



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 05:13 am GMT

PDB ID : 4GNK
Title : Crystal structure of Galphaq in complex with full-length human PLCbeta3
Authors : Lyon, A.M.; Tesmer, J.J.G.
Deposited on : 2012-08-17
Resolution : 4.00 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

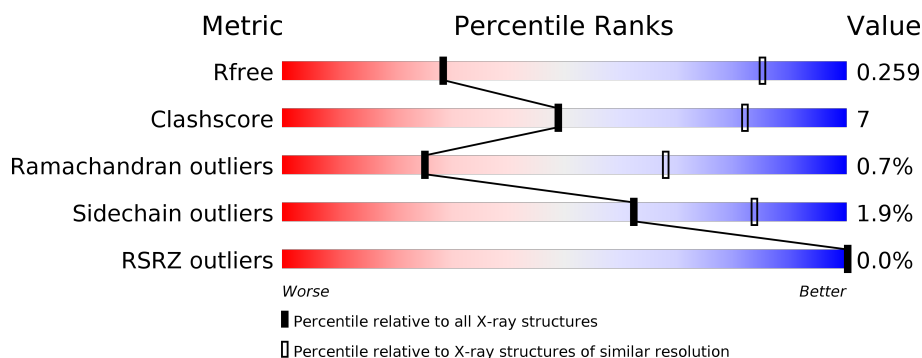
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.00 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	1088 (4.40-3.60)
Clashscore	112137	1187 (4.40-3.60)
Ramachandran outliers	110173	1139 (4.40-3.60)
Sidechain outliers	110143	1126 (4.40-3.60)
RSRZ outliers	101464	1099 (4.40-3.60)

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	353	
1	C	353	
2	B	1235	
2	D	1235	
3	E	259	

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 19764 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 Guanine nucleotide-binding protein G(q) subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	337	Total	C	N	O	S	0	0	0
			2794	1774	484	524	12			
1	C	321	Total	C	N	O	S	0	0	0
			2648	1689	447	500	12			

- Molecule 2 is a protein called 1-phosphatidylinositol 4,5-bisphosphate phosphodiesterase beta-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	769	Total	C	N	O	S	0	0	0
			6160	3926	1051	1152	31			
2	D	766	Total	C	N	O	S	0	0	0
			6132	3907	1047	1147	31			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	EXPRESSION TAG	UNP Q01970
B	1	ALA	-	EXPRESSION TAG	UNP Q01970
B	2	HIS	-	EXPRESSION TAG	UNP Q01970
B	3	HIS	-	EXPRESSION TAG	UNP Q01970
B	4	HIS	-	EXPRESSION TAG	UNP Q01970
B	5	HIS	-	EXPRESSION TAG	UNP Q01970
B	6	HIS	-	EXPRESSION TAG	UNP Q01970
B	7	HIS	-	EXPRESSION TAG	UNP Q01970
B	8	GLY	-	EXPRESSION TAG	UNP Q01970
B	9	THR	-	EXPRESSION TAG	UNP Q01970
D	0	MET	-	EXPRESSION TAG	UNP Q01970
D	1	ALA	-	EXPRESSION TAG	UNP Q01970
D	2	HIS	-	EXPRESSION TAG	UNP Q01970
D	3	HIS	-	EXPRESSION TAG	UNP Q01970
D	4	HIS	-	EXPRESSION TAG	UNP Q01970
D	5	HIS	-	EXPRESSION TAG	UNP Q01970

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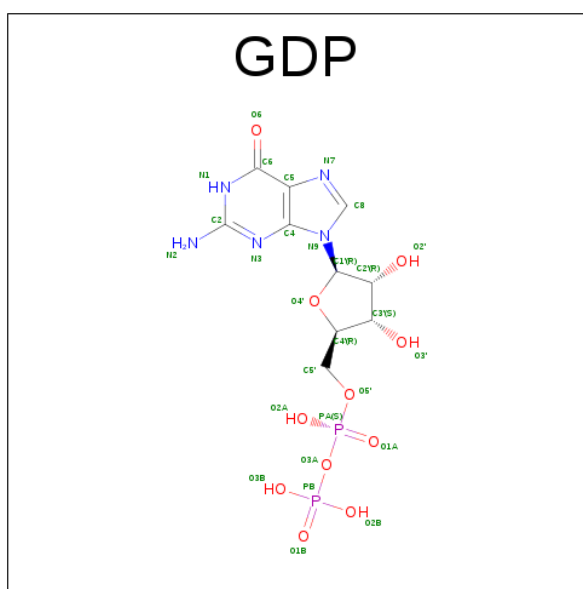
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Chain	Residue	Modelled	Actual	Comment	Reference
D	6	HIS	-	EXPRESSION TAG	UNP Q01970
D	7	HIS	-	EXPRESSION TAG	UNP Q01970
D	8	GLY	-	EXPRESSION TAG	UNP Q01970
D	9	THR	-	EXPRESSION TAG	UNP Q01970

- Molecule 3 is a protein called 1-phosphatidylinositol 4,5-bisphosphate phosphodiesterase beta-3.

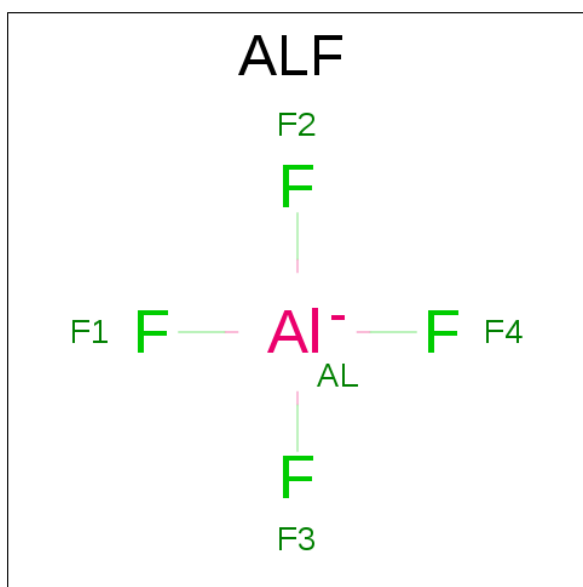
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	235	Total	C	N	O	S	0	0	0
			1954	1186	403	362	3			

- Molecule 4 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: $C_{10}H_{15}N_5O_{11}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
4	C	1	Total	C	N	O	P	0	0
			28	10	5	11	2		

- Molecule 5 is TETRAFLUOROALUMINATE ION (three-letter code: ALF) (formula: AlF_4).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	Al	F	0	0
			5	1	4		
5	C	1	Total	Al	F	0	0
			5	1	4		

- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Mg	0	0
			1	1		
6	C	1	Total	Mg	0	0
			1	1		

- Molecule 7 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	1	Total	Ca	0	0
			1	1		
7	D	1	Total	Ca	0	0
			1	1		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	3	Total	O	0	0
			3	3		

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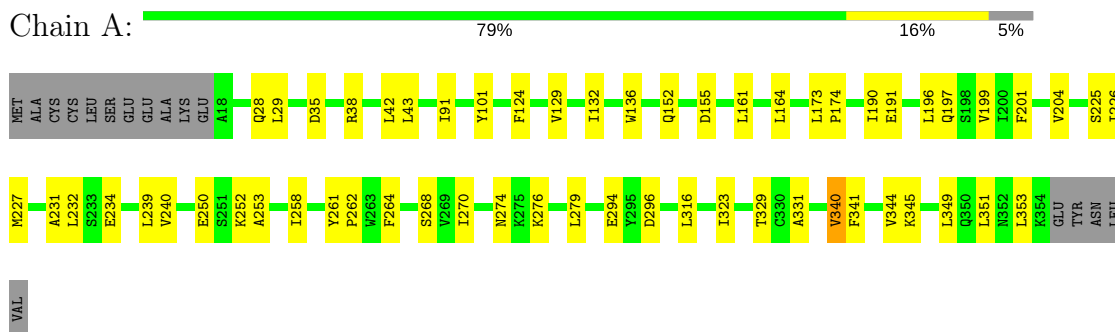
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	C	3	Total	O	0	0
			3	3		

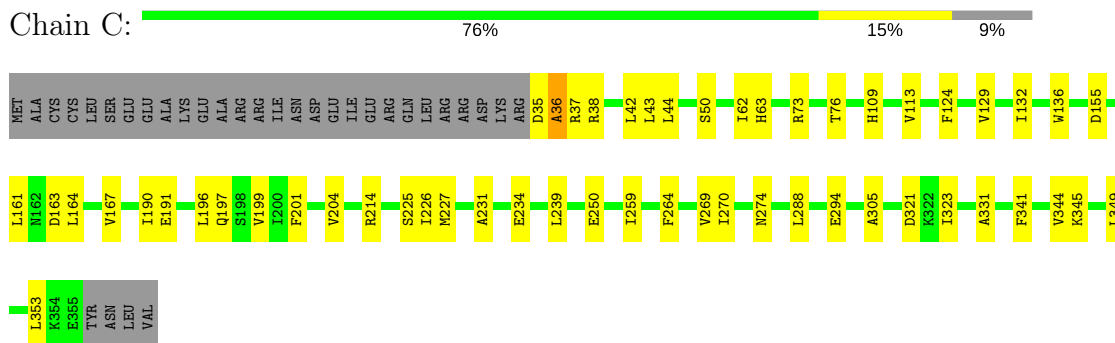
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

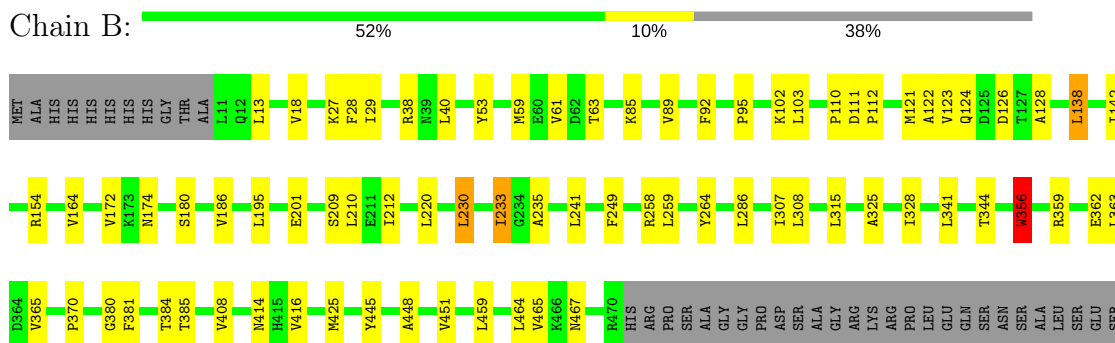
- Molecule 1: Guanine nucleotide-binding protein G(q) subunit alpha

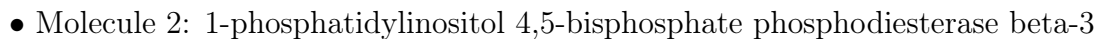


- Molecule 1: Guanine nucleotide-binding protein G(q) subunit alpha



- Molecule 2: 1-phosphatidylinositol 4,5-bisphosphate phosphodiesterase beta-3





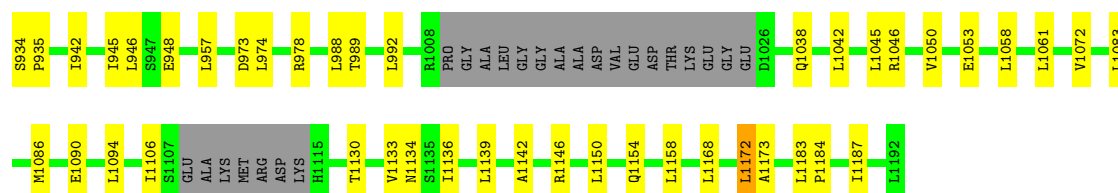
Device Type	Percentage
Smartphone	52%
Tablet	10%
Feature phone	38%



[illegible]

- Molecule 3: 1-phosphatidylinositol 4,5-bisphosphate phosphodiesterase beta-3

Chain E:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	91.42Å 188.84Å 293.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.99 – 4.00 19.99 – 4.00	Depositor EDS
% Data completeness (in resolution range)	77.8 (19.99-4.00) 78.5 (19.99-4.00)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.17 (at 4.07Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.214 , 0.255 0.222 , 0.259	Depositor DCC
R_{free} test set	1715 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	96.5	Xtriage
Anisotropy	0.126	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 25.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.37$, $\langle L^2 \rangle = 0.20$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	19764	wwPDB-VP
Average B, all atoms (Å ²)	131.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.62% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GDP, ALF, CA, 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	0.39	0/2850	0.49	0/3846
1	C	0.40	0/2704	0.50	0/3654
2	B	0.38	1/6295 (0.0%)	0.49	0/8523
2	D	0.38	0/6264	0.49	0/8479
3	E	0.37	0/1966	0.47	0/2622
All	All	0.38	1/20079 (0.0%)	0.49	0/27124

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	356	TRP	CD2-CE2	5.02	1.47	1.41

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	2794	0	2767	50	0
1	C	2648	0	2610	36	0
2	B	6160	0	6170	82	0
2	D	6132	0	6144	81	0
3	E	1954	0	2027	39	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	28	0	12	1	0
4	C	28	0	12	1	0
5	A	5	0	0	0	0
5	C	5	0	0	0	0
6	A	1	0	0	0	0
6	C	1	0	0	0	0
7	B	1	0	0	0	0
7	D	1	0	0	0	0
8	A	3	0	0	0	0
8	C	3	0	0	0	0
All	All	19764	0	19742	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 7.

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)
2:B:18:VAL:HG22	2:B:138:LEU:HD11	1.48	0.95
2:D:18:VAL:HG22	2:D:138:LEU:HD11	1.48	0.93
2:B:195:LEU:HD11	2:B:212:ILE:HG21	1.51	0.91
2:D:13:LEU:HD11	2:D:142:ILE:HD12	1.54	0.89
3:E:1058:LEU:HD21	3:E:1173:ALA:HB2	1.56	0.88
2:B:13:LEU:HD21	2:B:142:ILE:HD12	1.57	0.86
2:B:29:ILE:HD13	2:D:40:LEU:HD22	1.60	0.84
3:E:1061:LEU:HD22	3:E:1172:LEU:HD22	1.64	0.80
1:C:239:LEU:HD11	1:C:250:GLU:HG3	1.63	0.78
3:E:1094:LEU:HD22	3:E:1139:LEU:HD22	1.65	0.76
2:D:230:LEU:O	2:D:235:ALA:HB3	1.88	0.73
1:A:239:LEU:HD11	1:A:250:GLU:HG3	1.71	0.72
1:A:226:ILE:HD11	1:A:264:PHE:CD1	2.23	0.72
3:E:1042:LEU:HD23	3:E:1187:ILE:HG23	1.73	0.70
2:B:723:ILE:HD13	2:B:788:LEU:HD13	1.72	0.70
2:B:40:LEU:HD22	2:D:29:ILE:HD13	1.74	0.70
2:B:416:VAL:HG21	2:B:425:MET:HE1	1.74	0.70
1:A:29:LEU:HD11	3:E:1136:ILE:HD13	1.74	0.70
1:C:231:ALA:HB3	1:C:234:GLU:HG3	1.72	0.69
2:D:723:ILE:HD13	2:D:788:LEU:HD13	1.74	0.69
2:B:40:LEU:HD22	2:D:29:ILE:CD1	2.23	0.68
2:D:416:VAL:HG21	2:D:425:MET:HE1	1.77	0.67
2:B:230:LEU:O	2:B:235:ALA:HB3	1.94	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:191:GLU:HG2	1:C:204:VAL:HG22	1.77	0.66
2:B:629:LEU:HD12	2:B:667:VAL:HG11	1.77	0.66
1:A:231:ALA:HB3	1:A:234:GLU:HG3	1.78	0.66
2:D:629:LEU:HD12	2:D:667:VAL:HG11	1.76	0.66
2:B:195:LEU:HD11	2:B:212:ILE:CG2	2.23	0.65
2:D:416:VAL:HG21	2:D:425:MET:CE	2.26	0.65
2:B:620:PHE:HB3	2:B:625:ALA:HB2	1.78	0.65
3:E:1046:ARG:O	3:E:1050:VAL:HG23	1.98	0.64
2:B:713:ASP:HB3	2:B:716:THR:HG23	1.80	0.64
1:A:29:LEU:CD2	3:E:1133:VAL:HG13	2.27	0.64
2:D:620:PHE:HB3	2:D:625:ALA:HB2	1.80	0.63
2:B:344:THR:HG22	2:B:385:THR:HG23	1.81	0.63
2:D:713:ASP:HB3	2:D:716:THR:HG23	1.80	0.63
2:B:416:VAL:HG21	2:B:425:MET:CE	2.29	0.63
3:E:942:ILE:HG23	3:E:1139:LEU:HD13	1.79	0.63
1:A:43:LEU:HD23	1:A:204:VAL:HB	1.78	0.63
1:C:270:ILE:HD13	1:C:344:VAL:HG13	1.82	0.62
2:D:235:ALA:HB2	2:D:241:LEU:HD21	1.82	0.62
2:B:154:ARG:HG3	2:B:210:LEU:HD11	1.81	0.61
2:D:154:ARG:HG3	2:D:210:LEU:HD11	1.81	0.61
2:B:459:LEU:HD13	2:B:464:LEU:HD21	1.83	0.60
1:C:196:LEU:HD22	1:C:201:PHE:CE2	2.36	0.60
2:D:235:ALA:HB2	2:D:241:LEU:CD2	2.32	0.60
2:D:344:THR:HG22	2:D:385:THR:HG23	1.82	0.60
1:C:44:LEU:HD23	1:C:227:MET:HB2	1.85	0.59
2:B:235:ALA:HB2	2:B:241:LEU:HD21	1.85	0.59
2:B:38:ARG:NE	2:B:89:VAL:HG11	2.18	0.59
1:A:29:LEU:HD13	3:E:1136:ILE:HG21	1.84	0.59
3:E:988:LEU:HD22	3:E:1045:LEU:HD22	1.84	0.59
1:A:196:LEU:HD22	1:A:201:PHE:CE2	2.38	0.58
2:D:464:LEU:N	2:D:464:LEU:HD12	2.18	0.58
2:D:676:PHE:HA	2:D:683:MET:HE3	1.84	0.58
1:A:232:LEU:HB3	1:A:279:LEU:HD22	1.85	0.58
2:B:235:ALA:HB2	2:B:241:LEU:CD2	2.34	0.58
2:B:464:LEU:N	2:B:464:LEU:HD12	2.18	0.58
1:A:29:LEU:CD1	3:E:1136:ILE:HG21	2.34	0.58
2:D:459:LEU:HD13	2:D:464:LEU:HD21	1.85	0.58
2:B:61:VAL:HG21	2:B:164:VAL:HG22	1.86	0.58
2:D:465:VAL:HG23	2:D:594:VAL:HG22	1.86	0.58
2:B:356:TRP:CH2	2:B:687:ALA:HB1	2.39	0.58
1:A:258:ILE:HD11	2:B:859:LEU:HD21	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:85:LYS:O	2:B:89:VAL:HG23	2.04	0.57
2:D:728:LEU:HD23	2:D:729:ARG:N	2.18	0.57
1:A:253:ALA:CB	2:B:866:VAL:HG21	2.34	0.57
2:B:465:VAL:HG23	2:B:594:VAL:HG22	1.87	0.57
3:E:1083:LEU:HD12	3:E:1086:MET:CE	2.34	0.57
2:B:344:THR:HG22	2:B:385:THR:CG2	2.35	0.57
2:D:700:LEU:HD23	2:D:701:LYS:O	2.05	0.57
2:B:676:PHE:HA	2:B:683:MET:HE3	1.87	0.56
2:D:605:PHE:O	2:D:609:ARG:HB2	2.06	0.56
1:A:29:LEU:HD21	3:E:1133:VAL:HG13	1.87	0.56
1:A:331:ALA:HB3	4:A:401:GDP:N7	2.21	0.56
1:C:38:ARG:HH21	1:C:353:LEU:HD23	1.71	0.56
2:D:344:THR:HG22	2:D:385:THR:CG2	2.36	0.56
2:D:38:ARG:NE	2:D:89:VAL:HG11	2.21	0.56
2:B:700:LEU:HD23	2:B:701:LYS:O	2.06	0.56
2:B:728:LEU:HD22	2:B:730:VAL:HG23	1.89	0.55
2:B:605:PHE:O	2:B:609:ARG:HB2	2.06	0.55
3:E:1042:LEU:HD23	3:E:1187:ILE:CG2	2.37	0.55
1:A:38:ARG:HH21	1:A:353:LEU:HD23	1.72	0.55
2:D:61:VAL:HG21	2:D:164:VAL:HG22	1.88	0.55
2:D:356:TRP:CH2	2:D:687:ALA:HB1	2.41	0.55
1:C:42:LEU:HD22	1:C:225:SER:HB3	1.89	0.54
2:B:728:LEU:HD23	2:B:729:ARG:N	2.22	0.54
1:A:253:ALA:HB1	2:B:866:VAL:HG21	1.89	0.54
1:C:124:PHE:CE2	1:C:129:VAL:HG22	2.42	0.54
2:D:85:LYS:O	2:D:89:VAL:HG23	2.08	0.54
3:E:988:LEU:HD22	3:E:1045:LEU:CD2	2.38	0.54
3:E:1154:GLN:O	3:E:1158:LEU:HD13	2.09	0.53
2:B:172:VAL:HG23	2:B:201:GLU:C	2.28	0.53
1:A:196:LEU:HD21	1:A:341:PHE:HE2	1.73	0.53
1:C:196:LEU:HD13	1:C:201:PHE:HD2	1.74	0.53
2:B:341:LEU:O	2:B:384:THR:HG23	2.09	0.52
1:C:35:ASP:O	1:C:36:ALA:HB2	2.09	0.52
1:A:261:TYR:OH	2:B:855:TYR:O	2.28	0.52
2:D:111:ASP:HB2	2:D:112:PRO:HD2	1.91	0.52
2:B:465:VAL:HG21	2:B:593:LEU:HB2	1.91	0.52
1:C:43:LEU:HD23	1:C:204:VAL:HB	1.91	0.52
1:C:196:LEU:HD21	1:C:341:PHE:HE2	1.74	0.52
2:D:728:LEU:HD22	2:D:730:VAL:HG23	1.91	0.52
3:E:946:LEU:HD22	3:E:1146:ARG:NH1	2.25	0.52
2:D:465:VAL:HG21	2:D:593:LEU:HB2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:349:LEU:HD22	1:A:353:LEU:HD11	1.92	0.51
2:B:840:ILE:HD13	2:B:842:THR:HG23	1.92	0.51
2:D:233:ILE:HG21	2:D:249:PHE:CD1	2.45	0.51
3:E:1130:THR:HG22	3:E:1134:ASN:ND2	2.26	0.51
1:C:136:TRP:CE3	1:C:161:LEU:HD13	2.46	0.51
1:A:274:ASN:HD21	1:A:331:ALA:HB2	1.75	0.51
3:E:1061:LEU:HD11	3:E:1168:LEU:HD13	1.91	0.51
2:B:111:ASP:HB2	2:B:112:PRO:HD2	1.92	0.51
1:A:253:ALA:HB1	2:B:866:VAL:CG2	2.41	0.51
1:A:42:LEU:HD22	1:A:225:SER:HB3	1.93	0.50
2:D:414:ASN:O	2:D:584:VAL:HG22	2.10	0.50
1:C:196:LEU:HD21	1:C:341:PHE:CE2	2.47	0.50
3:E:957:LEU:HD12	3:E:1072:VAL:HG22	1.93	0.50
1:C:163:ASP:O	1:C:167:VAL:HG23	2.12	0.50
2:D:840:ILE:HD13	2:D:842:THR:HG23	1.93	0.50
2:B:467:ASN:HD22	2:B:590:MET:CE	2.25	0.50
3:E:1083:LEU:HD12	3:E:1086:MET:HE2	1.94	0.50
1:A:42:LEU:HD11	1:A:201:PHE:CD1	2.47	0.50
1:A:124:PHE:CE2	1:A:129:VAL:HG22	2.47	0.50
1:C:190:ILE:HD12	1:C:190:ILE:N	2.26	0.50
2:B:233:ILE:HG21	2:B:249:PHE:CD1	2.46	0.49
2:D:341:LEU:O	2:D:384:THR:HG23	2.11	0.49
3:E:1061:LEU:HD21	3:E:1168:LEU:HB3	1.94	0.49
1:A:152:GLN:HG3	1:A:240:VAL:HG23	1.94	0.49
3:E:992:LEU:HD22	3:E:1038:GLN:HB3	1.94	0.49
1:C:226:ILE:HD11	1:C:264:PHE:CD1	2.48	0.49
1:C:226:ILE:HD11	1:C:264:PHE:CE1	2.48	0.49
1:A:28:GLN:HG2	3:E:1133:VAL:HG21	1.94	0.49
1:A:323:ILE:HG23	1:A:323:ILE:O	2.12	0.48
1:A:329:THR:HG22	1:A:340:VAL:HG21	1.96	0.48
1:A:38:ARG:NH2	1:A:353:LEU:HD23	2.28	0.48
2:B:53:TYR:HB3	2:B:63:THR:HG22	1.96	0.48
1:A:91:ILE:HG23	1:A:101:TYR:OH	2.14	0.48
2:B:259:LEU:HD22	2:B:264:TYR:CE1	2.48	0.48
3:E:1083:LEU:HD22	3:E:1150:LEU:CD2	2.43	0.48
2:D:53:TYR:HB3	2:D:63:THR:HG22	1.96	0.48
2:B:380:GLY:HA3	2:B:579:THR:HG21	1.96	0.47
1:C:349:LEU:HD22	1:C:353:LEU:HD11	1.95	0.47
2:B:629:LEU:CD1	2:B:667:VAL:HG11	2.44	0.47
2:D:307:ILE:HG23	2:D:308:LEU:HG	1.96	0.47
2:B:414:ASN:O	2:B:584:VAL:HG22	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:268:SER:HB3	1:A:351:LEU:HD23	1.95	0.47
1:C:196:LEU:HD23	1:C:345:LYS:HD2	1.96	0.47
1:A:196:LEU:HD22	1:A:201:PHE:CD2	2.50	0.47
1:A:199:VAL:HG11	1:A:349:LEU:HD21	1.97	0.47
2:D:869:MET:HG3	2:D:870:ASP:N	2.30	0.47
1:A:227:MET:HE2	1:A:344:VAL:HG11	1.95	0.47
1:C:239:LEU:HD11	1:C:250:GLU:CG	2.40	0.47
3:E:989:THR:HA	3:E:992:LEU:HD12	1.97	0.47
1:A:132:ILE:HG22	1:A:164:LEU:HD11	1.97	0.46
2:B:448:ALA:HB3	2:B:451:VAL:CG2	2.45	0.46
1:C:323:ILE:HG23	1:C:323:ILE:O	2.15	0.46
2:D:259:LEU:HD22	2:D:264:TYR:CE1	2.50	0.46
2:D:380:GLY:HA3	2:D:579:THR:HG21	1.97	0.46
1:A:196:LEU:HD23	1:A:345:LYS:HD2	1.96	0.46
1:A:270:ILE:HD13	1:A:344:VAL:HG13	1.97	0.46
1:C:259:ILE:HD11	1:C:269:VAL:HG11	1.98	0.46
2:D:467:ASN:HD22	2:D:590:MET:CE	2.28	0.46
1:A:226:ILE:HD11	1:A:264:PHE:CG	2.50	0.46
1:C:196:LEU:HD22	1:C:201:PHE:CD2	2.51	0.45
2:D:629:LEU:CD1	2:D:667:VAL:HG11	2.44	0.45
1:A:42:LEU:HD11	1:A:201:PHE:CE1	2.51	0.45
1:C:38:ARG:NH2	1:C:353:LEU:HD23	2.31	0.45
2:B:28:PHE:O	2:B:40:LEU:HD12	2.17	0.45
2:D:325:ALA:HB1	2:D:706:ARG:CD	2.47	0.45
2:B:307:ILE:HG23	2:B:308:LEU:HG	1.99	0.45
1:A:191:GLU:HG2	1:A:204:VAL:HG22	1.99	0.45
2:D:362:GLU:C	2:D:363:LEU:HD12	2.38	0.45
2:D:467:ASN:HD22	2:D:590:MET:HE3	1.82	0.45
2:D:258:ARG:CZ	2:D:844:ALA:HB1	2.46	0.45
1:A:196:LEU:HD21	1:A:341:PHE:CE2	2.50	0.44
2:B:445:TYR:O	2:B:596:TYR:OH	2.27	0.44
2:B:745:ILE:HD11	2:B:771:SER:CB	2.47	0.44
2:D:686:ASN:HD22	2:D:686:ASN:C	2.19	0.44
2:D:325:ALA:HB1	2:D:706:ARG:HD2	1.99	0.44
3:E:942:ILE:HG12	3:E:1094:LEU:HD21	1.98	0.44
3:E:1046:ARG:HB3	3:E:1183:LEU:HD11	1.99	0.44
2:B:325:ALA:HB1	2:B:706:ARG:CD	2.47	0.44
2:B:448:ALA:HB3	2:B:451:VAL:HG23	1.99	0.44
2:D:122:ALA:HB2	2:D:128:ALA:HB2	2.00	0.44
1:A:232:LEU:HD12	1:A:276:LYS:HA	1.98	0.44
3:E:1083:LEU:HD12	3:E:1086:MET:HE1	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:445:TYR:O	2:D:596:TYR:OH	2.27	0.44
2:D:29:ILE:HD12	2:D:121:MET:CE	2.47	0.44
2:D:337:THR:HG22	2:D:349:MET:CE	2.48	0.44
1:A:226:ILE:HD11	1:A:264:PHE:CE1	2.53	0.43
1:A:252:LYS:HE2	1:A:316:LEU:HD11	1.99	0.43
1:C:199:VAL:HG11	1:C:349:LEU:HD21	2.00	0.43
1:C:288:LEU:HD22	1:C:305:ALA:HB1	2.00	0.43
2:D:102:LYS:HA	2:D:128:ALA:HB3	2.00	0.43
2:D:745:ILE:HD11	2:D:771:SER:CB	2.48	0.43
2:B:258:ARG:CZ	2:B:844:ALA:HB1	2.48	0.43
2:D:700:LEU:HD21	2:D:706:ARG:NH2	2.32	0.43
2:B:29:ILE:HD12	2:B:121:MET:HE2	2.00	0.43
2:B:325:ALA:HB1	2:B:706:ARG:HD2	1.99	0.43
2:B:700:LEU:HD21	2:B:706:ARG:NH2	2.33	0.43
2:B:29:ILE:HD12	2:B:121:MET:CE	2.48	0.43
2:D:448:ALA:HB3	2:D:451:VAL:CG2	2.49	0.43
2:B:686:ASN:HD22	2:B:686:ASN:C	2.21	0.43
1:A:173:LEU:HD12	1:A:174:PRO:HD2	2.00	0.43
1:C:132:ILE:HG22	1:C:164:LEU:HD11	2.00	0.43
2:D:869:MET:O	2:D:872:ARG:N	2.51	0.43
2:D:233:ILE:HD11	2:D:241:LEU:CD1	2.48	0.43
1:C:274:ASN:HD21	1:C:331:ALA:HB2	1.84	0.42
2:D:365:VAL:HG23	2:D:416:VAL:HG23	2.01	0.42
2:D:665:TRP:CD1	2:D:672:VAL:HG21	2.54	0.42
2:D:92:PHE:HZ	2:D:103:LEU:HD11	1.84	0.42
1:A:261:TYR:HA	1:A:262:PRO:HD3	1.95	0.42
2:D:123:VAL:HG13	2:D:124:GLN:N	2.34	0.42
2:D:233:ILE:HD11	2:D:241:LEU:HD11	2.01	0.42
2:D:626:MET:CE	2:D:663:LEU:HD13	2.50	0.42
3:E:1090:GLU:HB3	3:E:1139:LEU:HD21	2.01	0.42
3:E:934:SER:N	3:E:935:PRO:CD	2.83	0.42
2:D:379:HIS:ND1	2:D:382:THR:HG21	2.35	0.42
2:B:362:GLU:C	2:B:363:LEU:HD12	2.40	0.42
1:C:124:PHE:CZ	1:C:129:VAL:HG22	2.53	0.42
2:D:28:PHE:O	2:D:40:LEU:HD12	2.19	0.42
2:B:27:LYS:O	2:B:123:VAL:HG12	2.20	0.42
1:C:109:HIS:O	1:C:113:VAL:HG23	2.19	0.42
2:D:350:TYR:CD2	2:D:392:VAL:HG13	2.55	0.42
2:B:102:LYS:HA	2:B:128:ALA:HB3	2.01	0.42
2:B:123:VAL:HG23	2:D:123:VAL:HG23	2.01	0.42
2:D:27:LYS:O	2:D:123:VAL:HG12	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:337:THR:HG22	2:D:349:MET:HE3	2.01	0.42
3:E:978:ARG:NE	3:E:1053:GLU:OE2	2.52	0.42
3:E:1106:ILE:O	3:E:1106:ILE:HG22	2.19	0.42
1:A:136:TRP:CE3	1:A:161:LEU:HD13	2.54	0.42
1:A:190:ILE:N	1:A:190:ILE:HD12	2.35	0.42
2:B:209:SER:H	2:B:212:ILE:HD12	1.84	0.42
2:B:186:VAL:HG22	2:B:220:LEU:HD21	2.02	0.42
2:B:233:ILE:HD11	2:B:241:LEU:CD1	2.50	0.42
2:D:209:SER:H	2:D:212:ILE:HD12	1.85	0.42
2:B:123:VAL:HG13	2:B:124:GLN:N	2.36	0.41
2:D:448:ALA:HB3	2:D:451:VAL:HG23	2.01	0.41
2:D:647:ILE:HD12	2:D:664:PHE:HB3	2.02	0.41
3:E:974:LEU:HD22	3:E:978:ARG:HH21	1.86	0.41
1:A:274:ASN:HD21	1:A:331:ALA:CB	2.34	0.41
2:B:286:LEU:C	2:B:286:LEU:HD23	2.40	0.41
1:C:73:ARG:O	1:C:76:THR:HG22	2.21	0.41
1:C:62:ILE:HG22	1:C:63:HIS:CE1	2.54	0.41
3:E:945:ILE:HD13	3:E:1094:LEU:HD11	2.01	0.41
2:B:13:LEU:HD22	2:B:142:ILE:HB	2.02	0.41
1:A:258:ILE:CD1	2:B:859:LEU:HD21	2.49	0.41
2:B:92:PHE:HZ	2:B:103:LEU:HD11	1.85	0.41
2:D:186:VAL:HG22	2:D:220:LEU:HD21	2.02	0.41
2:B:365:VAL:HG23	2:B:416:VAL:HG23	2.02	0.41
2:B:647:ILE:HD12	2:B:664:PHE:HB3	2.01	0.41
2:B:665:TRP:CD1	2:B:672:VAL:HG21	2.55	0.41
3:E:1130:THR:HG22	3:E:1134:ASN:HD21	1.86	0.41
2:B:122:ALA:HB2	2:B:128:ALA:HB2	2.02	0.41
2:B:29:ILE:HG21	2:D:40:LEU:HD22	2.02	0.41
2:B:359:ARG:O	2:B:408:VAL:HG13	2.21	0.41
3:E:946:LEU:CD1	3:E:1142:ALA:HB1	2.51	0.41
2:D:328:ILE:HB	2:D:671:LEU:HD23	2.03	0.41
2:D:840:ILE:C	2:D:840:ILE:HD12	2.41	0.40
2:B:59:MET:CE	2:D:85:LYS:HB3	2.51	0.40
3:E:1042:LEU:CD2	3:E:1187:ILE:HG23	2.46	0.40
3:E:946:LEU:HD23	3:E:1090:GLU:OE2	2.20	0.40
2:B:328:ILE:HB	2:B:671:LEU:HD23	2.02	0.40
1:C:214:ARG:HD3	2:D:848:ILE:HD11	2.04	0.40
1:C:331:ALA:HB3	4:C:401:GDP:N7	2.36	0.40
2:D:286:LEU:HD23	2:D:286:LEU:C	2.42	0.40
2:D:629:LEU:CD1	2:D:667:VAL:CG1	3.00	0.40
2:B:840:ILE:HD12	2:B:840:ILE:C	2.42	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	335/353 (95%)	319 (95%)	15 (4%)	1 (0%)	44	80
1	C	319/353 (90%)	295 (92%)	22 (7%)	2 (1%)	28	70
2	B	765/1235 (62%)	704 (92%)	54 (7%)	7 (1%)	20	63
2	D	760/1235 (62%)	697 (92%)	56 (7%)	7 (1%)	20	63
3	E	229/259 (88%)	224 (98%)	4 (2%)	1 (0%)	38	77
All	All	2408/3435 (70%)	2239 (93%)	151 (6%)	18 (1%)	25	68

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	381	PHE
1	C	36	ALA
2	D	381	PHE
1	A	197	GLN
2	B	95	PRO
2	D	95	PRO
2	B	233	ILE
2	B	595	ASN
1	C	197	GLN
2	D	233	ILE
2	D	595	ASN
2	B	180	SER
2	D	180	SER
2	B	110	PRO
2	B	633	PRO
3	E	1184	PRO
2	D	110	PRO
2	D	633	PRO

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	308/322 (96%)	303 (98%)	5 (2%)	68	86
1	C	293/322 (91%)	288 (98%)	5 (2%)	66	86
2	B	682/1076 (63%)	667 (98%)	15 (2%)	57	81
2	D	678/1076 (63%)	664 (98%)	14 (2%)	59	82
3	E	210/226 (93%)	207 (99%)	3 (1%)	71	87
All	All	2171/3022 (72%)	2129 (98%)	42 (2%)	62	83

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	ASP
1	A	155	ASP
1	A	294	GLU
1	A	296	ASP
1	A	340	VAL
2	B	126	ASP
2	B	138	LEU
2	B	174	ASN
2	B	230	LEU
2	B	315	LEU
2	B	356	TRP
2	B	370	PRO
2	B	595	ASN
2	B	632	SER
2	B	636	PHE
2	B	686	ASN
2	B	736	GLN
2	B	790	THR
2	B	808	HIS
2	B	830	ASN
1	C	37	ARG
1	C	50	SER
1	C	155	ASP

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Mol	Chain	Res	Type
1	C	294	GLU
1	C	321	ASP
2	D	126	ASP
2	D	138	LEU
2	D	174	ASN
2	D	230	LEU
2	D	315	LEU
2	D	356	TRP
2	D	370	PRO
2	D	595	ASN
2	D	636	PHE
2	D	686	ASN
2	D	736	GLN
2	D	790	THR
2	D	808	HIS
2	D	830	ASN
3	E	948	GLU
3	E	973	ASP
3	E	1172	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	352	ASN
2	B	39	ASN
2	B	467	ASN
2	D	39	ASN
2	D	467	ASN
3	E	1000	GLN
3	E	1125	ASN
3	E	1134	ASN

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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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$
4	GDP	A	401	6	25,30,30	1.16	2 (8%)	26,47,47	2.07	6 (23%)
5	ALF	A	402	-	0,4,4	0.00	-	0,6,6	0.00	-
4	GDP	C	401	6	25,30,30	1.06	2 (8%)	26,47,47	1.94	6 (23%)
5	ALF	C	402	-	0,4,4	0.00	-	0,6,6	0.00	-

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GDP	A	401	6	-	0/12/32/32	0/3/3/3
5	ALF	A	402	-	-	0/0/0/0	0/0/0/0
4	GDP	C	401	6	-	0/12/32/32	0/3/3/3
5	ALF	C	402	-	-	0/0/0/0	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	401	GDP	C5-C4	2.71	1.46	1.40
4	A	401	GDP	C5-C4	2.96	1.47	1.40
4	C	401	GDP	C6-C5	3.41	1.47	1.41
4	A	401	GDP	C6-C5	3.99	1.48	1.41

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	401	GDP	C6-C5-C4	-3.80	117.06	120.84
4	C	401	GDP	C6-C5-C4	-3.76	117.10	120.84
4	C	401	GDP	C5-C6-N1	-3.59	118.37	123.48
4	A	401	GDP	C5-C6-N1	-3.56	118.42	123.48
4	A	401	GDP	N3-C2-N1	-3.34	122.59	127.46
4	A	401	GDP	C4-C5-N7	-3.19	106.33	109.41
4	C	401	GDP	N3-C2-N1	-2.95	123.15	127.46
4	C	401	GDP	C4-C5-N7	-2.69	106.81	109.41
4	C	401	GDP	C6-N1-C2	3.95	121.74	116.06
4	A	401	GDP	C6-N1-C2	4.13	122.00	116.06
4	C	401	GDP	C2-N3-C4	4.67	120.61	115.16
4	A	401	GDP	C2-N3-C4	5.27	121.31	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	401	GDP	1	0
4	C	401	GDP	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 [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	337/353 (95%)	-0.63	0 100 100	88, 128, 178, 232	0
1	C	321/353 (90%)	-0.67	0 100 100	86, 120, 157, 192	0
2	B	769/1235 (62%)	-0.63	0 100 100	76, 119, 169, 209	0
2	D	766/1235 (62%)	-0.60	1 (0%) 95 94	78, 130, 185, 219	0
3	E	235/259 (90%)	-0.41	0 100 100	136, 167, 202, 223	0
All	All	2428/3435 (70%)	-0.60	1 (0%) 100 100	76, 127, 182, 232	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	199	ARG	2.7

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(\AA^2)	Q<0.9
7	CA	D	1301	1/1	0.95	0.24	0.08	121,121,121,121	0
4	GDP	A	401	28/28	0.97	0.16	-0.41	93,98,114,115	0
4	GDP	C	401	28/28	0.98	0.15	-0.60	83,90,96,99	0
7	CA	B	1301	1/1	0.99	0.18	-0.72	89,89,89,89	0
5	ALF	C	402	5/5	0.99	0.12	-1.32	82,82,85,85	0
6	MG	C	403	1/1	0.99	0.09	-1.40	87,87,87,87	0
6	MG	A	403	1/1	0.99	0.10	-1.52	96,96,96,96	0
5	ALF	A	402	5/5	0.99	0.11	-1.54	93,93,95,95	0

6.5 Other polymers [i](#)

There are no such residues in this entry.