



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:03 AM UTC

PDB ID : 1Z2C / pdb_00001z2c
Title : Crystal structure of mDIA1 GBD-FH3 in complex with RhoC-GMPPNP
Authors : Rose, R.; Weyand, M.; Lammers, M.; Ishizaki, T.; Ahmadian, M.R.; Wittinghofer, A.
Deposited on : 2005-03-08
Resolution : 3.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<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 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

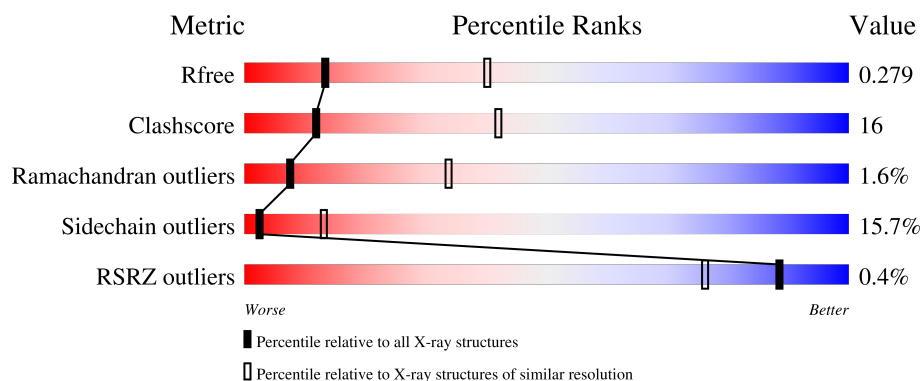
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	193	55%30%8%7%
1	C	193	% 58%28%6%7%
2	B	383	55%30%. . 10%
2	D	383	% 50%31%7%13%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8399 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 Rho-related GTP-binding protein RhoC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	179	Total	C	N	O	S	0	0	0
			1419	894	242	273	10			
1	C	179	Total	C	N	O	S	0	0	0
			1416	892	242	273	9			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	25	ASN	PHE	engineered mutation	UNP P08134
C	25	ASN	PHE	engineered mutation	UNP P08134

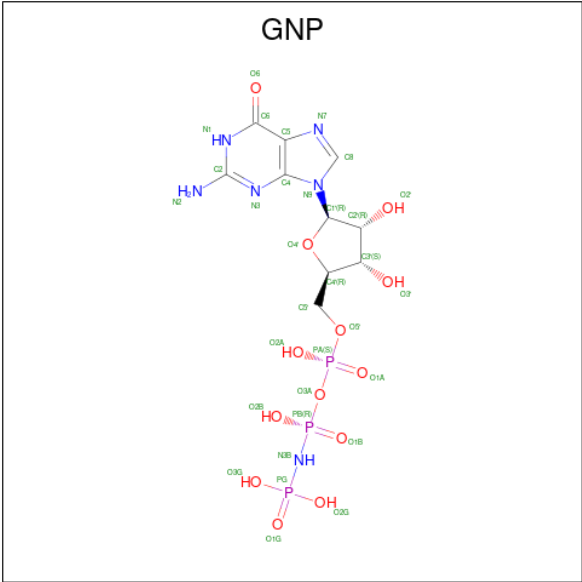
- Molecule 2 is a protein called Diaphanous protein homolog 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	346	Total	C	N	O	S	0	0	0
			2788	1756	477	531	24			
2	D	335	Total	C	N	O	S	0	0	0
			2709	1712	461	512	24			

- 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	B	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		

- Molecule 4 is PHOSPHOAMINOPHOSPHONIC ACID-GUANYLATE ESTER (CCD ID: GNP) (formula: C₁₀H₁₇N₆O₁₃P₃).

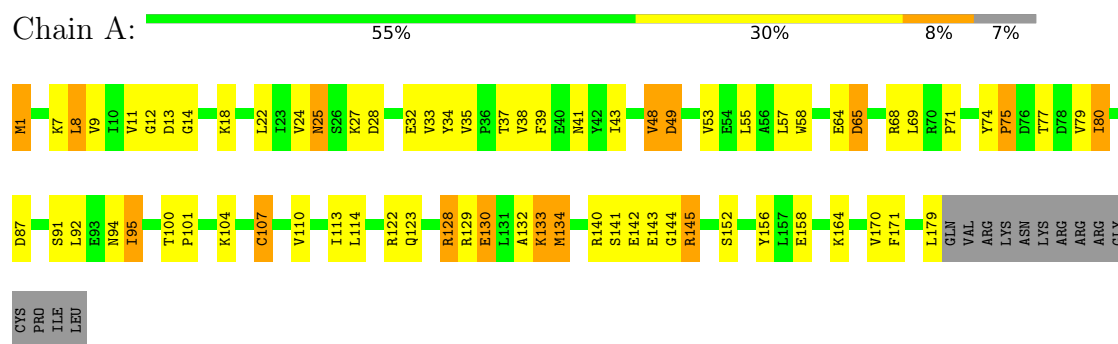


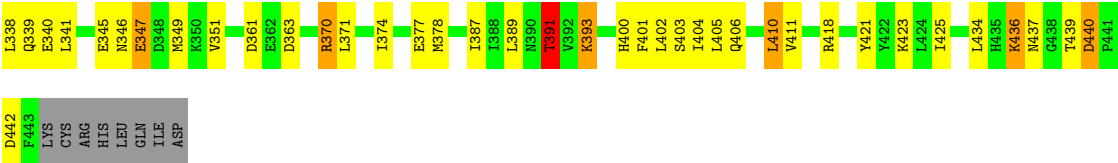
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			32	10	6	13	3		
4	C	1	Total	C	N	O	P	0	0
			32	10	6	13	3		

3 Residue-property plots [i](#)

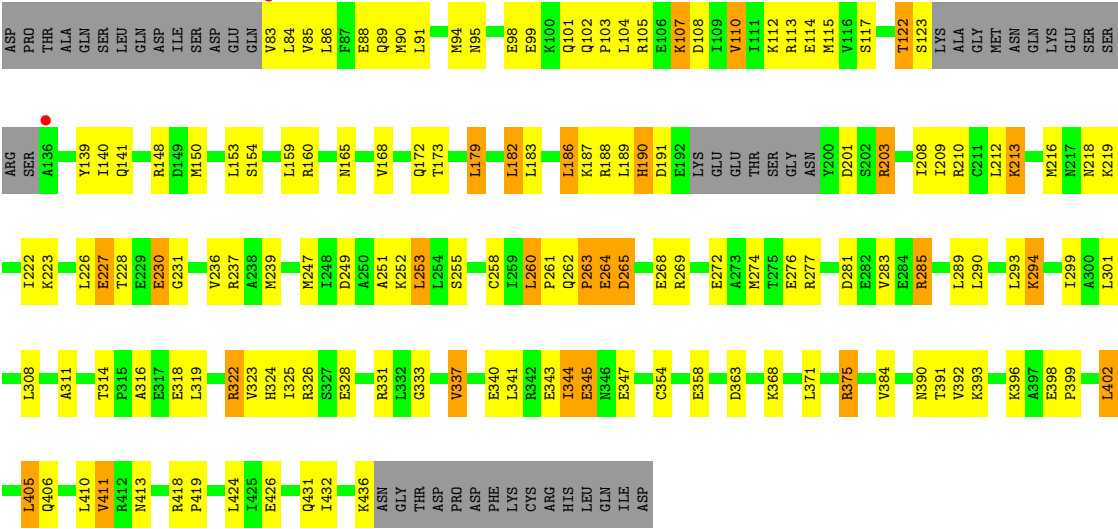
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Rho-related GTP-binding protein RhoC





● Molecule 2: Diaphanous protein homolog 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	148.64Å 85.35Å 123.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.00 20.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-3.00) 99.5 (20.00-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.95 (at 2.98Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.211 , 0.285 0.209 , 0.279	Depositor DCC
R_{free} test set	1605 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	57.5	Xtriage
Anisotropy	0.069	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 52.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8399	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.37% 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: GNP, 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	0.89	0/1445	1.12	5/1953 (0.3%)
1	C	0.80	1/1442 (0.1%)	1.08	2/1950 (0.1%)
2	B	0.91	1/2825 (0.0%)	1.18	12/3801 (0.3%)
2	D	0.96	3/2745 (0.1%)	1.17	14/3694 (0.4%)
All	All	0.90	5/8457 (0.1%)	1.15	33/11398 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	B	0	1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	107	LYS	CD-CE	16.96	2.03	1.52
2	D	107	LYS	CE-NZ	11.81	1.84	1.49
2	D	107	LYS	CG-CD	5.63	1.69	1.52
1	C	35	VAL	CA-CB	5.16	1.59	1.54
2	B	347	GLU	CG-CD	5.02	1.64	1.52

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	107	LYS	CG-CD-CE	-9.74	88.91	111.30
2	D	107	LYS	CD-CE-NZ	-8.56	84.51	111.90
2	D	263	PRO	CA-C-N	7.79	135.73	121.70
2	D	263	PRO	C-N-CA	7.79	135.73	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	442	ASP	N-CA-C	7.35	119.44	107.32
1	C	30	PHE	CA-C-N	7.31	128.97	119.84
1	C	30	PHE	C-N-CA	7.31	128.97	119.84
2	B	393	LYS	CA-C-N	-6.97	112.78	123.17
2	B	393	LYS	C-N-CA	-6.97	112.78	123.17
2	D	392	VAL	N-CA-C	-6.64	105.85	111.56
1	A	41	ASN	N-CA-C	6.34	119.92	108.69
2	B	253	LEU	N-CA-C	-6.33	104.46	111.36
1	A	35	VAL	CA-C-N	6.14	127.11	119.98
1	A	35	VAL	C-N-CA	6.14	127.11	119.98
2	B	391	THR	N-CA-C	-6.14	107.01	114.75
2	D	384	VAL	N-CA-C	-6.04	104.62	110.42
1	A	107	CYS	CA-C-N	-6.00	113.50	119.99
1	A	107	CYS	C-N-CA	-6.00	113.50	119.99
2	B	285	ARG	N-CA-C	5.88	118.47	111.71
2	D	393	LYS	CA-C-N	-5.83	112.40	122.14
2	D	393	LYS	C-N-CA	-5.83	112.40	122.14
2	D	141	GLN	N-CA-C	5.82	117.63	111.28
2	D	263	PRO	N-CA-C	5.78	124.37	112.47
2	B	339	GLN	N-CA-C	5.74	117.99	111.11
2	D	318	GLU	N-CA-C	5.63	118.14	109.07
2	B	149	ASP	CB-CA-C	-5.55	108.56	115.89
2	D	411	VAL	CB-CA-C	-5.52	103.10	111.33
2	D	344	ILE	N-CA-C	5.52	115.51	108.12
2	B	163	LEU	N-CA-C	-5.39	105.51	111.71
2	B	109	ILE	N-CA-CB	5.36	117.17	110.47
2	B	243	VAL	CA-C-N	5.23	124.67	119.24
2	B	243	VAL	C-N-CA	5.23	124.67	119.24
2	D	89	GLN	N-CA-C	-5.13	105.58	111.07

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	216	MET	Peptide

5.2 Too-close contacts

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	1419	0	1415	43	0
1	C	1416	0	1409	59	0
2	B	2788	0	2811	79	0
2	D	2709	0	2739	93	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
4	A	32	0	13	2	0
4	C	32	0	13	2	0
All	All	8399	0	8400	266	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 (266) 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:D:107:LYS:CE	2:D:107:LYS:NZ	1.84	1.38
2:D:107:LYS:CE	2:D:107:LYS:CD	2.03	1.35
2:B:262:GLN:HG3	2:B:263:PRO:HD3	1.20	1.12
2:D:239:MET:HE1	2:D:308:LEU:HD22	1.28	1.10
2:D:91:LEU:HD22	2:D:101:GLN:HG3	1.41	1.02
1:C:140:ARG:HB2	1:C:140:ARG:NH1	1.81	0.96
2:D:268:GLU:O	2:D:272:GLU:HG3	1.69	0.93
1:C:82:MET:HE3	1:C:95:ILE:HG23	1.50	0.92
2:D:264:GLU:O	2:D:265:ASP:HB2	1.70	0.90
2:B:214:ALA:O	2:B:217:ASN:HB2	1.73	0.89
2:B:262:GLN:CG	2:B:263:PRO:HD3	2.04	0.88
2:D:322:ARG:HG2	2:D:322:ARG:HH11	1.37	0.87
1:A:107:CYS:HB3	1:A:110:VAL:HG21	1.58	0.86
2:B:239:MET:HE1	2:B:308:LEU:HD12	1.56	0.86
1:A:128:ARG:HH11	1:A:128:ARG:HG3	1.40	0.85
2:B:277:ARG:NH2	2:B:280:MET:HG2	1.95	0.82
1:C:140:ARG:HH11	1:C:140:ARG:CB	1.93	0.81
1:C:113:ILE:CD1	1:C:174:ALA:HA	2.11	0.81
2:D:90:MET:O	2:D:94:MET:HG2	1.82	0.80
2:D:333:GLY:O	2:D:337:VAL:HG23	1.80	0.80
1:C:93:GLU:HA	1:C:93:GLU:OE1	1.83	0.79
2:D:239:MET:HE1	2:D:308:LEU:CD2	2.12	0.79
2:D:107:LYS:CE	2:D:107:LYS:CG	2.61	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:GLY:HA2	4:A:195:GNP:O3G	1.85	0.76
2:D:239:MET:CE	2:D:308:LEU:HD22	2.12	0.76
2:B:302:LYS:HB3	2:B:349:MET:HE1	1.68	0.76
1:C:82:MET:CE	1:C:95:ILE:HG23	2.17	0.75
1:C:140:ARG:NH1	1:C:140:ARG:CB	2.47	0.75
2:B:403:SER:HA	2:B:406:GLN:HE21	1.51	0.74
1:A:1:MET:HE2	2:D:114:GLU:OE1	1.88	0.73
1:A:130:GLU:OE2	1:A:133:LYS:HD2	1.88	0.73
1:A:134:MET:HE1	2:B:346:ASN:HD22	1.53	0.72
1:C:113:ILE:HD13	1:C:174:ALA:HA	1.71	0.72
1:A:25:ASN:OD1	1:A:171:PHE:CD1	2.44	0.70
2:D:260:LEU:HD23	2:D:263:PRO:HG2	1.73	0.70
1:A:8:LEU:HG	1:A:57:LEU:HD23	1.73	0.70
1:C:163:THR:O	1:C:164:LYS:HB2	1.92	0.70
2:B:90:MET:HG2	2:B:94:MET:HE2	1.72	0.70
1:C:124:ASP:HB3	1:C:127:THR:OG1	1.91	0.70
2:D:201:ASP:OD1	2:D:203:ARG:NH1	2.24	0.70
1:C:85:SER:OG	1:C:87:ASP:HB2	1.93	0.69
2:D:239:MET:HE3	2:D:251:ALA:HA	1.74	0.69
1:A:140:ARG:HB2	1:A:143:GLU:HG3	1.76	0.68
2:B:314:THR:HB	2:B:315:PRO:HD3	1.75	0.68
2:D:322:ARG:NH1	2:D:363:ASP:OD1	2.26	0.67
2:B:286:PHE:CD1	2:B:329:LEU:HD13	2.29	0.67
1:C:140:ARG:HB2	1:C:140:ARG:CZ	2.25	0.67
1:A:1:MET:HG2	2:D:114:GLU:OE2	1.95	0.66
2:D:322:ARG:HH11	2:D:322:ARG:CG	2.09	0.65
2:D:160:ARG:HD2	2:D:210:ARG:HH11	1.61	0.65
2:D:418:ARG:HB3	2:D:419:PRO:HD3	1.78	0.65
2:B:214:ALA:O	2:B:217:ASN:CB	2.44	0.65
1:C:140:ARG:HH11	1:C:140:ARG:HB3	1.60	0.65
1:A:123:GLN:HA	1:A:128:ARG:HH22	1.61	0.64
2:B:327:SER:HB3	2:B:439:THR:HG21	1.79	0.64
2:B:216:MET:HE1	2:B:225:MET:HE2	1.77	0.64
2:B:212:LEU:O	2:B:216:MET:HG3	1.97	0.63
1:C:113:ILE:HG23	1:C:173:MET:CE	2.28	0.63
2:D:107:LYS:NZ	2:D:107:LYS:CD	2.61	0.63
2:B:237:ARG:HG2	2:B:277:ARG:HD3	1.81	0.63
2:B:330:MET:HG2	2:B:437:ASN:HA	1.79	0.63
1:A:69:LEU:HA	2:B:94:MET:SD	2.39	0.63
1:C:123:GLN:HE21	1:C:123:GLN:N	1.97	0.62
1:C:125:GLU:HG3	1:C:129:ARG:NH1	2.14	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:371:LEU:HD21	2:D:375:ARG:NH1	2.15	0.62
2:B:216:MET:C	2:B:218:ASN:H	2.09	0.61
2:D:264:GLU:O	2:D:265:ASP:CB	2.45	0.61
2:B:403:SER:HA	2:B:406:GLN:NE2	2.15	0.60
1:C:125:GLU:HG3	1:C:129:ARG:HH11	1.66	0.60
2:B:313:ILE:HD11	2:B:325:ILE:HD11	1.83	0.60
2:B:262:GLN:HG3	2:B:263:PRO:CD	2.13	0.60
1:A:144:GLY:HA3	1:A:156:TYR:CZ	2.36	0.60
2:B:263:PRO:HB2	2:B:266:MET:HE3	1.83	0.60
2:D:277:ARG:HD2	2:D:281:ASP:OD2	2.02	0.59
2:B:91:LEU:HD22	2:B:101:GLN:HG2	1.84	0.59
2:D:294:LYS:HE3	2:D:294:LYS:HA	1.84	0.59
2:B:378:MET:CE	2:B:387:ILE:HG13	2.33	0.58
1:A:122:ARG:HH22	1:A:158:GLU:CD	2.11	0.58
2:D:247:MET:HE2	2:D:301:LEU:HD11	1.84	0.58
2:B:138:MET:O	2:B:142:GLU:HG3	2.03	0.58
2:B:322:ARG:NH1	2:B:363:ASP:OD1	2.36	0.58
2:B:216:MET:O	2:B:218:ASN:N	2.37	0.58
1:A:128:ARG:HG3	1:A:128:ARG:NH1	2.15	0.58
2:D:319:LEU:O	2:D:323:VAL:HG23	2.04	0.57
2:B:387:ILE:O	2:B:391:THR:HB	2.04	0.57
2:D:186:LEU:HG	2:D:208:ILE:HG21	1.86	0.57
2:B:212:LEU:HB3	2:B:216:MET:HE3	1.86	0.57
2:D:316:ALA:HB3	2:D:322:ARG:HD3	1.87	0.56
1:C:9:VAL:HG22	1:C:58:TRP:HB2	1.87	0.56
1:C:1:MET:O	1:C:2:ALA:HB3	2.04	0.56
2:B:334:LEU:HD12	2:B:337:VAL:CG2	2.35	0.56
2:B:320:ASP:OD1	2:B:370:ARG:NH2	2.39	0.56
2:D:160:ARG:CD	2:D:210:ARG:HH11	2.17	0.56
2:D:107:LYS:CE	2:D:107:LYS:HG2	2.35	0.55
2:D:189:LEU:C	2:D:191:ASP:H	2.15	0.55
2:B:102:GLN:HB3	2:B:103:PRO:HD3	1.88	0.55
2:D:341:LEU:C	2:D:343:GLU:H	2.15	0.55
2:D:212:LEU:HB3	2:D:216:MET:HE3	1.89	0.54
2:D:91:LEU:HD22	2:D:101:GLN:CG	2.28	0.54
2:B:235:LEU:O	2:B:238:ALA:HB3	2.07	0.54
1:C:6:LYS:HD2	1:C:179:LEU:HD13	1.89	0.54
1:C:93:GLU:O	1:C:96:PRO:HD2	2.08	0.54
1:C:113:ILE:HG23	1:C:173:MET:HE2	1.87	0.54
1:A:65:ASP:OD2	1:A:65:ASP:N	2.40	0.54
2:D:398:GLU:HB3	2:D:399:PRO:HD3	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:182:LEU:HD23	2:D:208:ILE:HG23	1.90	0.54
2:D:159:LEU:HD21	2:D:182:LEU:HD11	1.90	0.53
2:D:218:ASN:O	2:D:222:ILE:HG13	2.08	0.53
1:A:68:ARG:HH11	1:A:68:ARG:HG2	1.73	0.53
1:A:8:LEU:HG	1:A:57:LEU:CD2	2.38	0.53
2:D:91:LEU:CD2	2:D:101:GLN:HG3	2.29	0.53
1:C:113:ILE:HD11	1:C:174:ALA:HA	1.89	0.53
2:B:276:GLU:HG2	2:B:280:MET:HE1	1.91	0.53
1:C:1:MET:O	1:C:2:ALA:CB	2.57	0.53
1:C:13:ASP:OD2	1:C:91:SER:HA	2.09	0.53
2:D:95:ASN:HB2	2:D:168:VAL:HG21	1.91	0.53
2:D:213:LYS:HA	2:D:253:LEU:HD11	1.91	0.53
1:A:107:CYS:HB3	1:A:110:VAL:CG2	2.37	0.52
2:D:86:LEU:HD12	2:D:86:LEU:H	1.75	0.52
2:D:201:ASP:OD2	2:D:201:ASP:C	2.51	0.52
1:C:8:LEU:HD13	1:C:79:VAL:HG23	1.91	0.52
2:D:187:LYS:HE2	2:D:230:GLU:HG2	1.92	0.52
1:A:12:GLY:N	1:A:18:LYS:HD3	2.25	0.52
1:A:13:ASP:CG	1:A:94:ASN:HD22	2.18	0.51
1:C:126:HIS:HA	1:C:129:ARG:NH1	2.26	0.51
2:D:85:VAL:O	2:D:88:GLU:HB2	2.11	0.51
1:C:38:VAL:HG11	2:D:104:LEU:HG	1.92	0.51
2:D:285:ARG:NH1	2:D:328:GLU:OE1	2.44	0.51
1:C:87:ASP:OD1	1:C:127:THR:HG21	2.11	0.51
2:D:110:VAL:O	2:D:113:ARG:HB2	2.11	0.51
1:A:77:THR:HG21	1:A:80:ILE:HD12	1.93	0.50
1:C:113:ILE:HG12	1:C:173:MET:HE3	1.91	0.50
1:A:71:PRO:HA	1:A:74:TYR:CD2	2.46	0.50
2:B:216:MET:C	2:B:218:ASN:N	2.69	0.50
2:B:216:MET:CE	2:B:225:MET:HE2	2.42	0.50
1:C:85:SER:OG	1:C:118:LYS:HD2	2.12	0.50
1:C:160:SER:O	1:C:164:LYS:N	2.43	0.50
2:B:93:ASP:OD2	2:B:167:PRO:HB3	2.13	0.49
2:B:400:HIS:O	2:B:404:ILE:HG13	2.12	0.49
2:D:324:HIS:C	2:D:324:HIS:CD2	2.91	0.49
2:D:108:ASP:O	2:D:112:LYS:HG3	2.12	0.49
2:B:276:GLU:HG2	2:B:280:MET:CE	2.43	0.49
2:D:402:LEU:O	2:D:406:GLN:HG3	2.13	0.49
2:B:325:ILE:HD12	2:B:329:LEU:CD2	2.43	0.48
1:C:133:LYS:HB3	1:C:134:MET:HE2	1.96	0.48
1:C:85:SER:HG	1:C:118:LYS:HD2	1.77	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:237:ARG:HG2	2:B:277:ARG:CD	2.43	0.48
2:D:101:GLN:OE1	2:D:105:ARG:NH2	2.43	0.48
2:D:190:HIS:HB3	2:D:237:ARG:HH21	1.77	0.48
1:C:5:ARG:NE	1:C:54:GLU:OE1	2.46	0.48
2:D:262:GLN:N	2:D:263:PRO:HD3	2.28	0.48
2:D:289:LEU:HD21	2:D:308:LEU:HD23	1.96	0.48
2:B:88:GLU:O	2:B:92:VAL:HG23	2.14	0.48
1:C:122:ARG:C	1:C:123:GLN:HE21	2.22	0.48
1:C:174:ALA:O	1:C:177:ALA:HB3	2.14	0.48
1:A:8:LEU:C	1:A:8:LEU:HD12	2.39	0.47
2:B:97:ASN:OD1	2:B:97:ASN:C	2.57	0.47
1:C:79:VAL:HG11	1:C:177:ALA:HB3	1.96	0.47
2:D:371:LEU:O	2:D:375:ARG:HB2	2.15	0.47
2:B:325:ILE:O	2:B:329:LEU:HD22	2.15	0.47
1:A:122:ARG:NH2	1:A:158:GLU:OE1	2.48	0.47
2:B:118:GLN:OE1	2:B:118:GLN:HA	2.15	0.47
2:D:148:ARG:HD3	2:D:188:ARG:NH2	2.30	0.47
1:A:13:ASP:OD2	1:A:91:SER:HA	2.15	0.46
2:B:190:HIS:HB3	2:B:237:ARG:NH2	2.30	0.46
2:B:242:ALA:C	2:B:244:PRO:HD3	2.40	0.46
2:D:293:LEU:HA	2:D:293:LEU:HD23	1.67	0.46
2:D:337:VAL:O	2:D:341:LEU:HG	2.15	0.46
1:C:69:LEU:HD11	2:D:115:MET:HE1	1.98	0.46
2:D:139:TYR:O	2:D:140:ILE:C	2.58	0.46
2:B:389:LEU:HA	2:B:393:LYS:HB2	1.98	0.46
2:B:371:LEU:HD22	2:B:434:LEU:HD11	1.97	0.46
2:D:247:MET:HG2	2:D:301:LEU:HD11	1.97	0.46
2:B:177:GLU:O	2:B:178:GLY:C	2.59	0.46
1:C:28:ASP:OD2	1:C:29:GLN:HG3	2.16	0.46
2:B:134:ARG:HB3	2:B:138:MET:HB2	1.98	0.46
1:C:34:TYR:HB2	4:C:195:GNP:O3'	2.16	0.46
2:D:258:CYS:HB2	2:D:311:ALA:HB1	1.97	0.45
2:B:361:ASP:OD1	2:B:436:LYS:HE2	2.17	0.45
1:A:74:TYR:N	1:A:75:PRO:CD	2.80	0.45
2:D:122:THR:O	2:D:123:SER:C	2.60	0.45
2:D:368:LYS:O	2:D:371:LEU:HB3	2.16	0.45
1:A:9:VAL:HG22	1:A:58:TRP:HB2	1.98	0.45
2:B:121:HIS:O	2:B:123:SER:N	2.42	0.45
2:B:299:ILE:HD12	2:B:299:ILE:H	1.81	0.45
2:D:179:LEU:HD22	2:D:183:LEU:HG	1.97	0.45
1:A:34:TYR:HE1	4:A:195:GNP:HNB3	1.65	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:LEU:HD12	1:A:55:LEU:HA	1.88	0.45
1:C:74:TYR:O	1:C:75:PRO:C	2.60	0.45
1:C:122:ARG:O	1:C:122:ARG:HG3	2.17	0.45
1:A:87:ASP:OD1	1:A:87:ASP:C	2.59	0.44
2:B:239:MET:HE3	2:B:251:ALA:HB2	1.99	0.44
2:B:314:THR:HB	2:B:315:PRO:CD	2.44	0.44
2:D:249:ASP:O	2:D:253:LEU:HD22	2.17	0.44
2:D:262:GLN:HG3	2:D:262:GLN:O	2.16	0.44
2:B:234:LEU:HD23	2:B:234:LEU:HA	1.77	0.44
1:C:156:TYR:O	1:C:157:LEU:HD23	2.16	0.44
1:C:7:LYS:N	1:C:78:ASP:OD2	2.44	0.44
2:D:347:GLU:OE2	2:D:347:GLU:HA	2.18	0.44
2:D:223:LYS:O	2:D:227:GLU:HG2	2.18	0.44
2:B:302:LYS:CB	2:B:349:MET:HE1	2.43	0.44
1:C:67:ASP:O	2:D:165:ASN:ND2	2.49	0.43
2:B:410:LEU:HD23	2:B:410:LEU:HA	1.42	0.43
1:C:43:ILE:H	1:C:43:ILE:HG13	1.62	0.43
2:D:322:ARG:NH1	2:D:322:ARG:CG	2.75	0.43
2:D:354:CYS:O	2:D:358:GLU:HG3	2.18	0.43
2:B:208:ILE:O	2:B:211:CYS:HB2	2.18	0.43
2:B:334:LEU:HD12	2:B:337:VAL:HG22	2.00	0.43
1:C:40:GLU:HB3	1:C:59:ASP:HB3	2.01	0.43
1:A:145:ARG:HE	1:A:145:ARG:HB2	1.72	0.43
2:B:90:MET:HE3	2:B:94:MET:HE1	2.00	0.43
2:B:378:MET:HE1	2:B:387:ILE:HG13	1.97	0.43
1:A:48:VAL:O	1:A:49:ASP:C	2.61	0.43
2:B:401:PHE:O	2:B:402:LEU:C	2.62	0.43
2:B:322:ARG:HH11	2:B:363:ASP:CG	2.25	0.43
1:C:133:LYS:CB	1:C:134:MET:HE2	2.48	0.43
2:D:190:HIS:O	2:D:237:ARG:NH2	2.52	0.43
2:B:167:PRO:O	2:B:168:VAL:C	2.60	0.43
1:A:13:ASP:OD2	1:A:94:ASN:ND2	2.51	0.42
1:A:38:VAL:HG23	1:A:39:PHE:N	2.34	0.42
1:C:157:LEU:HD12	1:C:170:VAL:HA	2.00	0.42
2:D:94:MET:HA	2:D:94:MET:HE2	2.01	0.42
1:C:8:LEU:HD13	1:C:79:VAL:CG2	2.48	0.42
2:B:257:LEU:O	2:B:260:LEU:HB2	2.18	0.42
1:A:141:SER:HB3	1:A:145:ARG:HH21	1.84	0.42
1:A:114:LEU:O	1:A:156:TYR:HA	2.19	0.42
1:A:132:ALA:C	1:A:134:MET:H	2.27	0.42
2:B:242:ALA:O	2:B:244:PRO:HD3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:124:ASP:O	1:C:127:THR:N	2.53	0.42
2:D:262:GLN:H	2:D:263:PRO:HD3	1.84	0.42
2:D:285:ARG:HH11	2:D:328:GLU:CD	2.28	0.42
2:B:421:TYR:CZ	2:B:425:ILE:HD11	2.55	0.42
2:B:322:ARG:HB3	2:B:363:ASP:OD2	2.20	0.42
2:D:189:LEU:O	2:D:191:ASP:N	2.53	0.42
2:D:405:LEU:HD12	2:D:405:LEU:HA	1.86	0.42
2:D:418:ARG:HB3	2:D:419:PRO:CD	2.47	0.42
1:C:132:ALA:O	1:C:133:LYS:C	2.62	0.42
2:D:186:LEU:HD23	2:D:186:LEU:HA	1.89	0.42
2:D:236:VAL:HG11	2:D:274:MET:HA	2.00	0.42
1:A:8:LEU:C	1:A:8:LEU:CD1	2.93	0.41
1:A:100:THR:O	1:A:101:PRO:C	2.63	0.41
2:B:330:MET:CG	2:B:437:ASN:HA	2.47	0.41
2:D:102:GLN:N	2:D:103:PRO:CD	2.83	0.41
2:D:150:MET:HE2	2:D:150:MET:HB3	1.81	0.41
2:D:325:ILE:O	2:D:326:ARG:C	2.62	0.41
2:D:153:LEU:HD22	2:D:203:ARG:HG3	2.02	0.41
1:C:39:PHE:CD1	2:D:115:MET:HE3	2.55	0.41
1:C:118:LYS:HG2	4:C:195:GNP:C6	2.51	0.41
2:D:189:LEU:C	2:D:191:ASP:N	2.78	0.41
2:D:341:LEU:C	2:D:343:GLU:N	2.78	0.41
2:B:313:ILE:HD11	2:B:325:ILE:CD1	2.49	0.41
2:B:313:ILE:CD1	2:B:325:ILE:HD11	2.49	0.41
2:B:324:HIS:C	2:B:324:HIS:CD2	2.98	0.41
2:B:437:ASN:O	2:B:440:ASP:HB2	2.21	0.41
1:C:35:VAL:HA	1:C:36:PRO:HD2	1.92	0.41
1:C:123:GLN:HA	1:C:128:ARG:HH22	1.86	0.41
2:D:390:ASN:C	2:D:391:THR:CG2	2.94	0.41
2:B:167:PRO:O	2:B:169:SER:N	2.54	0.41
2:B:167:PRO:C	2:B:169:SER:N	2.79	0.40
1:A:7:LYS:HE3	1:A:58:TRP:CE2	2.57	0.40
1:A:11:VAL:C	1:A:18:LYS:HD3	2.46	0.40
1:A:92:LEU:HD12	1:A:95:ILE:HD13	2.03	0.40
2:B:302:LYS:HD3	2:B:341:LEU:HD22	2.02	0.40
1:C:5:ARG:NH2	1:C:54:GLU:OE1	2.54	0.40
1:C:28:ASP:HB3	1:C:29:GLN:NE2	2.37	0.40
2:D:260:LEU:HA	2:D:261:PRO:HD3	1.94	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	177/193 (92%)	156 (88%)	18 (10%)	3 (2%)	7	32
1	C	177/193 (92%)	160 (90%)	14 (8%)	3 (2%)	7	32
2	B	340/383 (89%)	307 (90%)	28 (8%)	5 (2%)	8	35
2	D	329/383 (86%)	292 (89%)	32 (10%)	5 (2%)	8	35
All	All	1023/1152 (89%)	915 (89%)	92 (9%)	16 (2%)	7	34

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	33	VAL
2	B	122	THR
2	B	217	ASN
1	C	28	ASP
2	D	190	HIS
2	D	265	ASP
2	D	345	GLU
1	A	133	LYS
1	C	2	ALA
2	D	122	THR
1	A	28	ASP
2	B	168	VAL
2	D	231	GLY
2	B	133	SER
1	C	133	LYS
2	B	261	PRO

5.3.2 Protein sidechains [i](#)

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	157/170 (92%)	127 (81%)	30 (19%)	1	8
1	C	156/170 (92%)	139 (89%)	17 (11%)	6	26
2	B	316/350 (90%)	267 (84%)	49 (16%)	2	13
2	D	306/350 (87%)	255 (83%)	51 (17%)	2	12
All	All	935/1040 (90%)	788 (84%)	147 (16%)	2	13

All (147) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	8	LEU
1	A	22	LEU
1	A	24	VAL
1	A	25	ASN
1	A	27	LYS
1	A	32	GLU
1	A	37	THR
1	A	43	ILE
1	A	48	VAL
1	A	49	ASP
1	A	53	VAL
1	A	64	GLU
1	A	65	ASP
1	A	75	PRO
1	A	79	VAL
1	A	80	ILE
1	A	95	ILE
1	A	104	LYS
1	A	113	ILE
1	A	128	ARG
1	A	129	ARG
1	A	130	GLU
1	A	134	MET
1	A	142	GLU
1	A	145	ARG
1	A	152	SER
1	A	164	LYS
1	A	170	VAL
1	A	179	LEU

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Mol	Chain	Res	Type
2	B	83	VAL
2	B	109	ILE
2	B	110	VAL
2	B	132	SER
2	B	147	LEU
2	B	152	LEU
2	B	158	SER
2	B	168	VAL
2	B	173	THR
2	B	179	LEU
2	B	182	LEU
2	B	201	ASP
2	B	203	ARG
2	B	210	ARG
2	B	222	ILE
2	B	228	THR
2	B	230	GLU
2	B	248	ILE
2	B	259	ILE
2	B	260	LEU
2	B	262	GLN
2	B	266	MET
2	B	280	MET
2	B	283	VAL
2	B	295	SER
2	B	306	LEU
2	B	308	LEU
2	B	317	GLU
2	B	319	LEU
2	B	323	VAL
2	B	325	ILE
2	B	329	LEU
2	B	331	ARG
2	B	338	LEU
2	B	340	GLU
2	B	345	GLU
2	B	347	GLU
2	B	351	VAL
2	B	370	ARG
2	B	374	ILE
2	B	377	GLU
2	B	391	THR

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Mol	Chain	Res	Type
2	B	405	LEU
2	B	410	LEU
2	B	411	VAL
2	B	418	ARG
2	B	423	LYS
2	B	436	LYS
2	B	440	ASP
1	C	6	LYS
1	C	16	CYS
1	C	32	GLU
1	C	43	ILE
1	C	60	THR
1	C	65	ASP
1	C	79	VAL
1	C	93	GLU
1	C	115	VAL
1	C	123	GLN
1	C	125	GLU
1	C	127	THR
1	C	137	GLU
1	C	142	GLU
1	C	168	ARG
1	C	176	ARG
1	C	179	LEU
2	D	83	VAL
2	D	84	LEU
2	D	98	GLU
2	D	99	GLU
2	D	110	VAL
2	D	117	SER
2	D	154	SER
2	D	172	GLN
2	D	173	THR
2	D	179	LEU
2	D	182	LEU
2	D	186	LEU
2	D	203	ARG
2	D	209	ILE
2	D	213	LYS
2	D	219	LYS
2	D	226	LEU
2	D	227	GLU

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Mol	Chain	Res	Type
2	D	228	THR
2	D	230	GLU
2	D	252	LYS
2	D	253	LEU
2	D	255	SER
2	D	260	LEU
2	D	264	GLU
2	D	269	ARG
2	D	276	GLU
2	D	283	VAL
2	D	285	ARG
2	D	290	LEU
2	D	294	LYS
2	D	299	ILE
2	D	314	THR
2	D	322	ARG
2	D	331	ARG
2	D	337	VAL
2	D	340	GLU
2	D	344	ILE
2	D	345	GLU
2	D	375	ARG
2	D	396	LYS
2	D	402	LEU
2	D	405	LEU
2	D	410	LEU
2	D	411	VAL
2	D	413	ASN
2	D	424	LEU
2	D	426	GLU
2	D	431	GLN
2	D	432	ILE
2	D	436	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	63	GLN
1	A	136	GLN
2	B	102	GLN
2	B	165	ASN
2	B	166	ASN

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Mol	Chain	Res	Type
2	B	204	ASN
2	B	262	GLN
2	B	324	HIS
2	B	335	HIS
2	B	339	GLN
2	B	406	GLN
2	B	431	GLN
2	B	437	ASN
1	C	123	GLN
2	D	165	ASN
2	D	262	GLN
2	D	310	ASN
2	D	324	HIS
2	D	335	HIS
2	D	339	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 3 are monoatomic - leaving 2 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	GNP	C	195	3	34,34,34	1.60	5 (14%)	47,54,54	1.43	7 (14%)
4	GNP	A	195	3	34,34,34	1.42	5 (14%)	47,54,54	1.17	4 (8%)

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	GNP	C	195	3	-	6/18/38/38	0/3/3/3
4	GNP	A	195	3	-	7/18/38/38	0/3/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	195	GNP	PA-O3A	-5.60	1.53	1.59
4	C	195	GNP	PB-O3A	-4.10	1.54	1.59
4	A	195	GNP	PA-O3A	-4.02	1.55	1.59
4	A	195	GNP	PG-O1G	3.95	1.52	1.46
4	A	195	GNP	PB-O3A	-3.39	1.54	1.59
4	A	195	GNP	PB-O2B	-3.18	1.48	1.56
4	C	195	GNP	PG-O1G	2.95	1.50	1.46
4	C	195	GNP	PB-O2B	-2.87	1.49	1.56
4	C	195	GNP	PG-O2G	-2.13	1.51	1.56
4	A	195	GNP	PG-O3G	-2.04	1.51	1.56

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	195	GNP	O3G-PG-O1G	-4.71	101.65	113.45
4	C	195	GNP	O2B-PB-O1B	4.26	119.01	109.87
4	C	195	GNP	O3G-PG-O1G	-3.81	103.89	113.45
4	C	195	GNP	O1B-PB-N3B	-3.69	106.34	111.77
4	C	195	GNP	O2G-PG-O3G	3.09	115.91	107.59
4	A	195	GNP	O2G-PG-O3G	3.02	115.71	107.59
4	A	195	GNP	O2B-PB-O1B	2.49	115.21	109.87
4	C	195	GNP	O2A-PA-O3A	2.27	113.41	107.27
4	A	195	GNP	O2B-PB-O3A	2.19	111.95	104.64
4	C	195	GNP	C2'-C3'-C4'	2.18	106.83	102.61
4	C	195	GNP	O3A-PB-N3B	2.15	112.56	106.59

There are no chirality outliers.

All (13) torsion outliers are listed below:

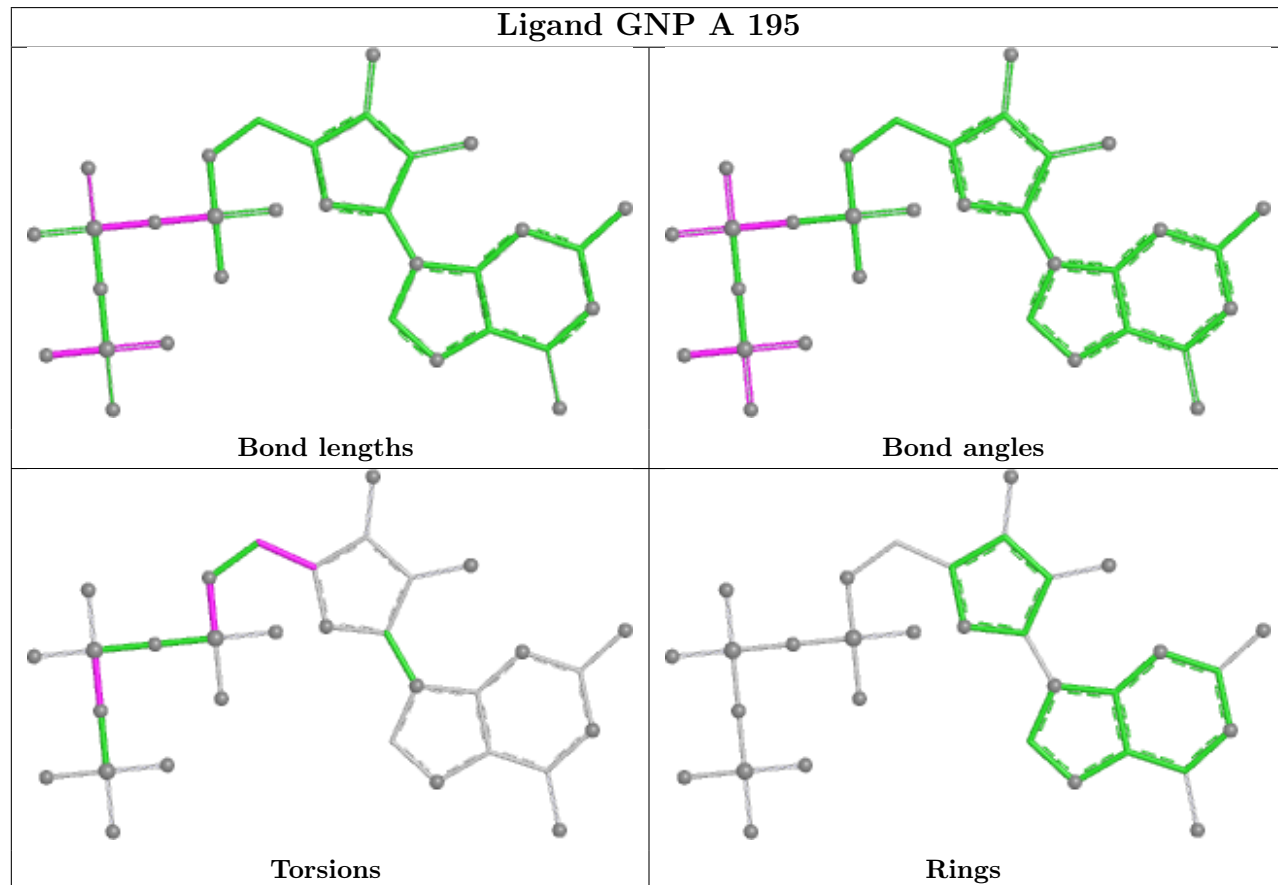
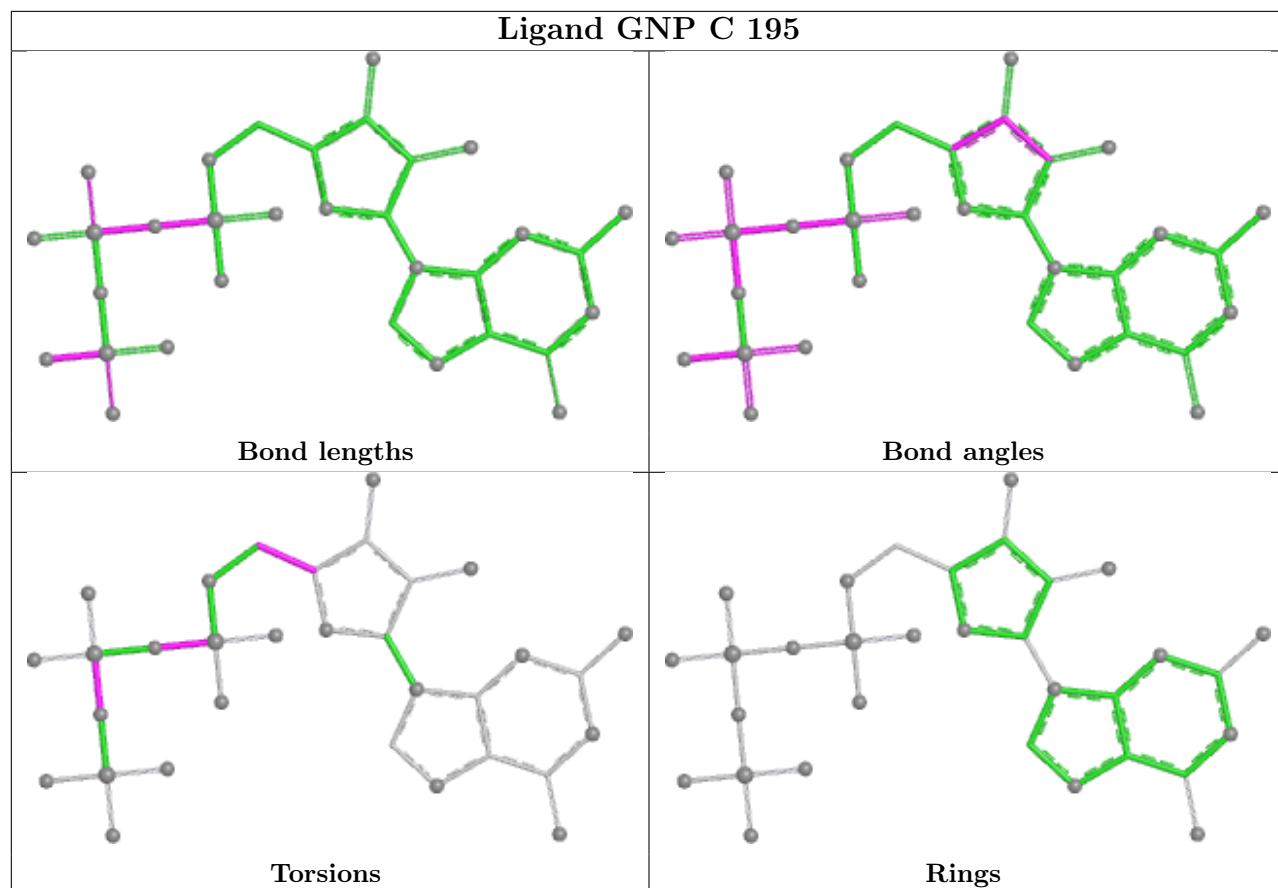
Mol	Chain	Res	Type	Atoms
4	A	195	GNP	PG-N3B-PB-O1B
4	A	195	GNP	PG-N3B-PB-O3A
4	C	195	GNP	PG-N3B-PB-O1B
4	C	195	GNP	PG-N3B-PB-O3A
4	A	195	GNP	O4'-C4'-C5'-O5'
4	A	195	GNP	C3'-C4'-C5'-O5'
4	C	195	GNP	O4'-C4'-C5'-O5'
4	C	195	GNP	C3'-C4'-C5'-O5'
4	A	195	GNP	C5'-O5'-PA-O3A
4	A	195	GNP	C5'-O5'-PA-O1A
4	A	195	GNP	C5'-O5'-PA-O2A
4	C	195	GNP	PB-O3A-PA-O2A
4	C	195	GNP	PB-O3A-PA-O1A

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	195	GNP	2	0
4	A	195	GNP	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	179/193 (92%)	-0.47	0 100 100	28, 52, 93, 108	0
1	C	179/193 (92%)	-0.40	1 (0%) 85 69	35, 56, 100, 119	0
2	B	346/383 (90%)	-0.55	1 (0%) 90 79	26, 44, 76, 100	0
2	D	335/383 (87%)	-0.49	2 (0%) 85 69	21, 49, 78, 95	0
All	All	1039/1152 (90%)	-0.49	4 (0%) 88 76	21, 48, 85, 119	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	83	VAL	3.7
2	D	136	ALA	2.8
2	B	83	VAL	2.5
1	C	29	GLN	2.2

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](#)

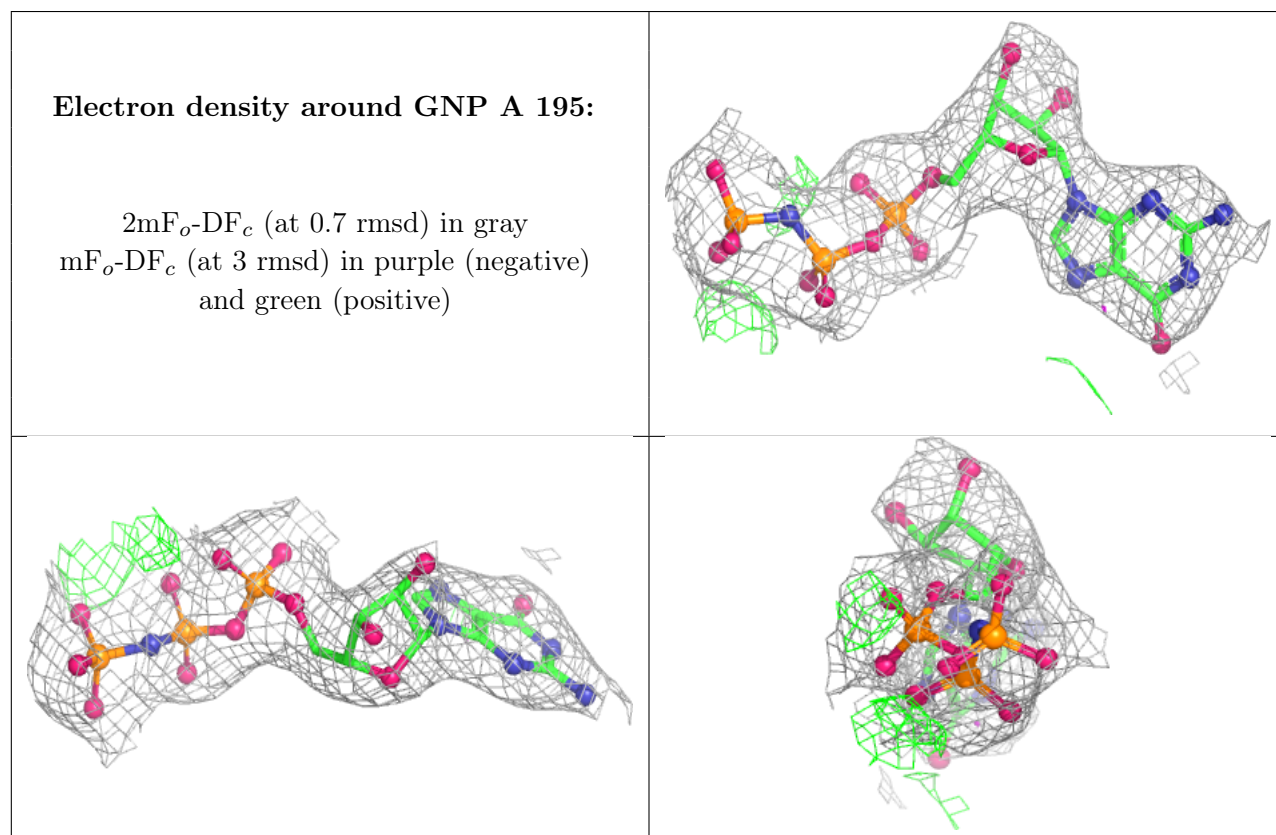
There are no oligosaccharides in this entry.

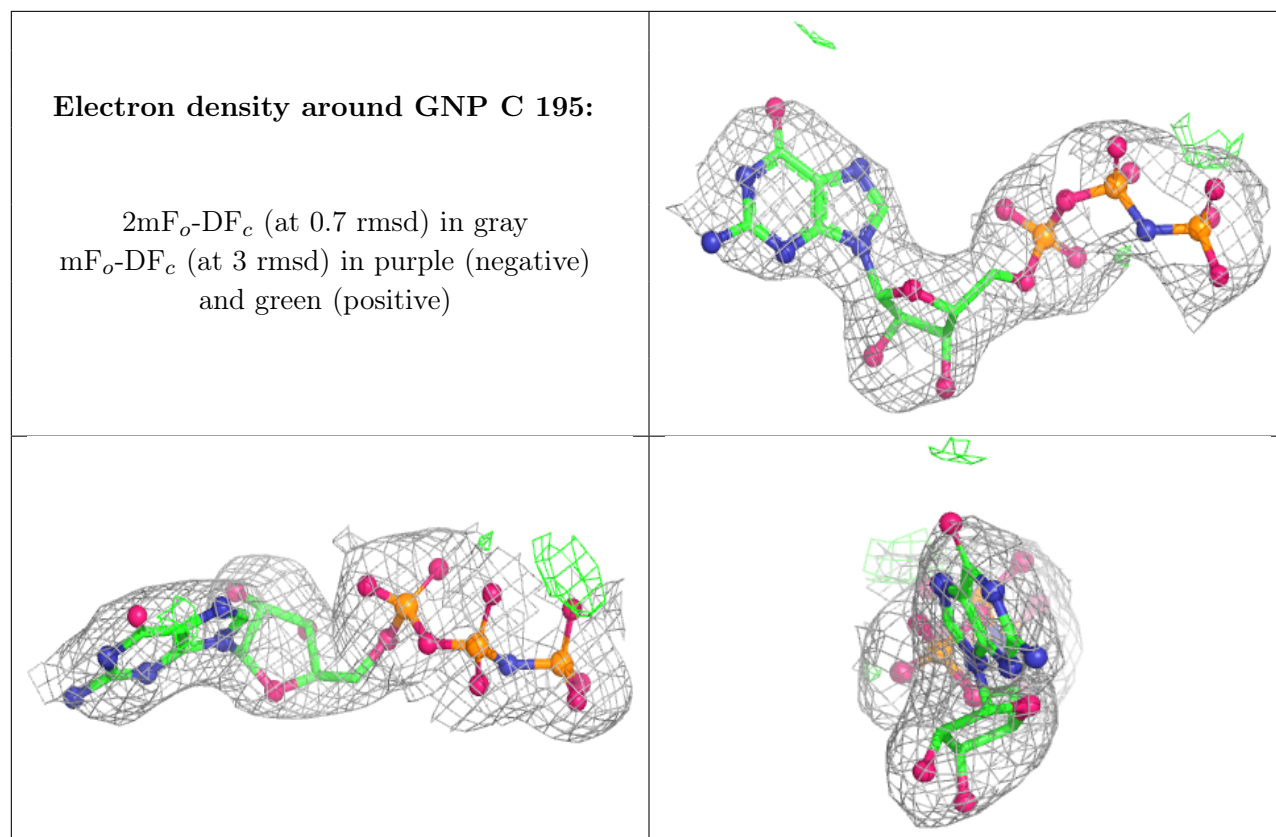
6.4 Ligands [i](#)

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
3	MG	B	5	1/1	0.90	0.52	66,66,66,66	0
3	MG	C	194	1/1	0.92	0.12	56,56,56,56	0
3	MG	A	194	1/1	0.94	0.11	31,31,31,31	0
4	GNP	A	195	32/32	0.96	0.07	42,54,56,57	0
4	GNP	C	195	32/32	0.97	0.06	49,57,61,61	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.