



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 12, 2016 – 12:27 PM EDT

PDB ID : 5B03
Title : Structure of MoeN5-Sso7d fusion protein in complex with geranyl pyrophosphate
Authors : Ko, T.-P.; Zhang, L.; Chen, C.-C.; Guo, R.-T.; Oldfield, E.O.
Deposited on : 2015-10-27
Resolution : 2.60 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-20027107
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-20027107

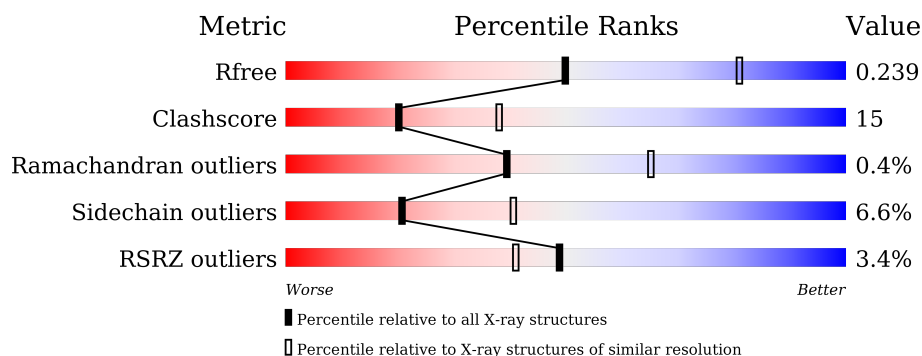
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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}	91344	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (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 ≥ 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	343	<div> <div>3%</div> <div> <div></div> <div>59%</div> <div>17%</div> <div>•</div> <div>23%</div> </div> </div>
1	B	343	<div> <div>3%</div> <div> <div></div> <div>68%</div> <div>27%</div> <div>• •</div> </div> </div>
1	C	343	<div> <div></div> <div> <div>57%</div> <div>17%</div> <div>•</div> <div>24%</div> </div> </div>
1	D	343	<div> <div>7%</div> <div> <div></div> <div>68%</div> <div>26%</div> <div>• •</div> </div> </div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GPP	B	500	-	-	-	X
2	GPP	C	500	-	-	-	X
2	GPP	D	500	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9801 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 MoeN5,DNA-binding protein 7d.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	264	Total	C	N	O	S	0	0	0
			2013	1243	374	385	11			
1	B	332	Total	C	N	O	S	0	0	0
			2533	1571	463	485	14			
1	C	261	Total	C	N	O	S	0	0	0
			1990	1230	371	378	11			
1	D	333	Total	C	N	O	S	0	0	0
			2544	1579	466	485	14			

There are 76 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	expression tag	UNP A0A010
A	-12	ALA	-	expression tag	UNP A0A010
A	-11	HIS	-	expression tag	UNP A0A010
A	-10	HIS	-	expression tag	UNP A0A010
A	-9	HIS	-	expression tag	UNP A0A010
A	-8	HIS	-	expression tag	UNP A0A010
A	-7	HIS	-	expression tag	UNP A0A010
A	-6	HIS	-	expression tag	UNP A0A010
A	-5	VAL	-	expression tag	UNP A0A010
A	-4	ASP	-	expression tag	UNP A0A010
A	-3	ASP	-	expression tag	UNP A0A010
A	-2	ASP	-	expression tag	UNP A0A010
A	-1	ASP	-	expression tag	UNP A0A010
A	0	LYS	-	expression tag	UNP A0A010
A	261	ALA	-	linker	UNP A0A010
A	262	GLY	-	linker	UNP A0A010
A	263	ALA	-	linker	UNP A0A010
A	264	GLY	-	linker	UNP A0A010
A	265	ALA	-	linker	UNP A0A010
B	-13	MET	-	expression tag	UNP A0A010
B	-12	ALA	-	expression tag	UNP A0A010

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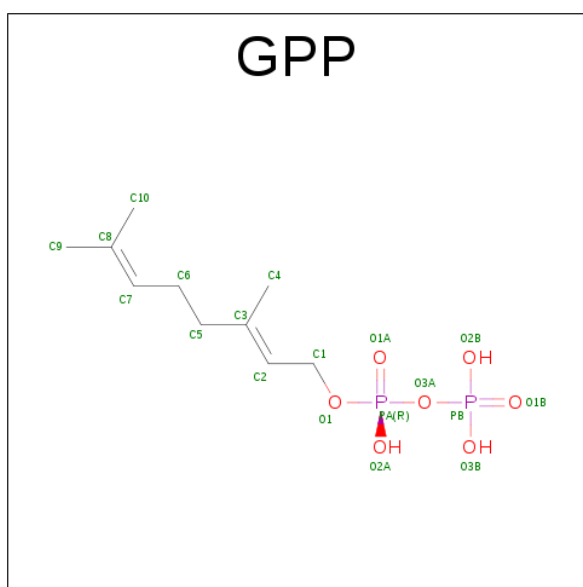
Chain	Residue	Modelled	Actual	Comment	Reference
B	-11	HIS	-	expression tag	UNP A0A010
B	-10	HIS	-	expression tag	UNP A0A010
B	-9	HIS	-	expression tag	UNP A0A010
B	-8	HIS	-	expression tag	UNP A0A010
B	-7	HIS	-	expression tag	UNP A0A010
B	-6	HIS	-	expression tag	UNP A0A010
B	-5	VAL	-	expression tag	UNP A0A010
B	-4	ASP	-	expression tag	UNP A0A010
B	-3	ASP	-	expression tag	UNP A0A010
B	-2	ASP	-	expression tag	UNP A0A010
B	-1	ASP	-	expression tag	UNP A0A010
B	0	LYS	-	expression tag	UNP A0A010
B	261	ALA	-	linker	UNP A0A010
B	262	GLY	-	linker	UNP A0A010
B	263	ALA	-	linker	UNP A0A010
B	264	GLY	-	linker	UNP A0A010
B	265	ALA	-	linker	UNP A0A010
C	-13	MET	-	expression tag	UNP A0A010
C	-12	ALA	-	expression tag	UNP A0A010
C	-11	HIS	-	expression tag	UNP A0A010
C	-10	HIS	-	expression tag	UNP A0A010
C	-9	HIS	-	expression tag	UNP A0A010
C	-8	HIS	-	expression tag	UNP A0A010
C	-7	HIS	-	expression tag	UNP A0A010
C	-6	HIS	-	expression tag	UNP A0A010
C	-5	VAL	-	expression tag	UNP A0A010
C	-4	ASP	-	expression tag	UNP A0A010
C	-3	ASP	-	expression tag	UNP A0A010
C	-2	ASP	-	expression tag	UNP A0A010
C	-1	ASP	-	expression tag	UNP A0A010
C	0	LYS	-	expression tag	UNP A0A010
C	261	ALA	-	linker	UNP A0A010
C	262	GLY	-	linker	UNP A0A010
C	263	ALA	-	linker	UNP A0A010
C	264	GLY	-	linker	UNP A0A010
C	265	ALA	-	linker	UNP A0A010
D	-13	MET	-	expression tag	UNP A0A010
D	-12	ALA	-	expression tag	UNP A0A010
D	-11	HIS	-	expression tag	UNP A0A010
D	-10	HIS	-	expression tag	UNP A0A010
D	-9	HIS	-	expression tag	UNP A0A010
D	-8	HIS	-	expression tag	UNP A0A010

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-7	HIS	-	expression tag	UNP A0A010
D	-6	HIS	-	expression tag	UNP A0A010
D	-5	VAL	-	expression tag	UNP A0A010
D	-4	ASP	-	expression tag	UNP A0A010
D	-3	ASP	-	expression tag	UNP A0A010
D	-2	ASP	-	expression tag	UNP A0A010
D	-1	ASP	-	expression tag	UNP A0A010
D	0	LYS	-	expression tag	UNP A0A010
D	261	ALA	-	linker	UNP A0A010
D	262	GLY	-	linker	UNP A0A010
D	263	ALA	-	linker	UNP A0A010
D	264	GLY	-	linker	UNP A0A010
D	265	ALA	-	linker	UNP A0A010

- Molecule 2 is GERANYL DIPHOSPHATE (three-letter code: GPP) (formula: $C_{10}H_{20}O_7P_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total 19	C 10	O 7	P 2	0	0
2	B	1	Total 19	C 10	O 7	P 2	0	0
2	C	1	Total 19	C 10	O 7	P 2	0	0
2	D	1	Total 19	C 10	O 7	P 2	0	0

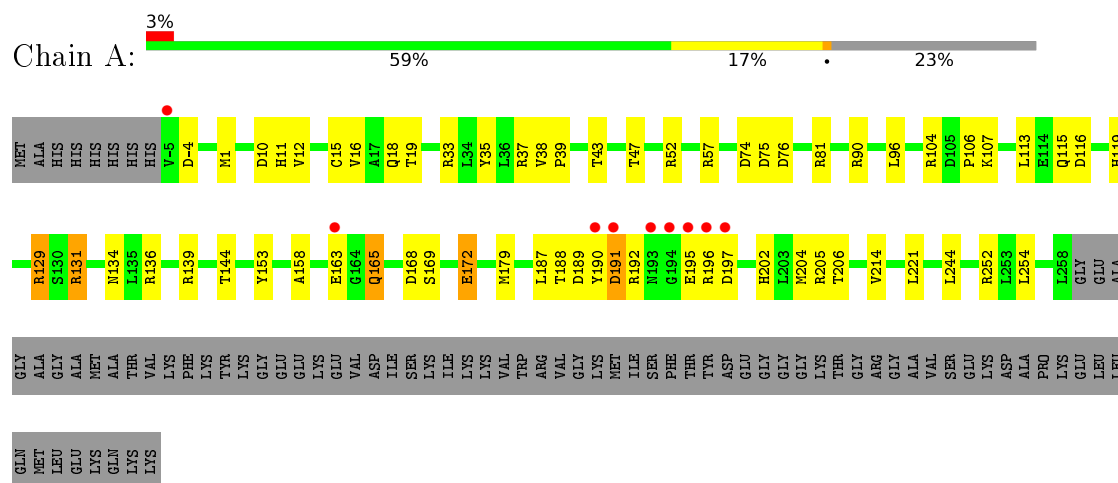
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	152	Total 152	O 152	0	0
3	B	197	Total 197	O 197	0	0
3	C	123	Total 123	O 123	0	0
3	D	173	Total 173	O 173	0	0

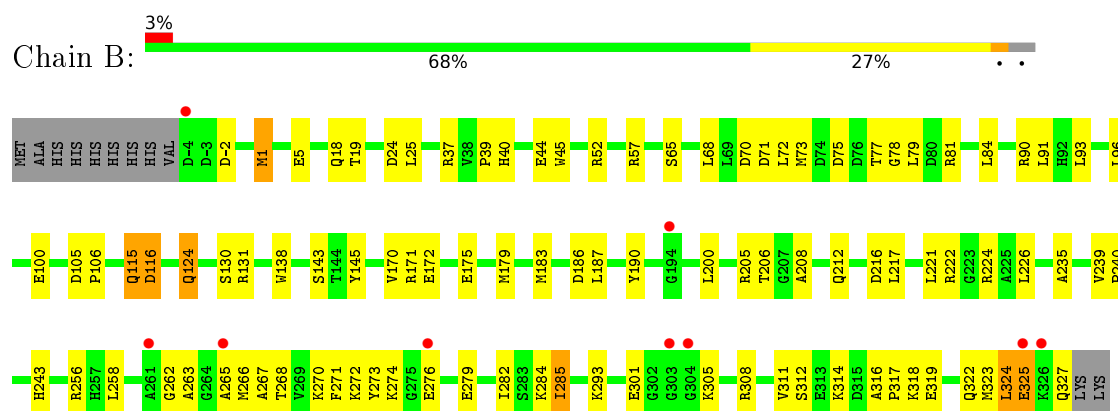
3 Residue-property plots [i](#)

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.

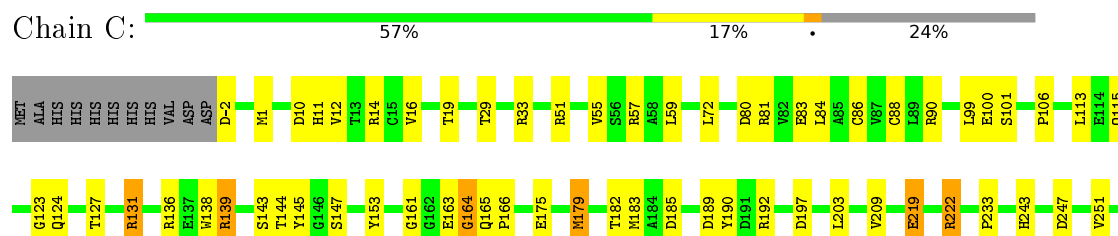
• Molecule 1: MoeN5,DNA-binding protein 7d

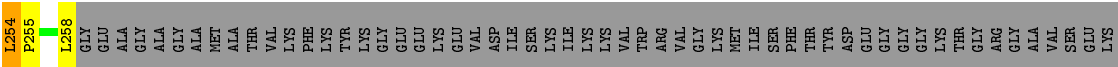


• Molecule 1: MoeN5,DNA-binding protein 7d

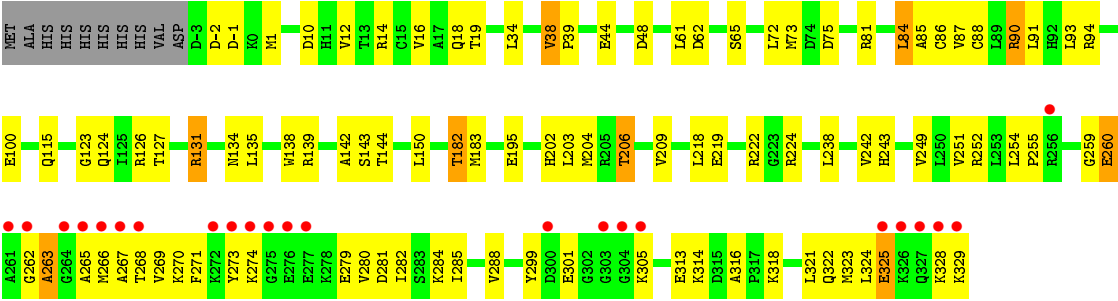


• Molecule 1: MoeN5,DNA-binding protein 7d





● Molecule 1: MoeN5,DNA-binding protein 7d



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	137.55Å 217.34Å 104.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.60 24.97 – 2.61	Depositor EDS
% Data completeness (in resolution range)	95.3 (25.00-2.60) 95.8 (24.97-2.61)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.31 (at 2.60Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.178 , 0.239 0.177 , 0.239	Depositor DCC
R_{free} test set	2296 reflections (4.98%)	DCC
Wilson B-factor (Å ²)	32.8	Xtriage
Anisotropy	0.237	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 53.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 46082 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9801	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.67% 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 ⓘ

5.1 Standard geometry ⓘ

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

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	0.82	1/2045 (0.0%)	0.87	0/2780
1	B	0.82	2/2572 (0.1%)	0.89	2/3478 (0.1%)
1	C	0.81	1/2022 (0.0%)	0.90	2/2748 (0.1%)
1	D	0.76	0/2583	0.85	1/3489 (0.0%)
All	All	0.80	4/9222 (0.0%)	0.88	5/12495 (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
1	A	0	1
1	C	0	1
All	All	0	2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1	MET	CG-SD	7.28	2.00	1.81
1	A	172	GLU	CG-CD	6.72	1.62	1.51
1	B	172	GLU	CG-CD	6.59	1.61	1.51
1	C	219	GLU	CG-CD	5.83	1.60	1.51

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	72	LEU	CA-CB-CG	6.19	129.55	115.30
1	C	222	ARG	NE-CZ-NH1	5.32	122.96	120.30
1	B	1	MET	CG-SD-CE	5.25	108.61	100.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	72	LEU	CA-CB-CG	5.15	127.14	115.30
1	B	216	ASP	CB-CG-OD1	5.03	122.83	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	153	TYR	Sidechain
1	C	153	TYR	Sidechain

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	2013	0	1996	65	0
1	B	2533	0	2532	82	0
1	C	1990	0	1979	55	0
1	D	2544	0	2554	85	0
2	A	19	0	17	1	0
2	B	19	0	17	0	0
2	C	19	0	17	0	0
2	D	19	0	17	0	0
3	A	152	0	0	8	0
3	B	197	0	0	8	0
3	C	123	0	0	6	0
3	D	173	0	0	7	0
All	All	9801	0	9129	269	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 15.

All (269) 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:D:267:ALA:HB1	1:D:282:ILE:HB	1.25	1.09

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1:MET:SD	1:B:37:ARG:HD3	2.02	1.00
1:A:129:ARG:HG3	1:A:129:ARG:HH11	1.28	0.99
1:C:222:ARG:NH1	1:C:243:HIS:HD2	1.64	0.96
1:D:202:HIS:O	1:D:206:THR:HG22	1.66	0.95
1:A:129:ARG:HH11	1:A:129:ARG:CG	1.85	0.88
1:C:222:ARG:NH1	1:C:243:HIS:CD2	2.42	0.88
1:D:267:ALA:HB1	1:D:282:ILE:CB	2.05	0.86
1:C:222:ARG:HH11	1:C:243:HIS:CD2	1.95	0.84
1:A:38:VAL:HG12	1:A:39:PRO:HD3	1.63	0.80
1:C:247:ASP:O	1:C:251:VAL:HG12	1.80	0.80
1:A:33:ARG:HH11	1:A:33:ARG:HG2	1.46	0.80
1:C:81:ARG:HD3	1:D:81:ARG:NH2	1.97	0.79
1:A:202:HIS:O	1:A:206:THR:HG23	1.82	0.79
1:B:272:LYS:HG3	1:B:276:GLU:O	1.84	0.78
1:C:131:ARG:HH11	1:C:131:ARG:CB	1.97	0.78
1:B:222:ARG:HH11	1:B:243:HIS:HD2	1.30	0.78
1:B:318:LYS:O	1:B:322:GLN:HG3	1.85	0.76
1:D:270:LYS:HG2	1:D:279:GLU:HG3	1.67	0.76
1:C:123:GLY:O	1:C:127:THR:HG23	1.85	0.75
1:D:224:ARG:NH1	3:D:602:HOH:O	2.18	0.75
1:D:204:MET:HG3	1:D:209:VAL:HG12	1.68	0.74
1:C:131:ARG:HH11	1:C:131:ARG:HB2	1.52	0.74
1:B:1:MET:SD	1:B:37:ARG:CD	2.77	0.73
1:D:10:ASP:OD1	1:D:14:ARG:HD2	1.88	0.73
1:D:38:VAL:HG12	1:D:39:PRO:HD3	1.69	0.73
1:A:214:VAL:HG11	1:A:254:LEU:HD21	1.70	0.72
1:A:205:ARG:HG2	1:D:131:ARG:NH1	2.04	0.72
1:A:35:TYR:OH	2:A:500:GPP:H12	1.89	0.72
1:C:14:ARG:HG3	3:C:672:HOH:O	1.90	0.72
1:B:314:LYS:HB2	1:B:314:LYS:NZ	2.06	0.71
1:D:222:ARG:HH11	1:D:243:HIS:HD2	1.36	0.71
1:C:131:ARG:NH1	1:C:131:ARG:HB3	2.06	0.71
1:B:224:ARG:HD2	3:B:660:HOH:O	1.90	0.70
1:B:68:LEU:HD22	1:B:84:LEU:CD2	2.22	0.70
1:C:166:PRO:HD3	1:C:233:PRO:HD2	1.73	0.70
1:D:123:GLY:O	1:D:127:THR:HG23	1.91	0.70
1:D:259:GLY:O	1:D:260:GLU:HG2	1.92	0.69
1:A:169:SER:HA	1:A:172:GLU:HG3	1.73	0.69
1:A:195:GLU:O	1:A:196:ARG:HD2	1.92	0.69
1:D:251:VAL:O	1:D:255:PRO:HG2	1.93	0.68
1:C:203:LEU:HB3	1:C:209:VAL:HG23	1.74	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:59:LEU:HD23	1:C:99:LEU:HD21	1.76	0.68
1:B:270:LYS:HE2	1:B:279:GLU:OE2	1.95	0.66
1:D:126:ARG:NH2	3:D:603:HOH:O	2.25	0.66
1:D:263:ALA:HB3	1:D:266:MET:HB2	1.77	0.66
1:A:129:ARG:NH1	1:A:129:ARG:HG3	2.08	0.66
1:D:218:LEU:HD21	1:D:249:VAL:HG11	1.78	0.66
1:C:131:ARG:NH1	1:C:131:ARG:CB	2.58	0.65
1:A:168:ASP:O	1:A:172:GLU:HG3	1.96	0.65
1:B:239:VAL:HB	1:B:240:PRO:HD3	1.80	0.63
1:B:319:GLU:HA	1:B:322:GLN:HE21	1.62	0.63
1:B:273:TYR:CE2	1:B:274:LYS:HG3	2.33	0.63
1:D:142:ALA:HB1	1:D:182:THR:HG21	1.80	0.63
1:D:18:GLN:HG2	1:D:18:GLN:O	1.99	0.63
1:D:252:ARG:O	1:D:255:PRO:HD2	1.98	0.63
1:A:33:ARG:NH1	1:A:33:ARG:HG2	2.10	0.63
1:D:318:LYS:O	1:D:322:GLN:HG2	1.98	0.63
1:A:163:GLU:HA	3:A:642:HOH:O	1.99	0.62
1:A:12:VAL:O	1:A:16:VAL:HG23	1.99	0.62
1:A:252:ARG:HD2	3:C:610:HOH:O	2.00	0.62
1:C:81:ARG:HD3	1:D:81:ARG:HH22	1.63	0.61
1:D:203:LEU:HA	1:D:206:THR:CG2	2.30	0.61
1:B:322:GLN:O	1:B:325:GLU:HG3	2.01	0.61
1:C:12:VAL:HG12	1:C:29:THR:HG21	1.83	0.61
1:C:131:ARG:HD2	1:C:197:ASP:HB3	1.83	0.60
1:D:135:LEU:HB2	1:D:209:VAL:HG22	1.84	0.60
1:A:129:ARG:NH1	1:A:129:ARG:CG	2.56	0.60
1:A:129:ARG:NH1	1:A:129:ARG:HB2	2.17	0.60
1:B:78:GLY:N	3:B:602:HOH:O	2.24	0.60
1:D:316:ALA:HB3	1:D:321:LEU:HD21	1.83	0.59
1:A:158:ALA:HA	1:A:165:GLN:HG2	1.84	0.59
1:C:14:ARG:NH2	1:C:14:ARG:HB2	2.17	0.59
1:C:138:TRP:CH2	1:C:182:THR:HG22	2.37	0.59
1:C:219:GLU:OE2	1:C:222:ARG:NE	2.31	0.59
1:A:192:ARG:HA	3:A:625:HOH:O	2.02	0.58
1:B:285:ILE:HG22	1:B:285:ILE:O	2.04	0.58
1:D:267:ALA:CB	1:D:282:ILE:HB	2.17	0.57
1:D:329:LYS:OXT	1:D:329:LYS:HG2	2.04	0.57
1:A:188:THR:CG2	1:A:192:ARG:HD3	2.34	0.57
1:D:321:LEU:O	3:D:601:HOH:O	2.17	0.57
1:B:267:ALA:HB2	1:B:319:GLU:OE2	2.05	0.56
1:C:254:LEU:HD22	1:C:258:LEU:HG	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:75:ASP:HB2	1:D:81:ARG:NH1	2.21	0.56
1:B:105:ASP:CG	1:B:106:PRO:HD2	2.25	0.56
1:B:314:LYS:CB	1:B:314:LYS:NZ	2.70	0.55
1:D:285:ILE:HG21	1:D:323:MET:SD	2.46	0.55
1:A:81:ARG:HE	1:B:75:ASP:HB2	1.71	0.55
1:A:15:CYS:O	1:A:19:THR:HG22	2.07	0.55
1:B:266:MET:HE2	1:B:267:ALA:N	2.22	0.54
1:A:10:ASP:CG	1:C:192:ARG:HD2	2.28	0.54
1:A:129:ARG:HH11	1:A:129:ARG:CB	2.20	0.54
1:A:47:THR:HG23	3:A:613:HOH:O	2.06	0.54
1:A:205:ARG:NH1	1:D:131:ARG:HD2	2.23	0.54
1:D:12:VAL:HG22	1:D:61:LEU:HD23	1.90	0.53
1:C:51:ARG:HG3	1:C:51:ARG:NH2	2.23	0.53
1:A:38:VAL:CG1	1:A:39:PRO:HD3	2.38	0.53
1:D:62:ASP:O	1:D:65:SER:HB3	2.09	0.53
1:B:1:MET:HB2	3:B:738:HOH:O	2.08	0.53
1:B:314:LYS:HB2	1:B:314:LYS:HZ2	1.74	0.53
1:D:44:GLU:HG2	1:D:238:LEU:HG	1.90	0.53
1:B:77:THR:OG1	1:B:79:LEU:HB2	2.09	0.53
1:D:259:GLY:O	1:D:260:GLU:CG	2.57	0.53
1:B:115:GLN:HA	1:B:115:GLN:OE1	2.10	0.52
1:A:187:LEU:HD21	1:A:204:MET:CE	2.39	0.52
1:B:96:LEU:O	1:B:100:GLU:HG3	2.09	0.52
1:D:328:LYS:NZ	3:D:604:HOH:O	2.38	0.52
1:D:301:GLU:O	1:D:305:LYS:HB2	2.09	0.52
1:A:-4:ASP:HB2	3:A:657:HOH:O	2.10	0.52
1:B:45:TRP:CE3	1:B:170:VAL:HG22	2.44	0.52
1:B:179:MET:HE2	1:B:221:LEU:HD21	1.91	0.52
1:A:191:ASP:HB2	3:D:646:HOH:O	2.08	0.52
1:A:187:LEU:CD2	1:A:204:MET:HE1	2.39	0.52
1:D:16:VAL:HG13	1:D:87:VAL:HG11	1.91	0.52
1:B:190:TYR:OH	1:B:205:ARG:HD2	2.09	0.52
1:B:124:GLN:NE2	1:B:145:TYR:HB3	2.25	0.52
1:D:263:ALA:CB	1:D:266:MET:HB2	2.39	0.52
1:C:51:ARG:HG3	1:C:51:ARG:HH21	1.74	0.52
1:D:280:VAL:HG12	1:D:281:ASP:N	2.24	0.52
1:B:171:ARG:O	1:B:175:GLU:HG3	2.10	0.51
1:B:273:TYR:CD2	1:B:274:LYS:HG3	2.45	0.51
1:C:11:HIS:CD2	1:C:57:ARG:HD2	2.45	0.51
1:C:161:GLY:HA3	1:C:165:GLN:OE1	2.10	0.51
1:B:138:TRP:CH2	1:B:183:MET:HG2	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:101:SER:HB2	3:D:653:HOH:O	2.11	0.51
1:A:113:LEU:HB3	1:B:93:LEU:HD22	1.92	0.51
1:D:273:TYR:O	1:D:274:LYS:HB2	2.11	0.51
1:D:85:ALA:O	1:D:88:CYS:HB3	2.11	0.51
1:C:113:LEU:HB3	1:D:93:LEU:HD22	1.93	0.50
1:A:190:TYR:CZ	1:A:196:ARG:NH2	2.79	0.50
1:A:129:ARG:NH1	1:A:129:ARG:CB	2.75	0.50
1:C:179:MET:O	1:C:183:MET:HG3	2.12	0.50
1:B:266:MET:CE	1:B:267:ALA:N	2.75	0.50
1:C:100:GLU:OE1	1:D:100:GLU:OE1	2.30	0.50
1:A:33:ARG:CG	1:A:33:ARG:NH1	2.75	0.50
1:D:280:VAL:HG21	1:D:299:TYR:CE1	2.46	0.50
1:C:14:ARG:HD3	3:C:651:HOH:O	2.12	0.49
1:C:12:VAL:CG1	1:C:29:THR:HG21	2.41	0.49
1:B:116:ASP:OD1	1:B:116:ASP:N	2.42	0.49
1:A:81:ARG:NH2	1:B:72:LEU:O	2.45	0.49
1:B:267:ALA:HB3	1:B:282:ILE:CG1	2.43	0.49
1:C:136:ARG:HB2	3:C:646:HOH:O	2.13	0.48
1:B:190:TYR:CZ	1:B:205:ARG:HD2	2.48	0.48
1:B:293:LYS:HE2	3:B:638:HOH:O	2.12	0.48
1:B:258:LEU:HD13	1:B:262:GLY:HA3	1.96	0.48
1:B:323:MET:O	1:B:327:GLN:HB2	2.14	0.48
1:A:195:GLU:C	1:A:196:ARG:HD2	2.33	0.48
1:B:222:ARG:NH1	1:B:243:HIS:HD2	2.05	0.48
1:A:96:LEU:HD11	1:B:96:LEU:HD11	1.95	0.47
1:B:71:ASP:HB3	1:B:77:THR:HG21	1.95	0.47
1:D:265:ALA:C	1:D:267:ALA:H	2.18	0.47
1:C:55:VAL:O	1:C:59:LEU:HG	2.14	0.47
1:A:81:ARG:NE	1:B:75:ASP:HB2	2.30	0.47
1:B:5:GLU:OE1	1:B:37:ARG:NH2	2.39	0.47
1:D:238:LEU:O	1:D:242:VAL:HG23	2.15	0.47
1:D:271:PHE:HE2	1:D:273:TYR:HB2	1.79	0.47
1:A:119:HIS:HD2	3:A:713:HOH:O	1.98	0.47
1:B:1:MET:CE	1:B:37:ARG:HD3	2.44	0.47
1:D:288:VAL:HG12	1:D:324:LEU:HD13	1.97	0.47
1:B:314:LYS:CB	1:B:314:LYS:HZ3	2.28	0.47
1:D:124:GLN:OE1	1:D:124:GLN:HA	2.14	0.46
1:D:203:LEU:O	1:D:206:THR:HG23	2.16	0.46
1:A:119:HIS:CE1	1:A:144:THR:HG22	2.50	0.46
1:B:301:GLU:CD	1:B:305:LYS:HE3	2.35	0.46
1:C:251:VAL:O	1:C:255:PRO:HG2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:24:ASP:HB3	1:B:79:LEU:CD2	2.46	0.46
1:D:219:GLU:OE1	1:D:219:GLU:HA	2.16	0.46
1:A:129:ARG:HB2	1:A:129:ARG:HH11	1.79	0.46
1:A:81:ARG:HH11	1:A:81:ARG:HG2	1.81	0.46
1:B:265:ALA:O	1:B:266:MET:C	2.54	0.46
1:B:25:LEU:HD13	1:B:79:LEU:HD13	1.98	0.46
1:D:150:LEU:HD12	1:D:150:LEU:HA	1.76	0.46
1:C:10:ASP:CG	1:C:14:ARG:HH22	2.20	0.46
1:B:57:ARG:HD3	3:B:736:HOH:O	2.16	0.46
1:B:267:ALA:O	1:B:282:ILE:HG12	2.17	0.45
1:B:267:ALA:HB3	1:B:282:ILE:HG13	1.97	0.45
1:C:11:HIS:CE1	1:C:57:ARG:HH11	2.34	0.45
1:D:270:LYS:HG2	1:D:279:GLU:CG	2.42	0.45
1:D:134:ASN:C	1:D:134:ASN:OD1	2.54	0.45
1:B:271:PHE:CE2	1:B:273:TYR:HB2	2.52	0.45
1:A:43:THR:HB	1:A:52:ARG:HG3	1.99	0.45
1:C:124:GLN:HB3	1:D:86:CYS:SG	2.57	0.45
1:C:14:ARG:HH21	1:C:14:ARG:HB2	1.82	0.45
1:C:1:MET:CE	3:C:690:HOH:O	2.65	0.45
1:D:12:VAL:O	1:D:16:VAL:HG23	2.17	0.45
1:B:1:MET:CE	1:B:37:ARG:CD	2.95	0.44
1:D:267:ALA:O	1:D:281:ASP:HA	2.17	0.44
1:A:158:ALA:CA	1:A:165:GLN:HG2	2.46	0.44
1:D:271:PHE:CE2	1:D:273:TYR:HB2	2.53	0.44
1:D:267:ALA:C	1:D:282:ILE:HG12	2.38	0.44
1:D:87:VAL:O	1:D:91:LEU:HG	2.18	0.44
1:A:81:ARG:HD2	3:B:621:HOH:O	2.17	0.44
1:B:212:GLN:HB2	1:B:263:ALA:O	2.18	0.44
1:B:221:LEU:HD23	1:B:221:LEU:HA	1.84	0.44
1:D:281:ASP:HB3	1:D:284:LYS:HG2	1.98	0.44
1:D:204:MET:HG3	1:D:209:VAL:CG1	2.41	0.44
1:A:104:ARG:HB2	3:A:651:HOH:O	2.16	0.44
1:C:147:SER:O	1:C:175:GLU:HG2	2.17	0.43
1:A:11:HIS:CE1	1:A:57:ARG:CZ	3.02	0.43
1:A:74:ASP:HB3	1:A:76:ASP:OD2	2.18	0.43
1:B:25:LEU:HD11	1:B:84:LEU:HD23	1.99	0.43
1:A:192:ARG:HD2	1:C:80:ASP:HB2	2.00	0.43
1:A:244:LEU:HD12	3:A:641:HOH:O	2.18	0.43
1:A:75:ASP:OD1	1:B:81:ARG:NH2	2.50	0.43
1:C:138:TRP:CZ2	1:C:182:THR:HG22	2.53	0.43
1:D:285:ILE:CG2	1:D:323:MET:SD	3.06	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:190:TYR:CZ	1:A:196:ARG:CZ	3.02	0.43
1:A:33:ARG:NH2	3:A:609:HOH:O	2.50	0.43
1:B:179:MET:CE	1:B:217:LEU:HD11	2.47	0.43
1:A:179:MET:HG2	1:A:221:LEU:HD11	2.00	0.43
1:A:187:LEU:HD21	1:A:204:MET:HE1	1.99	0.43
1:A:1:MET:CE	1:A:37:ARG:HD3	2.49	0.43
1:C:219:GLU:CD	1:C:222:ARG:HE	2.19	0.43
1:D:48:ASP:C	1:D:48:ASP:OD1	2.56	0.43
1:D:19:THR:CG2	1:D:90:ARG:HG3	2.49	0.43
1:B:40:HIS:CE1	1:B:52:ARG:NH2	2.87	0.43
1:D:268:THR:HA	1:D:280:VAL:O	2.19	0.43
1:A:188:THR:HG23	1:A:192:ARG:HD3	2.00	0.43
1:D:238:LEU:HD23	1:D:238:LEU:HA	1.69	0.43
1:A:179:MET:CG	1:A:221:LEU:HD11	2.49	0.43
1:B:282:ILE:HG21	1:B:319:GLU:HB3	2.01	0.43
1:C:145:TYR:C	1:C:145:TYR:CD1	2.92	0.42
1:B:68:LEU:HD22	1:B:84:LEU:HD22	1.99	0.42
1:B:68:LEU:HD23	1:B:68:LEU:HA	1.65	0.42
1:C:1:MET:HE1	3:C:690:HOH:O	2.18	0.42
1:B:206:THR:HG22	1:B:208:ALA:H	1.84	0.42
1:A:205:ARG:CZ	1:D:131:ARG:HD2	2.50	0.42
1:D:218:LEU:HD11	1:D:249:VAL:CG1	2.49	0.42
1:B:301:GLU:OE1	1:B:305:LYS:HE3	2.20	0.42
1:B:266:MET:HE3	1:B:267:ALA:H	1.85	0.42
1:A:205:ARG:HG2	1:D:131:ARG:HH12	1.81	0.42
1:B:266:MET:CE	1:B:267:ALA:H	2.32	0.42
1:D:138:TRP:CH2	1:D:183:MET:HG2	2.55	0.42
1:D:254:LEU:HD12	1:D:254:LEU:O	2.19	0.42
1:B:187:LEU:HD21	1:B:200:LEU:HD23	2.01	0.41
1:B:256:ARG:NH1	3:B:603:HOH:O	2.32	0.41
1:D:84:LEU:HA	1:D:84:LEU:HD23	1.74	0.41
1:A:131:ARG:N	1:A:131:ARG:CD	2.83	0.41
1:B:235:ALA:HA	3:B:645:HOH:O	2.20	0.41
1:B:44:GLU:HG3	1:B:44:GLU:O	2.21	0.41
1:C:101:SER:CB	3:D:653:HOH:O	2.68	0.41
1:C:86:CYS:SG	1:D:124:GLN:HB3	2.60	0.41
1:D:288:VAL:HG12	1:D:324:LEU:CD1	2.51	0.41
1:C:165:GLN:HE21	1:C:165:GLN:HB3	1.60	0.41
1:C:80:ASP:HB3	1:C:83:GLU:HG2	2.02	0.41
1:A:134:ASN:OD1	1:A:136:ARG:N	2.54	0.41
1:A:188:THR:CG2	1:A:192:ARG:HH11	2.34	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:268:THR:HB	1:B:279:GLU:HG3	2.03	0.41
1:B:311:VAL:HG23	1:B:312:SER:O	2.21	0.41
1:C:189:ASP:O	1:C:190:TYR:C	2.58	0.41
1:C:139:ARG:HG2	1:C:179:MET:HE3	2.02	0.41
1:C:12:VAL:O	1:C:16:VAL:HG23	2.21	0.41
1:B:179:MET:HE3	1:B:217:LEU:HD11	2.01	0.41
1:B:65:SER:HB2	1:B:91:LEU:HB2	2.02	0.41
1:B:273:TYR:O	1:B:276:GLU:HB2	2.20	0.41
1:D:38:VAL:CG1	1:D:39:PRO:HD3	2.45	0.41
1:B:70:ASP:HA	1:B:73:MET:HG2	2.03	0.40
1:D:249:VAL:O	1:D:254:LEU:HB2	2.21	0.40
1:D:90:ARG:NH1	1:D:94:ARG:HB2	2.35	0.40
1:D:75:ASP:HB2	1:D:81:ARG:HH12	1.84	0.40
1:B:316:ALA:HA	1:B:317:PRO:HD2	1.89	0.40
1:C:84:LEU:HA	1:C:84:LEU:HD23	1.92	0.40
1:D:280:VAL:HG21	1:D:299:TYR:CD1	2.57	0.40
1:B:226:LEU:HD23	1:B:226:LEU:HA	1.96	0.40
1:C:163:GLU:O	1:C:164:GLY:O	2.39	0.40
1:D:34:LEU:HD12	1:D:34:LEU:HA	1.94	0.40
1:D:206:THR:OG1	1:D:206:THR:O	2.39	0.40
1:D:269:VAL:HG23	1:D:282:ILE:HD13	2.03	0.40
1:D:325:GLU:HA	1:D:329:LYS:HB3	2.03	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	262/343 (76%)	252 (96%)	10 (4%)	0	100	100
1	B	330/343 (96%)	304 (92%)	25 (8%)	1 (0%)	46	72
1	C	259/343 (76%)	248 (96%)	10 (4%)	1 (0%)	39	65

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	331/343 (96%)	305 (92%)	23 (7%)	3 (1%)	21	42
All	All	1182/1372 (86%)	1109 (94%)	68 (6%)	5 (0%)	39	65

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	263	ALA
1	C	164	GLY
1	D	260	GLU
1	B	324	LEU
1	D	262	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	208/270 (77%)	195 (94%)	13 (6%)	22	44
1	B	260/270 (96%)	243 (94%)	17 (6%)	21	42
1	C	205/270 (76%)	191 (93%)	14 (7%)	20	39
1	D	261/270 (97%)	243 (93%)	18 (7%)	19	38
All	All	934/1080 (86%)	872 (93%)	62 (7%)	21	40

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	90	ARG
1	A	106	PRO
1	A	107	LYS
1	A	115	GLN
1	A	116	ASP
1	A	129	ARG
1	A	131	ARG
1	A	139	ARG

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Mol	Chain	Res	Type
1	A	165	GLN
1	A	189	ASP
1	A	191	ASP
1	A	197	ASP
1	B	-2	ASP
1	B	18	GLN
1	B	19	THR
1	B	39	PRO
1	B	90	ARG
1	B	115	GLN
1	B	116	ASP
1	B	124	GLN
1	B	130	SER
1	B	131	ARG
1	B	143	SER
1	B	186	ASP
1	B	284	LYS
1	B	285	ILE
1	B	308	ARG
1	B	324	LEU
1	B	325	GLU
1	C	-2	ASP
1	C	19	THR
1	C	33	ARG
1	C	88	CYS
1	C	90	ARG
1	C	106	PRO
1	C	115	GLN
1	C	131	ARG
1	C	139	ARG
1	C	143	SER
1	C	144	THR
1	C	179	MET
1	C	185	ASP
1	C	254	LEU
1	D	-2	ASP
1	D	-1	ASP
1	D	1	MET
1	D	38	VAL
1	D	73	MET
1	D	84	LEU
1	D	90	ARG

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Mol	Chain	Res	Type
1	D	115	GLN
1	D	131	ARG
1	D	139	ARG
1	D	143	SER
1	D	144	THR
1	D	182	THR
1	D	195	GLU
1	D	206	THR
1	D	313	GLU
1	D	314	LYS
1	D	325	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	119	HIS
1	A	165	GLN
1	A	193	ASN
1	B	18	GLN
1	B	124	GLN
1	B	243	HIS
1	B	322	GLN
1	C	115	GLN
1	C	165	GLN
1	C	202	HIS
1	C	243	HIS
1	D	165	GLN
1	D	243	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

4 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$
2	GPP	A	500	-	16,18,18	1.03	2 (12%)	17,25,25	0.87	0
2	GPP	B	500	-	16,18,18	0.79	0	17,25,25	1.47	2 (11%)
2	GPP	C	500	-	16,18,18	0.70	0	17,25,25	1.19	2 (11%)
2	GPP	D	500	-	16,18,18	0.89	0	17,25,25	1.02	1 (5%)

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	GPP	A	500	-	-	0/19/19/19	0/0/0/0
2	GPP	B	500	-	-	0/19/19/19	0/0/0/0
2	GPP	C	500	-	-	0/19/19/19	0/0/0/0
2	GPP	D	500	-	-	0/19/19/19	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	GPP	O1-C1	2.10	1.46	1.43
2	A	500	GPP	C1-C2	2.57	1.57	1.49

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	B	500	GPP	C5-C3-C2	-3.03	115.33	120.98
2	C	500	GPP	C5-C3-C2	-2.31	116.69	120.98
2	D	500	GPP	C4-C3-C5	2.38	118.99	115.37
2	C	500	GPP	C4-C3-C5	2.60	119.34	115.37
2	B	500	GPP	C4-C3-C5	3.99	121.45	115.37

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	GPP	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	264/343 (76%)	-0.49	9 (3%)	49	41	17, 32, 66, 104	0
1	B	332/343 (96%)	-0.40	9 (2%)	58	51	17, 33, 76, 84	0
1	C	261/343 (76%)	-0.61	0	100	100	17, 33, 50, 70	0
1	D	333/343 (97%)	-0.15	23 (6%)	20	14	16, 35, 115, 134	0
All	All	1190/1372 (86%)	-0.40	41 (3%)	49	41	16, 33, 85, 134	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	266	MET	7.3
1	D	265	ALA	6.1
1	D	328	LYS	5.4
1	D	275	GLY	5.3
1	D	303	GLY	5.1
1	A	194	GLY	5.0
1	D	268	THR	4.4
1	D	276	GLU	4.3
1	D	325	GLU	4.2
1	D	272	LYS	4.2
1	B	304	GLY	4.1
1	A	191	ASP	3.9
1	D	264	GLY	3.8
1	A	195	GLU	3.7
1	D	267	ALA	3.6
1	D	305	LYS	3.5
1	D	329	LYS	3.5
1	A	190	TYR	3.4
1	D	326	LYS	3.3
1	D	327	GLN	3.3
1	B	194	GLY	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	325	GLU	3.1
1	A	193	ASN	3.1
1	D	273	TYR	3.1
1	B	303	GLY	3.0
1	A	196	ARG	3.0
1	B	261	ALA	2.9
1	D	256	ARG	2.9
1	D	304	GLY	2.8
1	D	277	GLU	2.8
1	B	326	LYS	2.6
1	B	276	GLU	2.5
1	D	262	GLY	2.5
1	A	-5	VAL	2.5
1	B	-4	ASP	2.4
1	D	300	ASP	2.4
1	D	261	ALA	2.3
1	A	163	GLU	2.2
1	A	197	ASP	2.2
1	B	265	ALA	2.2
1	D	274	LYS	2.1

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
2	GPP	B	500	19/19	0.75	0.34	6.01	96,106,118,119	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	GPP	D	500	19/19	0.88	0.21	4.98	69,83,102,102	0
2	GPP	C	500	19/19	0.83	0.27	4.16	93,99,112,112	0
2	GPP	A	500	19/19	0.71	0.34	1.91	96,111,125,125	0

6.5 Other polymers [i](#)

There are no such residues in this entry.