



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 4, 2026 – 10:29 AM EDT

PDB ID : 1SHZ / pdb_00001shz
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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

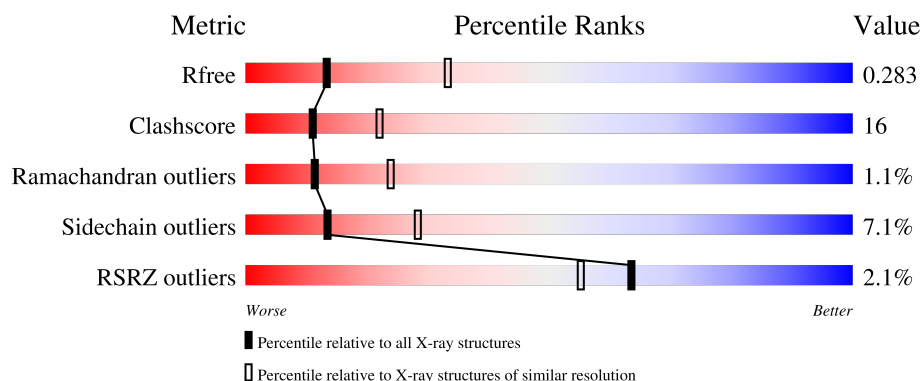
1 Overall quality at a glance

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}	180053	1407 (2.88-2.84)
Clashscore	190562	1446 (2.88-2.84)
Ramachandran outliers	187476	1406 (2.88-2.84)
Sidechain outliers	187428	1407 (2.88-2.84)
RSRZ outliers	180081	1408 (2.88-2.84)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	 61% 29% 5% .
1	D	340	 57% 36% . .
2	C	233	 4% 52% 31% . 16%
2	F	233	 3% 47% 32% . 19%

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			

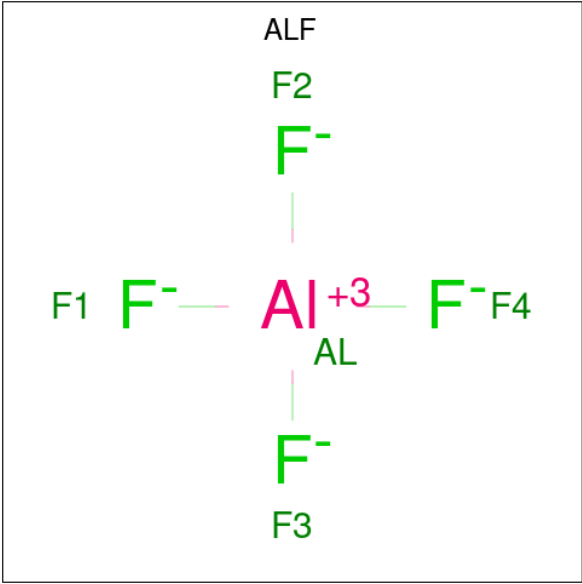
- Molecule 2 is a protein called Rho guanine nucleotide exchange factor 1.

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 (CCD ID: MG) (formula: Mg).

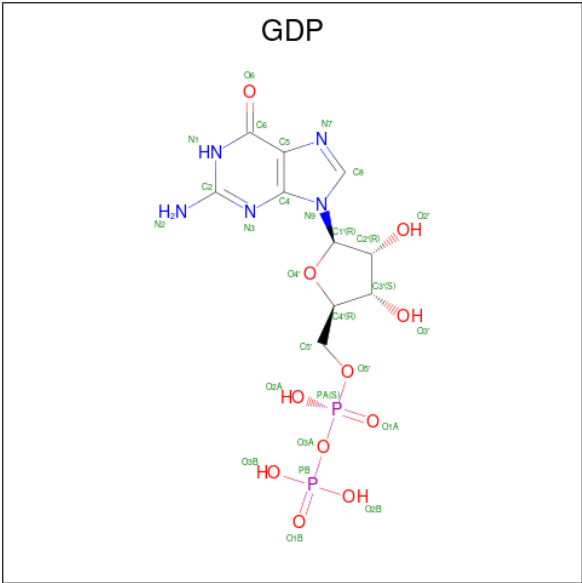
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		

- Molecule 4 is TETRAFLUOROALUMINATE ION (CCD ID: ALF) (formula: 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 (CCD ID: GDP) (formula: C₁₀H₁₅N₅O₁₁P₂).



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

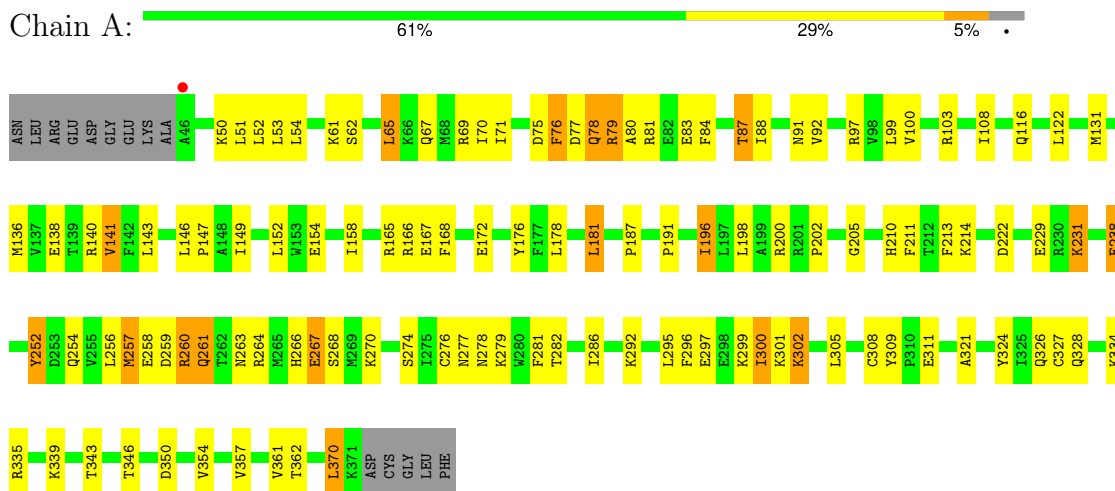
- Molecule 6 is water.

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

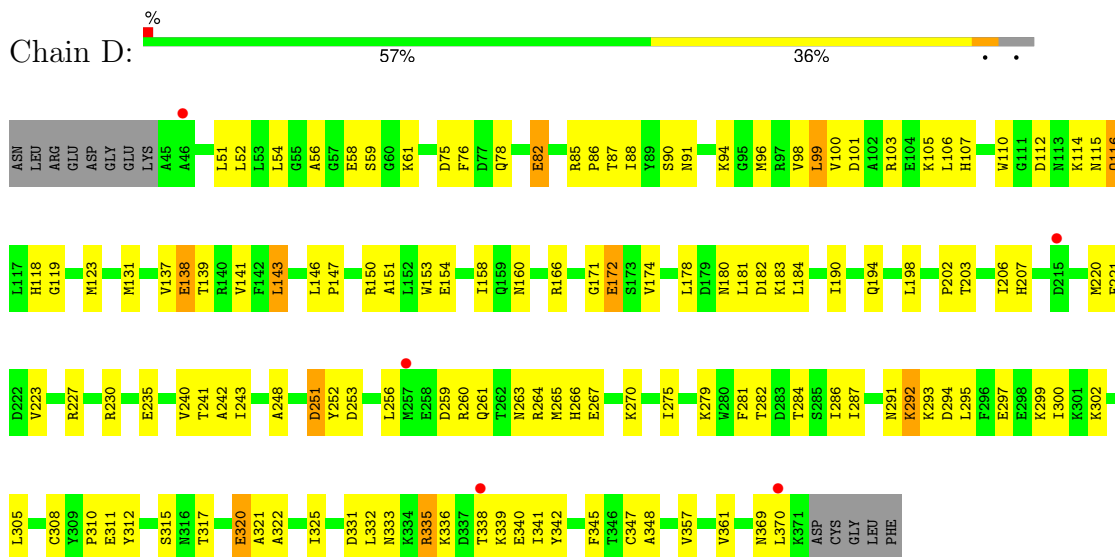
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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 Galpha(13):Galpha(i1) Chimera

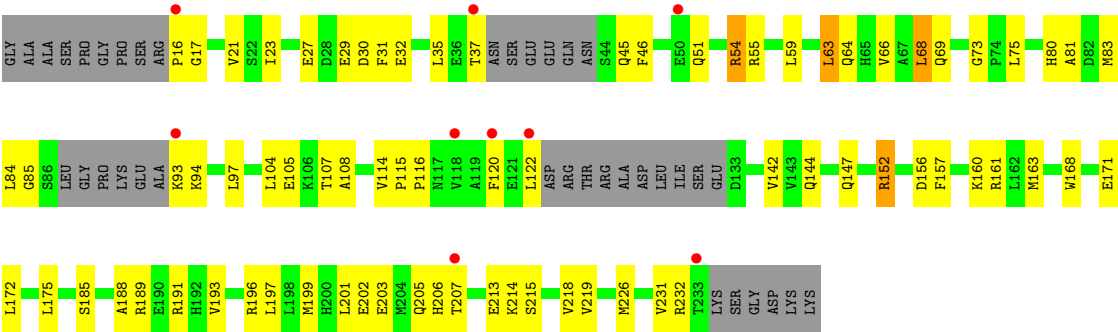


• Molecule 1: Guanine Nucleotide-Binding Protein Galpha(13):Galpha(i1) Chimera

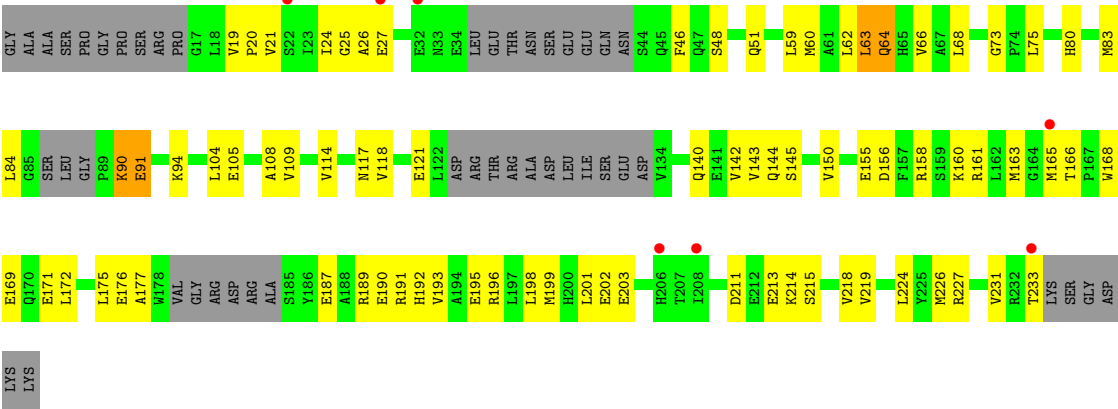


• Molecule 2: Rho guanine nucleotide exchange factor 1





● 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, α , β , γ	199.75Å 105.27Å 71.75Å 90.00° 96.91° 90.00°	Depositor
Resolution (Å)	46.49 – 2.85 46.49 – 2.85	Depositor EDS
% Data completeness (in resolution range)	90.7 (46.49-2.85) 90.7 (46.49-2.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.19	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 2.77Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.229 , 0.297 0.216 , 0.283	Depositor DCC
R_{free} test set	2203 reflections (6.63%)	wwPDB-VP
Wilson B-factor (Å ²)	58.2	Xtriage
Anisotropy	0.099	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 53.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8602	wwPDB-VP
Average B, all atoms (Å ²)	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.*

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: ALF, GDP, 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	0.52	0/2734	0.89	2/3680 (0.1%)
1	D	0.53	0/2739	0.88	0/3687
2	C	0.43	0/1605	0.84	2/2164 (0.1%)
2	F	0.44	0/1543	0.83	3/2078 (0.1%)
All	All	0.49	0/8621	0.87	7/11609 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	252	TYR	N-CA-C	-7.61	103.58	113.17
2	F	114	VAL	CA-C-N	5.23	123.48	119.66
2	F	114	VAL	C-N-CA	5.23	123.48	119.66
2	F	155	GLU	N-CA-C	-5.19	105.51	111.07
1	A	76	PHE	N-CA-C	5.04	121.53	110.80
2	C	114	VAL	CA-C-N	5.01	123.32	119.66
2	C	114	VAL	C-N-CA	5.01	123.32	119.66

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	85	0
1	D	2682	0	2655	94	0
2	C	1576	0	1561	60	0
2	F	1515	0	1504	58	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
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	278	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 16.

All (278) 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:230:ARG:HD3	2:F:165:MET:HE1	1.34	1.08
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:252:TYR:HB3	1:D:265:MET:HE2	1.62	0.82
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:HG2	2:F:161:ARG:HH11	1.57	0.67
2:C:189:ARG:O	2:C:193:VAL:HG23	1.93	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:279:LYS:HA	1:D:282:THR:HG23	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:75:ASP:CG	1:D:76:PHE:H	2.07	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:NH1	1:A:103:ARG:HB3	2.14	0.62
2:F:226:MET:HE3	2:F:231:VAL:HG11	1.79	0.62
1:A:279:LYS:HG3	2:C:69:GLN:HA	1.82	0.61
2:C:64:GLN:HE22	2:C:168:TRP:NE1	1.98	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:257:MET:HE1	2:C:27:GLU:O	2.01	0.61
1:D:51:LEU:HD23	1:D:220:MET:HE2	1.82	0.60
1:A:103:ARG:HB3	1:A:103:ARG:HH11	1.66	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:A:311:GLU:CD	1:A:311:GLU:H	2.11	0.59
1:D:178:LEU:HA	1:D:181:LEU:HD13	1.84	0.59
2:F:59:LEU:HG	2:F:63:LEU:CD2	2.32	0.58
1:A:172:GLU:H	1:A:172:GLU:CD	2.11	0.58
1:D:270:LYS:HB3	2:F:163:MET:HE1	1.86	0.58
1:D:243:ILE:HD11	1:D:281:PHE:CE1	2.38	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:D:227:ARG:HG2	1:D:230:ARG:NH1	2.19	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:HG2	2:F:161:ARG:NH1	2.20	0.56
1:A:256:LEU:HD11	1:A:267:GLU:HG3	1.87	0.56
1:D:287:ILE:HG12	1:D:361:VAL:HG13	1.86	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:A:70:ILE:HG23	1:A:75:ASP:OD2	2.06	0.55
1:A:80:ALA:HA	1:A:83:GLU:OE2	2.06	0.55
2:C:16:PRO:CB	2:F:24:ILE:HD11	2.35	0.55
2:F:156:ASP:O	2:F:160:LYS:HG2	2.07	0.55
1:D:150:ARG:O	1:D:154:GLU:HG2	2.06	0.55
2:F:201:LEU:HD23	2:F:201:LEU:O	2.06	0.55
1:A:274:SER:HB2	2:C:160:LYS:HZ2	1.72	0.54
1:A:300:ILE:HD12	1:A:301:LYS:N	2.22	0.54
1:D:256:LEU:HD11	1:D:267:GLU:CG	2.38	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:D:52:LEU:HD23	1:D:221:PHE:HB2	1.88	0.54
2:F:175:LEU:HD13	2:F:189:ARG:HD3	1.89	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
2:F:60:MET:HE1	2:F:201:LEU:HD13	1.88	0.53
1:A:252:TYR:O	1:A:308:CYS:HB2	2.09	0.53
2:C:55:ARG:HB3	6:C:504:HOH:O	2.08	0.53
1:D:253:ASP:OD2	1:D:299:LYS:HE2	2.09	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
1:D:320:GLU:H	1:D:320:GLU:CD	2.16	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
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:A:238:GLU:HG2	1:A:238:GLU:O	2.10	0.52
1:D:139:THR:HG22	1:D:143:LEU:HD22	1.92	0.52
2:C:168:TRP:O	2:C:172:LEU:HB2	2.09	0.52
2:C:191:ARG:NH1	2:C:232:ARG:HH12	2.07	0.52
1:D:101:ASP:HB2	1:D:123:MET:HE1	1.90	0.51
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:60:MET:HE1	2:F:201:LEU:CD1	2.41	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
2:C:80:HIS:O	2:C:83:MET:HB3	2.12	0.50
1:D:293:LYS:O	1:D:297:GLU:HG3	2.11	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
1:A:97:ARG:O	1:A:100:VAL:HG22	2.11	0.50
2:C:68:LEU:HD21	2:C:207:THR:HB	1.94	0.50
1:A:296:PHE:O	1:A:300:ILE:HG13	2.11	0.50
2:F:211:ASP:C	2:F:213:GLU:H	2.19	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:54:LEU:C	1:D:61:LYS:HD3	2.38	0.48
1:D:331:ASP:C	1:D:333:ASN:H	2.22	0.48
1:A:122:LEU:HD12	1:A:152:LEU:CD1	2.43	0.48
2:F:26:ALA:O	2:F:27:GLU:HB2	2.14	0.47
2:F:175:LEU:HD21	2:F:190:GLU:HA	1.96	0.47
1:A:277:ASN:HD21	1:A:334:LYS:HG3	1.79	0.47
1:A:370:LEU:H	1:A:370:LEU:HD22	1.79	0.47
1:D:322:ALA:HB1	1:D:345:PHE:CE1	2.48	0.47
1:A:166:ARG:NH1	1:A:167:GLU:HG3	2.29	0.47
1:A:176:TYR:CD2	1:A:198:LEU:HD13	2.48	0.47
1:A:191:PRO:HB2	1:A:196:ILE:CD1	2.44	0.47
2:F:143:VAL:C	2:F:145:SER:H	2.22	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:165:ARG:C	1:A:167:GLU:H	2.23	0.47
1:D:52:LEU:CD2	1:D:221:PHE:HB2	2.44	0.47
1:D:146:LEU:HB3	1:D:147:PRO:HD3	1.96	0.47
1:D:240:VAL:O	1:D:284:THR:HG23	2.14	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
2:F:46:PHE:O	2:F:158:ARG:HD3	2.15	0.47
1:D:317:THR:HB	1:D:320:GLU:CD	2.39	0.47
1:A:335:ARG:HG2	1:A:335:ARG:HH11	1.80	0.47
1:A:305:LEU:HD22	1:A:321:ALA:HB1	1.95	0.47
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
2:F:59:LEU:O	2:F:63:LEU:HD22	2.16	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
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
1:A:259:ASP:C	1:A:261:GLN:H	2.22	0.46
1:A:259:ASP:HA	2:F:19:VAL:CG1	2.45	0.46
2:C:68:LEU:HD11	2:C:207:THR:CG2	2.45	0.46
2:F:117:ASN:OD1	2:F:118:VAL:HG23	2.16	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
1:D:259:ASP:C	1:D:261:GLN:H	2.23	0.46
1:D:265:MET:HB3	1:D:308:CYS:SG	2.56	0.46
2:C:185:SER:O	2:C:188:ALA:HB3	2.16	0.45
2:C:196:ARG:HA	2:C:199:MET:HE3	1.98	0.45
1:D:100:VAL:CG2	1:D:101:ASP:N	2.79	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:F:201:LEU:HD23	2:F:201:LEU:C	2.42	0.45
2:C:161:ARG:HG2	2:C:161:ARG:HH11	1.81	0.45
1:D:153:TRP:CE3	1:D:153:TRP:HA	2.52	0.45
1:D:265:MET:HE1	1:D:325:ILE:HD13	1.98	0.45
1:A:191:PRO:HB2	1:A:196:ILE:HD11	1.98	0.45
1:D:357:VAL:O	1:D:361:VAL:HG23	2.17	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:226:MET:HB3	2:C:231:VAL:HB	1.99	0.45
2:C:115:PRO:HA	2:C:116:PRO:HD3	1.92	0.44
2:C:191:ARG:NH1	2:C:232:ARG:HH22	2.03	0.44
2:F:195:GLU:O	2:F:199:MET:HE3	2.16	0.44
2:C:152:ARG:HD3	1:D:116:GLN:HG3	1.98	0.44
1:A:324:TYR:O	1:A:328:GLN:HG2	2.17	0.44
2:C:120:PHE:C	2:C:122:LEU:H	2.25	0.44
1:D:369:ASN:HD21	1:D:370:LEU:HG	1.79	0.44
1:A:88:ILE:O	1:A:92:VAL:HG23	2.17	0.44
1:D:292:LYS:HG3	5:D:476:GDP:C6	2.52	0.44
1:A:50:LYS:HZ2	1:A:238:GLU:HG2	1.82	0.44
1:D:90:SER:OG	1:D:94:LYS:HE2	2.17	0.44
2:F:60:MET:HE2	2:F:198:LEU:HA	1.97	0.44
1:D:160:ASN:N	1:D:160:ASN:HD22	2.15	0.44
2:C:54:ARG:HE	2:C:54:ARG:HB2	1.70	0.44
1:D:100:VAL:HG23	1:D:101:ASP:N	2.32	0.44
1:A:213:PHE:CE1	1:A:214:LYS:HG3	2.53	0.44
2:C:64:GLN:NE2	2:C:168:TRP:NE1	2.66	0.44
2:C:226:MET:HE2	2:C:226:MET:HA	2.00	0.44
2:F:211:ASP:C	2:F:213:GLU:N	2.76	0.43
1:A:53:LEU:O	1:A:54:LEU:HD23	2.17	0.43
1:D:138:GLU:CD	1:D:138:GLU:H	2.25	0.43
1:D:235:GLU:OE2	1:D:235:GLU:N	2.47	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:51:LEU:C	1:A:51:LEU:HD23	2.43	0.43
1:A:165:ARG:HA	1:A:168:PHE:CE2	2.52	0.43
1:A:279:LYS:HG3	2:C:68:LEU:O	2.19	0.43
1:A:300:ILE:HG13	1:A:300:ILE:H	1.52	0.43
1:D:305:LEU:HD22	1:D:321:ALA:HB1	1.99	0.43
1:A:65:LEU:HD11	1:A:222:ASP:HB2	2.01	0.43
2:C:29:GLU:HA	2:C:32:GLU:CG	2.43	0.43
1:D:56:ALA:O	1:D:59:SER:HB2	2.19	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
2:F:59:LEU:HG	2:F:63:LEU:HD22	2.01	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:140:ARG:HG3	1:A:140:ARG:HH11	1.82	0.43
1:A:302:LYS:HB2	1:A:302:LYS:NZ	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:293:LYS:HG3	1:D:294:ASP:N	2.34	0.42
2:F:168:TRP:O	2:F:172:LEU:HB2	2.19	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:231:LYS:NZ	1:A:231:LYS:HB3	2.34	0.42
1:D:115:ASN:OD1	1:D:118:HIS:HD2	2.02	0.42
1:D:308:CYS:O	1:D:310:PRO:HD3	2.20	0.42
2:C:168:TRP:CD1	2:C:197:LEU:HD22	2.54	0.42
1:D:96:MET:O	1:D:100:VAL:HG13	2.19	0.42
1:D:291:ASN:ND2	5:D:476:GDP:N7	2.66	0.42
1:A:172:GLU:CD	1:A:172:GLU:N	2.76	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
1:A:270:LYS:HB3	2:C:163:MET:SD	2.60	0.42
1:A:278:ASN:HB3	1:A:281:PHE:CD2	2.55	0.42
2:C:175:LEU:HD12	2:C:189:ARG:NH2	2.35	0.42
1:D:51:LEU:HD12	1:D:242:ALA:HB3	2.01	0.42
1:D:78:GLN:H	1:D:78:GLN:CD	2.27	0.42
2:F:224:LEU:HA	2:F:227:ARG:NH1	2.35	0.42
1:A:309:TYR:HB3	1:A:311:GLU:OE2	2.20	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:D:158:ILE:HD12	1:D:158:ILE:N	2.34	0.42
1:A:131:MET:HE3	1:A:136:MET:SD	2.60	0.41
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
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:160:LYS:HA	2:C:160:LYS:HD3	1.89	0.41
2:C:202:GLU:O	2:C:205:GLN:HG3	2.20	0.41
1:A:54:LEU:C	1:A:61:LYS:HD3	2.46	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:A:51:LEU:HD23	1:A:52:LEU:N	2.36	0.41
1:A:205:GLY:H	1:A:229:GLU:CD	2.28	0.41
1:D:293:LYS:HG2	1:D:347:CYS:HB2	2.02	0.41
2:C:202:GLU:O	2:C:203:GLU:C	2.64	0.41
1:D:263:ASN:HB3	1:D:266:HIS:CD2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:161:ARG:HE	2:F:169:GLU:CD	2.29	0.40
1:A:210:HIS:O	1:A:211:PHE:HB3	2.21	0.40
2:C:168:TRP:HA	2:C:168:TRP:HE3	1.86	0.40
2:C:171:GLU:H	2:C:171:GLU:CD	2.30	0.40
2:C:161:ARG:HG2	2:C:161:ARG:NH1	2.36	0.40
1:D:112:ASP:C	1:D:114:LYS:N	2.78	0.40
1:D:287:ILE:HG13	1:D:342:TYR:HB2	2.03	0.40
2:F:75:LEU:HD13	2:F:150:VAL:HG13	2.03	0.40
1:D:100:VAL:CG2	1:D:123:MET:SD	3.10	0.40
1:D:281:PHE:O	1:D:339:LYS:HE3	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	324/340 (95%)	300 (93%)	22 (7%)	2 (1%)	21	38
1	D	325/340 (96%)	293 (90%)	28 (9%)	4 (1%)	10	22
2	C	188/233 (81%)	163 (87%)	23 (12%)	2 (1%)	11	24
2	F	178/233 (76%)	155 (87%)	20 (11%)	3 (2%)	7	16
All	All	1015/1146 (89%)	911 (90%)	93 (9%)	11 (1%)	11	24

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	260	ARG
1	A	187	PRO
1	D	348	ALA

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Mol	Chain	Res	Type
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%)	263 (91%)	27 (9%)	8	18
1	D	290/301 (96%)	266 (92%)	24 (8%)	10	22
2	C	170/199 (85%)	163 (96%)	7 (4%)	27	52
2	F	163/199 (82%)	156 (96%)	7 (4%)	26	50
All	All	913/1000 (91%)	848 (93%)	65 (7%)	13	28

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	62	SER
1	A	65	LEU
1	A	77	ASP
1	A	78	GLN
1	A	79	ARG
1	A	87	THR
1	A	99	LEU
1	A	116	GLN
1	A	141	VAL
1	A	143	LEU
1	A	154	GLU
1	A	158	ILE
1	A	181	LEU

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Mol	Chain	Res	Type
1	A	196	ILE
1	A	231	LYS
1	A	238	GLU
1	A	257	MET
1	A	258	GLU
1	A	261	GLN
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	339	LYS
1	A	370	LEU
2	C	37	THR
2	C	54	ARG
2	C	63	LEU
2	C	68	LEU
2	C	152	ARG
2	C	206	HIS
2	C	213	GLU
1	D	58	GLU
1	D	82	GLU
1	D	87	THR
1	D	99	LEU
1	D	103	ARG
1	D	105	LYS
1	D	116	GLN
1	D	131	MET
1	D	138	GLU
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	241	THR
1	D	251	ASP
1	D	275	ILE
1	D	292	LYS
1	D	300	ILE
1	D	311	GLU

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Mol	Chain	Res	Type
1	D	320	GLU
1	D	335	ARG
1	D	340	GLU
2	F	63	LEU
2	F	64	GLN
2	F	91	GLU
2	F	94	LYS
2	F	109	VAL
2	F	166	THR
2	F	203	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	78	GLN
1	A	91	ASN
1	A	116	GLN
1	A	160	ASN
1	A	169	GLN
1	A	210	HIS
1	A	266	HIS
1	A	277	ASN
1	A	316	ASN
1	A	355	GLN
2	C	45	GLN
2	C	47	GLN
2	C	51	GLN
2	C	58	HIS
2	C	64	GLN
2	C	192	HIS
2	C	205	GLN
1	D	118	HIS
1	D	160	ASN
1	D	193	GLN
1	D	194	GLN
1	D	254	GLN
1	D	316	ASN
1	D	326	GLN
1	D	368	ASN
2	F	64	GLN
2	F	135	GLN
2	F	220	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

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	$\# Z > 2$
5	GDP	D	476	4,3	29,30,30	1.32	3 (10%)	45,47,47	1.57	10 (22%)
4	ALF	D	378	5,6	4,4,4	0.87	0	-		
5	GDP	A	475	4,3	29,30,30	1.32	3 (10%)	45,47,47	1.57	10 (22%)
4	ALF	A	378	5,6	4,4,4	0.99	0	-		

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	GDP	D	476	4,3	-	2/16/32/32	0/3/3/3
5	GDP	A	475	4,3	-	2/16/32/32	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	476	GDP	C1'-N9	2.32	1.54	1.47
5	A	475	GDP	C1'-N9	2.31	1.54	1.47
5	A	475	GDP	O5'-C5'	-2.27	1.36	1.44
5	D	476	GDP	O5'-C5'	-2.25	1.36	1.44
5	A	475	GDP	C3'-C4'	-2.17	1.47	1.53
5	D	476	GDP	C3'-C4'	-2.16	1.47	1.53

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	476	GDP	O2A-PA-O3A	4.73	120.06	107.27
5	A	475	GDP	O2A-PA-O3A	4.72	120.04	107.27
5	A	475	GDP	O2'-C2'-C1'	-3.26	98.88	110.10
5	D	476	GDP	O2'-C2'-C1'	-3.24	98.93	110.10
5	D	476	GDP	C2'-C1'-N9	2.72	120.82	113.25
5	A	475	GDP	C2'-C1'-N9	2.72	120.81	113.25
5	A	475	GDP	C3'-C2'-C1'	2.70	106.58	101.46
5	D	476	GDP	C3'-C2'-C1'	2.70	106.57	101.46
5	D	476	GDP	O3A-PA-O1A	-2.58	102.93	110.70
5	A	475	GDP	O3A-PA-O1A	-2.58	102.94	110.70
5	D	476	GDP	O3'-C3'-C2'	-2.41	104.09	111.82
5	A	475	GDP	O3'-C3'-C2'	-2.40	104.11	111.82
5	A	475	GDP	O4'-C1'-C2'	-2.31	101.68	106.62
5	D	476	GDP	N9-C8-N7	2.31	117.68	113.40
5	A	475	GDP	N9-C8-N7	2.30	117.66	113.40
5	D	476	GDP	O4'-C1'-C2'	-2.30	101.70	106.62
5	A	475	GDP	C1'-N9-C8	2.21	133.01	126.73
5	D	476	GDP	C1'-N9-C8	2.20	132.99	126.73
5	D	476	GDP	O2B-PB-O1B	2.14	119.19	110.83
5	A	475	GDP	O2B-PB-O1B	2.14	119.17	110.83

There are no chirality outliers.

All (4) torsion outliers are listed below:

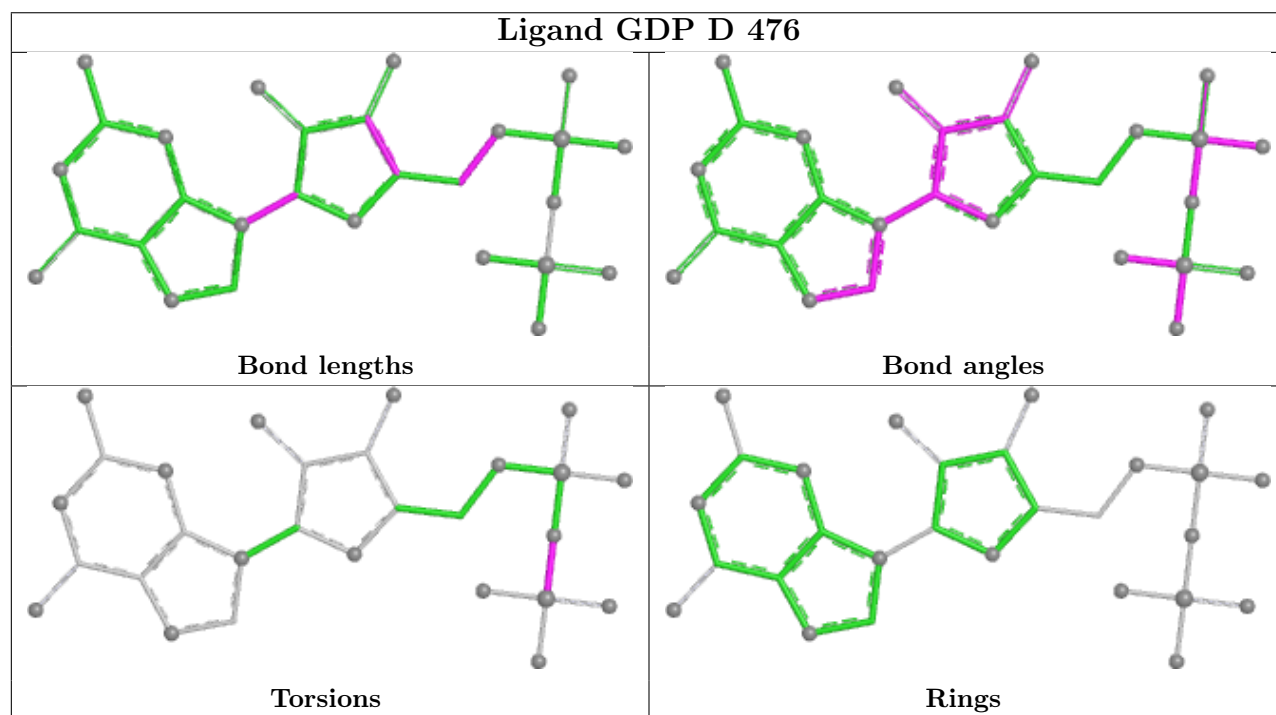
Mol	Chain	Res	Type	Atoms
5	A	475	GDP	PA-O3A-PB-O2B
5	D	476	GDP	PA-O3A-PB-O2B
5	A	475	GDP	PA-O3A-PB-O1B
5	D	476	GDP	PA-O3A-PB-O1B

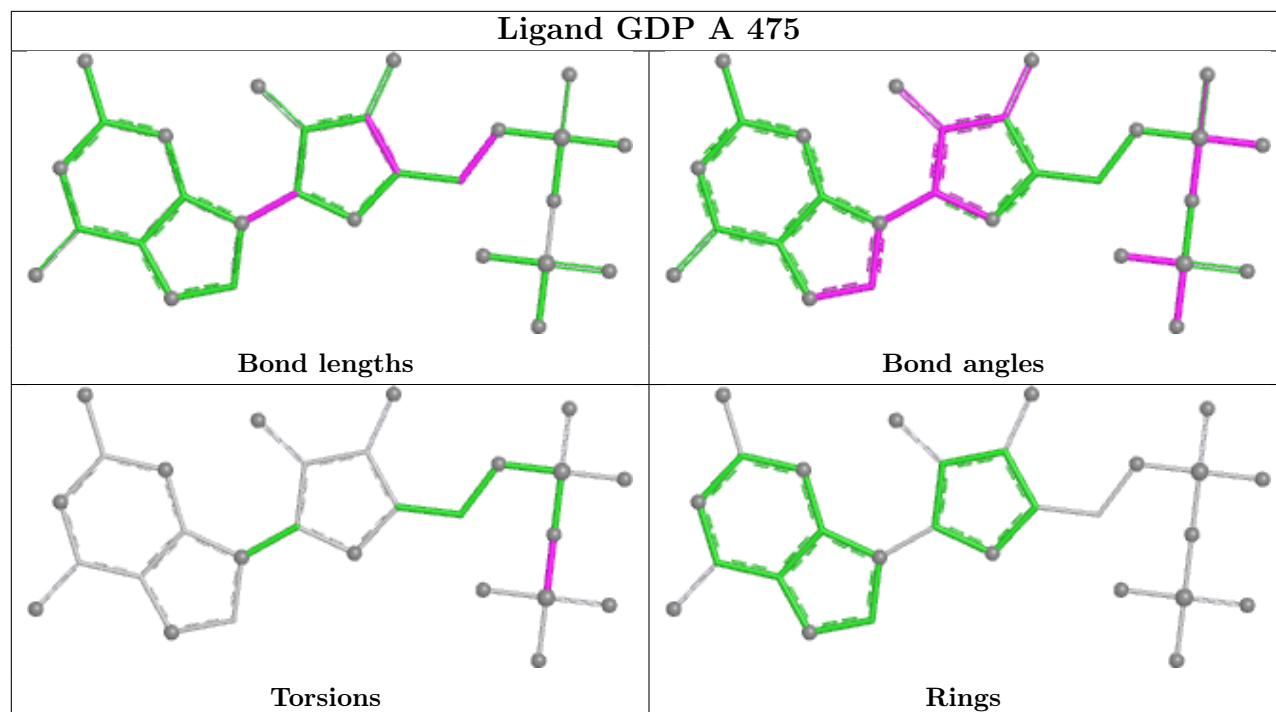
There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	476	GDP	2	0
5	A	475	GDP	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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	326/340 (95%)	-0.07	1 (0%) 90 89	22, 48, 73, 98	0
1	D	327/340 (96%)	-0.08	5 (1%) 72 65	22, 44, 76, 97	0
2	C	196/233 (84%)	0.38	9 (4%) 37 29	33, 72, 102, 103	1 (0%)
2	F	188/233 (80%)	0.44	7 (3%) 45 35	36, 73, 99, 103	1 (0%)
All	All	1037/1146 (90%)	0.10	22 (2%) 63 55	22, 53, 96, 103	2 (0%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	207	THR	3.9
2	C	37	THR	3.3
2	C	16	PRO	3.0
2	C	93	LYS	3.0
2	F	208	ILE	2.9
2	F	27	GLU	2.7
1	D	370	LEU	2.6
2	F	206	HIS	2.4
2	C	233	THR	2.4
1	D	257	MET	2.3
2	C	50	GLU	2.3
1	A	46	ALA	2.2
1	D	215	ASP	2.2
2	F	233	THR	2.2
2	F	22	SER	2.2
2	C	120	PHE	2.1
1	D	46	ALA	2.1
2	C	122	LEU	2.1
2	F	165	MET	2.1
2	F	32	GLU	2.1
1	D	338	THR	2.1
2	C	118	VAL	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

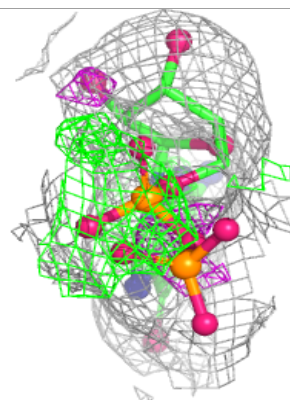
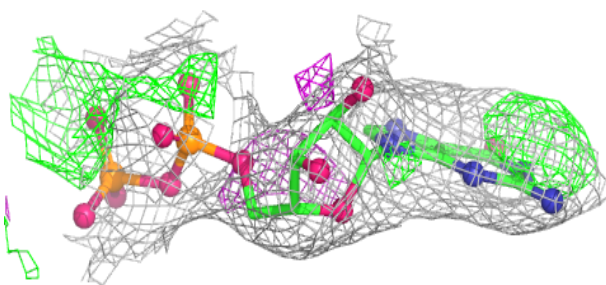
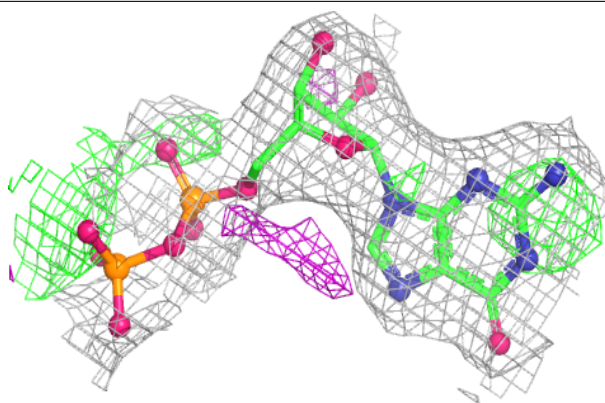
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	GDP	A	475	28/28	0.93	0.11	22,42,55,57	0
5	GDP	D	476	28/28	0.93	0.10	22,36,44,52	0
4	ALF	A	378	5/5	0.94	0.10	21,22,29,36	0
3	MG	D	377	1/1	0.96	0.17	9,9,9,9	0
3	MG	A	377	1/1	0.97	0.14	8,8,8,8	0
4	ALF	D	378	5/5	0.97	0.06	20,32,41,43	0

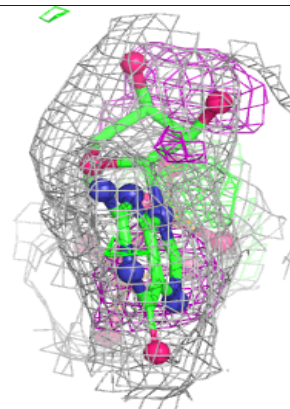
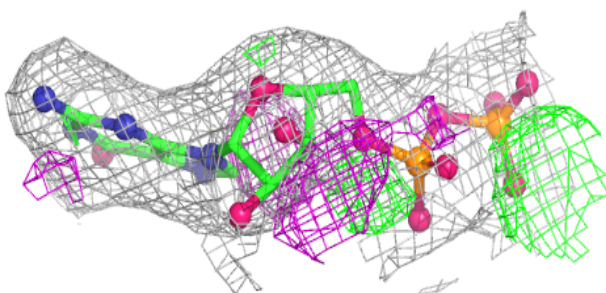
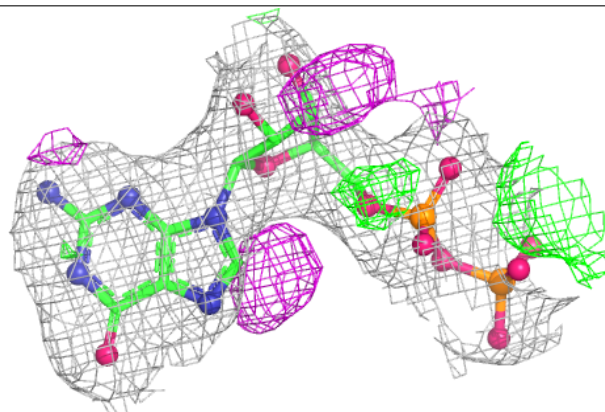
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around GDP A 475:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around GDP D 476:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers ⓘ

There are no such residues in this entry.