



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:39 PM GMT

PDB ID : 1Y3A
Title : Structure of G-Alpha-I1 bound to a GDP-selective peptide provides insight into guanine nucleotide exchange
Authors : Johnston, C.A.; Willard, F.S.; Jezyk, M.R.; Fredericks, Z.; Bodor, E.T.; Jones, M.B.; Blaesius, R.; Harden, T.K.; Sondek, J.; Watts, V.J.; Ramer, J.K.; Siderovski, D.P.
Deposited on : 2004-11-24
Resolution : 2.50 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 (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

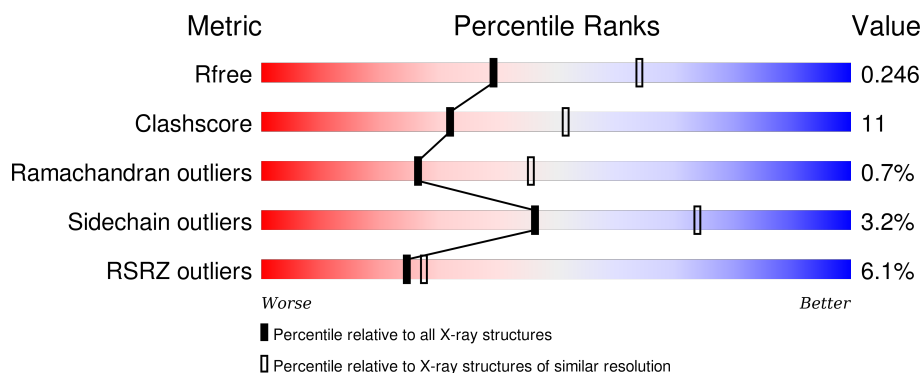
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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	329	<div> <div>5%</div> <div> <div>71%</div> <div>19%</div> <div>9%</div> </div> </div>
1	B	329	<div> <div>6%</div> <div> <div>73%</div> <div>19%</div> <div>7%</div> </div> </div>
1	C	329	<div> <div>4%</div> <div> <div>70%</div> <div>19%</div> <div>9%</div> </div> </div>
1	D	329	<div> <div>6%</div> <div> <div>70%</div> <div>20%</div> <div>8%</div> </div> </div>
2	E	16	<div> <div>13%</div> <div> <div>31%</div> <div>44%</div> <div>25%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
2	F	16	
2	G	16	
2	H	16	

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	GDP	B	356	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10424 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.

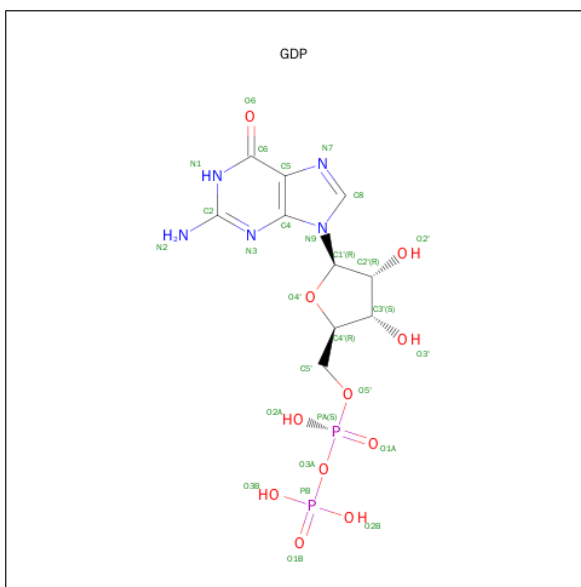
- Molecule 1 is a protein called Guanine nucleotide-binding protein G(i), alpha-1 subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	299	Total	C	N	O	S	0	0	0
			2413	1540	407	451	15			
1	B	307	Total	C	N	O	S	0	0	0
			2471	1575	414	467	15			
1	C	298	Total	C	N	O	S	0	0	0
			2404	1534	405	450	15			
1	D	304	Total	C	N	O	S	0	0	0
			2452	1562	411	464	15			

- Molecule 2 is a protein called KB752 peptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	12	Total	C	N	O	S	0	0	0
			109	71	16	21	1			
2	F	12	Total	C	N	O	S	0	0	0
			109	71	16	21	1			
2	G	12	Total	C	N	O	S	0	0	0
			109	71	16	21	1			
2	H	12	Total	C	N	O	S	0	0	0
			109	71	16	21	1			

- Molecule 3 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: C₁₀H₁₅N₅O₁₁P₂).



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

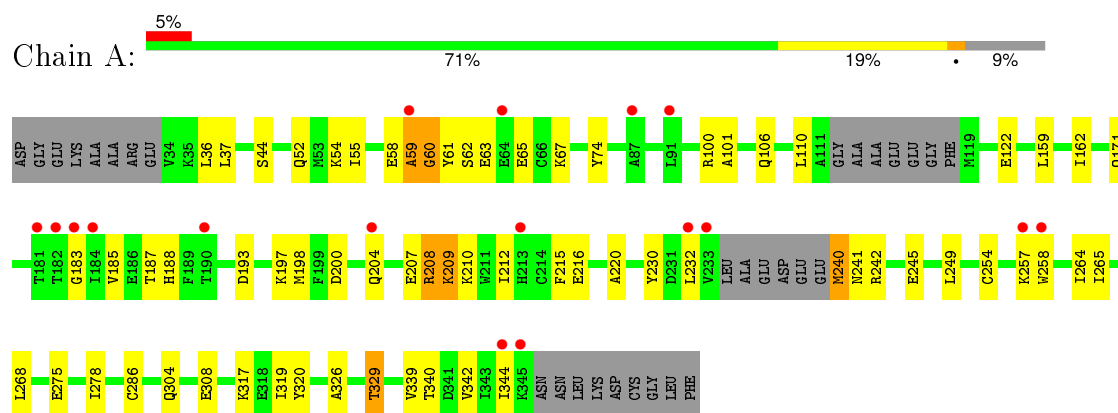
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	46	Total O 46 46	0	0
4	B	29	Total O 29 29	0	0
4	C	30	Total O 30 30	0	0
4	D	25	Total O 25 25	0	0
4	E	2	Total O 2 2	0	0
4	F	1	Total O 1 1	0	0
4	G	2	Total O 2 2	0	0
4	H	1	Total O 1 1	0	0

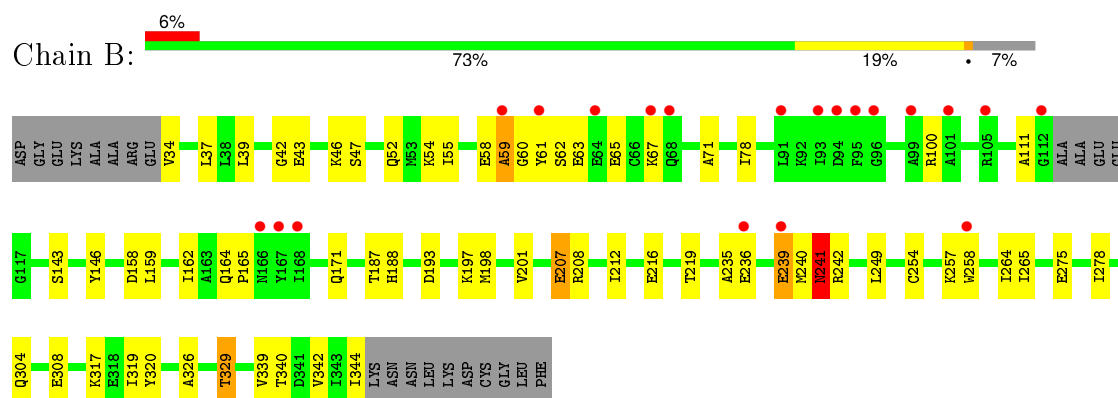
3 Residue-property plots [i](#)

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

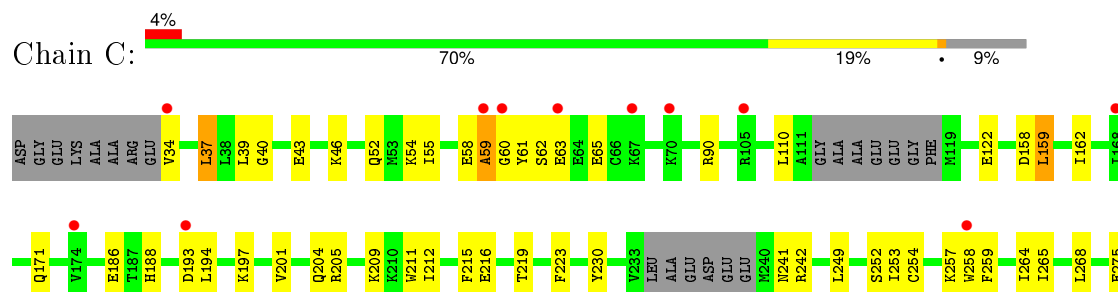
- Molecule 1: Guanine nucleotide-binding protein G(i), alpha-1 subunit



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

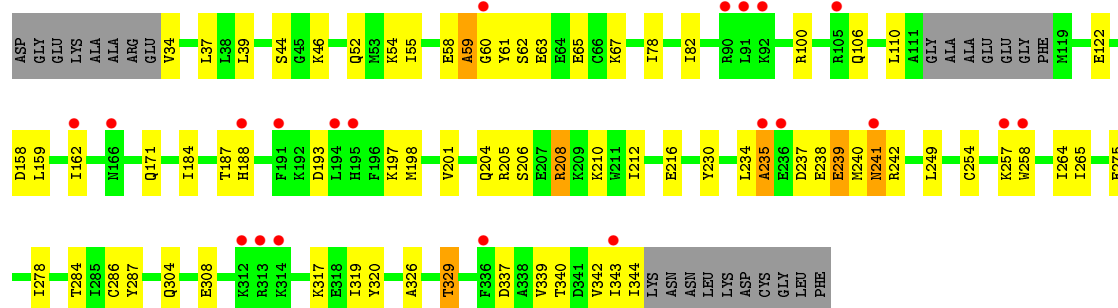


- Molecule 1: Guanine nucleotide-binding protein G(i), alpha-1 subunit

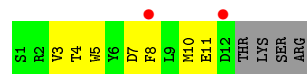




- Molecule 1: Guanine nucleotide-binding protein G(i), alpha-1 subunit



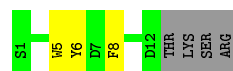
- Molecule 2: KB752 peptide



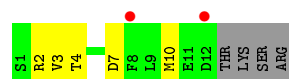
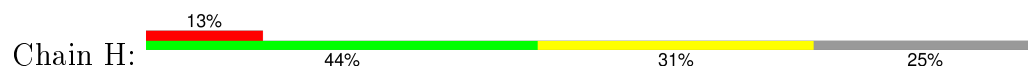
- Molecule 2: KB752 peptide



- Molecule 2: KB752 peptide



- Molecule 2: KB752 peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	72.94Å 112.78Å 109.49Å 90.00° 93.75° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50 37.59 – 2.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.50) 96.1 (37.59-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.49 (at 2.51Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.240 , 0.280 0.246 , 0.246	Depositor DCC
R_{free} test set	2987 reflections (5.34%)	DCC
Wilson B-factor (Å ²)	37.0	Xtriage
Anisotropy	0.107	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 42.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 59131 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	10424	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.09% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	1/2456 (0.0%)	0.59	0/3306
1	B	0.45	1/2516 (0.0%)	0.61	0/3389
1	C	0.43	0/2447	0.60	0/3295
1	D	0.43	1/2496 (0.0%)	0.60	1/3363 (0.0%)
2	E	0.35	0/112	0.51	0/151
2	F	0.41	0/112	0.51	0/151
2	G	0.37	0/112	0.48	0/151
2	H	0.38	0/112	0.50	0/151
All	All	0.43	3/10363 (0.0%)	0.60	1/13957 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	60	GLY	N-CA	5.80	1.54	1.46
1	B	241	ASN	N-CA	5.66	1.57	1.46
1	D	241	ASN	N-CA	5.52	1.57	1.46

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	241	ASN	N-CA-C	-7.67	90.29	111.00

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	2413	0	2411	52	0
1	B	2471	0	2452	43	0
1	C	2404	0	2398	54	1
1	D	2452	0	2437	58	1
2	E	109	0	98	8	0
2	F	109	0	98	5	0
2	G	109	0	98	4	0
2	H	109	0	98	4	0
3	A	28	0	12	2	0
3	B	28	0	12	2	0
3	C	28	0	12	3	0
3	D	28	0	12	3	0
4	A	46	0	0	8	0
4	B	29	0	0	2	0
4	C	30	0	0	3	0
4	D	25	0	0	1	0
4	E	2	0	0	0	0
4	F	1	0	0	0	0
4	G	2	0	0	1	0
4	H	1	0	0	0	0
All	All	10424	0	10138	216	1

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 11.

All (216) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:208:ARG:H	1:D:208:ARG:HD2	1.15	1.09
1:D:55:ILE:HA	1:D:60:GLY:HA2	1.39	1.03
1:C:55:ILE:HA	1:C:60:GLY:HA2	1.39	1.03
1:A:55:ILE:HA	1:A:60:GLY:HA2	1.38	1.03
1:B:55:ILE:HA	1:B:60:GLY:HA2	1.38	1.02
1:A:240:MET:HG2	1:A:241:ASN:H	1.29	0.94
1:A:240:MET:HG2	1:A:241:ASN:N	1.89	0.86
1:D:339:VAL:O	1:D:342:VAL:HG12	1.74	0.85
1:C:339:VAL:O	1:C:342:VAL:HG12	1.77	0.83
1:D:235:ALA:HB3	1:D:238:GLU:HG2	1.61	0.83
1:B:339:VAL:O	1:B:342:VAL:HG12	1.77	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:245:GLU:HB3	4:A:360:HOH:O	1.81	0.80
1:D:264:ILE:HG12	1:D:317:LYS:HE3	1.64	0.79
1:B:264:ILE:HG12	1:B:317:LYS:HE3	1.65	0.79
1:A:339:VAL:O	1:A:342:VAL:HG12	1.81	0.79
1:D:240:MET:HG2	1:D:241:ASN:H	1.51	0.74
1:A:264:ILE:HG12	1:A:317:LYS:HE3	1.70	0.74
1:B:208:ARG:HH12	2:F:3:VAL:HG21	1.53	0.74
1:A:230:TYR:O	1:A:286:CYS:HB2	1.88	0.73
1:C:60:GLY:HA3	4:C:387:HOH:O	1.89	0.72
1:C:264:ILE:HG12	1:C:317:LYS:HE3	1.72	0.72
1:C:253:ILE:HD11	2:G:5:TRP:CZ3	2.27	0.70
1:C:90:ARG:NH2	4:C:369:HOH:O	2.22	0.70
1:D:212:ILE:O	1:D:216:GLU:HG3	1.92	0.70
2:F:3:VAL:HG13	2:F:7:ASP:HB2	1.75	0.69
1:A:232:LEU:O	4:A:357:HOH:O	2.10	0.69
1:C:110:LEU:HD11	1:C:122:GLU:HG2	1.74	0.69
2:E:3:VAL:CG1	2:E:7:ASP:HB2	2.24	0.68
1:D:240:MET:HG2	1:D:241:ASN:N	2.09	0.68
1:A:183:GLY:HA2	1:A:204:GLN:CD	2.14	0.67
1:B:39:LEU:HD23	1:B:201:VAL:HB	1.77	0.67
1:D:337:ASP:HB2	4:D:371:HOH:O	1.94	0.66
1:A:52:GLN:OE1	1:A:329:THR:HB	1.96	0.66
1:C:280:LYS:HG2	4:C:365:HOH:O	1.94	0.66
1:C:212:ILE:O	1:C:216:GLU:HG3	1.94	0.66
1:A:101:ALA:HB3	4:A:369:HOH:O	1.96	0.66
1:D:240:MET:CG	1:D:241:ASN:H	2.10	0.65
1:D:234:LEU:HD23	1:D:234:LEU:H	1.63	0.63
1:D:208:ARG:HD2	1:D:208:ARG:N	2.00	0.63
1:C:340:THR:O	1:C:344:ILE:HG12	1.99	0.63
1:C:58:GLU:O	1:C:59:ALA:CB	2.47	0.63
1:D:264:ILE:CG1	1:D:317:LYS:HE3	2.29	0.62
1:D:58:GLU:O	1:D:59:ALA:CB	2.47	0.62
1:B:239:GLU:O	1:B:240:MET:HB3	1.99	0.62
1:C:264:ILE:CG1	1:C:317:LYS:HE3	2.30	0.62
1:A:183:GLY:HA2	1:A:204:GLN:NE2	2.14	0.62
1:B:58:GLU:O	1:B:59:ALA:CB	2.48	0.61
1:A:212:ILE:O	1:A:216:GLU:HG3	2.01	0.61
1:B:264:ILE:CG1	1:B:317:LYS:HE3	2.30	0.61
1:A:58:GLU:O	1:A:59:ALA:CB	2.48	0.61
1:D:208:ARG:O	1:D:212:ILE:HG12	2.01	0.61
1:B:239:GLU:HG3	1:B:239:GLU:O	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:GLN:OE1	4:A:361:HOH:O	2.16	0.61
1:A:264:ILE:CG1	1:A:317:LYS:HE3	2.30	0.61
1:B:241:ASN:HB2	4:B:360:HOH:O	2.00	0.61
1:B:52:GLN:OE1	1:B:329:THR:HB	2.00	0.60
1:C:58:GLU:O	1:C:59:ALA:HB3	2.02	0.60
1:D:254:CYS:SG	1:D:319:ILE:HD11	2.42	0.60
1:A:265:ILE:CD1	1:A:339:VAL:HG13	2.31	0.59
1:B:58:GLU:O	1:B:59:ALA:HB3	2.03	0.59
1:C:62:SER:OG	1:C:65:GLU:HG3	2.03	0.59
1:B:326:ALA:HB3	3:B:356:GDP:N7	2.18	0.58
1:A:183:GLY:HA2	1:A:204:GLN:OE1	2.03	0.58
1:D:58:GLU:O	1:D:59:ALA:HB3	2.02	0.58
1:A:58:GLU:O	1:A:59:ALA:HB3	2.03	0.58
1:D:62:SER:OG	1:D:65:GLU:HG3	2.03	0.58
1:B:207:GLU:HA	2:F:1:SER:O	2.04	0.58
1:D:52:GLN:OE1	1:D:329:THR:HB	2.03	0.58
1:A:304:GLN:O	1:A:308:GLU:HG3	2.04	0.58
1:D:340:THR:O	1:D:344:ILE:HG12	2.03	0.58
1:B:304:GLN:O	1:B:308:GLU:HG3	2.03	0.57
1:C:205:ARG:HD2	4:G:63:HOH:O	2.04	0.57
1:C:209:LYS:HB2	1:D:237:ASP:HB3	1.87	0.57
1:B:212:ILE:O	1:B:216:GLU:HG3	2.05	0.56
1:B:254:CYS:SG	1:B:319:ILE:HD11	2.46	0.56
1:D:39:LEU:HD23	1:D:201:VAL:HB	1.87	0.56
1:D:188:HIS:CE1	1:D:197:LYS:HD2	2.39	0.56
1:A:208:ARG:O	1:A:212:ILE:HG12	2.05	0.56
1:A:212:ILE:HD13	2:E:8:PHE:HZ	1.70	0.56
1:C:188:HIS:CE1	1:C:197:LYS:HD2	2.41	0.56
1:A:254:CYS:SG	1:A:319:ILE:HD11	2.45	0.56
1:B:54:LYS:HA	1:B:58:GLU:HB2	1.87	0.56
1:A:188:HIS:CE1	1:A:197:LYS:HD2	2.40	0.56
1:C:52:GLN:OE1	1:C:329:THR:HB	2.06	0.55
1:B:62:SER:OG	1:B:65:GLU:HG3	2.06	0.55
1:C:215:PHE:CD2	1:C:259:PHE:HZ	2.24	0.55
1:D:230:TYR:O	1:D:286:CYS:HB2	2.07	0.55
1:D:240:MET:CG	1:D:241:ASN:N	2.69	0.55
2:E:3:VAL:HG12	2:E:4:THR:N	2.22	0.55
1:A:265:ILE:HD13	1:A:339:VAL:HG13	1.89	0.55
2:H:3:VAL:HG12	2:H:4:THR:O	2.06	0.55
1:C:39:LEU:HD23	1:C:201:VAL:HB	1.89	0.55
1:D:265:ILE:CD1	1:D:339:VAL:HG13	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:7:ASP:O	2:E:11:GLU:HG3	2.08	0.54
1:A:54:LYS:HA	1:A:58:GLU:HB2	1.89	0.54
1:D:205:ARG:O	2:H:2:ARG:HD2	2.07	0.53
1:D:206:SER:HB3	1:D:210:LYS:HD2	1.89	0.53
1:C:254:CYS:SG	1:C:319:ILE:HD11	2.48	0.53
1:A:62:SER:OG	1:A:65:GLU:HG3	2.09	0.53
2:F:3:VAL:HG12	2:F:4:THR:O	2.09	0.53
1:C:54:LYS:HA	1:C:58:GLU:HB2	1.91	0.53
1:B:265:ILE:HD13	1:B:339:VAL:HG13	1.91	0.52
1:C:204:GLN:H	1:C:204:GLN:CD	2.13	0.52
1:C:110:LEU:CD1	1:C:122:GLU:HG2	2.39	0.52
1:B:340:THR:O	1:B:344:ILE:HG12	2.09	0.52
2:E:3:VAL:HG11	2:E:7:ASP:HB2	1.92	0.52
1:B:39:LEU:CD2	1:B:201:VAL:HB	2.40	0.51
2:E:3:VAL:HG12	2:E:7:ASP:HB2	1.91	0.51
1:B:188:HIS:CE1	1:B:197:LYS:HD2	2.46	0.51
1:D:158:ASP:O	1:D:162:ILE:HG12	2.11	0.51
1:D:54:LYS:HA	1:D:58:GLU:HB2	1.91	0.51
1:C:326:ALA:HB3	3:C:357:GDP:N7	2.26	0.51
1:C:34:VAL:HG21	1:C:194:LEU:HD13	1.93	0.51
1:A:36:LEU:HD11	1:A:220:ALA:HB3	1.94	0.50
1:B:265:ILE:CD1	1:B:339:VAL:HG13	2.41	0.50
1:C:39:LEU:CD2	1:C:201:VAL:HB	2.42	0.50
1:A:61:TYR:H	1:A:171:GLN:NE2	2.09	0.50
1:D:265:ILE:HD13	1:D:339:VAL:HG13	1.93	0.50
1:D:257:LYS:HG3	1:D:258:TRP:CD1	2.47	0.50
1:A:257:LYS:HG3	1:A:258:TRP:CD1	2.47	0.50
1:D:304:GLN:O	1:D:308:GLU:HG3	2.12	0.50
1:A:326:ALA:HB3	3:A:355:GDP:N7	2.27	0.49
1:C:55:ILE:HA	1:C:60:GLY:CA	2.28	0.49
1:D:234:LEU:O	1:D:235:ALA:O	2.29	0.49
1:D:239:GLU:O	1:D:240:MET:HB3	2.11	0.49
1:D:184:ILE:HG13	1:D:204:GLN:HE21	1.76	0.49
1:C:320:TYR:CE1	1:C:342:VAL:HG21	2.47	0.49
1:B:208:ARG:NH1	2:F:3:VAL:HG21	2.25	0.49
1:D:326:ALA:HB3	3:D:358:GDP:N7	2.28	0.49
1:A:187:THR:HG22	1:A:198:MET:HB3	1.94	0.48
1:B:257:LYS:HG3	1:B:258:TRP:CD1	2.49	0.48
1:C:34:VAL:HG13	1:C:219:THR:HG21	1.96	0.48
1:D:110:LEU:CD1	1:D:122:GLU:HG2	2.43	0.48
1:C:304:GLN:O	1:C:308:GLU:HG3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:340:THR:O	1:A:344:ILE:HG12	2.14	0.48
1:A:209:LYS:HE3	1:A:209:LYS:HA	1.95	0.48
1:B:78:ILE:HG21	1:B:111:ALA:HB1	1.96	0.48
1:C:275:GLU:O	1:C:278:ILE:HG22	2.13	0.48
1:C:257:LYS:HG3	1:C:258:TRP:CD1	2.48	0.48
1:D:235:ALA:HB3	1:D:238:GLU:CG	2.39	0.47
1:D:61:TYR:H	1:D:171:GLN:NE2	2.11	0.47
1:C:40:GLY:HA2	1:C:223:PHE:CE1	2.49	0.47
1:A:36:LEU:CD1	1:A:220:ALA:HB3	2.45	0.47
1:D:44:SER:N	3:D:358:GDP:O3B	2.48	0.47
1:D:238:GLU:O	1:D:238:GLU:HG3	2.15	0.47
1:A:210:LYS:HG3	4:A:396:HOH:O	2.14	0.47
1:D:275:GLU:O	1:D:278:ILE:HG22	2.15	0.47
1:D:46:LYS:HB2	3:D:358:GDP:O1B	2.15	0.46
1:C:46:LYS:HB2	3:C:357:GDP:O1B	2.15	0.46
1:D:55:ILE:HA	1:D:60:GLY:CA	2.28	0.46
1:C:265:ILE:HD13	1:C:339:VAL:HG13	1.98	0.46
1:B:320:TYR:CE1	1:B:342:VAL:HG21	2.51	0.46
1:D:39:LEU:CD2	1:D:201:VAL:HB	2.46	0.46
2:H:3:VAL:HG13	2:H:7:ASP:HB2	1.97	0.46
1:B:158:ASP:O	1:B:162:ILE:HG12	2.16	0.46
1:A:106:GLN:HB2	4:A:361:HOH:O	2.15	0.45
1:B:34:VAL:HG13	1:B:219:THR:HG21	1.98	0.45
1:B:187:THR:HG22	1:B:198:MET:HB3	1.97	0.45
1:B:63:GLU:O	1:B:67:LYS:HG3	2.16	0.45
1:A:100:ARG:HA	1:A:100:ARG:NE	2.31	0.45
1:C:230:TYR:O	1:C:286:CYS:HB2	2.16	0.45
1:C:265:ILE:CD1	1:C:339:VAL:HG13	2.46	0.45
1:C:215:PHE:CZ	2:G:5:TRP:HB2	2.52	0.45
1:A:44:SER:N	3:A:355:GDP:O3B	2.50	0.44
1:C:158:ASP:O	1:C:162:ILE:HG12	2.17	0.44
1:B:275:GLU:O	1:B:278:ILE:HG22	2.18	0.44
1:B:55:ILE:HA	1:B:60:GLY:CA	2.27	0.44
1:B:239:GLU:O	1:B:240:MET:CB	2.66	0.44
1:A:63:GLU:O	1:A:67:LYS:HG3	2.18	0.44
1:A:110:LEU:HD13	1:A:122:GLU:HG2	1.99	0.44
1:B:100:ARG:HA	1:B:100:ARG:NE	2.33	0.44
1:C:211:TRP:O	1:C:215:PHE:HD1	2.01	0.44
1:C:212:ILE:HD13	2:G:8:PHE:HZ	1.83	0.44
1:A:215:PHE:HZ	2:E:5:TRP:HB2	1.82	0.43
1:D:63:GLU:O	1:D:67:LYS:HG3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:78:ILE:O	1:D:82:ILE:HG13	2.18	0.43
1:C:43:GLU:HA	3:C:357:GDP:H5"	2.00	0.43
1:D:187:THR:HG22	1:D:198:MET:HB3	2.00	0.43
1:B:71:ALA:HB3	4:B:369:HOH:O	2.18	0.43
1:A:240:MET:N	4:A:375:HOH:O	2.52	0.43
1:A:106:GLN:HG3	1:A:110:LEU:CD1	2.49	0.43
1:A:185:VAL:HB	1:A:200:ASP:HB3	2.00	0.43
1:D:100:ARG:NE	1:D:100:ARG:HA	2.34	0.43
1:B:61:TYR:H	1:B:171:GLN:NE2	2.16	0.43
1:A:320:TYR:CE1	1:A:342:VAL:HG21	2.53	0.42
1:B:46:LYS:HB2	3:B:356:GDP:O1B	2.19	0.42
2:H:3:VAL:HG12	2:H:4:THR:N	2.34	0.42
1:C:61:TYR:H	1:C:171:GLN:NE2	2.16	0.42
1:A:275:GLU:O	1:A:278:ILE:HG22	2.20	0.42
1:A:55:ILE:HA	1:A:60:GLY:CA	2.28	0.42
1:C:268:LEU:HD23	1:C:268:LEU:HA	1.84	0.42
1:D:184:ILE:N	1:D:204:GLN:NE2	2.68	0.42
2:E:3:VAL:HG12	2:E:4:THR:H	1.83	0.42
1:D:110:LEU:HD13	1:D:122:GLU:HG2	2.02	0.42
1:D:284:THR:HA	1:D:287:TYR:O	2.20	0.42
1:D:234:LEU:CD2	1:D:234:LEU:H	2.31	0.41
1:D:34:VAL:HG11	1:D:343:ILE:HD13	2.02	0.41
1:C:252:SER:HB3	2:G:6:TYR:CE1	2.55	0.41
1:C:209:LYS:HD2	1:D:237:ASP:HA	2.02	0.41
1:A:207:GLU:HB2	1:A:210:LYS:HE2	2.02	0.41
1:B:143:SER:HA	1:B:146:TYR:CE1	2.56	0.41
1:D:106:GLN:HG3	1:D:110:LEU:HG	2.02	0.41
1:A:268:LEU:HD23	1:A:268:LEU:HA	1.86	0.41
1:C:215:PHE:CD2	1:C:259:PHE:CZ	3.07	0.41
1:C:194:LEU:HD11	1:C:344:ILE:HD13	2.03	0.41
1:D:320:TYR:CE1	1:D:342:VAL:HG21	2.56	0.41
1:A:216:GLU:HB2	4:A:384:HOH:O	2.20	0.41
1:C:159:LEU:HD23	1:C:159:LEU:HA	1.87	0.41
1:B:42:GLY:O	1:B:43:GLU:HB2	2.21	0.41
1:C:186:GLU:OE2	1:C:188:HIS:CE1	2.73	0.41
1:C:278:ILE:HD12	1:C:278:ILE:HA	1.96	0.40
1:B:143:SER:HA	1:B:146:TYR:CZ	2.57	0.40
1:C:37:LEU:HD12	1:C:37:LEU:HA	1.90	0.40
1:A:74:TYR:HE1	1:A:162:ILE:HG22	1.86	0.40
1:C:284:THR:HA	1:C:287:TYR:O	2.21	0.40
1:B:164:GLN:HA	1:B:165:PRO:HD3	1.93	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:63:GLU:OE1	1:D:257:LYS:NZ[2_656]	2.18	0.02

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	293/329 (89%)	291 (99%)	1 (0%)	1 (0%)	46	68
1	B	303/329 (92%)	297 (98%)	2 (1%)	4 (1%)	15	26
1	C	292/329 (89%)	288 (99%)	2 (1%)	2 (1%)	26	46
1	D	300/329 (91%)	292 (97%)	6 (2%)	2 (1%)	26	46
2	E	10/16 (62%)	10 (100%)	0	0	100	100
2	F	10/16 (62%)	10 (100%)	0	0	100	100
2	G	10/16 (62%)	10 (100%)	0	0	100	100
2	H	10/16 (62%)	9 (90%)	1 (10%)	0	100	100
All	All	1228/1380 (89%)	1207 (98%)	12 (1%)	9 (1%)	26	46

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	241	ASN
1	C	241	ASN
1	D	235	ALA
1	A	59	ALA
1	B	59	ALA
1	C	59	ALA
1	D	59	ALA
1	B	236	GLU
1	B	235	ALA

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	263/284 (93%)	254 (97%)	9 (3%)	44	72
1	B	268/284 (94%)	258 (96%)	10 (4%)	41	68
1	C	262/284 (92%)	256 (98%)	6 (2%)	58	83
1	D	267/284 (94%)	259 (97%)	8 (3%)	48	76
2	E	12/16 (75%)	11 (92%)	1 (8%)	14	26
2	F	12/16 (75%)	12 (100%)	0	100	100
2	G	12/16 (75%)	12 (100%)	0	100	100
2	H	12/16 (75%)	11 (92%)	1 (8%)	14	26
All	All	1108/1200 (92%)	1073 (97%)	35 (3%)	46	74

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	LEU
1	A	159	LEU
1	A	193	ASP
1	A	208	ARG
1	A	209	LYS
1	A	240	MET
1	A	242	ARG
1	A	249	LEU
1	A	329	THR
1	B	37	LEU
1	B	47	SER
1	B	159	LEU
1	B	193	ASP
1	B	207	GLU
1	B	239	GLU
1	B	241	ASN
1	B	242	ARG
1	B	249	LEU
1	B	329	THR

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Mol	Chain	Res	Type
1	C	37	LEU
1	C	159	LEU
1	C	193	ASP
1	C	242	ARG
1	C	249	LEU
1	C	329	THR
1	D	37	LEU
1	D	159	LEU
1	D	193	ASP
1	D	208	ARG
1	D	239	GLU
1	D	242	ARG
1	D	249	LEU
1	D	329	THR
2	E	10	MET
2	H	10	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	GLN
1	A	106	GLN
1	A	147	GLN
1	A	171	GLN
1	A	241	ASN
1	A	333	GLN
1	B	79	GLN
1	B	147	GLN
1	B	171	GLN
1	B	241	ASN
1	B	333	GLN
1	C	79	GLN
1	C	147	GLN
1	C	171	GLN
1	C	204	GLN
1	C	241	ASN
1	C	333	GLN
1	D	79	GLN
1	D	147	GLN
1	D	171	GLN
1	D	213	HIS
1	D	241	ASN

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Mol	Chain	Res	Type
1	D	333	GLN

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](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	GDP	A	355	-	23,30,30	1.83	7 (30%)	30,47,47	3.34	17 (56%)
3	GDP	B	356	-	23,30,30	1.91	8 (34%)	30,47,47	3.32	16 (53%)
3	GDP	C	357	-	23,30,30	1.68	5 (21%)	30,47,47	3.30	15 (50%)
3	GDP	D	358	-	23,30,30	1.72	6 (26%)	30,47,47	3.33	15 (50%)

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	GDP	A	355	-	-	0/12/32/32	0/3/3/3
3	GDP	B	356	-	-	0/12/32/32	0/3/3/3
3	GDP	C	357	-	-	0/12/32/32	0/3/3/3
3	GDP	D	358	-	-	0/12/32/32	0/3/3/3

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	358	GDP	C8-N7	-2.42	1.30	1.34
3	A	355	GDP	C8-N7	-2.42	1.30	1.34
3	B	356	GDP	C8-N7	-2.35	1.30	1.34
3	C	357	GDP	C8-N7	-2.22	1.30	1.34
3	B	356	GDP	PA-O2A	-2.14	1.45	1.54
3	B	356	GDP	PB-O1B	-2.12	1.44	1.51
3	A	355	GDP	PB-O1B	-2.05	1.44	1.51
3	D	358	GDP	C2-N1	2.03	1.39	1.35
3	D	358	GDP	C3'-C4'	2.14	1.58	1.53
3	C	357	GDP	C3'-C4'	2.24	1.59	1.53
3	D	358	GDP	O4'-C1'	2.27	1.44	1.41
3	B	356	GDP	C3'-C4'	2.28	1.59	1.53
3	C	357	GDP	C2-N1	2.29	1.39	1.35
3	C	357	GDP	C5'-C4'	2.32	1.59	1.51
3	A	355	GDP	C3'-C4'	2.34	1.59	1.53
3	B	356	GDP	C2-N1	2.50	1.39	1.35
3	A	355	GDP	C5'-C4'	2.51	1.59	1.51
3	B	356	GDP	C5'-C4'	2.52	1.59	1.51
3	B	356	GDP	O4'-C1'	2.64	1.44	1.41
3	D	358	GDP	C5'-C4'	2.66	1.60	1.51
3	A	355	GDP	C2-N1	2.71	1.40	1.35
3	A	355	GDP	O4'-C1'	3.01	1.45	1.41
3	A	355	GDP	C6-N1	5.04	1.42	1.33
3	C	357	GDP	C6-N1	5.09	1.42	1.33
3	D	358	GDP	C6-N1	5.25	1.42	1.33
3	B	356	GDP	C6-N1	5.78	1.43	1.33

All (63) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	355	GDP	C5-C6-N1	-8.60	111.84	123.59
3	D	358	GDP	C5-C6-N1	-8.40	112.10	123.59
3	B	356	GDP	C5-C6-N1	-8.34	112.19	123.59
3	C	357	GDP	C5-C6-N1	-8.17	112.42	123.59
3	C	357	GDP	O5'-PA-O1A	-7.76	79.48	109.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	356	GDP	O5'-PA-O1A	-7.69	79.75	109.62
3	D	358	GDP	O5'-PA-O1A	-7.68	79.80	109.62
3	A	355	GDP	O5'-PA-O1A	-7.63	80.01	109.62
3	A	355	GDP	O3A-PA-O5'	-4.60	90.74	102.94
3	D	358	GDP	O3A-PA-O5'	-4.34	91.42	102.94
3	B	356	GDP	O3A-PA-O5'	-4.14	91.96	102.94
3	C	357	GDP	O3A-PA-O5'	-4.10	92.06	102.94
3	C	357	GDP	C1'-N9-C4	-3.23	122.08	126.94
3	B	356	GDP	O4'-C4'-C3'	-3.10	98.89	105.15
3	A	355	GDP	O4'-C4'-C3'	-3.08	98.94	105.15
3	C	357	GDP	O4'-C4'-C3'	-3.04	99.01	105.15
3	D	358	GDP	C1'-N9-C4	-3.04	122.36	126.94
3	D	358	GDP	O4'-C4'-C3'	-2.99	99.12	105.15
3	B	356	GDP	C1'-N9-C4	-2.94	122.50	126.94
3	A	355	GDP	PA-O3A-PB	-2.84	123.15	132.67
3	C	357	GDP	N3-C2-N1	-2.80	123.18	127.44
3	D	358	GDP	N3-C2-N1	-2.78	123.21	127.44
3	B	356	GDP	N3-C2-N1	-2.76	123.24	127.44
3	C	357	GDP	PA-O3A-PB	-2.66	123.74	132.67
3	A	355	GDP	C1'-N9-C4	-2.66	122.93	126.94
3	A	355	GDP	N3-C2-N1	-2.57	123.53	127.44
3	D	358	GDP	C6-C5-C4	-2.54	117.86	120.90
3	C	357	GDP	O2A-PA-O5'	-2.53	95.72	108.46
3	D	358	GDP	PA-O3A-PB	-2.51	124.27	132.67
3	D	358	GDP	O2A-PA-O5'	-2.51	95.83	108.46
3	B	356	GDP	PA-O3A-PB	-2.50	124.27	132.67
3	B	356	GDP	C6-C5-C4	-2.46	117.96	120.90
3	C	357	GDP	C6-C5-C4	-2.46	117.96	120.90
3	A	355	GDP	C6-C5-C4	-2.43	118.00	120.90
3	B	356	GDP	O2A-PA-O5'	-2.36	96.55	108.46
3	A	355	GDP	O3B-PB-O2B	-2.11	99.33	107.38
3	A	355	GDP	O5'-C5'-C4'	-2.09	101.41	109.12
3	B	356	GDP	O5'-C5'-C4'	-2.07	101.47	109.12
3	D	358	GDP	O3B-PB-O2B	-2.04	99.60	107.38
3	C	357	GDP	O5'-C5'-C4'	-2.03	101.64	109.12
3	A	355	GDP	O2A-PA-O5'	-2.01	98.34	108.46
3	B	356	GDP	O3B-PB-O1B	2.01	117.05	110.58
3	A	355	GDP	O3B-PB-O1B	2.02	117.07	110.58
3	A	355	GDP	O2A-PA-O1A	3.07	129.14	112.53
3	B	356	GDP	O3B-PB-O3A	3.11	119.18	105.09
3	D	358	GDP	C4'-O4'-C1'	3.12	113.15	109.72
3	A	355	GDP	O3B-PB-O3A	3.17	119.48	105.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	358	GDP	O3B-PB-O3A	3.18	119.54	105.09
3	D	358	GDP	O2A-PA-O1A	3.19	129.79	112.53
3	B	356	GDP	O2A-PA-O1A	3.19	129.82	112.53
3	C	357	GDP	O2A-PA-O1A	3.28	130.28	112.53
3	C	357	GDP	O3B-PB-O3A	3.34	120.24	105.09
3	B	356	GDP	C4'-O4'-C1'	3.50	113.57	109.72
3	C	357	GDP	C4'-O4'-C1'	3.57	113.64	109.72
3	A	355	GDP	C4'-O4'-C1'	3.68	113.77	109.72
3	C	357	GDP	O2A-PA-O3A	5.73	131.08	105.09
3	D	358	GDP	O2A-PA-O3A	5.88	131.75	105.09
3	A	355	GDP	O2A-PA-O3A	5.92	131.94	105.09
3	B	356	GDP	O2A-PA-O3A	6.05	132.52	105.09
3	C	357	GDP	C6-N1-C2	6.48	124.94	115.94
3	A	355	GDP	C6-N1-C2	6.57	125.06	115.94
3	B	356	GDP	C6-N1-C2	6.70	125.24	115.94
3	D	358	GDP	C6-N1-C2	6.79	125.37	115.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	355	GDP	2	0
3	B	356	GDP	2	0
3	C	357	GDP	3	0
3	D	358	GDP	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	299/329 (90%)	0.35	17 (5%) 27 31	19, 40, 85, 154	0
1	B	307/329 (93%)	0.24	20 (6%) 22 25	16, 37, 78, 136	0
1	C	298/329 (90%)	0.21	14 (4%) 35 40	21, 38, 80, 104	0
1	D	304/329 (92%)	0.37	21 (6%) 20 22	22, 43, 96, 171	0
2	E	12/16 (75%)	1.29	2 (16%) 2 2	43, 64, 84, 142	0
2	F	12/16 (75%)	0.02	0 100 100	24, 44, 69, 110	0
2	G	12/16 (75%)	0.28	0 100 100	34, 48, 91, 130	0
2	H	12/16 (75%)	1.08	2 (16%) 2 2	43, 66, 100, 148	0
All	All	1256/1380 (91%)	0.31	76 (6%) 25 27	16, 40, 85, 171	0

All (76) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	257	LYS	6.3
1	D	258	TRP	5.6
1	C	59	ALA	5.0
1	A	345	LYS	4.9
2	E	12	ASP	4.8
1	C	67	LYS	4.4
1	C	258	TRP	4.4
2	H	8	PHE	4.0
1	A	258	TRP	3.8
1	A	233	VAL	3.8
1	D	194	LEU	3.6
1	B	168	ILE	3.6
1	A	344	ILE	3.6
1	D	60	GLY	3.6
1	C	60	GLY	3.5
1	B	64	GLU	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	257	LYS	3.4
1	C	168	ILE	3.3
1	B	95	PHE	3.2
1	C	105	ARG	3.2
1	B	94	ASP	3.1
1	D	312	LYS	3.1
1	B	96	GLY	3.1
1	B	167	TYR	3.0
1	C	34	VAL	3.0
1	D	92	LYS	3.0
1	A	181	THR	3.0
1	B	236	GLU	3.0
1	A	91	LEU	2.9
1	A	204	GLN	2.9
1	D	188	HIS	2.9
1	B	166	ASN	2.8
1	D	166	ASN	2.8
1	A	190	THR	2.8
1	C	63	GLU	2.8
1	D	191	PHE	2.8
1	D	314	LYS	2.8
1	A	232	LEU	2.7
1	B	105	ARG	2.7
1	B	101	ALA	2.7
1	D	91	LEU	2.5
1	B	61	TYR	2.5
1	C	344	ILE	2.5
1	B	99	ALA	2.5
1	D	343	ILE	2.4
1	D	235	ALA	2.4
1	B	258	TRP	2.4
1	B	68	GLN	2.3
1	C	342	VAL	2.3
1	A	183	GLY	2.3
1	B	93	ILE	2.3
1	A	64	GLU	2.3
1	B	239	GLU	2.3
2	H	12	ASP	2.3
1	B	67	LYS	2.2
1	D	90	ARG	2.2
1	A	184	ILE	2.2
1	A	213	HIS	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	313	ARG	2.2
1	D	336	PHE	2.2
1	B	59	ALA	2.1
2	E	8	PHE	2.1
1	A	59	ALA	2.1
1	A	87	ALA	2.1
1	C	193	ASP	2.1
1	C	70	LYS	2.1
1	D	105	ARG	2.1
1	C	174	VAL	2.1
1	D	195	HIS	2.1
1	D	162	ILE	2.0
1	D	236	GLU	2.0
1	A	182	THR	2.0
1	B	112	GLY	2.0
1	C	343	ILE	2.0
1	B	91	LEU	2.0
1	D	241	ASN	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	GDP	B	356	28/28	0.94	0.18	2.90	23,30,48,51	0
3	GDP	A	355	28/28	0.94	0.18	1.57	23,31,49,50	0
3	GDP	D	358	28/28	0.95	0.17	1.12	28,35,56,59	0
3	GDP	C	357	28/28	0.93	0.18	1.00	25,31,58,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.