



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 08:40 pm GMT

PDB ID : 4MDT
Title : Structure of LpxC bound to the reaction product UDP-(3-O-(R-3-hydroxymyristoyl))-glucosamine
Authors : Clayton, G.M.; Klein, D.J.; Rickert, K.W.; Patel, S.B.; Kornienko, M.; Zugay-Murphy, J.; Reid, J.C.; Tummala, S.; Sharma, S.; Singh, S.B.; Miesel, L.; Lumb, K.J.; Soisson, S.M.
Deposited on : 2013-08-23
Resolution : 2.59 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 : trunk28620
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) : recalc28949

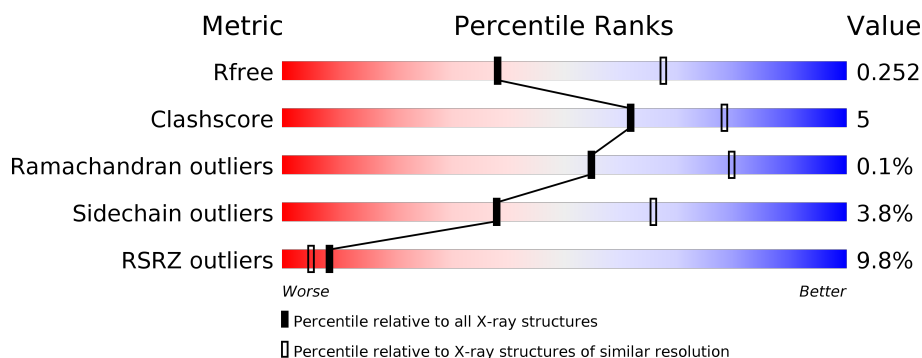
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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	2542 (2.60-2.60)
Clashscore	112137	2895 (2.60-2.60)
Ramachandran outliers	110173	2848 (2.60-2.60)
Sidechain outliers	110143	2848 (2.60-2.60)
RSRZ outliers	101464	2550 (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	305	<div> <div>87%</div> <div>11%</div> <div>•</div> </div>
1	B	305	<div> <div>83%</div> <div>14%</div> <div>••</div> </div>
1	C	305	<div> <div>16%</div> <div>82%</div> <div>16%</div> <div>•</div> </div>
1	D	305	<div> <div>20%</div> <div>80%</div> <div>15%</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
3	24G	A	402	-	-	-	X
4	PO4	D	403	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9746 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 UDP-3-O-[3-hydroxymyristoyl] N-acetylglucosamine deacetylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	300	Total	C	N	O	S	0	0	0
			2356	1497	408	438	13			
1	B	300	Total	C	N	O	S	0	0	0
			2356	1497	408	438	13			
1	C	299	Total	C	N	O	S	0	0	0
			2350	1494	407	436	13			
1	D	293	Total	C	N	O	S	0	0	0
			2299	1457	399	431	12			

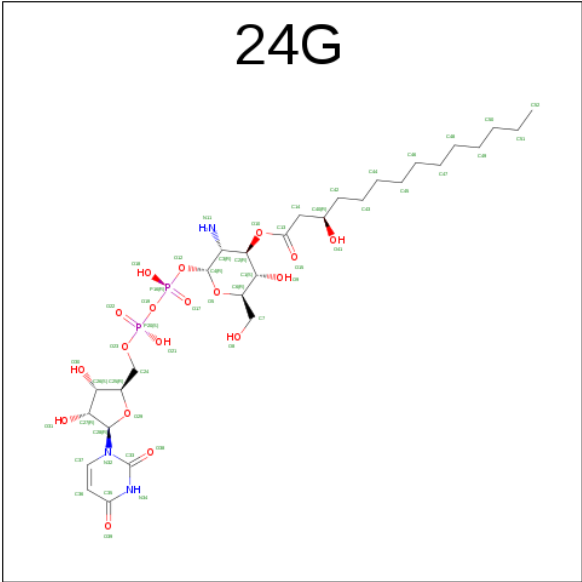
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	125	SER	CYS	ENGINEERED MUTATION	UNP K0BGQ2
B	125	SER	CYS	ENGINEERED MUTATION	UNP K0BGQ2
C	125	SER	CYS	ENGINEERED MUTATION	UNP K0BGQ2
D	125	SER	CYS	ENGINEERED MUTATION	UNP K0BGQ2

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

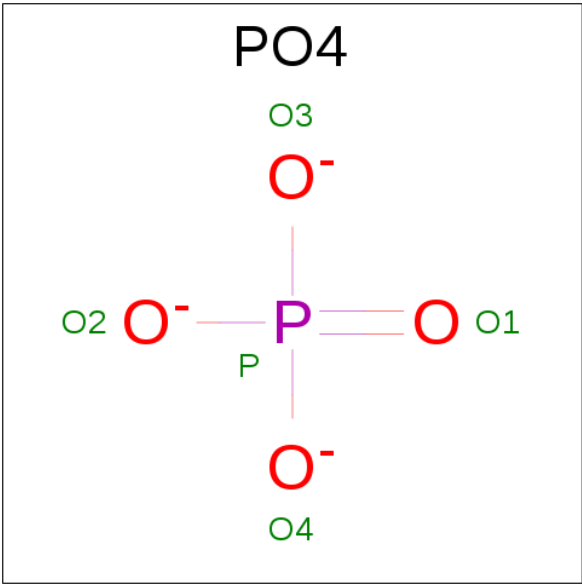
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		
2	D	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		

- Molecule 3 is URIDINE-5'-DIPHOSPHATE-3-O-(R-3-HYDROXYMYRISTOYL)-GLUCOSAMINE (three-letter code: 24G) (formula: C₂₉H₅₁N₃O₁₈P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			52	29	3	18	2		
3	B	1	Total	C	N	O	P	0	0
			52	29	3	18	2		
3	C	1	Total	C	N	O	P	0	0
			52	29	3	18	2		
3	D	1	Total	C	N	O	P	0	0
			52	29	3	18	2		

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O P 5 4 1	0	0
4	B	1	Total O P 5 4 1	0	0
4	C	1	Total O P 5 4 1	0	0
4	D	1	Total O P 5 4 1	0	0

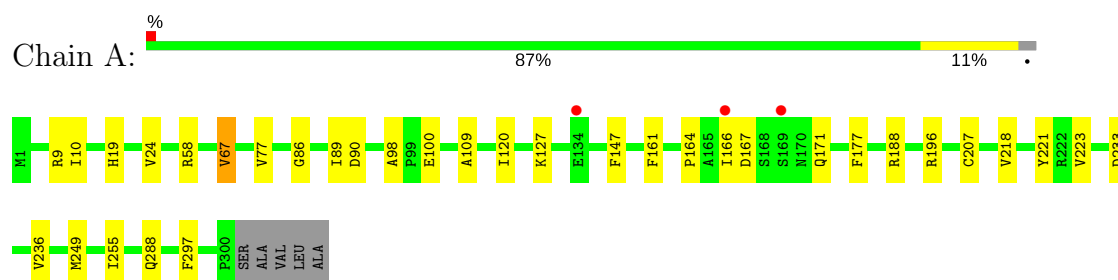
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	61	Total O 61 61	0	0
5	B	55	Total O 55 55	0	0
5	C	17	Total O 17 17	0	0
5	D	20	Total O 20 20	0	0

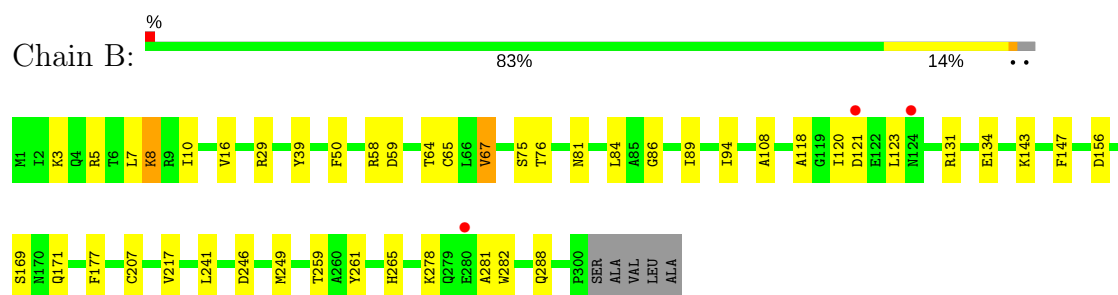
3 Residue-property plots

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

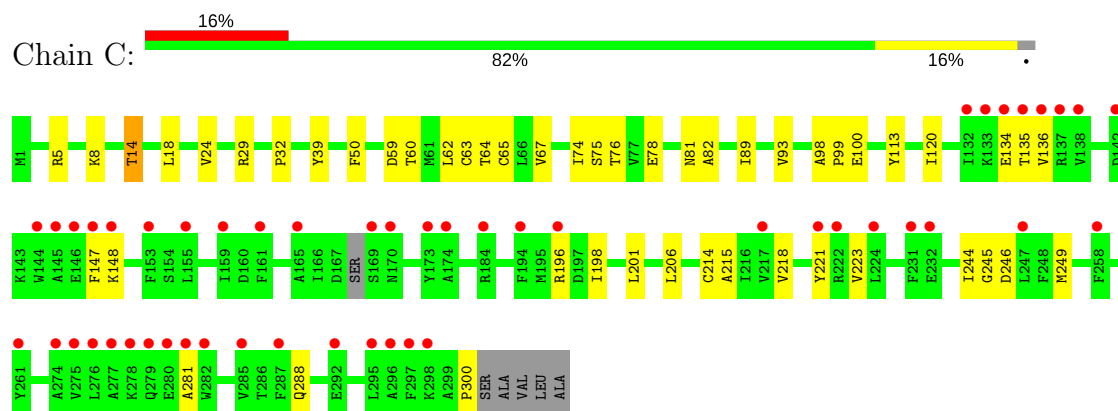
- Molecule 1: UDP-3-O-[3-hydroxymyristoyl] N-acetylglucosamine deacetylase



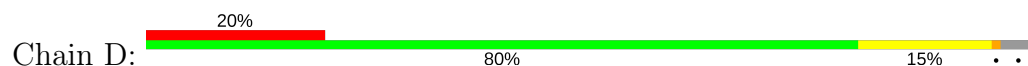
- Molecule 1: UDP-3-O-[3-hydroxymyristoyl] N-acetylglucosamine deacetylase

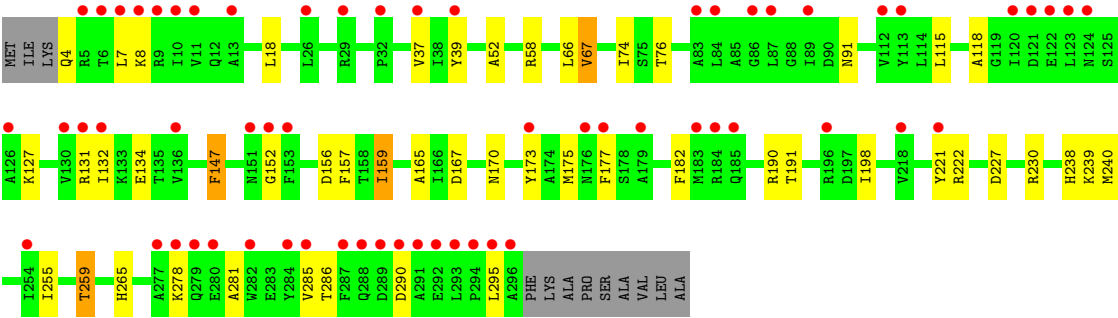


- Molecule 1: UDP-3-O-[3-hydroxymyristoyl] N-acetylglucosamine deacetylase



- Molecule 1: UDP-3-O-[3-hydroxymyristoyl] N-acetylglucosamine deacetylase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	168.97Å 103.52Å 103.97Å 90.00° 103.96° 90.00°	Depositor
Resolution (Å)	50.45 – 2.59 50.45 – 2.59	Depositor EDS
% Data completeness (in resolution range)	95.7 (50.45-2.59) 95.9 (50.45-2.59)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 2.58Å)	Xtriage
Refinement program	BUSTER 2.9.7	Depositor
R, R_{free}	0.197 , 0.240 0.205 , 0.252	Depositor DCC
R_{free} test set	2646 reflections (5.38%)	DCC
Wilson B-factor (Å ²)	54.7	Xtriage
Anisotropy	0.290	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 65.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9746	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.89% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, PO4, 24G

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	0.53	0/2404	0.79	0/3255
1	B	0.52	0/2404	0.79	0/3255
1	C	0.48	0/2397	0.75	0/3244
1	D	0.48	0/2345	0.76	0/3177
All	All	0.50	0/9550	0.77	0/12931

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

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	2356	0	2337	20	0
1	B	2356	0	2337	23	0
1	C	2350	0	2331	28	0
1	D	2299	0	2267	29	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	52	0	49	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	52	0	49	1	0
3	C	52	0	49	3	0
3	D	52	0	49	3	0
4	A	5	0	0	0	0
4	B	5	0	0	0	0
4	C	5	0	0	0	0
4	D	5	0	0	2	0
5	A	61	0	0	0	0
5	B	55	0	0	0	0
5	C	17	0	0	0	0
5	D	20	0	0	0	0
All	All	9746	0	9468	97	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (97) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:39:TYR:HB2	1:B:50:PHE:HB2	1.69	0.75
1:D:134:GLU:HB2	1:D:281:ALA:HA	1.73	0.71
1:C:81:ASN:HD21	1:C:249:MET:CE	2.05	0.69
1:D:4:GLN:HB3	1:D:91:ASN:HD21	1.58	0.69
1:B:64:THR:HG21	1:B:246:ASP:OD1	1.94	0.68
1:A:89:ILE:HD11	1:A:120:ILE:HD12	1.74	0.67
1:B:8:LYS:HG2	1:B:121:ASP:HB2	1.77	0.65
1:A:77:VAL:HG21	1:A:249:MET:HE3	1.77	0.65
1:B:81:ASN:HD21	1:B:249:MET:HE2	1.66	0.61
1:B:58:ARG:HB2	1:B:67:VAL:HG13	1.82	0.61
1:A:24:VAL:HG11	1:A:100:GLU:O	2.01	0.60
1:C:134:GLU:HB2	1:C:281:ALA:HA	1.84	0.60
1:C:81:ASN:HD21	1:C:249:MET:HE3	1.66	0.59
1:A:166:ILE:HG22	1:A:171:GLN:HE21	1.68	0.58
1:B:84:LEU:HD11	1:B:94:ILE:HD11	1.85	0.58
1:C:64:THR:OG1	1:C:246:ASP:OD1	2.18	0.58
1:C:135:THR:HG23	1:C:148:LYS:HG2	1.85	0.57
1:A:58:ARG:HB2	1:A:67:VAL:HG13	1.87	0.57
1:D:191:THR:HA	1:D:238:HIS:CD2	2.39	0.57
1:B:39:TYR:OH	1:B:81:ASN:OD1	2.23	0.57
1:A:10:ILE:CD1	1:D:167:ASP:HB2	2.36	0.56
1:C:39:TYR:HB2	1:C:50:PHE:HB2	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:156:ASP:O	1:D:259:THR:HA	2.07	0.54
1:C:5:ARG:HH22	1:C:300:PRO:HD3	1.71	0.54
1:B:134:GLU:HB2	1:B:281:ALA:HA	1.89	0.54
1:C:29:ARG:HB2	1:C:93:VAL:HB	1.90	0.53
1:C:81:ASN:HD21	1:C:249:MET:HE2	1.71	0.53
1:C:59:ASP:HB3	1:C:65:CYS:HB3	1.91	0.53
1:D:265:HIS:NE2	4:D:403:PO4:O3	2.35	0.53
1:B:64:THR:HG22	1:B:265:HIS:HB3	1.91	0.53
1:A:10:ILE:HD12	1:D:167:ASP:HB2	1.91	0.53
1:A:218:VAL:HG12	1:A:223:VAL:HA	1.91	0.52
1:D:58:ARG:HB2	1:D:67:VAL:HG13	1.92	0.52
1:C:82:ALA:HA	1:C:244:ILE:HG22	1.91	0.51
1:C:24:VAL:HG11	1:C:100:GLU:O	2.10	0.51
1:B:81:ASN:HD21	1:B:249:MET:CE	2.22	0.51
1:A:164:PRO:HG2	1:A:221:TYR:HA	1.93	0.50
1:C:24:VAL:HG13	1:C:98:ALA:HB3	1.94	0.50
1:D:66:LEU:HB2	1:D:74:ILE:HG23	1.94	0.50
1:B:169:SER:O	1:B:171:GLN:NE2	2.41	0.49
1:C:201:LEU:HB3	1:C:206:LEU:HB2	1.93	0.49
3:D:402:24G:N11	4:D:403:PO4:O4	2.46	0.49
1:C:218:VAL:HG12	1:C:223:VAL:HA	1.94	0.48
1:D:131:ARG:HD3	1:D:285:VAL:HG11	1.95	0.48
1:B:86:GLY:HA2	1:B:177:PHE:HZ	1.78	0.48
1:B:108:ALA:HB2	1:B:241:LEU:HD21	1.96	0.48
1:D:7:LEU:HD21	1:D:115:LEU:HD23	1.96	0.48
1:D:157:PHE:CE2	1:D:159:ILE:HG23	2.49	0.47
1:D:165:ALA:HB2	1:D:222:ARG:HA	1.97	0.47
1:C:198:ILE:HG21	3:C:402:24G:H27	1.97	0.47
1:B:7:LEU:HD22	1:B:118:ALA:HB3	1.97	0.46
1:D:198:ILE:HG21	3:D:402:24G:H27	1.96	0.46
1:A:19:HIS:CE1	1:A:207:CYS:HG	2.33	0.46
1:A:161:PHE:H	1:A:171:GLN:HE22	1.62	0.46
1:B:3:LYS:HB3	1:B:123:LEU:O	2.16	0.46
1:A:109:ALA:HB2	1:A:188:ARG:HG2	1.97	0.46
1:B:89:ILE:HD11	1:B:120:ILE:HD12	1.98	0.46
1:C:62:LEU:HD11	1:C:206:LEU:HD13	1.98	0.45
1:A:86:GLY:HA2	1:A:177:PHE:HZ	1.81	0.45
1:B:278:LYS:HD3	1:B:281:ALA:HB2	1.98	0.45
1:D:152:GLY:H	1:D:255:ILE:HG23	1.81	0.45
1:D:4:GLN:HG3	1:D:127:LYS:HA	1.97	0.44
1:B:59:ASP:HB3	1:B:65:CYS:HB3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:VAL:HG13	1:A:98:ALA:HB3	1.99	0.44
1:D:18:LEU:HD21	3:D:402:24G:H40	2.00	0.44
1:C:78:GLU:O	1:C:245:GLY:HA3	2.18	0.44
1:C:89:ILE:HD11	1:C:120:ILE:HD12	1.99	0.44
1:D:157:PHE:HE2	1:D:159:ILE:HG23	1.83	0.44
1:C:39:TYR:OH	1:C:81:ASN:OD1	2.31	0.43
1:A:24:VAL:HG22	1:A:98:ALA:HB3	1.99	0.43
1:C:18:LEU:HD11	3:C:402:24G:H36	1.99	0.43
1:D:177:PHE:CE1	1:D:182:PHE:HB2	2.53	0.43
1:C:81:ASN:ND2	1:C:249:MET:HE3	2.34	0.42
1:D:132:ILE:HG13	1:D:147:PHE:CD1	2.54	0.42
1:A:77:VAL:CG2	1:A:249:MET:HE3	2.48	0.42
1:B:10:ILE:HG12	1:B:29:ARG:HG2	2.00	0.42
1:A:233:ASP:HB2	1:A:236:VAL:HG12	2.01	0.42
1:B:156:ASP:O	1:B:259:THR:HA	2.19	0.42
1:B:131:ARG:O	1:B:282:TRP:HA	2.20	0.42
1:D:278:LYS:HD3	1:D:281:ALA:HB2	2.00	0.42
1:D:37:VAL:HB	1:D:52:ALA:HB2	2.02	0.42
1:D:173:TYR:O	1:D:240:MET:HE1	2.19	0.42
1:D:39:TYR:HB3	1:D:74:ILE:HD13	2.02	0.42
1:C:14:THR:HG23	1:D:227:ASP:OD1	2.20	0.41
1:D:159:ILE:HG21	1:D:239:LYS:HG3	2.02	0.41
1:B:217:VAL:HG21	3:B:402:24G:H30	2.01	0.41
1:C:74:ILE:HA	1:C:99:PRO:O	2.21	0.41
1:C:215:ALA:HB3	3:C:402:24G:H33	2.03	0.41
1:A:90:ASP:OD1	1:A:127:LYS:HG3	2.19	0.41
1:C:39:TYR:HB3	1:C:74:ILE:HD13	2.03	0.41
1:B:261:TYR:CG	1:C:32:PRO:HG3	2.55	0.41
1:D:190:ARG:HE	1:D:230:ARG:HG3	1.85	0.41
1:A:255:ILE:HG12	1:A:297:PHE:CE2	2.56	0.41
1:A:24:VAL:HG22	1:A:98:ALA:CB	2.50	0.41
1:C:63:CYS:SG	1:C:75:SER:HB2	2.61	0.41
1:D:173:TYR:HB3	1:D:240:MET:HE3	2.02	0.40
1:D:8:LYS:HB2	1:D:118:ALA:O	2.21	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	298/305 (98%)	286 (96%)	12 (4%)	0	100	100
1	B	298/305 (98%)	282 (95%)	16 (5%)	0	100	100
1	C	295/305 (97%)	271 (92%)	24 (8%)	0	100	100
1	D	291/305 (95%)	273 (94%)	17 (6%)	1 (0%)	44	70
All	All	1182/1220 (97%)	1112 (94%)	69 (6%)	1 (0%)	55	79

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	295	LEU

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	251/254 (99%)	245 (98%)	6 (2%)	54	80
1	B	251/254 (99%)	241 (96%)	10 (4%)	36	64
1	C	250/254 (98%)	238 (95%)	12 (5%)	30	55
1	D	245/254 (96%)	235 (96%)	10 (4%)	35	63
All	All	997/1016 (98%)	959 (96%)	38 (4%)	38	66

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ARG
1	A	67	VAL
1	A	147	PHE
1	A	167	ASP
1	A	196	ARG
1	A	288	GLN
1	B	5	ARG
1	B	8	LYS
1	B	16	VAL
1	B	67	VAL
1	B	75	SER
1	B	76	THR
1	B	143	LYS
1	B	147	PHE
1	B	207	CYS
1	B	288	GLN
1	C	8	LYS
1	C	14	THR
1	C	60	THR
1	C	67	VAL
1	C	76	THR
1	C	113	TYR
1	C	136	VAL
1	C	147	PHE
1	C	196	ARG
1	C	214	CYS
1	C	221	TYR
1	C	288	GLN
1	D	67	VAL
1	D	76	THR
1	D	147	PHE
1	D	159	ILE
1	D	170	ASN
1	D	175	MET
1	D	221	TYR
1	D	259	THR
1	D	286	THR
1	D	290	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	ASN

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Mol	Chain	Res	Type
1	A	171	GLN
1	A	176	ASN
1	B	12	GLN
1	B	176	ASN
1	B	185	GLN
1	C	81	ASN
1	C	176	ASN
1	D	91	ASN

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](#)

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 for Mogul analysis.

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
3	24G	A	402	-	47,54,54	1.11	4 (8%)	57,76,76	2.28	11 (19%)
4	PO4	A	403	2	4,4,4	2.24	1 (25%)	6,6,6	0.98	0
3	24G	B	402	-	47,54,54	0.99	3 (6%)	57,76,76	2.28	3 (5%)
4	PO4	B	403	2	4,4,4	1.83	1 (25%)	6,6,6	0.78	0
3	24G	C	402	-	47,54,54	1.10	3 (6%)	57,76,76	2.26	5 (8%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	PO4	C	403	2	4,4,4	2.15	1 (25%)	6,6,6	1.31	1 (16%)
3	24G	D	402	-	47,54,54	1.11	3 (6%)	57,76,76	2.26	4 (7%)
4	PO4	D	403	2	4,4,4	4.08	1 (25%)	6,6,6	1.43	1 (16%)

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
3	24G	A	402	-	-	0/37/78/78	0/3/3/3
4	PO4	A	403	2	-	0/0/0/0	0/0/0/0
3	24G	B	402	-	-	0/37/78/78	0/3/3/3
4	PO4	B	403	2	-	0/0/0/0	0/0/0/0
3	24G	C	402	-	-	0/37/78/78	0/3/3/3
4	PO4	C	403	2	-	0/0/0/0	0/0/0/0
3	24G	D	402	-	-	0/37/78/78	0/3/3/3
4	PO4	D	403	2	-	0/0/0/0	0/0/0/0

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	24G	O29-C28	2.30	1.44	1.41
3	B	402	24G	C35-N34	2.47	1.37	1.33
3	A	402	24G	C35-N34	2.82	1.38	1.33
3	C	402	24G	C35-N34	3.13	1.38	1.33
4	B	403	PO4	P-O1	3.17	1.57	1.50
3	B	402	24G	P20-O22	3.19	1.62	1.50
3	B	402	24G	P16-O17	3.19	1.62	1.50
3	D	402	24G	C35-N34	3.34	1.39	1.33
3	C	402	24G	P16-O17	3.41	1.63	1.50
3	A	402	24G	P16-O17	3.41	1.63	1.50
3	C	402	24G	P20-O22	3.46	1.63	1.50
3	A	402	24G	P20-O22	3.48	1.64	1.50
3	D	402	24G	P20-O22	3.51	1.64	1.50
3	D	402	24G	P16-O17	3.58	1.64	1.50
4	C	403	PO4	P-O1	3.83	1.58	1.50
4	A	403	PO4	P-O1	4.42	1.60	1.50
4	D	403	PO4	P-O1	7.94	1.67	1.50

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	402	24G	C36-C35-N34	-3.15	115.60	123.12
3	C	402	24G	C36-C35-N34	-3.08	115.78	123.12
3	B	402	24G	C36-C35-N34	-3.02	115.90	123.12
3	A	402	24G	C36-C35-N34	-2.63	116.83	123.12
4	C	403	PO4	O4-P-O1	-2.61	99.86	110.97
3	A	402	24G	C2-O10-C13	-2.50	113.37	117.62
3	B	402	24G	C2-C3-N11	-2.22	106.30	110.87
4	D	403	PO4	O4-P-O3	-2.16	99.97	107.90
3	A	402	24G	C36-C37-N32	2.02	125.30	120.67
3	C	402	24G	C36-C37-N32	2.04	125.35	120.67
3	C	402	24G	O18-P16-O12	2.04	114.71	106.49
3	A	402	24G	O29-C28-N32	2.08	112.24	108.08
3	A	402	24G	C1-C2-C3	2.16	114.87	111.09
3	A	402	24G	O12-C4-C3	2.33	112.61	108.40
3	A	402	24G	O21-P20-O23	2.34	119.19	108.14
3	A	402	24G	O18-P16-O12	2.55	116.72	106.49
3	D	402	24G	O5-C4-O12	2.58	114.74	111.36
3	A	402	24G	O10-C2-C1	2.60	113.47	107.75
3	C	402	24G	O5-C4-O12	2.89	115.15	111.36
3	D	402	24G	O29-C28-N32	3.23	114.55	108.08
3	A	402	24G	O5-C4-O12	3.70	116.20	111.36
3	A	402	24G	C35-N34-C33	13.98	126.14	114.13
3	D	402	24G	C35-N34-C33	14.89	126.92	114.13
3	C	402	24G	C35-N34-C33	15.02	127.03	114.13
3	B	402	24G	C35-N34-C33	15.27	127.24	114.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	402	24G	1	0
3	C	402	24G	3	0
3	D	402	24G	3	0
4	D	403	PO4	2	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	300/305 (98%)	0.16	3 (1%) 82 79	32, 46, 69, 90	0
1	B	300/305 (98%)	0.26	3 (1%) 82 79	32, 51, 75, 95	0
1	C	299/305 (98%)	1.01	50 (16%) 2 1	42, 72, 111, 148	0
1	D	293/305 (96%)	1.26	61 (20%) 1 0	44, 74, 117, 156	0
All	All	1192/1220 (97%)	0.67	117 (9%) 8 5	32, 59, 105, 156	0

All (117) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	293	LEU	13.6
1	D	123	LEU	7.3
1	D	296	ALA	7.2
1	C	281	ALA	6.3
1	D	152	GLY	6.2
1	C	144	TRP	6.2
1	D	287	PHE	6.1
1	D	291	ALA	6.1
1	D	289	ASP	5.9
1	C	280	GLU	5.7
1	D	87	LEU	5.6
1	C	278	LYS	5.4
1	C	134	GLU	5.4
1	C	296	ALA	5.3
1	D	288	GLN	5.2
1	C	221	TYR	5.2
1	D	177	PHE	5.2
1	D	120	ILE	5.0
1	C	287	PHE	4.8
1	C	222	ARG	4.7
1	D	278	LYS	4.7

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Mol	Chain	Res	Type	RSRZ
1	D	10	ILE	4.7
1	D	285	VAL	4.6
1	C	169	SER	4.6
1	D	130	VAL	4.5
1	D	196	ARG	4.5
1	D	290	ASP	4.5
1	C	138	VAL	4.1
1	D	124	ASN	4.1
1	D	294	PRO	4.1
1	C	136	VAL	4.1
1	C	161	PHE	4.0
1	C	135	THR	4.0
1	D	6	THR	4.0
1	C	276	LEU	3.9
1	D	9	ARG	3.9
1	C	170	ASN	3.8
1	C	137	ARG	3.8
1	C	297	PHE	3.7
1	C	261	TYR	3.7
1	D	151	ASN	3.7
1	D	284	TYR	3.6
1	C	133	LYS	3.6
1	D	254	ILE	3.6
1	D	279	GLN	3.6
1	C	285	VAL	3.5
1	C	277	ALA	3.5
1	C	159	ILE	3.4
1	C	292	GLU	3.4
1	D	176	ASN	3.4
1	C	217	VAL	3.3
1	C	173	TYR	3.3
1	C	298	LYS	3.3
1	C	232	GLU	3.2
1	D	292	GLU	3.2
1	D	39	TYR	3.2
1	D	8	LYS	3.2
1	D	7	LEU	3.2
1	D	126	ALA	3.1
1	C	184	ARG	3.1
1	D	113	TYR	3.0
1	D	5	ARG	3.0
1	D	89	ILE	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	124	ASN	3.0
1	D	122	GLU	3.0
1	D	32	PRO	3.0
1	D	280	GLU	3.0
1	C	279	GLN	3.0
1	C	295	LEU	2.9
1	D	277	ALA	2.9
1	D	83	ALA	2.9
1	C	155	LEU	2.9
1	D	131	ARG	2.9
1	D	295	LEU	2.8
1	D	184	ARG	2.8
1	C	153	PHE	2.8
1	A	169	SER	2.8
1	C	194	PHE	2.8
1	C	147	PHE	2.7
1	D	179	ALA	2.7
1	C	145	ALA	2.6
1	C	231	PHE	2.6
1	D	132	ILE	2.6
1	D	153	PHE	2.6
1	A	166	ILE	2.6
1	D	185	GLN	2.6
1	C	258	PHE	2.6
1	D	29	ARG	2.6
1	B	280	GLU	2.5
1	D	37	VAL	2.5
1	C	132	ILE	2.5
1	D	13	ALA	2.5
1	C	275	VAL	2.4
1	D	221	TYR	2.4
1	C	146	GLU	2.4
1	D	218	VAL	2.4
1	C	224	LEU	2.4
1	C	274	ALA	2.4
1	C	148	LYS	2.4
1	D	136	VAL	2.3
1	C	165	ALA	2.3
1	C	247	LEU	2.3
1	C	282	TRP	2.3
1	D	282	TRP	2.3
1	C	196	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	174	ALA	2.2
1	D	86	GLY	2.2
1	C	142	ASP	2.2
1	D	84	LEU	2.2
1	A	134	GLU	2.2
1	B	121	ASP	2.2
1	D	11	VAL	2.1
1	D	112	VAL	2.1
1	D	121	ASP	2.1
1	D	26	LEU	2.1
1	D	173	TYR	2.0
1	D	183	MET	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
3	24G	A	402	52/52	0.91	0.26	2.56	59,70,88,89	0
3	24G	C	402	52/52	0.81	0.32	1.39	97,105,132,132	0
2	ZN	A	401	1/1	1.00	0.17	0.98	39,39,39,39	0
3	24G	D	402	52/52	0.92	0.21	0.72	55,83,90,92	0
3	24G	B	402	52/52	0.97	0.18	0.30	38,48,60,63	0
4	PO4	D	403	5/5	0.97	0.16	0.03	42,48,52,57	0
4	PO4	B	403	5/5	0.99	0.16	0.02	43,46,50,50	0
4	PO4	A	403	5/5	0.99	0.16	-0.16	43,47,50,53	0
2	ZN	D	401	1/1	0.98	0.14	-0.28	55,55,55,55	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	ZN	B	401	1/1	1.00	0.16	-0.45	40,40,40,40	0
2	ZN	C	401	1/1	0.99	0.16	-0.55	63,63,63,63	0
4	PO4	C	403	5/5	0.98	0.14	-1.33	62,67,68,70	0

6.5 Other polymers [i](#)

There are no such residues in this entry.