Full wwPDB X-ray Structure Validation Report (i)

Jun 16, 2024 – 01:32 PM EDT

PDB ID	:	4X9E
Title	:	DEOXYGUANOSINETRIPHOSPHATE TRIPHOSPHOHYDROLASE from Escherichia coli with two DNA effector molecules
Authors	:	Singh, D.; Gawel, D.; Itsko, M.; Krahn, J.M.; London, R.E.; Schaaper, R.M.
Deposited on	:	2014-12-11
Resolution	:	3.10 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

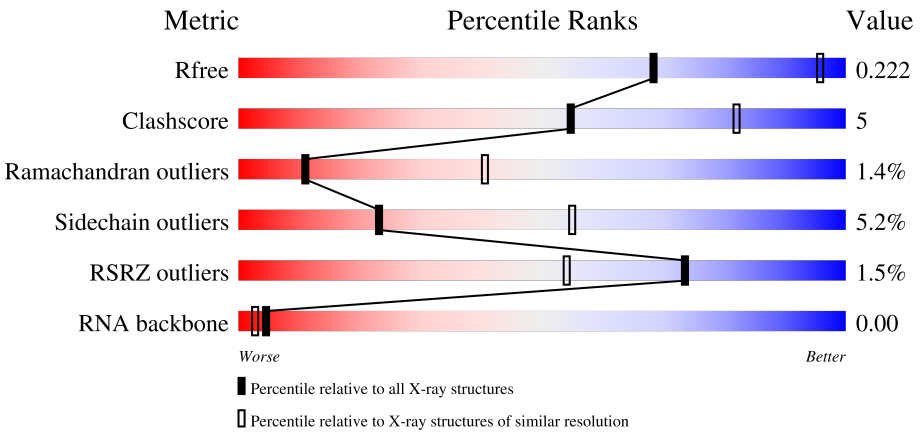
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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}	130704	1094 (3.10-3.10)
Clashscore	141614	1184 (3.10-3.10)
Ramachandran outliers	138981	1141 (3.10-3.10)
Sidechain outliers	138945	1141 (3.10-3.10)
RSRZ outliers	127900	1067 (3.10-3.10)
RNA backbone	3102	1116 (3.40-2.80)

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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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 <=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	505	2% 78% 19% • •
1	B	505	3% 81% 16% • •
1	C	505	.% 82% 16% • •
1	D	505	2% 82% 16% • •

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Mol	Chain	Length	Quality of chain
1	E	505	82% 15%
1	F	505	% 84% 13%
2	G	3	33% 67%
2	H	3	67% 33%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 24471 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 Deoxyguanosinetriphosphate triphosphohydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	499	4092	2618	728	730	16	0	0	0
1	B	495	3983	2562	709	696	16	0	0	0
1	C	500	4107	2627	730	733	17	0	0	0
1	D	501	4067	2606	713	732	16	0	0	0
1	E	496	4075	2609	715	735	16	0	0	0
1	F	497	4034	2588	715	715	16	0	0	0

- Molecule 2 is a RNA chain called RNA (5'-R(P*CP*CP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	G	3	58	27	9	19	3	0	0	0
2	H	3	53	27	9	15	2	0	0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
3	A	1	1	1	0	0
3	B	1	1	1	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Deoxyguanosinetriphosphate triphosphohydrolase

Chain A: 2% 78% 19% ••

M1 A2 F6 T24 E25 H26 E27 I31 S34 R38 I45 R46 R47 L48 Q49 T52 Q53 V54 F55 P56 L57 GLU ARG N60 A61 T65 R66 L67 T68 V73 K82 E83 I84 R87 E95 E107 C114 L115 M116 I119 A145 E146 S147 Q148

T151 V157 A158 A159 L160 R161 L162 E166 L169 N170 E171 L172 R173 R174 K175 L192 V193 L196 M197 R198 L210 W217 Y228 L229 I243 R257 M264 E265 I270 S271 Y272 C273 V274 I285 F286 E289 L295 G300 Q301 H302 E303 K304 G305

S309 L310 V311 V312 E313 N314 A315 W316 E317 K318 S319 R320 S321 N322 SER LEU SER ARG S327 T328 L341 L344 F352 L356 P357 F360 A361 G362 D371 A372 S373 E374 C375 N384 V385 K388 H389 H393 E400 I413 S423 R436 F437 P438 R442

S448 T449 R450 H451 L471 W472 E473 Y474 R479 I485 L500 Q505

• Molecule 1: Deoxyguanosinetriphosphate triphosphohydrolase

Chain B: 3% 81% 16% ••

MET A2 Q3 T24 E27 I31 R38 I45 R46 R47 L48 Q49 T52 Q53 V54 F55 P56 LEU GLU ARG N60 L67 T68 I84 F106 M116 I119 L140 A145 GLU SER Q148 T151 D152 V157 A158 A159 L160 R161 L162 G165 E166 E167 P168

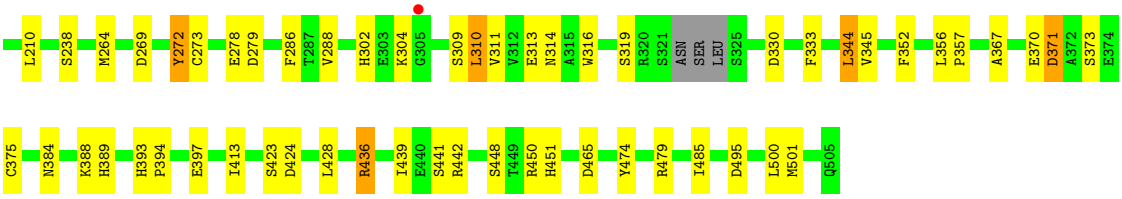
L169 N170 E171 L172 K175 L192 V193 R198 L210 K211 A216 W217 Y228 L229 P233 G234 Y235 I243 R257 L260 M264 E265 I270 S271 Y272 C273 V274 I285 F286 L295 W299 G300 Q301 H302 S309 L310 V311 V312 E313 N314 S321 ASN SER

LEU SER R326 L341 L344 F352 L356 P357 D371 A372 S373 E374 C375 N384 V385 K388 H389 H393 E400 I413 S423 V434 P438 R442 S448 H451 K461 L471 W472 E473 Y474 R479 I485 L500 Q505

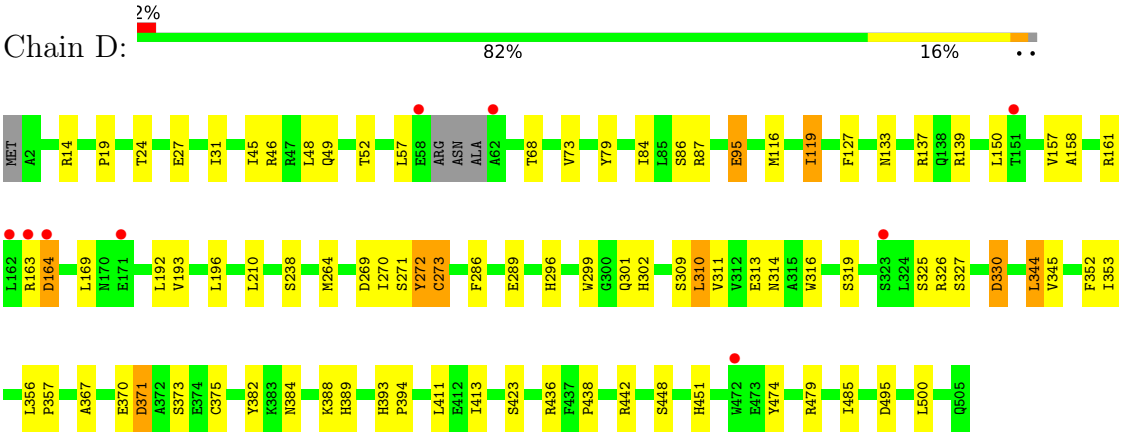
• Molecule 1: Deoxyguanosinetriphosphate triphosphohydrolase

Chain C: % 82% 16% ••

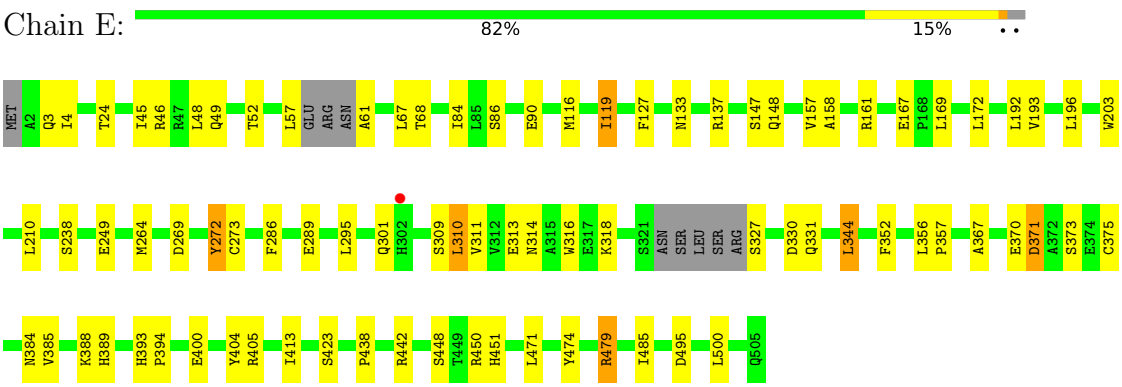
M1 T24 I45 R46 R47 L48 Q49 T52 L57 E58 R59 ASN ALA A62 V63 R64 Y79 I84 L85 S86 R87 K92 M116 I119 N133 R137 S147 Q148 P149 V157 A158 R161 L162 R163 D164 L169 Q178 H182 L192 V193 L196



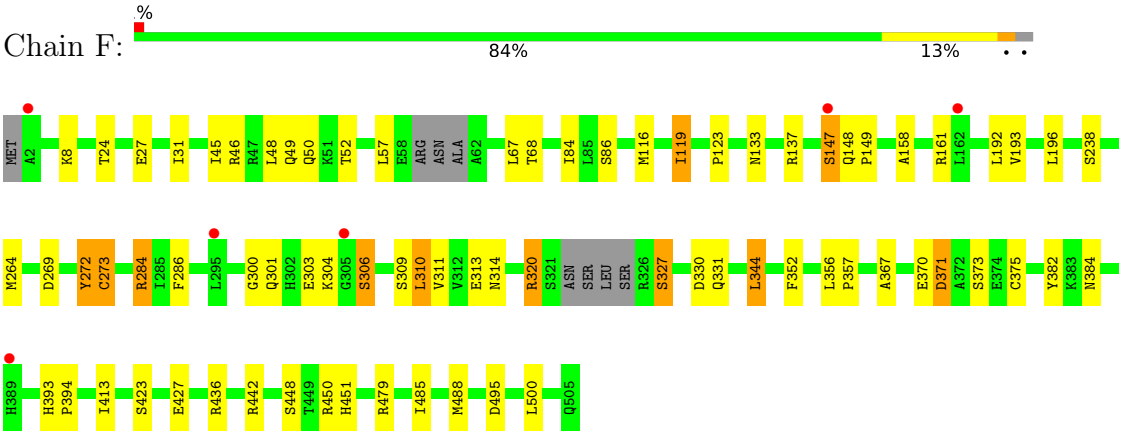
• Molecule 1: Deoxyguanosinetriphosphate triphosphohydrolase



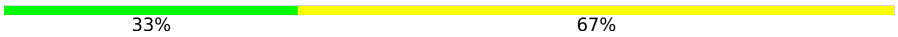
• Molecule 1: Deoxyguanosinetriphosphate triphosphohydrolase



• Molecule 1: Deoxyguanosinetriphosphate triphosphohydrolase



• Molecule 2: RNA (5'-R(P*CP*CP*C)-3')

Chain G: 

- Molecule 2: RNA (5'-R(P*CP*CP*C)-3')

Chain H:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	136.80Å 159.80Å 190.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.63 – 3.10 49.63 – 3.09	Depositor EDS
% Data completeness (in resolution range)	99.7 (49.63-3.10) 95.2 (49.63-3.09)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.29 (at 3.07Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.167 , 0.223 0.168 , 0.222	Depositor DCC
R_{free} test set	3881 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å^2)	87.0	Xtriage
Anisotropy	0.573	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 69.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	24471	wwPDB-VP
Average B, all atoms (Å^2)	108.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.46% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: MG

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	0.47	0/4193	0.60	0/5674
1	B	0.42	0/4083	0.57	0/5532
1	C	0.53	0/4208	0.61	0/5692
1	D	0.47	0/4169	0.59	0/5649
1	E	0.53	0/4176	0.61	0/5651
1	F	0.48	0/4134	0.59	0/5599
2	G	1.61	1/63 (1.6%)	0.71	0/92
2	H	0.78	0/58	0.71	0/87
All	All	0.49	1/25084 (0.0%)	0.60	0/33976

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	1	C	OP3-P	-10.19	1.49	1.61

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	4092	0	4021	57	1
1	B	3983	0	3877	44	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	4107	0	4033	41	1
1	D	4067	0	3945	45	1
1	E	4075	0	3997	43	0
1	F	4034	0	3941	34	1
2	G	58	0	34	1	0
2	H	53	0	32	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
All	All	24471	0	23880	249	2

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 (249) 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:E:61:ALA:HB2	1:F:50:GLN:HG3	1.52	0.91
1:E:371:ASP:HB3	1:E:373:SER:H	1.48	0.79
1:F:371:ASP:HB3	1:F:373:SER:H	1.48	0.79
1:B:371:ASP:HB3	1:B:373:SER:H	1.49	0.78
1:D:371:ASP:HB3	1:D:373:SER:H	1.48	0.78
1:C:371:ASP:HB3	1:C:373:SER:H	1.48	0.76
1:A:371:ASP:HB3	1:A:373:SER:H	1.51	0.74
1:C:442:ARG:CZ	1:E:127:PHE:HZ	2.01	0.73
1:C:310:LEU:O	1:C:314:ASN:HB3	1.90	0.72
1:D:310:LEU:O	1:D:314:ASN:HB3	1.90	0.72
1:A:289:GLU:HG3	1:A:316:TRP:HH2	1.55	0.72
1:E:310:LEU:O	1:E:314:ASN:HB3	1.91	0.71
1:F:310:LEU:O	1:F:314:ASN:HB3	1.91	0.71
1:F:284:ARG:HA	1:F:284:ARG:HH11	1.56	0.70
1:A:57:LEU:HD22	1:A:65:THR:HG22	1.73	0.69
1:B:413:ILE:HG21	1:B:500:LEU:HD13	1.75	0.69
1:D:119:ILE:HG23	1:D:192:LEU:HD23	1.76	0.68
1:A:310:LEU:O	1:A:314:ASN:HB3	1.93	0.68
1:F:367:ALA:HB3	1:F:370:GLU:HB3	1.75	0.68
1:B:310:LEU:O	1:B:314:ASN:HB3	1.93	0.67
1:F:119:ILE:HG23	1:F:192:LEU:HD23	1.77	0.67
1:E:367:ALA:HB3	1:E:370:GLU:HB3	1.76	0.67
1:A:68:THR:HG21	1:B:46:ARG:HG2	1.77	0.67
1:C:367:ALA:HB3	1:C:370:GLU:HB3	1.76	0.67
1:E:119:ILE:HG23	1:E:192:LEU:HD23	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:119:ILE:HG23	1:C:192:LEU:HD23	1.78	0.65
1:D:299:TRP:HD1	1:D:302:HIS:HD1	1.44	0.65
1:D:367:ALA:HB3	1:D:370:GLU:HB3	1.78	0.65
1:A:119:ILE:HG23	1:A:192:LEU:HD23	1.78	0.64
1:D:161:ARG:HD2	1:D:163:ARG:HH22	1.61	0.64
1:B:119:ILE:HG23	1:B:192:LEU:HD23	1.79	0.64
1:A:229:LEU:HD22	1:A:257:ARG:HD3	1.78	0.64
1:D:299:TRP:HD1	1:D:302:HIS:ND1	1.96	0.64
1:A:413:ILE:HG21	1:A:500:LEU:HD13	1.80	0.63
1:A:45:ILE:O	1:A:48:LEU:HB2	2.00	0.62
1:A:46:ARG:HG2	1:B:68:THR:HG21	1.81	0.62
1:D:164:ASP:OD1	1:D:164:ASP:N	2.32	0.62
1:B:400:GLU:OE1	1:D:442:ARG:NE	2.32	0.62
1:B:229:LEU:HD22	1:B:257:ARG:HD3	1.81	0.62
1:C:448:SER:HB2	1:C:451:HIS:CD2	2.37	0.60
1:A:82:LYS:HE2	1:A:107:GLU:OE1	2.01	0.60
1:C:288:VAL:HG12	1:C:316:TRP:HZ3	1.67	0.60
1:B:45:ILE:O	1:B:48:LEU:HB2	2.01	0.59
1:C:393:HIS:CD2	1:C:394:PRO:HD2	2.37	0.59
1:D:45:ILE:O	1:D:48:LEU:HB2	2.03	0.59
1:D:95:GLU:CD	1:D:95:GLU:H	2.05	0.59
1:D:289:GLU:HG3	1:D:316:TRP:HH2	1.67	0.58
1:E:158:ALA:HA	1:E:161:ARG:HH12	1.69	0.58
1:E:413:ILE:HG21	1:E:500:LEU:HD13	1.85	0.58
1:E:448:SER:HB2	1:E:451:HIS:CD2	2.39	0.58
1:D:87:ARG:HD3	1:D:353:ILE:HD12	1.86	0.57
1:F:413:ILE:HG21	1:F:500:LEU:HD13	1.86	0.57
1:A:158:ALA:HA	1:A:161:ARG:NH1	2.20	0.56
1:D:413:ILE:HG21	1:D:500:LEU:HD13	1.87	0.56
1:F:393:HIS:CD2	1:F:394:PRO:HD2	2.41	0.56
1:C:439:ILE:HD11	1:E:404:TYR:CD1	2.41	0.55
1:A:438:PRO:O	1:A:442:ARG:HG3	2.06	0.55
1:B:49:GLN:HB3	1:B:67:LEU:HD22	1.88	0.55
1:E:3:GLN:OE1	1:E:4:ILE:N	2.38	0.55
1:C:413:ILE:HG21	1:C:500:LEU:HD13	1.88	0.55
1:E:344:LEU:HD22	1:E:375:CYS:HB3	1.90	0.54
1:F:45:ILE:O	1:F:48:LEU:HB2	2.07	0.54
1:A:49:GLN:HB3	1:A:67:LEU:HD22	1.89	0.54
1:A:160:LEU:HD23	1:A:471:LEU:HD22	1.89	0.54
1:D:84:ILE:HD13	1:D:352:PHE:CD2	2.43	0.53
1:C:84:ILE:HD13	1:C:352:PHE:CD2	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:448:SER:HB2	1:D:451:HIS:CD2	2.44	0.53
1:C:424:ASP:OD1	1:C:436:ARG:NH2	2.42	0.53
1:E:84:ILE:HD13	1:E:352:PHE:CD2	2.44	0.53
1:E:45:ILE:O	1:E:48:LEU:HB2	2.08	0.53
1:C:64:ARG:NH1	1:C:279:ASP:OD1	2.42	0.52
1:A:145:ALA:O	1:A:174:ARG:HG2	2.10	0.52
1:A:175:LYS:HG2	1:A:217:TRP:CD1	2.44	0.52
1:F:84:ILE:HD13	1:F:352:PHE:CD2	2.44	0.52
1:A:303:GLU:O	1:A:305:GLY:N	2.42	0.52
1:C:45:ILE:O	1:C:48:LEU:HB2	2.07	0.52
1:B:448:SER:HB2	1:B:451:HIS:CD2	2.45	0.52
1:E:133:ASN:O	1:E:137:ARG:HB2	2.09	0.52
1:E:46:ARG:O	1:E:49:GLN:HG2	2.10	0.52
1:E:158:ALA:HA	1:E:161:ARG:NH1	2.25	0.52
1:C:158:ALA:HA	1:C:161:ARG:HH12	1.75	0.51
1:A:270:ILE:HG21	1:A:341:LEU:CD2	2.40	0.51
1:A:448:SER:HB2	1:A:451:HIS:CD2	2.46	0.51
1:E:314:ASN:O	1:E:318:LYS:HB2	2.09	0.51
1:B:160:LEU:HD23	1:B:471:LEU:HD22	1.91	0.51
1:B:175:LYS:HG2	1:B:217:TRP:CD1	2.46	0.51
1:C:393:HIS:CG	1:C:394:PRO:HD2	2.46	0.51
1:C:442:ARG:NH1	1:E:127:PHE:HZ	2.08	0.51
1:F:448:SER:HB2	1:F:451:HIS:CD2	2.44	0.51
1:B:158:ALA:HA	1:B:161:ARG:NH1	2.26	0.51
1:C:269:ASP:O	1:C:272:TYR:HB3	2.10	0.50
1:B:285:ILE:HG23	1:B:393:HIS:HD2	1.74	0.50
1:F:344:LEU:HD22	1:F:375:CYS:HB3	1.93	0.50
1:A:228:TYR:OH	1:A:265:GLU:OE1	2.28	0.50
1:B:270:ILE:HG21	1:B:341:LEU:CD2	2.41	0.50
1:F:269:ASP:O	1:F:272:TYR:HB3	2.11	0.50
1:F:303:GLU:O	1:F:306:SER:HB3	2.12	0.50
1:A:116:MET:HB2	1:A:119:ILE:HD12	1.94	0.50
1:C:147:SER:OG	1:C:148:GLN:N	2.43	0.50
1:E:393:HIS:CD2	1:E:394:PRO:HD2	2.47	0.49
1:A:27:GLU:O	1:A:31:ILE:HG13	2.12	0.49
1:A:289:GLU:HG3	1:A:316:TRP:CH2	2.42	0.49
1:D:46:ARG:O	1:D:49:GLN:HG2	2.12	0.49
1:F:427:GLU:CD	1:F:436:ARG:HH22	2.16	0.49
1:C:442:ARG:CZ	1:E:127:PHE:CZ	2.90	0.49
1:C:442:ARG:NE	1:E:400:GLU:OE1	2.39	0.49
1:D:158:ALA:HA	1:D:161:ARG:HH12	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:344:LEU:HD22	1:D:375:CYS:HB3	1.95	0.49
1:A:67:LEU:HD23	1:B:67:LEU:HD23	1.94	0.48
1:B:228:TYR:OH	1:B:265:GLU:OE1	2.30	0.48
1:B:84:ILE:HD13	1:B:352:PHE:CD2	2.48	0.48
1:C:46:ARG:O	1:C:49:GLN:HG2	2.13	0.48
1:C:344:LEU:HD22	1:C:375:CYS:HB3	1.95	0.48
1:D:158:ALA:HA	1:D:161:ARG:NH1	2.29	0.48
1:E:147:SER:OG	1:E:148:GLN:N	2.46	0.48
1:F:123:PRO:HG3	1:F:488:MET:O	2.14	0.48
1:B:270:ILE:HG21	1:B:341:LEU:HD23	1.95	0.48
1:C:158:ALA:HA	1:C:161:ARG:NH1	2.28	0.48
1:E:388:LYS:HB3	1:E:389:HIS:CD2	2.49	0.48
1:B:388:LYS:HB3	1:B:389:HIS:CD2	2.48	0.47
1:F:133:ASN:O	1:F:137:ARG:HB2	2.15	0.47
1:C:133:ASN:O	1:C:137:ARG:HB2	2.15	0.47
1:F:158:ALA:HA	1:F:161:ARG:HH12	1.79	0.47
1:E:269:ASP:O	1:E:272:TYR:HB3	2.14	0.47
1:D:269:ASP:O	1:D:272:TYR:HB3	2.13	0.47
1:D:393:HIS:CD2	1:D:394:PRO:HD2	2.50	0.47
1:B:27:GLU:O	1:B:31:ILE:HG13	2.15	0.47
1:B:169:LEU:HD23	1:B:169:LEU:HA	1.74	0.46
1:F:46:ARG:O	1:F:49:GLN:HG2	2.15	0.46
1:F:147:SER:HB3	1:F:148:GLN:H	1.50	0.46
1:A:172:LEU:HD23	1:A:471:LEU:HD12	1.96	0.46
1:A:273:CYS:SG	1:A:274:VAL:N	2.88	0.46
1:B:116:MET:HB2	1:B:119:ILE:HD12	1.96	0.46
1:E:289:GLU:HG3	1:E:316:TRP:HH2	1.81	0.46
1:A:73:VAL:HG12	1:A:114:CYS:HB3	1.98	0.46
1:D:192:LEU:HD12	1:D:196:LEU:HB2	1.98	0.46
1:D:388:LYS:HB3	1:D:389:HIS:CD2	2.51	0.46
1:F:393:HIS:CG	1:F:394:PRO:HD2	2.50	0.46
1:B:438:PRO:O	1:B:442:ARG:HG3	2.16	0.46
1:A:344:LEU:HD22	1:A:375:CYS:HB3	1.97	0.46
1:C:309:SER:O	1:C:313:GLU:HB3	2.16	0.46
1:A:84:ILE:HD13	1:A:352:PHE:CD2	2.52	0.45
1:D:289:GLU:HG3	1:D:316:TRP:CH2	2.49	0.45
1:C:157:VAL:HG21	1:C:474:TYR:CE2	2.52	0.45
1:A:388:LYS:HB3	1:A:389:HIS:CD2	2.51	0.45
1:F:116:MET:HB2	1:F:119:ILE:HD12	1.98	0.45
1:F:309:SER:O	1:F:313:GLU:HB3	2.17	0.45
1:D:133:ASN:O	1:D:137:ARG:HB2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:116:MET:HB2	1:E:119:ILE:HD12	1.98	0.45
1:A:356:LEU:N	1:A:357:PRO:HD2	2.32	0.45
1:F:158:ALA:HA	1:F:161:ARG:NH1	2.32	0.45
1:B:172:LEU:HD23	1:B:471:LEU:HD12	2.00	0.44
1:A:53:GLN:HG3	1:A:54:VAL:HG23	1.99	0.44
1:C:169:LEU:HD23	1:C:169:LEU:HA	1.76	0.44
1:C:178:GLN:O	1:C:182:HIS:HD2	2.00	0.44
1:D:116:MET:HB2	1:D:119:ILE:HD12	1.98	0.44
1:D:356:LEU:N	1:D:357:PRO:HD2	2.32	0.44
1:E:393:HIS:CG	1:E:394:PRO:HD2	2.52	0.44
1:C:501:MET:HA	1:E:405:ARG:HD2	1.99	0.44
1:A:158:ALA:HA	1:A:161:ARG:HH12	1.81	0.44
1:D:52:THR:HG22	1:D:57:LEU:HG	2.00	0.44
1:C:388:LYS:HB3	1:C:389:HIS:CD2	2.52	0.44
1:E:327:SER:O	1:E:331:GLN:HG2	2.18	0.44
1:A:166:GLU:O	1:A:170:ASN:ND2	2.50	0.44
1:B:344:LEU:HD22	1:B:375:CYS:HB3	1.99	0.44
1:D:27:GLU:O	1:D:31:ILE:HG13	2.17	0.44
1:E:309:SER:O	1:E:313:GLU:HB3	2.18	0.44
1:F:356:LEU:N	1:F:357:PRO:HD2	2.32	0.44
1:B:157:VAL:HG21	1:B:474:TYR:CE2	2.53	0.44
1:D:296:HIS:CD2	1:D:302:HIS:CE1	3.05	0.44
1:A:172:LEU:HD11	1:A:472:TRP:CZ2	2.53	0.44
1:B:309:SER:O	1:B:313:GLU:HB3	2.18	0.44
1:E:169:LEU:HD23	1:E:169:LEU:HA	1.84	0.44
1:A:243:ILE:HD13	1:A:243:ILE:HA	1.80	0.43
1:B:53:GLN:HG3	1:B:54:VAL:HG23	1.98	0.43
1:C:116:MET:HB2	1:C:119:ILE:HD12	1.99	0.43
1:C:356:LEU:N	1:C:357:PRO:HD2	2.32	0.43
1:A:320:ARG:HE	1:A:320:ARG:HB3	1.66	0.43
1:C:192:LEU:HD12	1:C:196:LEU:HB2	1.99	0.43
1:C:79:TYR:CD1	1:C:345:VAL:HG21	2.53	0.43
1:D:79:TYR:CD1	1:D:345:VAL:HG21	2.54	0.43
1:D:411:LEU:HD23	1:D:411:LEU:HA	1.89	0.43
1:A:270:ILE:HG21	1:A:341:LEU:HD23	2.00	0.43
1:A:400:GLU:OE1	1:E:442:ARG:NE	2.51	0.43
1:B:166:GLU:O	1:B:170:ASN:ND2	2.51	0.43
1:F:192:LEU:HD12	1:F:196:LEU:HB2	1.99	0.43
1:A:83:GLU:O	1:A:87:ARG:HB2	2.19	0.43
1:C:52:THR:HG22	1:C:57:LEU:HG	2.01	0.43
1:A:46:ARG:O	1:A:49:GLN:HG2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:295:LEU:HD23	1:B:385:VAL:HG21	2.01	0.43
1:E:67:LEU:HD23	1:F:67:LEU:HD23	2.00	0.43
1:A:46:ARG:O	1:A:49:GLN:N	2.52	0.43
1:E:203:TRP:CE2	1:E:249:GLU:HG3	2.54	0.43
1:F:27:GLU:O	1:F:31:ILE:HG13	2.19	0.43
1:E:356:LEU:N	1:E:357:PRO:HD2	2.34	0.43
1:A:38:ARG:CZ	2:G:2:C:H2'	2.49	0.42
1:A:309:SER:O	1:A:313:GLU:HB3	2.19	0.42
1:B:172:LEU:HD11	1:B:472:TRP:CZ2	2.54	0.42
1:A:169:LEU:HD23	1:A:169:LEU:HA	1.69	0.42
1:A:442:ARG:HE	1:C:397:GLU:HG2	1.84	0.42
1:D:316:TRP:O	1:D:319:SER:N	2.52	0.42
1:A:295:LEU:HD23	1:A:385:VAL:HG21	2.01	0.42
1:B:273:CYS:SG	1:B:274:VAL:N	2.93	0.42
1:F:327:SER:O	1:F:331:GLN:HG2	2.19	0.42
1:B:106:PHE:CE1	1:B:260:LEU:HD21	2.54	0.42
1:C:149:PRO:HB3	1:C:162:LEU:HB3	2.01	0.42
1:A:52:THR:HB	1:A:55:PHE:O	2.20	0.42
1:A:285:ILE:HG23	1:A:393:HIS:HD2	1.84	0.42
1:B:356:LEU:N	1:B:357:PRO:HD2	2.35	0.42
1:E:438:PRO:O	1:E:442:ARG:HG3	2.19	0.42
1:A:157:VAL:HG21	1:A:474:TYR:CE2	2.55	0.42
1:E:295:LEU:HD23	1:E:385:VAL:HG21	2.01	0.42
1:F:46:ARG:O	1:F:49:GLN:N	2.51	0.42
1:C:428:LEU:HD11	1:C:441:SER:HA	2.02	0.41
1:F:320:ARG:HD2	1:F:320:ARG:HA	1.43	0.41
1:B:46:ARG:O	1:B:49:GLN:HG2	2.19	0.41
1:D:161:ARG:HB3	1:D:163:ARG:HH12	1.85	0.41
1:D:393:HIS:CG	1:D:394:PRO:HD2	2.54	0.41
1:B:264:MET:HG3	1:B:265:GLU:N	2.35	0.41
1:B:461:LYS:HB2	1:B:461:LYS:HE3	1.57	0.41
1:A:169:LEU:O	1:A:173:ARG:HG3	2.21	0.41
1:B:211:LYS:O	1:B:257:ARG:HD2	2.20	0.41
1:E:172:LEU:HD23	1:E:471:LEU:HD12	2.03	0.41
1:E:192:LEU:HD12	1:E:196:LEU:HB2	2.02	0.41
1:A:31:ILE:O	1:A:34:SER:HB3	2.21	0.41
1:A:192:LEU:HD12	1:A:196:LEU:HB2	2.03	0.41
1:C:163:ARG:HA	1:C:163:ARG:HD3	1.91	0.41
1:D:157:VAL:HG21	1:D:474:TYR:CE2	2.55	0.41
1:F:273:CYS:HB2	1:F:382:TYR:HB3	2.02	0.41
1:A:229:LEU:CD2	1:A:257:ARG:HD3	2.47	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:MET:HG3	1:A:265:GLU:N	2.36	0.41
1:B:216:ALA:HA	1:B:233:PRO:HB3	2.02	0.41
1:D:273:CYS:HB2	1:D:382:TYR:HB3	2.02	0.41
1:E:479:ARG:HD3	1:E:479:ARG:HA	1.91	0.41
1:A:270:ILE:HG21	1:A:341:LEU:HD21	2.02	0.41
1:B:243:ILE:HA	1:B:243:ILE:HD13	1.82	0.41
1:E:157:VAL:HG21	1:E:474:TYR:CE2	2.56	0.41
1:B:52:THR:HB	1:B:55:PHE:O	2.21	0.40
1:C:278:GLU:HA	1:C:333:PHE:CZ	2.56	0.40
1:F:52:THR:HG22	1:F:57:LEU:HG	2.02	0.40
1:B:38:ARG:CZ	2:H:2:C:H2'	2.50	0.40
1:D:73:VAL:HG22	1:D:271:SER:HB2	2.02	0.40
1:D:309:SER:O	1:D:313:GLU:HB3	2.21	0.40
1:D:327:SER:OG	1:D:330:ASP:HB2	2.22	0.40
1:E:52:THR:HG22	1:E:57:LEU:HG	2.03	0.40
1:D:14:ARG:HG3	1:D:19:PRO:HD2	2.02	0.40
1:D:127:PHE:HZ	1:F:442:ARG:CZ	2.35	0.40
1:D:169:LEU:HA	1:D:169:LEU:HD23	1.81	0.40
1:D:438:PRO:O	1:D:442:ARG:HG3	2.21	0.40

All (2) 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:D:164:ASP:O	1:F:8:LYS:NZ[3_644]	2.03	0.17
1:A:26:HIS:NE2	1:C:465:ASP:OD2[3_544]	2.15	0.05

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	493/505 (98%)	454 (92%)	32 (6%)	7 (1%)	11 40

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	487/505 (96%)	453 (93%)	28 (6%)	6 (1%)	13	44
1	C	494/505 (98%)	468 (95%)	20 (4%)	6 (1%)	13	44
1	D	497/505 (98%)	465 (94%)	24 (5%)	8 (2%)	9	37
1	E	490/505 (97%)	466 (95%)	19 (4%)	5 (1%)	15	49
1	F	491/505 (97%)	462 (94%)	21 (4%)	8 (2%)	9	37
All	All	2952/3030 (97%)	2768 (94%)	144 (5%)	40 (1%)	11	40

All (40) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	148	GLN
1	A	304	LYS
1	B	301	GLN
1	D	326	ARG
1	F	301	GLN
1	A	310	LEU
1	A	311	VAL
1	B	310	LEU
1	B	311	VAL
1	C	272	TYR
1	C	304	LYS
1	C	310	LEU
1	D	272	TYR
1	D	325	SER
1	E	272	TYR
1	E	301	GLN
1	E	311	VAL
1	F	272	TYR
1	A	146	GLU
1	A	210	LEU
1	B	210	LEU
1	B	302	HIS
1	C	311	VAL
1	D	310	LEU
1	D	311	VAL
1	E	310	LEU
1	F	149	PRO
1	F	306	SER
1	F	310	LEU
1	C	319	SER

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Mol	Chain	Res	Type
1	D	210	LEU
1	F	300	GLY
1	F	304	LYS
1	F	311	VAL
1	A	61	ALA
1	C	210	LEU
1	D	301	GLN
1	B	299	TRP
1	E	210	LEU
1	D	270	ILE

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	430/450 (96%)	406 (94%)	24 (6%)	21	52
1	B	405/450 (90%)	386 (95%)	19 (5%)	26	59
1	C	432/450 (96%)	408 (94%)	24 (6%)	21	52
1	D	422/450 (94%)	400 (95%)	22 (5%)	23	55
1	E	431/450 (96%)	411 (95%)	20 (5%)	27	59
1	F	418/450 (93%)	396 (95%)	22 (5%)	22	54
All	All	2538/2700 (94%)	2407 (95%)	131 (5%)	23	55

All (131) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	THR
1	A	57	LEU
1	A	68	THR
1	A	87	ARG
1	A	95	GLU
1	A	119	ILE
1	A	147	SER
1	A	151	THR

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Mol	Chain	Res	Type
1	A	193	VAL
1	A	198	ARG
1	A	264	MET
1	A	271	SER
1	A	272	TYR
1	A	286	PHE
1	A	318	LYS
1	A	328	THR
1	A	344	LEU
1	A	371	ASP
1	A	384	ASN
1	A	423	SER
1	A	436	ARG
1	A	450	ARG
1	A	479	ARG
1	A	485	ILE
1	B	3	GLN
1	B	24	THR
1	B	68	THR
1	B	119	ILE
1	B	151	THR
1	B	152	ASP
1	B	193	VAL
1	B	198	ARG
1	B	264	MET
1	B	271	SER
1	B	272	TYR
1	B	286	PHE
1	B	302	HIS
1	B	344	LEU
1	B	371	ASP
1	B	384	ASN
1	B	423	SER
1	B	479	ARG
1	B	485	ILE
1	C	24	THR
1	C	63	VAL
1	C	86	SER
1	C	87	ARG
1	C	92	LYS
1	C	119	ILE
1	C	148	GLN

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Mol	Chain	Res	Type
1	C	164	ASP
1	C	193	VAL
1	C	238	SER
1	C	264	MET
1	C	273	CYS
1	C	286	PHE
1	C	302	HIS
1	C	330	ASP
1	C	344	LEU
1	C	371	ASP
1	C	384	ASN
1	C	423	SER
1	C	436	ARG
1	C	450	ARG
1	C	479	ARG
1	C	485	ILE
1	C	495	ASP
1	D	24	THR
1	D	68	THR
1	D	86	SER
1	D	95	GLU
1	D	119	ILE
1	D	139	ARG
1	D	150	LEU
1	D	164	ASP
1	D	193	VAL
1	D	238	SER
1	D	264	MET
1	D	273	CYS
1	D	286	PHE
1	D	330	ASP
1	D	344	LEU
1	D	371	ASP
1	D	384	ASN
1	D	423	SER
1	D	436	ARG
1	D	479	ARG
1	D	485	ILE
1	D	495	ASP
1	E	24	THR
1	E	68	THR
1	E	86	SER

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Mol	Chain	Res	Type
1	E	90	GLU
1	E	119	ILE
1	E	167	GLU
1	E	193	VAL
1	E	238	SER
1	E	264	MET
1	E	273	CYS
1	E	286	PHE
1	E	330	ASP
1	E	344	LEU
1	E	371	ASP
1	E	384	ASN
1	E	423	SER
1	E	450	ARG
1	E	479	ARG
1	E	485	ILE
1	E	495	ASP
1	F	24	THR
1	F	68	THR
1	F	86	SER
1	F	119	ILE
1	F	147	SER
1	F	193	VAL
1	F	238	SER
1	F	264	MET
1	F	273	CYS
1	F	284	ARG
1	F	286	PHE
1	F	320	ARG
1	F	327	SER
1	F	330	ASP
1	F	344	LEU
1	F	371	ASP
1	F	384	ASN
1	F	423	SER
1	F	450	ARG
1	F	479	ARG
1	F	485	ILE
1	F	495	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	302	HIS
1	A	384	ASN
1	B	302	HIS
1	B	384	ASN
1	C	13	HIS
1	C	141	HIS
1	C	182	HIS
1	C	384	ASN
1	C	393	HIS
1	C	505	GLN
1	D	13	HIS
1	D	141	HIS
1	D	182	HIS
1	D	227	HIS
1	D	296	HIS
1	D	384	ASN
1	E	13	HIS
1	E	141	HIS
1	E	182	HIS
1	E	384	ASN
1	E	505	GLN
1	F	13	HIS
1	F	141	HIS
1	F	182	HIS
1	F	227	HIS
1	F	384	ASN
1	F	393	HIS

5.3.3 RNA (i)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	G	0/3	-	-
2	H	0/3	-	-
All	All	0/6	-	-

There are no RNA backbone outliers to report.

There are no RNA pucker outliers to report.

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

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 (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	499/505 (98%)	-0.21	11 (2%)	62	41	61, 104, 156, 212	0
1	B	495/505 (98%)	-0.02	16 (3%)	47	25	73, 125, 176, 213	0
1	C	500/505 (99%)	-0.34	3 (0%)	89	78	55, 89, 140, 204	0
1	D	501/505 (99%)	-0.08	9 (1%)	68	47	59, 113, 180, 217	0
1	E	496/505 (98%)	-0.34	1 (0%)	95	90	56, 84, 139, 193	0
1	F	497/505 (98%)	-0.17	6 (1%)	79	61	65, 110, 175, 219	0
2	G	3/3 (100%)	0.01	0	100	100	95, 95, 109, 155	0
2	H	3/3 (100%)	0.45	0	100	100	117, 117, 119, 204	0
All	All	2994/3036 (98%)	-0.19	46 (1%)	73	54	55, 104, 166, 219	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	1	MET	7.4
1	D	323	SER	4.4
1	B	168	PRO	4.4
1	B	162	LEU	3.7
1	D	164	ASP	3.6
1	A	1	MET	3.6
1	B	60	ASN	3.5
1	A	2	ALA	3.5
1	A	302	HIS	3.4
1	F	162	LEU	3.4
1	A	322	ASN	3.1
1	D	162	LEU	3.0
1	B	160	LEU	3.0
1	A	6	PHE	2.8
1	F	305	GLY	2.7
1	C	305	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	151	THR	2.7
1	D	62	ALA	2.6
1	F	2	ALA	2.6
1	B	169	LEU	2.6
1	B	172	LEU	2.6
1	D	472	TRP	2.5
1	D	58	GLU	2.5
1	B	434	VAL	2.5
1	B	474	TYR	2.5
1	F	389	HIS	2.5
1	A	362	GLY	2.4
1	A	162	LEU	2.4
1	B	165	GLY	2.4
1	D	163	ARG	2.3
1	B	235	TYR	2.3
1	B	157	VAL	2.3
1	F	147	SER	2.3
1	A	361	ALA	2.3
1	B	56	PRO	2.3
1	A	360	PHE	2.3
1	E	302	HIS	2.2
1	B	356	LEU	2.2
1	B	140	LEU	2.2
1	C	62	ALA	2.1
1	A	148	GLN	2.1
1	B	471	LEU	2.1
1	F	295	LEU	2.0
1	A	300	GLY	2.0
1	D	171	GLU	2.0
1	B	159	ALA	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 monosaccharides 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. 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	B-factors(Å^2)	Q<0.9
3	MG	B	1000	1/1	0.77	0.29	134,134,134,134	0
3	MG	A	1000	1/1	0.93	0.24	120,120,120,120	0

6.5 Other polymers (i)

There are no such residues in this entry.