



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:07 PM GMT

PDB ID : 4BKM  
Title : Crystal structure of the murine AUM (phosphoglycolate phosphatase) capping domain as a fusion protein with the catalytic core domain of murine chronophin (pyridoxal phosphate phosphatase)  
Authors : Knobloch, G.; Seifried, A.; Gohla, A.; Schindelin, H.  
Deposited on : 2013-04-26  
Resolution : 2.65 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
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) : trunk26865

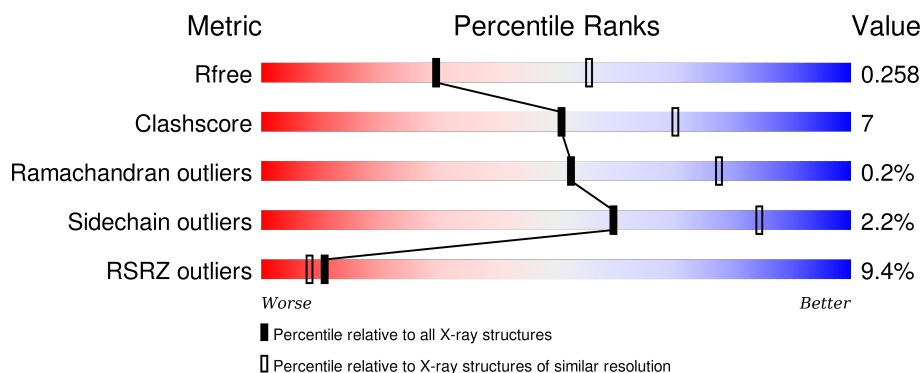
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.65 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	3152 (2.70-2.62)
Clashscore	102246	3524 (2.70-2.62)
Ramachandran outliers	100387	3469 (2.70-2.62)
Sidechain outliers	100360	3469 (2.70-2.62)
RSRZ outliers	91569	3161 (2.70-2.62)

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  $\geq 3$ , 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	309	<div> <div>10%</div> <div>82%</div> <div>16%</div> <div>••</div> </div>
1	B	309	<div> <div>10%</div> <div>81%</div> <div>17%</div> <div>•</div> </div>
1	C	309	<div> <div>10%</div> <div>86%</div> <div>12%</div> <div>•</div> </div>
1	D	309	<div> <div>7%</div> <div>84%</div> <div>14%</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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NO3	A	1305	-	X	-	-
3	NO3	A	1306	-	X	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 18782 atoms, of which 9332 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 PYRIDOXAL PHOSPHATE PHOSPHATASE, PHOSPHO-GLYCOLATE PHOSPHATASE, PYRIDOXAL PHOSPHATE PHOSPHATASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	304	Total	C	H	N	O	S	0	0	0
			4642	1454	2333	416	424	15			
1	B	304	Total	C	H	N	O	S	0	0	0
			4642	1454	2333	416	424	15			
1	C	304	Total	C	H	N	O	S	0	0	0
			4642	1454	2333	416	424	15			
1	D	304	Total	C	H	N	O	S	0	0	0
			4642	1454	2333	416	424	15			

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 3 is NITRATE ION (three-letter code: NO3) (formula: NO<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	N	O	0	0
			4	1	3		
3	A	1	Total	N	O	0	0
			4	1	3		

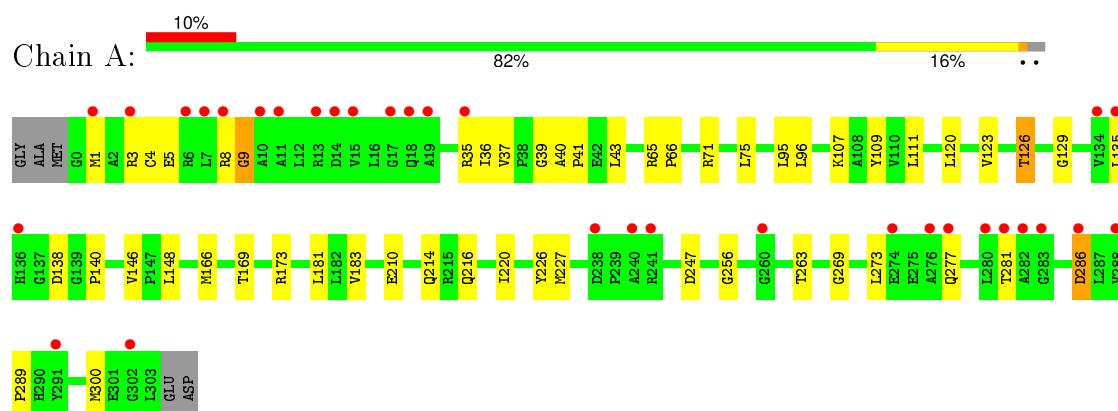
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	58	Total	O	0	0
			58	58		
4	B	52	Total	O	0	0
			52	52		
4	C	43	Total	O	0	0
			43	43		
4	D	49	Total	O	0	0
			49	49		

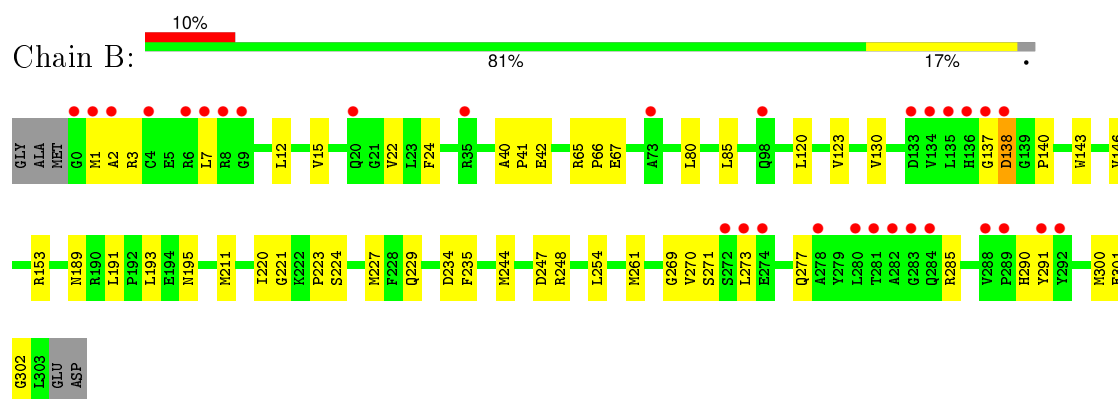
### 3 Residue-property plots

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

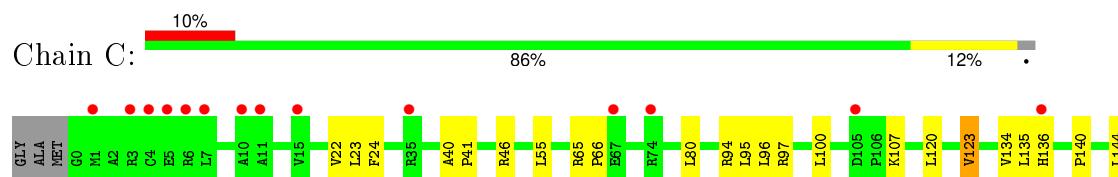
- Molecule 1: PYRIDOXAL PHOSPHATE PHOSPHATASE, PHOSPHOGLYCOLATE PHOSPHATASE, PYRIDOXAL PHOSPHATE PHOSPHATASE

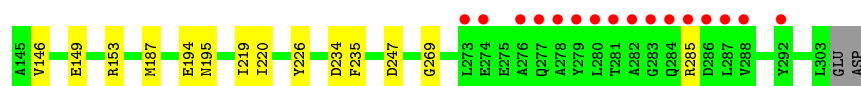


- Molecule 1: PYRIDOXAL PHOSPHATE PHOSPHATASE, PHOSPHOGLYCOLATE PHOSPHATASE, PYRIDOXAL PHOSPHATE PHOSPHATASE

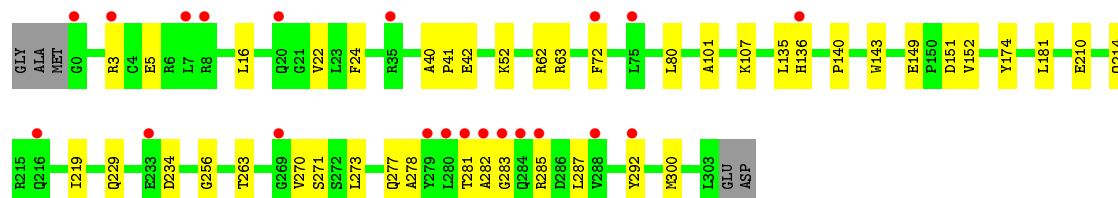
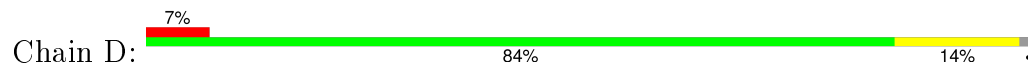


- Molecule 1: PYRIDOXAL PHOSPHATE PHOSPHATASE, PHOSPHOGLYCOLATE PHOSPHATASE, PYRIDOXAL PHOSPHATE PHOSPHATASE





- Molecule 1: PYRIDOXAL PHOSPHATE PHOSPHATASE, PHOSPHOGLYCOLATE PHOSPHATASE, PYRIDOXAL PHOSPHATE PHOSPHATASE



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	67.50Å 91.96Å 105.92Å 90.00° 90.20° 90.00°	Depositor
Resolution (Å)	41.74 – 2.65 41.74 – 2.65	Depositor EDS
% Data completeness (in resolution range)	95.3 (41.74-2.65) 95.8 (41.74-2.65)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.97 (at 2.65Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.198 , 0.256 0.206 , 0.258	Depositor DCC
$R_{free}$ test set	3512 reflections (5.07%)	DCC
Wilson B-factor (Å <sup>2</sup> )	50.8	Xtriage
Anisotropy	0.828	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.41 , 68.4	EDS
Estimated twinning fraction	0.095 for h,-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	2 of 36203 reflections (0.006%)	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	18782	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	95.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 40.32 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 2.7852e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.



## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, NO3

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	0.25	0/2351	0.45	0/3190
1	B	0.26	0/2351	0.46	0/3190
1	C	0.25	0/2351	0.43	0/3190
1	D	0.25	0/2351	0.48	0/3190
All	All	0.25	0/9404	0.46	0/12760

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2309	2333	2329	34	0
1	B	2309	2333	2329	40	0
1	C	2309	2333	2329	24	0
1	D	2309	2333	2329	30	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	8	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	58	0	0	4	0
4	B	52	0	0	5	0
4	C	43	0	0	4	0
4	D	49	0	0	0	0
All	All	9450	9332	9316	128	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 7.

All (128) 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:1:MET:HE1	1:B:285:ARG:HD2	1.42	0.96
1:B:1:MET:HE1	1:B:285:ARG:CD	2.15	0.77
1:A:35:ARG:NH2	4:A:2008:HOH:O	2.16	0.77
1:A:216:GLN:O	4:A:2044:HOH:O	2.15	0.64
1:B:244:MET:HB2	1:B:261:MET:HE2	1.80	0.63
1:D:278:ALA:HA	1:D:281:THR:HB	1.81	0.62
1:C:195:ASN:ND2	4:C:2034:HOH:O	2.33	0.61
1:D:277:GLN:O	1:D:281:THR:N	2.32	0.60
1:B:1:MET:HG3	1:B:2:ALA:N	2.19	0.58
1:D:3:ARG:HD3	1:D:5:GLU:OE2	2.04	0.58
1:D:277:GLN:O	1:D:281:THR:HB	2.05	0.57
1:D:285:ARG:HD3	1:D:285:ARG:N	2.19	0.57
1:A:3:ARG:HD3	1:A:5:GLU:OE2	2.04	0.56
1:D:149:GLU:HB2	1:D:152:VAL:HG23	1.87	0.56
1:A:286:ASP:OD1	1:A:286:ASP:N	2.38	0.56
1:A:71:ARG:O	1:A:75:LEU:HD23	2.05	0.56
1:B:2:ALA:HA	1:B:290:HIS:HA	1.89	0.55
1:B:1:MET:CG	1:B:2:ALA:N	2.71	0.54
1:C:195:ASN:N	4:C:2032:HOH:O	2.37	0.54
1:B:195:ASN:ND2	4:B:2041:HOH:O	2.41	0.54
1:B:234:ASP:HB3	1:B:235:PHE:CD1	2.43	0.53
1:A:273:LEU:O	1:A:277:GLN:HG3	2.09	0.53
1:D:283:GLY:HA2	1:D:285:ARG:CZ	2.39	0.53
1:B:247:ASP:HB2	1:B:269:GLY:HA3	1.91	0.52
1:B:270:VAL:CG1	1:B:271:SER:N	2.72	0.52
1:B:1:MET:CE	1:B:285:ARG:HH11	2.22	0.52
1:C:247:ASP:HB2	1:C:269:GLY:HA3	1.91	0.52
1:A:183:VAL:CG1	1:A:220:ILE:HD11	2.39	0.52
3:A:1305:NO3:O2	4:A:2036:HOH:O	2.18	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1:MET:HE1	1:B:285:ARG:CG	2.40	0.51
1:B:270:VAL:HG13	1:B:271:SER:N	2.25	0.51
1:D:22:VAL:HG11	1:D:24:PHE:CZ	2.46	0.51
1:D:278:ALA:O	1:D:282:ALA:N	2.43	0.51
1:C:123:VAL:HG13	1:C:123:VAL:O	2.09	0.51
1:B:211:MET:O	4:B:2044:HOH:O	2.18	0.51
1:B:130:VAL:O	4:B:2023:HOH:O	2.18	0.51
1:B:1:MET:CE	1:B:285:ARG:CG	2.89	0.50
1:C:40:ALA:HB3	1:C:41:PRO:HD3	1.92	0.50
1:C:187:MET:HG2	1:C:219:ILE:HG13	1.94	0.50
1:D:277:GLN:O	1:D:281:THR:CB	2.60	0.50
1:C:65:ARG:N	1:C:66:PRO:CD	2.75	0.49
1:B:1:MET:HE1	1:B:285:ARG:HH11	1.76	0.49
1:B:301:GLU:N	1:B:301:GLU:OE1	2.45	0.49
1:B:220:ILE:O	1:B:227:MET:HG2	2.13	0.48
1:C:46:ARG:NH1	4:C:2008:HOH:O	2.47	0.48
1:D:62:ARG:C	1:D:63:ARG:HG3	2.35	0.47
1:B:273:LEU:O	1:B:277:GLN:HG3	2.14	0.47
1:B:189:ASN:ND2	4:B:2038:HOH:O	2.40	0.47
1:A:256:GLY:HA3	1:A:263:THR:HG22	1.95	0.47
1:A:181:LEU:N	1:A:181:LEU:HD12	2.30	0.47
1:D:107:LYS:HD2	1:D:151:ASP:HB2	1.96	0.47
1:D:278:ALA:CA	1:D:281:THR:HB	2.43	0.47
1:C:107:LYS:NZ	1:C:149:GLU:OE1	2.48	0.47
1:A:169:THR:HG23	4:A:2028:HOH:O	2.15	0.46
1:A:247:ASP:HB2	1:A:269:GLY:HA3	1.97	0.46
1:A:40:ALA:HB3	1:A:41:PRO:HD3	1.97	0.46
1:D:256:GLY:HA3	1:D:263:THR:HG22	1.97	0.46
1:C:134:VAL:HG11	1:C:136:HIS:CE1	2.51	0.46
1:B:22:VAL:HG11	1:B:24:PHE:CZ	2.51	0.46
1:D:270:VAL:HG12	1:D:271:SER:N	2.30	0.45
1:C:95:LEU:HD13	1:C:226:TYR:CE1	2.52	0.45
1:B:7:LEU:HB2	1:B:291:TYR:HB3	1.98	0.45
1:B:12:LEU:HG	1:B:302:GLY:HA3	1.99	0.45
1:A:109:TYR:CE1	1:A:129:GLY:HA3	2.51	0.45
1:D:16:LEU:O	1:D:52:LYS:HE2	2.17	0.45
1:D:219:ILE:N	1:D:219:ILE:HD12	2.32	0.44
1:A:36:ILE:HG22	1:A:37:VAL:N	2.32	0.44
1:D:40:ALA:N	1:D:41:PRO:CD	2.81	0.44
1:D:40:ALA:HB3	1:D:41:PRO:HD3	2.00	0.44
1:B:42:GLU:HB3	1:B:300:MET:HE1	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:120:LEU:O	1:A:123:VAL:HG12	2.17	0.44
1:D:273:LEU:HD11	1:D:292:TYR:CE2	2.52	0.44
1:A:4:CYS:HA	1:A:289:PRO:O	2.17	0.44
1:B:80:LEU:HA	1:B:80:LEU:HD23	1.93	0.44
1:C:23:LEU:HD13	1:C:55:LEU:HB2	1.99	0.44
1:A:65:ARG:N	1:A:66:PRO:CD	2.81	0.44
1:A:1:MET:O	1:A:3:ARG:HG3	2.18	0.44
1:C:247:ASP:OD1	4:C:2004:HOH:O	2.21	0.44
1:A:183:VAL:CG1	1:A:220:ILE:CD1	2.95	0.44
1:A:40:ALA:N	1:A:41:PRO:CD	2.80	0.44
1:D:3:ARG:CD	1:D:5:GLU:OE2	2.65	0.44
1:B:221:GLY:O	1:B:224:SER:HB3	2.17	0.43
1:C:123:VAL:CG1	1:C:123:VAL:O	2.65	0.43
1:C:94:ARG:HD2	1:C:97:ARG:NH2	2.34	0.43
1:C:22:VAL:HG11	1:C:24:PHE:CZ	2.53	0.43
1:C:96:LEU:HD12	1:C:120:LEU:HD22	2.00	0.43
1:C:40:ALA:N	1:C:41:PRO:CD	2.81	0.43
1:D:219:ILE:CD1	1:D:219:ILE:N	2.80	0.43
1:C:220:ILE:N	1:C:220:ILE:CD1	2.82	0.43
1:B:248:ARG:HD2	1:B:270:VAL:HG11	2.00	0.43
1:A:183:VAL:HG13	1:A:220:ILE:HD11	2.00	0.43
1:A:140:PRO:HB3	1:B:193:LEU:HD21	1.99	0.43
1:A:43:LEU:N	1:A:300:MET:HE2	2.33	0.43
1:B:67:GLU:OE2	4:B:2015:HOH:O	2.21	0.42
1:C:140:PRO:O	1:C:144:LEU:HD13	2.18	0.42
1:B:120:LEU:O	1:B:123:VAL:HG12	2.19	0.42
1:A:216:GLN:CG	1:A:216:GLN:O	2.68	0.42
1:D:140:PRO:HA	1:D:143:TRP:CE3	2.54	0.42
1:B:40:ALA:N	1:B:41:PRO:CD	2.82	0.42
1:C:146:VAL:O	1:C:146:VAL:HG23	2.19	0.42
1:D:152:VAL:HG21	1:D:174:TYR:CE1	2.55	0.42
1:D:72:PHE:CD1	1:D:80:LEU:CD1	3.02	0.42
1:B:40:ALA:HB3	1:B:41:PRO:HD3	2.02	0.42
1:A:8:ARG:O	1:A:9:GLY:C	2.58	0.42
1:B:12:LEU:HG	1:B:302:GLY:CA	2.49	0.42
1:B:223:PRO:HG3	1:B:254:LEU:HD23	2.00	0.42
1:A:109:TYR:CE2	1:A:111:LEU:HD11	2.55	0.42
1:D:283:GLY:HA2	1:D:285:ARG:NE	2.35	0.42
1:A:39:GLY:O	1:A:300:MET:CE	2.68	0.42
1:B:65:ARG:NE	1:B:85:LEU:O	2.53	0.42
1:A:220:ILE:O	1:A:227:MET:HG2	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:137:GLY:O	1:B:138:ASP:HB2	2.20	0.41
1:C:134:VAL:CG1	1:C:136:HIS:CE1	3.03	0.41
1:A:107:LYS:HG2	1:A:126:THR:HB	2.02	0.41
1:D:42:GLU:HB2	1:D:300:MET:HE1	2.03	0.41
1:A:148:LEU:CD1	1:A:173:ARG:NH1	2.83	0.41
1:A:96:LEU:HD12	1:A:120:LEU:HD22	2.02	0.41
1:B:65:ARG:N	1:B:66:PRO:CD	2.83	0.41
1:C:220:ILE:N	1:C:220:ILE:HD12	2.36	0.41
1:C:234:ASP:HB3	1:C:235:PHE:CD1	2.56	0.41
1:A:210:GLU:O	1:A:214:GLN:N	2.54	0.41
1:D:181:LEU:N	1:D:181:LEU:HD12	2.36	0.41
1:B:146:VAL:O	1:B:146:VAL:HG13	2.20	0.41
1:A:146:VAL:HG21	1:A:166:MET:HE1	2.03	0.40
1:A:95:LEU:HD13	1:A:226:TYR:CE1	2.57	0.40
1:B:140:PRO:HA	1:B:143:TRP:CE3	2.56	0.40
1:D:210:GLU:O	1:D:214:GLN:N	2.54	0.40
1:D:287:LEU:HD12	1:D:287:LEU:N	2.36	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	302/309 (98%)	295 (98%)	6 (2%)	1 (0%)	46	72
1	B	302/309 (98%)	295 (98%)	6 (2%)	1 (0%)	46	72
1	C	302/309 (98%)	295 (98%)	7 (2%)	0	100	100
1	D	302/309 (98%)	295 (98%)	6 (2%)	1 (0%)	46	72
All	All	1208/1236 (98%)	1180 (98%)	25 (2%)	3 (0%)	52	77

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	101	ALA
1	A	9	GLY
1	B	138	ASP

### 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	241/244 (99%)	236 (98%)	5 (2%)	61	85
1	B	241/244 (99%)	236 (98%)	5 (2%)	61	85
1	C	241/244 (99%)	234 (97%)	7 (3%)	50	77
1	D	241/244 (99%)	237 (98%)	4 (2%)	68	88
All	All	964/976 (99%)	943 (98%)	21 (2%)	60	84

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	126	THR
1	A	135	LEU
1	A	138	ASP
1	A	281	THR
1	A	286	ASP
1	B	3	ARG
1	B	15	VAL
1	B	153	ARG
1	B	191	LEU
1	B	229	GLN
1	C	80	LEU
1	C	100	LEU
1	C	123	VAL
1	C	135	LEU
1	C	153	ARG
1	C	194	GLU
1	C	285	ARG
1	D	135	LEU
1	D	136	HIS

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Mol	Chain	Res	Type
1	D	229	GLN
1	D	234	ASP

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

Mol	Chain	Res	Type
1	C	136	HIS
1	D	136	HIS

### 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 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	NO3	A	1305	-	3,3,3	3.25	3 (100%)	3,3,3	0.13	0
3	NO3	A	1306	-	3,3,3	3.29	3 (100%)	3,3,3	0.10	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NO3	A	1305	-	-	0/0/0/0	0/0/0/0
3	NO3	A	1306	-	-	0/0/0/0	0/0/0/0

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1305	NO3	O3-N	2.86	1.40	1.25
3	A	1305	NO3	O2-N	2.90	1.40	1.25
3	A	1306	NO3	O2-N	2.91	1.40	1.25
3	A	1306	NO3	O3-N	2.91	1.40	1.25
3	A	1305	NO3	O1-N	3.90	1.40	1.24
3	A	1306	NO3	O1-N	3.94	1.40	1.24

There are no bond angle outliers.

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
3	A	1305	NO3	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	304/309 (98%)	0.67	32 (10%) <b>8</b> <b>6</b>	53, 79, 118, 190	0
1	B	304/309 (98%)	0.69	31 (10%) <b>9</b> <b>6</b>	49, 83, 132, 168	0
1	C	304/309 (98%)	0.61	30 (9%) <b>9</b> <b>7</b>	50, 80, 130, 148	0
1	D	304/309 (98%)	0.61	21 (6%) <b>20</b> <b>17</b>	52, 85, 135, 182	0
All	All	1216/1236 (98%)	0.65	114 (9%) <b>11</b> <b>8</b>	49, 82, 131, 190	0

All (114) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	280	LEU	8.1
1	D	283	GLY	6.6
1	D	282	ALA	6.5
1	C	281	THR	6.0
1	B	289	PRO	5.7
1	B	282	ALA	5.5
1	D	284	GLN	5.3
1	B	7	LEU	5.2
1	C	285	ARG	5.0
1	A	18	GLN	4.9
1	B	136	HIS	4.8
1	C	276	ALA	4.7
1	A	8	ARG	4.7
1	C	273	LEU	4.7
1	A	136	HIS	4.7
1	B	284	GLN	4.6
1	C	277	GLN	4.6
1	B	137	GLY	4.6
1	C	288	VAL	4.4
1	A	14	ASP	4.4
1	A	35	ARG	4.3

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Mol	Chain	Res	Type	RSRZ
1	C	282	ALA	4.3
1	B	283	GLY	4.3
1	A	282	ALA	4.3
1	A	276	ALA	4.2
1	A	134	VAL	4.2
1	A	241	ARG	4.2
1	D	285	ARG	4.1
1	A	288	VAL	4.1
1	C	283	GLY	4.1
1	B	73	ALA	4.0
1	A	11	ALA	3.9
1	A	17	GLY	3.9
1	B	134	VAL	3.9
1	A	1	MET	3.8
1	C	3	ARG	3.7
1	B	8	ARG	3.7
1	C	11	ALA	3.5
1	A	7	LEU	3.5
1	B	4	CYS	3.5
1	B	0	GLY	3.5
1	B	138	ASP	3.5
1	C	74	ARG	3.5
1	A	13	ARG	3.5
1	D	75	LEU	3.4
1	B	281	THR	3.3
1	C	105	ASP	3.2
1	D	281	THR	3.1
1	C	279	TYR	3.1
1	A	240	ALA	3.1
1	A	280	LEU	3.1
1	A	10	ALA	3.0
1	C	10	ALA	3.0
1	A	274	GLU	3.0
1	C	278	ALA	3.0
1	D	7	LEU	2.9
1	B	292	TYR	2.9
1	C	35	ARG	2.8
1	A	15	VAL	2.8
1	C	274	GLU	2.8
1	D	216	GLN	2.8
1	B	273	LEU	2.7
1	B	9	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	286	ASP	2.7
1	B	1	MET	2.7
1	C	1	MET	2.7
1	B	288	VAL	2.6
1	B	35	ARG	2.6
1	B	274	GLU	2.6
1	D	279	TYR	2.6
1	D	0	GLY	2.6
1	B	135	LEU	2.6
1	B	6	ARG	2.6
1	D	3	ARG	2.6
1	A	238	ASP	2.5
1	D	292	TYR	2.5
1	D	35	ARG	2.5
1	A	19	ALA	2.5
1	A	6	ARG	2.5
1	B	98	GLN	2.5
1	D	288	VAL	2.5
1	A	3	ARG	2.5
1	C	284	GLN	2.4
1	C	4	CYS	2.4
1	D	233	GLU	2.4
1	A	135	LEU	2.4
1	B	278	ALA	2.4
1	C	6	ARG	2.3
1	A	277	GLN	2.3
1	C	5	GLU	2.3
1	A	302	GLY	2.3
1	C	136	HIS	2.2
1	B	20	GLN	2.2
1	A	291	TYR	2.2
1	C	292	TYR	2.2
1	B	280	LEU	2.2
1	A	260	GLY	2.2
1	D	136	HIS	2.2
1	D	269	GLY	2.2
1	D	280	LEU	2.2
1	B	2	ALA	2.2
1	A	286	ASP	2.2
1	B	133	ASP	2.2
1	B	291	TYR	2.1
1	D	8	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	7	LEU	2.1
1	D	72	PHE	2.1
1	D	20	GLN	2.1
1	A	283	GLY	2.1
1	A	281	THR	2.1
1	C	15	VAL	2.1
1	B	272	SER	2.1
1	C	67	GLU	2.1
1	C	287	LEU	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	NO3	A	1306	4/4	0.94	0.27	3.66	64,73,81,96	0
2	MG	A	1304	1/1	0.93	0.27	1.54	88,88,88,88	0
3	NO3	A	1305	4/4	0.91	0.24	1.47	77,84,93,95	0
2	MG	D	1304	1/1	0.97	0.22	0.44	67,67,67,67	0
2	MG	C	1304	1/1	0.96	0.18	-0.73	61,61,61,61	0
2	MG	B	1304	1/1	0.96	0.18	-0.87	59,59,59,59	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.