



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:11 PM GMT

PDB ID : 1SHZ  
Title : Crystal Structure of the p115RhoGEF rgRGS Domain in A Complex with G  
alpha(13):Galpha(i1) Chimera  
Authors : Chen, Z.; Singer, W.D.; Sternweis, P.C.; Sprang, S.R.  
Deposited on : 2004-02-26  
Resolution : 2.85 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

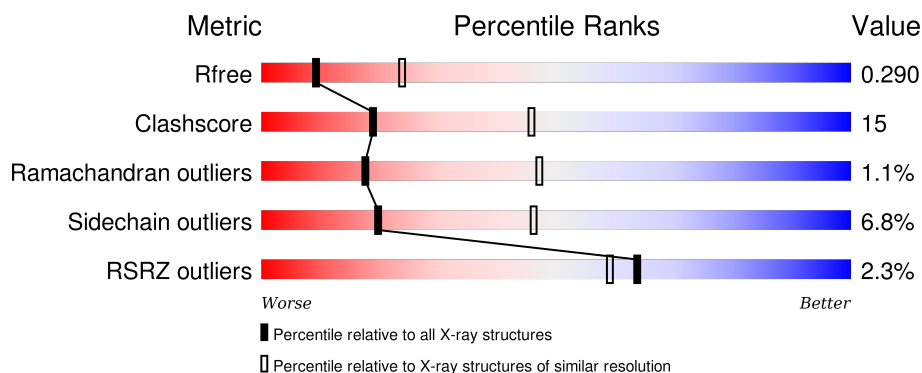
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.85 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	2228 (2.90-2.82)
Clashscore	102246	2499 (2.90-2.82)
Ramachandran outliers	100387	2439 (2.90-2.82)
Sidechain outliers	100360	2442 (2.90-2.82)
RSRZ outliers	91569	2236 (2.90-2.82)

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	340	 63% 28% 5% .
1	D	340	 59% 35% . .
2	C	233	 55% 27% . 16%
2	F	233	 51% 28% . 19%

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	MG	A	377	-	-	-	X
3	MG	D	377	-	-	-	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8602 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 Galpha(13):Galpha(i1) Chimera.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	326	Total	C	N	O	S	0	0	0
			2677	1706	463	493	15			
1	D	327	Total	C	N	O	S	0	0	0
			2682	1709	464	494	15			

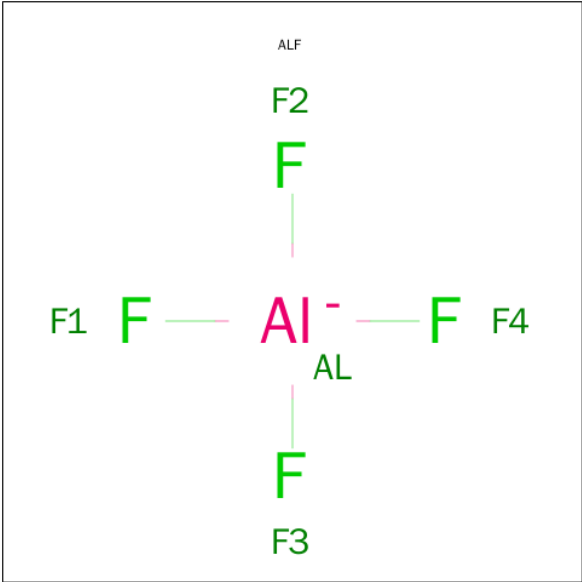
- Molecule 2 is a protein called Rho guanine nucleotide exchange factor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	196	Total	C	N	O	S	0	0	0
			1576	995	282	290	9			
2	F	188	Total	C	N	O	S	0	0	0
			1515	961	269	276	9			

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

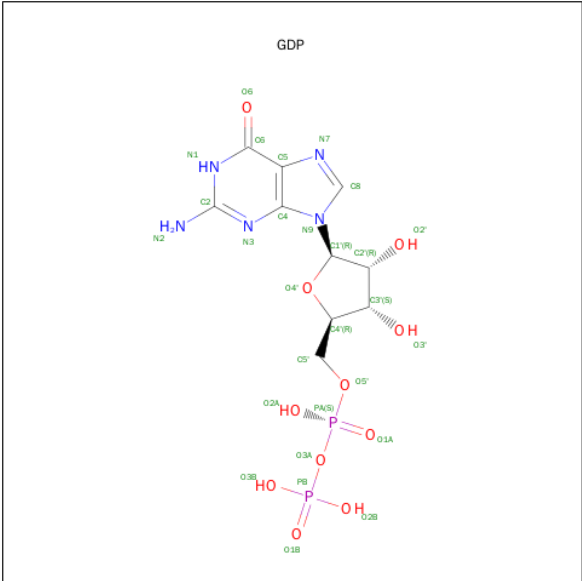
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		

- Molecule 4 is TETRAFLUOROALUMINATE ION (three-letter code: ALF) (formula:  $\text{AlF}_4$ ).



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

- Molecule 5 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: C<sub>10</sub>H<sub>15</sub>N<sub>5</sub>O<sub>11</sub>P<sub>2</sub>).



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

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

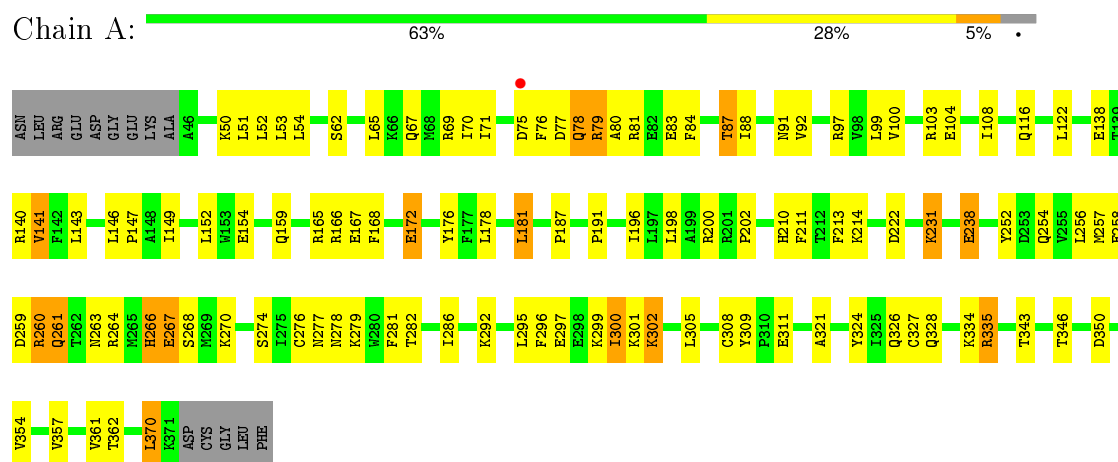
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	25	Total	O	0	0
			25	25		
6	C	20	Total	O	0	0
			20	20		
6	D	29	Total	O	0	0
			29	29		
6	F	10	Total	O	0	0
			10	10		

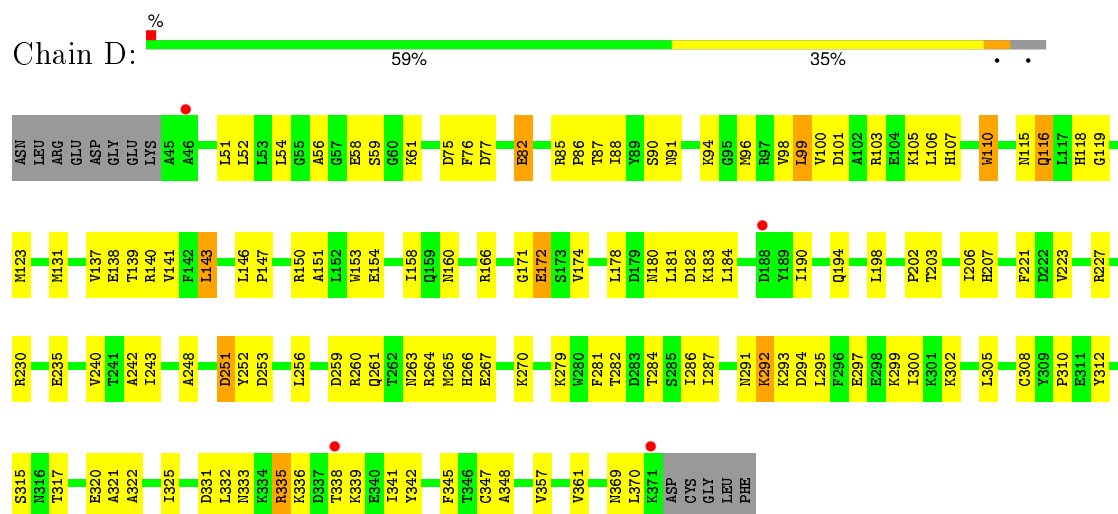
### 3 Residue-property plots

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: Guanine Nucleotide-Binding Protein Galpha(13):Galpha(i1) Chimera

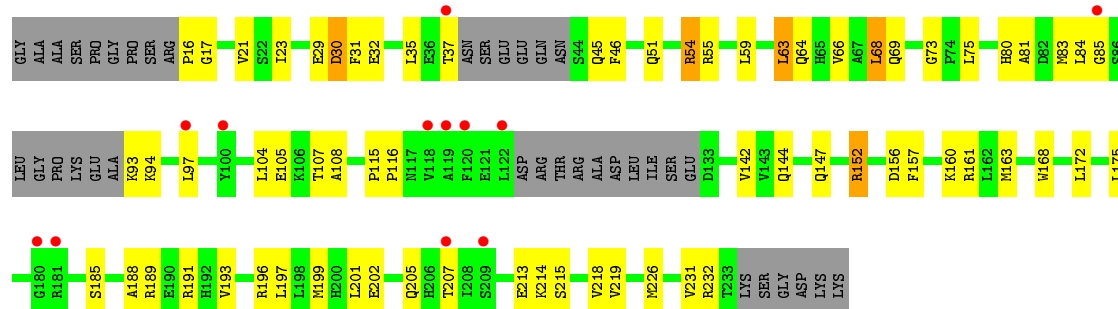


- Molecule 1: Guanine Nucleotide-Binding Protein Galpha(13):Galpha(i1) Chimera

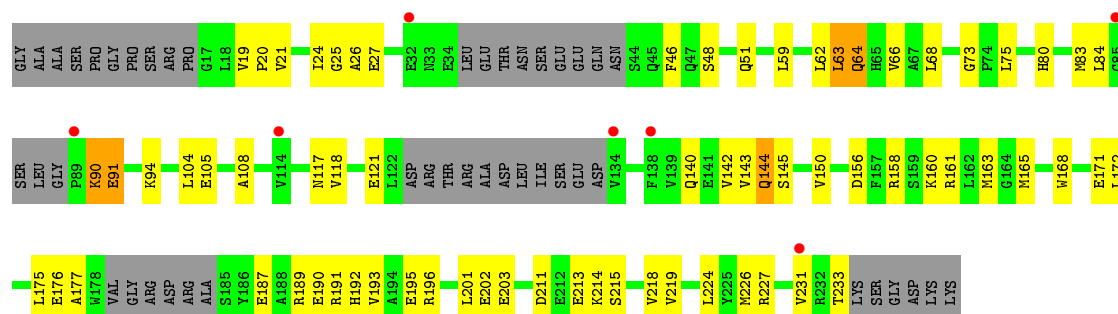


- Molecule 2: Rho guanine nucleotide exchange factor 1





- Molecule 2: Rho guanine nucleotide exchange factor 1





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	199.75Å 105.27Å 71.75Å 90.00° 96.91° 90.00°	Depositor
Resolution (Å)	46.49 – 2.85 46.49 – 2.75	Depositor EDS
% Data completeness (in resolution range)	90.7 (46.49-2.85) 86.3 (46.49-2.75)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.19	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.14 (at 2.77Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.229 , 0.297 0.225 , 0.290	Depositor DCC
$R_{free}$ test set	2203 reflections (7.56%)	DCC
Wilson B-factor (Å <sup>2</sup> )	58.2	Xtriage
Anisotropy	0.099	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 54.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 33204 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	8602	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.02% of the height of the origin peak. No significant pseudotranslation is detected.*

<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: GDP, ALF, 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	0.43	0/2734	0.60	0/3680
1	D	0.44	0/2739	0.58	0/3687
2	C	0.36	0/1605	0.51	0/2164
2	F	0.36	0/1543	0.51	0/2078
All	All	0.41	0/8621	0.56	0/11609

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	2677	0	2650	79	0
1	D	2682	0	2655	90	0
2	C	1576	0	1561	55	0
2	F	1515	0	1504	54	0
3	A	1	0	0	0	0
3	D	1	0	0	0	0
4	A	5	0	0	0	0
4	D	5	0	0	0	0
5	A	28	0	12	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	D	28	0	12	2	0
6	A	25	0	0	0	0
6	C	20	0	0	3	0
6	D	29	0	0	2	0
6	F	10	0	0	1	0
All	All	8602	0	8394	259	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 (259) 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:F:105:GLU:HB3	2:F:108:ALA:HB2	1.35	1.07
1:D:286:ILE:HD12	1:D:339:LYS:HD3	1.50	0.92
1:D:138:GLU:HG2	1:D:141:VAL:HG12	1.50	0.92
1:D:256:LEU:HD21	1:D:267:GLU:HG3	1.58	0.84
2:C:191:ARG:HH12	2:C:232:ARG:HH22	1.22	0.83
2:C:105:GLU:HB3	2:C:108:ALA:HB2	1.58	0.83
1:D:230:ARG:HD3	2:F:165:MET:HE1	1.62	0.81
1:A:91:ASN:HD21	1:A:200:ARG:H	1.24	0.81
2:C:16:PRO:HB2	2:F:24:ILE:HD11	1.62	0.79
2:C:29:GLU:HA	2:C:32:GLU:HG3	1.66	0.77
1:D:206:ILE:HG13	1:D:223:VAL:HG12	1.69	0.73
2:F:48:SER:OG	2:F:51:GLN:HG2	1.88	0.73
2:C:59:LEU:HG	2:C:63:LEU:HD23	1.70	0.72
2:F:161:ARG:HH11	2:F:161:ARG:HG2	1.57	0.67
2:C:189:ARG:O	2:C:193:VAL:HG23	1.93	0.67
1:A:50:LYS:HZ3	1:A:238:GLU:HG2	1.59	0.67
1:A:300:ILE:HD12	1:A:301:LYS:H	1.61	0.66
1:A:78:GLN:HG2	1:A:79:ARG:NH1	2.11	0.66
2:C:94:LYS:O	2:C:97:LEU:HB3	1.95	0.65
1:D:265:MET:HE3	1:D:325:ILE:HD13	1.78	0.65
1:D:279:LYS:HA	1:D:282:THR:HG23	1.80	0.64
2:C:191:ARG:HH12	2:C:232:ARG:NH2	1.93	0.64
1:A:149:ILE:HG22	1:A:181:LEU:HD21	1.81	0.63
1:D:137:VAL:HG13	1:D:141:VAL:HG13	1.81	0.63
2:C:45:GLN:HA	2:C:51:GLN:HE22	1.63	0.63
1:D:82:GLU:OE1	1:D:190:ILE:HG12	1.99	0.63
1:D:259:ASP:O	1:D:261:GLN:N	2.30	0.62
1:A:103:ARG:HB3	1:A:103:ARG:NH1	2.14	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:230:ARG:HD3	2:F:165:MET:CE	2.30	0.61
1:A:279:LYS:HG3	2:C:69:GLN:HA	1.82	0.61
1:D:369:ASN:ND2	1:D:370:LEU:HG	2.15	0.61
1:D:98:VAL:HG22	2:F:25:GLY:HA2	1.83	0.61
1:A:103:ARG:HB3	1:A:103:ARG:HH11	1.66	0.60
2:F:226:MET:HE3	2:F:231:VAL:HG21	1.83	0.60
2:C:152:ARG:HG2	1:D:116:GLN:HG3	1.83	0.60
1:A:78:GLN:HG2	1:A:79:ARG:HH12	1.66	0.60
1:A:346:THR:HG22	1:A:357:VAL:HG21	1.83	0.59
2:C:35:LEU:HD11	6:C:580:HOH:O	2.01	0.59
1:D:178:LEU:HA	1:D:181:LEU:HD13	1.84	0.59
1:D:75:ASP:CG	1:D:76:PHE:H	2.07	0.59
2:F:59:LEU:HG	2:F:63:LEU:CD2	2.32	0.58
1:D:243:ILE:HD11	1:D:281:PHE:CE1	2.38	0.58
1:D:227:ARG:HG2	1:D:230:ARG:NH1	2.19	0.58
1:A:50:LYS:NZ	1:A:238:GLU:HG2	2.19	0.58
1:A:67:GLN:HG3	1:A:354:VAL:HG21	1.86	0.58
1:A:146:LEU:HB3	1:A:147:PRO:HD3	1.86	0.58
1:A:326:GLN:HG3	1:A:343:THR:HG21	1.85	0.58
2:C:93:LYS:HB3	6:C:536:HOH:O	2.04	0.57
1:A:84:PHE:O	1:A:87:THR:HG23	2.05	0.57
2:F:19:VAL:HG13	2:F:20:PRO:HD2	1.87	0.57
2:F:171:GLU:HG2	2:F:196:ARG:HH21	1.70	0.57
2:F:161:ARG:NH1	2:F:161:ARG:HG2	2.20	0.56
1:D:287:ILE:HG12	1:D:361:VAL:HG13	1.86	0.56
1:A:256:LEU:HD11	1:A:267:GLU:HG3	1.87	0.56
1:A:292:LYS:HD3	1:A:295:LEU:HD12	1.87	0.56
1:D:333:ASN:HD22	1:D:341:ILE:HD11	1.70	0.56
2:F:90:LYS:HG2	2:F:91:GLU:HG3	1.88	0.56
1:A:69:ARG:HD3	1:A:76:PHE:CZ	2.41	0.56
2:F:192:HIS:O	2:F:195:GLU:HG2	2.05	0.56
1:D:333:ASN:O	1:D:336:LYS:HG3	2.04	0.56
1:D:335:ARG:HB3	1:D:338:THR:HB	1.88	0.56
2:C:16:PRO:CB	2:F:24:ILE:HD11	2.35	0.55
1:A:80:ALA:HA	1:A:83:GLU:OE2	2.06	0.55
2:F:156:ASP:O	2:F:160:LYS:HG2	2.07	0.55
1:A:70:ILE:HG23	1:A:75:ASP:OD2	2.06	0.55
1:D:150:ARG:O	1:D:154:GLU:HG2	2.06	0.55
2:F:201:LEU:O	2:F:201:LEU:HD23	2.06	0.55
1:A:274:SER:HB2	2:C:160:LYS:HZ2	1.72	0.54
1:D:256:LEU:HD11	1:D:267:GLU:CG	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:300:ILE:HD12	1:A:301:LYS:N	2.22	0.54
1:A:350:ASP:O	1:A:354:VAL:HG23	2.08	0.54
2:F:104:LEU:HD21	2:F:142:VAL:HG21	1.90	0.54
1:A:311:GLU:H	1:A:311:GLU:CD	2.11	0.54
2:F:175:LEU:HD13	2:F:189:ARG:HD3	1.89	0.54
1:D:52:LEU:HD23	1:D:221:PHE:HB2	1.88	0.54
2:F:175:LEU:HD22	2:F:193:VAL:HG21	1.90	0.54
2:C:214:LYS:O	2:C:218:VAL:HG23	2.07	0.54
1:D:252:TYR:HB3	1:D:265:MET:HE2	1.90	0.54
1:A:172:GLU:CD	1:A:172:GLU:H	2.11	0.54
1:A:252:TYR:O	1:A:308:CYS:HB2	2.09	0.53
1:D:253:ASP:OD2	1:D:299:LYS:HE2	2.09	0.53
2:C:55:ARG:HB3	6:C:504:HOH:O	2.08	0.53
2:F:117:ASN:O	2:F:121:GLU:HG3	2.08	0.53
1:D:287:ILE:HG23	1:D:342:TYR:O	2.09	0.52
2:F:64:GLN:O	2:F:68:LEU:HD23	2.09	0.52
2:F:80:HIS:O	2:F:83:MET:HB2	2.09	0.52
1:A:263:ASN:HB3	1:A:266:HIS:CG	2.45	0.52
1:D:203:THR:HG23	1:D:207:HIS:NE2	2.24	0.52
1:A:238:GLU:O	1:A:238:GLU:HG2	2.10	0.52
2:C:66:VAL:HG21	2:C:75:LEU:HD22	1.92	0.52
1:D:119:GLY:O	1:D:123:MET:HG2	2.09	0.52
1:A:202:PRO:HG3	2:C:30:ASP:OD2	2.09	0.52
1:D:139:THR:HG22	1:D:143:LEU:HD22	1.92	0.52
2:C:191:ARG:NH1	2:C:232:ARG:HH12	2.07	0.52
2:C:168:TRP:O	2:C:172:LEU:HB2	2.09	0.52
2:C:144:GLN:HA	2:C:147:GLN:HG2	1.91	0.51
1:A:178:LEU:HA	1:A:181:LEU:HD13	1.92	0.51
2:C:201:LEU:HD23	2:C:201:LEU:O	2.11	0.51
1:D:305:LEU:HD23	1:D:312:TYR:CE2	2.45	0.51
1:A:297:GLU:O	1:A:300:ILE:HD11	2.11	0.51
1:D:85:ARG:N	1:D:86:PRO:HD2	2.25	0.51
2:F:172:LEU:O	2:F:176:GLU:HG3	2.10	0.51
1:D:251:ASP:OD2	1:D:264:ARG:HD3	2.11	0.50
1:D:293:LYS:O	1:D:297:GLU:HG3	2.11	0.50
2:C:80:HIS:O	2:C:83:MET:HB3	2.12	0.50
2:C:105:GLU:HG2	2:C:107:THR:H	1.76	0.50
1:D:302:LYS:HB2	6:D:570:HOH:O	2.10	0.50
2:F:140:GLN:O	2:F:144:GLN:HB2	2.12	0.50
2:C:68:LEU:HD21	2:C:207:THR:HB	1.94	0.50
1:A:97:ARG:O	1:A:100:VAL:HG22	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:296:PHE:O	1:A:300:ILE:HG13	2.11	0.50
1:D:171:GLY:O	1:D:174:VAL:HG13	2.12	0.50
1:D:183:LYS:HG3	1:D:184:LEU:HD12	1.94	0.49
1:A:214:LYS:HD2	1:A:362:THR:HG21	1.94	0.49
1:A:67:GLN:O	1:A:71:ILE:HG12	2.13	0.49
1:A:266:HIS:O	1:A:270:LYS:HD3	2.13	0.49
1:D:99:LEU:HG	1:D:158:ILE:HG23	1.95	0.49
2:C:17:GLY:O	2:F:21:VAL:HG23	2.12	0.49
2:F:160:LYS:HE2	6:F:553:HOH:O	2.12	0.49
1:D:292:LYS:HG2	1:D:295:LEU:HD12	1.95	0.49
2:C:31:PHE:O	2:C:35:LEU:HG	2.13	0.49
2:F:171:GLU:CG	2:F:196:ARG:HH21	2.26	0.48
1:D:180:ASN:O	1:D:184:LEU:HD13	2.13	0.48
1:D:320:GLU:CD	1:D:320:GLU:H	2.16	0.48
2:C:196:ARG:HA	2:C:199:MET:HE3	1.96	0.48
1:A:122:LEU:HD12	1:A:152:LEU:CD1	2.43	0.48
2:F:175:LEU:HD21	2:F:190:GLU:HA	1.96	0.47
2:F:26:ALA:O	2:F:27:GLU:HB2	2.14	0.47
1:D:322:ALA:HB1	1:D:345:PHE:CE1	2.48	0.47
1:A:370:LEU:H	1:A:370:LEU:HD22	1.79	0.47
1:A:277:ASN:HD21	1:A:334:LYS:HG3	1.79	0.47
1:A:166:ARG:NH1	1:A:167:GLU:HG3	2.29	0.47
1:A:191:PRO:HB2	1:A:196:ILE:CD1	2.44	0.47
1:A:176:TYR:CD2	1:A:198:LEU:HD13	2.48	0.47
2:F:187:GLU:HB3	2:F:191:ARG:HE	1.80	0.47
1:D:82:GLU:HG3	1:D:85:ARG:NH2	2.29	0.47
1:D:270:LYS:HE2	2:F:163:MET:HE2	1.95	0.47
1:D:240:VAL:O	1:D:284:THR:HG23	2.14	0.47
1:D:52:LEU:CD2	1:D:221:PHE:HB2	2.44	0.47
1:D:248:ALA:O	1:D:251:ASP:HB2	2.15	0.47
1:D:317:THR:HB	1:D:320:GLU:OE2	2.15	0.47
1:D:146:LEU:HB3	1:D:147:PRO:HD3	1.96	0.47
2:F:46:PHE:O	2:F:158:ARG:HD3	2.15	0.47
1:A:335:ARG:HH11	1:A:335:ARG:HG2	1.80	0.47
1:A:305:LEU:HD22	1:A:321:ALA:HB1	1.95	0.47
2:F:59:LEU:O	2:F:63:LEU:HD22	2.16	0.46
1:A:168:PHE:HB3	2:C:23:ILE:HD12	1.97	0.46
2:C:104:LEU:HD21	2:C:142:VAL:HG11	1.97	0.46
1:D:110:TRP:NE1	1:D:115:ASN:HB2	2.31	0.46
1:A:259:ASP:O	1:A:261:GLN:N	2.41	0.46
2:F:215:SER:O	2:F:219:VAL:HG23	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:215:SER:O	2:C:219:VAL:HG23	2.16	0.46
1:D:248:ALA:HB1	1:D:292:LYS:HD3	1.96	0.46
2:F:211:ASP:HB3	2:F:214:LYS:HB3	1.97	0.46
2:C:68:LEU:HD11	2:C:207:THR:CG2	2.45	0.46
1:A:259:ASP:HA	2:F:19:VAL:CG1	2.45	0.46
2:F:117:ASN:OD1	2:F:118:VAL:HG23	2.16	0.46
1:D:265:MET:HB3	1:D:308:CYS:SG	2.56	0.46
1:A:103:ARG:HG2	1:A:108:ILE:HB	1.98	0.46
1:A:259:ASP:HA	2:F:19:VAL:HG11	1.98	0.46
2:C:46:PHE:CZ	2:C:157:PHE:HE2	2.34	0.46
2:C:185:SER:O	2:C:188:ALA:HB3	2.16	0.45
1:D:100:VAL:CG2	1:D:101:ASP:N	2.79	0.45
2:F:211:ASP:C	2:F:213:GLU:H	2.19	0.45
2:C:156:ASP:O	2:C:160:LYS:HG2	2.16	0.45
2:F:62:LEU:O	2:F:66:VAL:HG23	2.17	0.45
2:C:161:ARG:HH11	2:C:161:ARG:HG2	1.81	0.45
1:D:153:TRP:HA	1:D:153:TRP:CE3	2.52	0.45
1:D:357:VAL:O	1:D:361:VAL:HG23	2.17	0.45
1:A:191:PRO:HB2	1:A:196:ILE:HD11	1.98	0.45
1:A:138:GLU:HB3	1:A:141:VAL:HG12	1.98	0.45
1:A:295:LEU:O	1:A:299:LYS:HB2	2.17	0.45
2:C:226:MET:HB3	2:C:231:VAL:HB	1.99	0.45
2:C:191:ARG:NH1	2:C:232:ARG:HH22	2.03	0.44
2:C:115:PRO:HA	2:C:116:PRO:HD3	1.92	0.44
2:C:152:ARG:HD3	1:D:116:GLN:HG3	1.98	0.44
1:D:369:ASN:HD21	1:D:370:LEU:HG	1.79	0.44
1:A:324:TYR:O	1:A:328:GLN:HG2	2.17	0.44
1:D:292:LYS:HG3	5:D:476:GDP:C6	2.52	0.44
1:A:88:ILE:O	1:A:92:VAL:HG23	2.17	0.44
1:D:90:SER:OG	1:D:94:LYS:HE2	2.17	0.44
1:D:160:ASN:N	1:D:160:ASN:HD22	2.15	0.44
1:D:100:VAL:HG23	1:D:101:ASP:N	2.32	0.44
2:C:54:ARG:HE	2:C:54:ARG:HB2	1.70	0.44
2:C:64:GLN:NE2	2:C:168:TRP:NE1	2.66	0.44
1:A:213:PHE:CE1	1:A:214:LYS:HG3	2.53	0.44
1:D:54:LEU:C	1:D:61:LYS:HD3	2.38	0.43
1:D:331:ASP:C	1:D:333:ASN:H	2.22	0.43
1:D:315:SER:OG	1:D:320:GLU:HG3	2.19	0.43
2:F:214:LYS:O	2:F:218:VAL:HG23	2.19	0.43
1:A:53:LEU:O	1:A:54:LEU:HD23	2.17	0.43
1:D:235:GLU:N	1:D:235:GLU:OE2	2.47	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:300:ILE:H	1:A:300:ILE:HG13	1.52	0.43
1:A:279:LYS:HG3	2:C:68:LEU:O	2.19	0.43
1:D:305:LEU:HD22	1:D:321:ALA:HB1	1.99	0.43
1:A:165:ARG:HA	1:A:168:PHE:CE2	2.52	0.43
2:C:29:GLU:HA	2:C:32:GLU:CG	2.43	0.43
1:A:65:LEU:HD11	1:A:222:ASP:HB2	2.01	0.43
1:D:56:ALA:O	1:D:59:SER:HB2	2.19	0.43
2:F:59:LEU:HG	2:F:63:LEU:HD22	2.01	0.43
2:C:226:MET:HE2	2:C:226:MET:HA	2.01	0.43
1:A:260:ARG:HG3	2:C:21:VAL:HG13	2.01	0.43
1:D:106:LEU:O	1:D:107:HIS:HB2	2.17	0.43
1:A:79:ARG:H	1:A:79:ARG:HG2	1.64	0.43
1:D:279:LYS:HE3	6:D:519:HOH:O	2.18	0.43
2:F:224:LEU:HA	2:F:227:ARG:HH12	1.83	0.43
1:A:302:LYS:HB2	1:A:302:LYS:NZ	2.33	0.43
1:A:140:ARG:HH11	1:A:140:ARG:HG3	1.82	0.43
2:F:143:VAL:C	2:F:145:SER:H	2.22	0.43
2:F:168:TRP:O	2:F:172:LEU:HB2	2.19	0.42
1:D:293:LYS:HG3	1:D:294:ASP:N	2.34	0.42
1:A:292:LYS:HA	5:A:475:GDP:O6	2.19	0.42
1:D:88:ILE:O	1:D:91:ASN:HB2	2.19	0.42
1:A:165:ARG:C	1:A:167:GLU:H	2.23	0.42
1:A:231:LYS:HB3	1:A:231:LYS:NZ	2.34	0.42
1:D:308:CYS:O	1:D:310:PRO:HD3	2.20	0.42
1:D:115:ASN:OD1	1:D:118:HIS:HD2	2.02	0.42
1:D:96:MET:O	1:D:100:VAL:HG13	2.19	0.42
2:C:168:TRP:CD1	2:C:197:LEU:HD22	2.54	0.42
1:D:291:ASN:ND2	5:D:476:GDP:N7	2.66	0.42
1:D:317:THR:HB	1:D:320:GLU:CD	2.39	0.42
1:D:100:VAL:HG23	1:D:123:MET:HE2	2.02	0.42
1:D:317:THR:HB	1:D:320:GLU:CG	2.49	0.42
2:F:84:LEU:N	2:F:84:LEU:HD12	2.35	0.42
2:C:175:LEU:HD12	2:C:189:ARG:NH2	2.35	0.42
1:A:270:LYS:HB3	2:C:163:MET:SD	2.60	0.42
2:F:224:LEU:HA	2:F:227:ARG:NH1	2.35	0.42
1:D:51:LEU:HD12	1:D:242:ALA:HB3	2.01	0.42
1:A:278:ASN:HB3	1:A:281:PHE:CD2	2.55	0.42
1:A:357:VAL:O	1:A:361:VAL:HG23	2.19	0.42
1:D:151:ALA:O	1:D:154:GLU:HB2	2.20	0.42
1:A:309:TYR:HB3	1:A:311:GLU:OE2	2.20	0.42
1:D:158:ILE:N	1:D:158:ILE:HD12	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:292:LYS:HG2	5:A:475:GDP:C6	2.55	0.41
1:D:202:PRO:HB3	2:F:27:GLU:OE2	2.20	0.41
2:F:202:GLU:OE1	2:F:233:THR:HG21	2.20	0.41
1:D:137:VAL:HG13	1:D:141:VAL:CG1	2.49	0.41
2:C:160:LYS:HA	2:C:160:LYS:HD3	1.89	0.41
1:A:276:CYS:HA	1:A:286:ILE:HD13	2.02	0.41
2:C:81:ALA:O	2:C:84:LEU:HD13	2.20	0.41
2:C:202:GLU:O	2:C:205:GLN:HG3	2.20	0.41
1:A:259:ASP:C	1:A:261:GLN:H	2.22	0.41
1:D:281:PHE:HB3	1:D:284:THR:HB	2.01	0.41
1:A:263:ASN:HB3	1:A:266:HIS:HB2	2.02	0.41
2:F:187:GLU:CB	2:F:191:ARG:HH21	2.33	0.41
1:D:293:LYS:HG2	1:D:347:CYS:HB2	2.02	0.41
1:A:51:LEU:HD23	1:A:52:LEU:N	2.36	0.41
1:D:263:ASN:HB3	1:D:266:HIS:CD2	2.55	0.41
1:A:81:ARG:HB3	1:A:196:ILE:HD11	2.03	0.40
1:A:254:GLN:HB2	1:A:264:ARG:HD2	2.04	0.40
2:F:201:LEU:C	2:F:201:LEU:HD23	2.42	0.40
2:C:168:TRP:HA	2:C:168:TRP:HE3	1.86	0.40
1:A:210:HIS:O	1:A:211:PHE:HB3	2.21	0.40
1:D:287:ILE:HG13	1:D:342:TYR:HB2	2.03	0.40
2:C:161:ARG:HG2	2:C:161:ARG:NH1	2.36	0.40
2:F:75:LEU:HD13	2:F:150:VAL:HG13	2.03	0.40
1:D:281:PHE:O	1:D:339:LYS:HE3	2.22	0.40
2:F:226:MET:HE3	2:F:231:VAL:HG11	2.03	0.40
1:D:100:VAL:CG2	1:D:123:MET:SD	3.10	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	324/340 (95%)	300 (93%)	22 (7%)	2 (1%)	30	63
1	D	325/340 (96%)	293 (90%)	28 (9%)	4 (1%)	16	44
2	C	188/233 (81%)	163 (87%)	23 (12%)	2 (1%)	17	47
2	F	178/233 (76%)	155 (87%)	20 (11%)	3 (2%)	11	35
All	All	1015/1146 (89%)	911 (90%)	93 (9%)	11 (1%)	17	47

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	260	ARG
1	A	187	PRO
1	D	348	ALA
2	C	73	GLY
2	C	85	GLY
1	D	172	GLU
2	F	177	ALA
1	D	332	LEU
2	F	90	LYS
1	A	260	ARG
2	F	73	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	290/301 (96%)	262 (90%)	28 (10%)	10	27
1	D	290/301 (96%)	269 (93%)	21 (7%)	18	43
2	C	170/199 (85%)	163 (96%)	7 (4%)	37	71
2	F	163/199 (82%)	157 (96%)	6 (4%)	41	74
All	All	913/1000 (91%)	851 (93%)	62 (7%)	20	46

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

Mol	Chain	Res	Type
1	A	62	SER
1	A	77	ASP
1	A	78	GLN
1	A	79	ARG
1	A	87	THR
1	A	99	LEU
1	A	104	GLU
1	A	116	GLN
1	A	141	VAL
1	A	143	LEU
1	A	154	GLU
1	A	159	GLN
1	A	172	GLU
1	A	181	LEU
1	A	231	LYS
1	A	238	GLU
1	A	257	MET
1	A	258	GLU
1	A	261	GLN
1	A	266	HIS
1	A	267	GLU
1	A	268	SER
1	A	282	THR
1	A	300	ILE
1	A	302	LYS
1	A	327	CYS
1	A	335	ARG
1	A	370	LEU
2	C	30	ASP
2	C	37	THR
2	C	54	ARG
2	C	63	LEU
2	C	68	LEU
2	C	152	ARG
2	C	213	GLU
1	D	58	GLU
1	D	77	ASP
1	D	82	GLU
1	D	87	THR
1	D	99	LEU
1	D	103	ARG
1	D	105	LYS
1	D	110	TRP

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Mol	Chain	Res	Type
1	D	116	GLN
1	D	131	MET
1	D	140	ARG
1	D	143	LEU
1	D	166	ARG
1	D	172	GLU
1	D	182	ASP
1	D	194	GLN
1	D	198	LEU
1	D	251	ASP
1	D	292	LYS
1	D	300	ILE
1	D	335	ARG
2	F	63	LEU
2	F	64	GLN
2	F	91	GLU
2	F	94	LYS
2	F	144	GLN
2	F	203	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	HIS
1	A	78	GLN
1	A	91	ASN
1	A	107	HIS
1	A	116	GLN
1	A	160	ASN
1	A	169	GLN
1	A	210	HIS
1	A	261	GLN
1	A	266	HIS
1	A	277	ASN
1	A	355	GLN
2	C	45	GLN
2	C	51	GLN
2	C	58	HIS
2	C	153	GLN
2	C	192	HIS
2	C	205	GLN
1	D	72	HIS

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Mol	Chain	Res	Type
1	D	115	ASN
1	D	118	HIS
1	D	160	ASN
1	D	169	GLN
1	D	193	GLN
1	D	194	GLN
1	D	266	HIS
1	D	326	GLN
1	D	368	ASN
1	D	369	ASN
2	F	64	GLN
2	F	153	GLN
2	F	220	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 6 ligands modelled in this entry, 2 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	ALF	A	378	3,5,6	0,4,4	0.00	-	0,6,6	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	GDP	A	475	3,4	23,30,30	1.90	6 (26%)	30,47,47	2.82	12 (40%)
4	ALF	D	378	3,5,6	0,4,4	0.00	-	0,6,6	0.00	-
5	GDP	D	476	3,4	23,30,30	1.89	6 (26%)	30,47,47	2.82	12 (40%)

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	ALF	A	378	3,5,6	-	0/0/0/0	0/0/0/0
5	GDP	A	475	3,4	-	0/12/32/32	0/3/3/3
4	ALF	D	378	3,5,6	-	0/0/0/0	0/0/0/0
5	GDP	D	476	3,4	-	0/12/32/32	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	475	GDP	C8-N7	-2.87	1.29	1.34
5	D	476	GDP	C8-N7	-2.84	1.29	1.34
5	A	475	GDP	O5'-C5'	-2.15	1.36	1.44
5	D	476	GDP	O5'-C5'	-2.13	1.36	1.44
5	A	475	GDP	C3'-C4'	-2.04	1.47	1.53
5	D	476	GDP	C3'-C4'	-2.04	1.47	1.53
5	D	476	GDP	C2-N1	2.72	1.40	1.35
5	A	475	GDP	C2-N1	2.76	1.40	1.35
5	D	476	GDP	O4'-C1'	3.59	1.45	1.41
5	A	475	GDP	O4'-C1'	3.61	1.45	1.41
5	A	475	GDP	C6-N1	4.63	1.41	1.33
5	D	476	GDP	C6-N1	4.63	1.41	1.33

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	476	GDP	C5-C6-N1	-9.29	110.88	123.59
5	A	475	GDP	C5-C6-N1	-9.29	110.89	123.59
5	D	476	GDP	PA-O3A-PB	-4.41	117.88	132.67
5	A	475	GDP	PA-O3A-PB	-4.40	117.92	132.67
5	A	475	GDP	N3-C2-N1	-2.67	123.37	127.44
5	D	476	GDP	N3-C2-N1	-2.65	123.40	127.44
5	D	476	GDP	O3'-C3'-C2'	-2.38	104.09	111.83
5	A	475	GDP	O3'-C3'-C2'	-2.37	104.11	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	476	GDP	C1'-N9-C4	-2.12	123.74	126.94
5	A	475	GDP	C1'-N9-C4	-2.11	123.75	126.94
5	A	475	GDP	N2-C2-N1	2.03	120.57	117.20
5	D	476	GDP	N2-C2-N1	2.04	120.58	117.20
5	A	475	GDP	C4-C5-N7	2.10	111.41	109.48
5	D	476	GDP	C4-C5-N7	2.10	111.41	109.48
5	A	475	GDP	O2B-PB-O1B	2.67	119.17	110.58
5	D	476	GDP	O2B-PB-O1B	2.67	119.19	110.58
5	A	475	GDP	O2A-PA-O3A	3.29	120.04	105.09
5	D	476	GDP	O2A-PA-O3A	3.30	120.06	105.09
5	D	476	GDP	C4'-O4'-C1'	3.69	113.77	109.72
5	A	475	GDP	C4'-O4'-C1'	3.69	113.77	109.72
5	A	475	GDP	C2'-C1'-N9	4.27	120.81	114.29
5	D	476	GDP	C2'-C1'-N9	4.27	120.82	114.29
5	A	475	GDP	C6-N1-C2	6.84	125.44	115.94
5	D	476	GDP	C6-N1-C2	6.86	125.46	115.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	475	GDP	2	0
5	D	476	GDP	2	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	326/340 (95%)	-0.20	1 (0%) 94 93	22, 48, 73, 98	0
1	D	327/340 (96%)	-0.12	4 (1%) 81 78	22, 44, 76, 97	0
2	C	196/233 (84%)	0.28	12 (6%) 25 18	33, 72, 102, 103	1 (0%)
2	F	188/233 (80%)	0.24	7 (3%) 45 38	36, 73, 99, 103	1 (0%)
All	All	1037/1146 (90%)	-0.00	24 (2%) 64 59	22, 53, 96, 103	2 (0%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	120	PHE	5.2
2	C	118	VAL	4.4
2	C	37	THR	4.3
2	F	89	PRO	4.1
2	C	100	TYR	3.9
2	F	85	GLY	3.8
2	C	122	LEU	3.8
2	C	207	THR	3.6
2	C	209	SER	3.4
2	F	114	VAL	3.1
2	C	119	ALA	3.1
2	F	231	VAL	2.9
1	D	188	ASP	2.9
2	F	138	PHE	2.9
1	D	46	ALA	2.8
2	C	180	GLY	2.8
1	D	338	THR	2.6
2	C	181	ARG	2.6
2	F	134	VAL	2.3
1	A	75	ASP	2.2
2	C	97	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
2	F	32	GLU	2.2
1	D	371	LYS	2.1
2	C	85	GLY	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, 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	MG	A	377	1/1	0.97	0.28	9.83	8,8,8,8	0
3	MG	D	377	1/1	0.98	0.31	5.02	9,9,9,9	0
5	GDP	A	475	28/28	0.94	0.24	1.69	22,42,55,57	0
5	GDP	D	476	28/28	0.95	0.21	1.03	22,36,44,52	0
4	ALF	A	378	5/5	0.96	0.22	0.81	21,22,29,36	0
4	ALF	D	378	5/5	0.98	0.18	-0.17	20,32,41,43	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.