	402	-	-	0/7/7/7	0/0/0/0
7	PGE	E	403	-	-	0/7/7/7	0/0/0/0
6	PUT	F	401	-	-	0/3/3/3	0/0/0/0
8	PG4	F	402	-	-	0/10/10/10	0/0/0/0

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	402	DY7	C07-C06	2.38	1.43	1.39
4	B	401	B3P	C6-C4	2.43	1.56	1.53
5	B	402	DY7	C07-C06	2.45	1.43	1.39
5	D	402	DY7	C04-N05	2.49	1.43	1.38
4	B	401	B3P	C10-C8	2.88	1.56	1.53
5	B	402	DY7	C04-N05	3.03	1.43	1.38
5	D	402	DY7	C06-N05	3.23	1.47	1.40
5	D	402	DY7	BR2-C11	3.28	1.97	1.90
5	D	402	DY7	BR9-C08	3.37	1.97	1.90
5	B	402	DY7	C15-N17	3.52	1.41	1.35
5	B	402	DY7	C06-N05	3.53	1.48	1.40
5	B	402	DY7	C02-N03	3.61	1.37	1.33
5	B	402	DY7	BR2-C11	3.61	1.97	1.90
5	D	402	DY7	C02-N03	3.72	1.38	1.33
5	D	402	DY7	C15-N17	3.78	1.41	1.35
5	B	402	DY7	BR9-C08	3.78	1.98	1.90
4	D	401	B3P	C10-C8	3.81	1.57	1.53
4	D	401	B3P	C6-C4	4.04	1.58	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	401	B3P	C7-C4	5.45	1.59	1.53
5	B	402	DY7	C02-N17	5.58	1.44	1.36
5	D	402	DY7	C02-N17	5.67	1.44	1.36
4	B	401	B3P	C11-C8	6.69	1.61	1.53
4	D	401	B3P	C11-C8	6.86	1.61	1.53
4	D	401	B3P	C7-C4	8.09	1.62	1.53

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	402	DY7	C02-N17-C15	-5.78	115.79	121.10
5	D	402	DY7	C02-N17-C15	-5.58	115.97	121.10
5	D	402	DY7	C16-C15-N17	-4.81	110.43	117.18
5	B	402	DY7	C16-C15-N17	-4.76	110.49	117.18
5	B	402	DY7	C14-C04-N03	-4.72	112.97	123.41
5	D	402	DY7	C14-C04-N03	-4.56	113.32	123.41
5	B	402	DY7	N01-C02-N17	-3.22	114.41	117.84
4	D	401	B3P	C3-N1-C4	-3.01	111.75	116.12
4	D	401	B3P	C2-N2-C8	-2.93	111.87	116.12
5	D	402	DY7	N01-C02-N17	-2.77	114.89	117.84
4	B	401	B3P	C2-N2-C8	-2.51	112.48	116.12
4	D	401	B3P	O5-C6-C4	-2.36	106.66	111.54
4	B	401	B3P	C3-N1-C4	-2.35	112.71	116.12
4	B	401	B3P	O4-C5-C4	-2.30	106.78	111.54
4	D	401	B3P	O1-C9-C8	-2.28	106.83	111.54
4	B	401	B3P	O2-C10-C8	-2.23	106.93	111.54
4	B	401	B3P	O5-C6-C4	-2.23	106.93	111.54
4	D	401	B3P	O2-C10-C8	-2.22	106.95	111.54
4	B	401	B3P	C7-C4-C5	-2.19	105.58	110.09
4	D	401	B3P	O4-C5-C4	-2.01	107.38	111.54
5	B	402	DY7	N01-C02-N03	3.11	124.06	120.26
5	D	402	DY7	N01-C02-N03	3.21	124.19	120.26
5	B	402	DY7	C02-N03-C04	3.43	123.53	116.58
5	D	402	DY7	C02-N03-C04	3.54	123.75	116.58
5	D	402	DY7	C16-C15-C14	5.28	130.25	124.55
5	B	402	DY7	C16-C15-C14	5.43	130.42	124.55

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	401	B3P	1	0
5	D	402	DY7	1	0
6	F	401	PUT	3	0
8	F	402	PG4	1	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 ⓘ

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	77/85 (90%)	0.50	1 (1%) 77 57	32, 61, 82, 93	0
1	C	79/85 (92%)	0.62	5 (6%) 21 11	35, 63, 91, 101	0
2	B	270/285 (94%)	0.48	17 (6%) 21 11	29, 50, 75, 100	0
2	D	270/285 (94%)	0.47	10 (3%) 42 25	28, 54, 77, 86	0
3	E	291/325 (89%)	0.52	18 (6%) 21 11	25, 50, 86, 98	0
3	F	289/325 (88%)	0.46	17 (5%) 23 12	25, 44, 79, 106	0
All	All	1276/1390 (91%)	0.49	68 (5%) 27 15	25, 51, 82, 106	0

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	F	24	GLY	5.5
1	C	74	PHE	4.4
3	F	22	CYS	4.2
2	B	120	CYS	4.1
3	E	206	GLN	4.1
3	F	134	TYR	4.0
3	E	205	VAL	3.9
3	F	37	LEU	3.5
2	D	292	ARG	3.4
2	D	120	CYS	3.4
3	E	37	LEU	3.3
3	F	238	THR	3.3
2	B	208	TYR	3.3
3	E	134	TYR	3.2
3	F	206	GLN	3.1
3	F	245	PHE	3.1
3	F	205	VAL	3.0
1	C	23	LYS	2.9
3	E	190	CYS	2.9

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Mol	Chain	Res	Type	RSRZ
3	E	24	GLY	2.9
1	C	83	LEU	2.8
3	F	259	ILE	2.8
2	B	226	THR	2.8
2	B	209	SER	2.8
2	D	192	LEU	2.8
2	B	213	LEU	2.7
3	E	259	ILE	2.7
3	E	135	GLY	2.7
3	F	86	ARG	2.7
2	D	182	VAL	2.6
3	F	253	VAL	2.6
2	B	207	TRP	2.6
2	D	119	VAL	2.6
3	F	243	HIS	2.6
3	E	84	PHE	2.6
2	D	237	ILE	2.5
2	D	279	CYS	2.5
1	A	82	VAL	2.5
2	B	210	ASP	2.5
2	B	242	GLN	2.5
1	C	57	VAL	2.5
2	B	292	ARG	2.5
1	C	76	GLU	2.5
3	E	245	PHE	2.4
3	E	285	ALA	2.4
3	F	298	GLU	2.4
3	E	70	LEU	2.4
2	B	121	GLY	2.4
3	E	298	GLU	2.3
2	B	239	HIS	2.3
2	D	285	LYS	2.3
3	E	219	GLU	2.3
3	F	177	TYR	2.3
3	F	52	GLY	2.2
3	F	254	SER	2.2
2	D	253	ILE	2.2
3	E	4	THR	2.2
2	B	227	GLY	2.2
2	B	230	GLU	2.1
2	B	249	SER	2.1
3	E	68	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
2	B	250	ILE	2.1
3	F	135	GLY	2.1
3	E	176	PHE	2.1
2	B	247	GLY	2.1
2	B	93	ASP	2.1
2	D	211	LYS	2.0
3	E	88	LEU	2.0

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
7	PGE	E	402	10/10	0.64	0.37	3.22	46,69,77,82	0
8	PG4	F	402	13/13	0.68	0.33	2.65	41,58,72,72	0
5	DY7	B	402	17/17	0.86	0.30	0.97	57,76,99,122	0
5	DY7	D	402	17/17	0.81	0.32	0.91	60,85,107,133	0
6	PUT	E	401	6/6	0.95	0.25	0.91	35,46,57,57	0
6	PUT	F	401	6/6	0.97	0.22	0.77	25,32,48,48	0
4	B3P	B	401	19/19	0.93	0.21	0.19	35,55,72,87	0
7	PGE	E	403	10/10	0.61	0.20	0.18	65,88,103,103	0
4	B3P	D	401	19/19	0.85	0.21	-0.49	49,71,93,100	0

6.5 Other polymers [i](#)

There are no such residues in this entry.