



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:17 PM UTC

PDB ID : 2GA9 / pdb_00002ga9
Title : Crystal Structure of the Heterodimeric Vaccinia Virus Polyadenylate Polymerase with Bound ATP-gamma-S
Authors : Moure, C.M.; Bowman, B.R.; Gershon, P.D.; Quiocho, F.A.
Deposited on : 2006-03-08
Resolution : 2.30 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

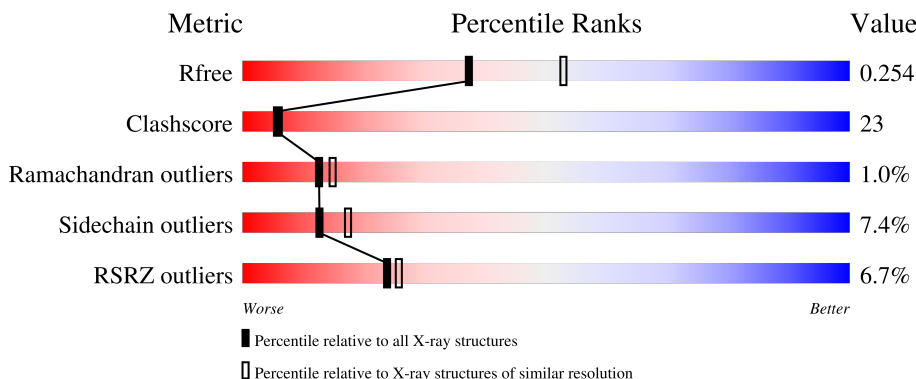
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	297	7% 70% 23% • •
2	D	469	6% 53% 34% 6% • 5%

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
4	AGS	D	483	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6333 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 Cap-specific mRNA (nucleoside-2'-O-)-methyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	288	Total	C	N	O	S	0	0	0
			2382	1552	394	424	12			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	140	ALA	ARG	engineered mutation	UNP P07617
A	142	ALA	LYS	engineered mutation	UNP P07617
A	143	ALA	ARG	engineered mutation	UNP P07617

- Molecule 2 is a protein called Poly(A) polymerase catalytic subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	445	Total	C	N	O	S	0	0	0
			3627	2318	605	680	24			

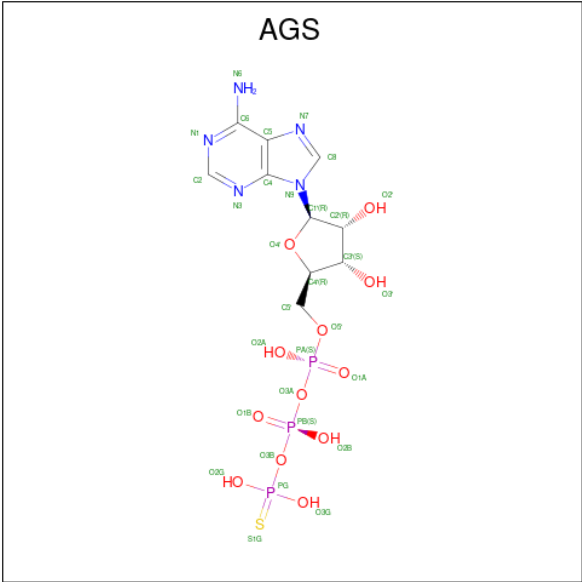
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	36	SER	LEU	variant	UNP P23371

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	2	Total	Ca	0	0
			2	2		

- Molecule 4 is PHOSPHOTHIOPHOSPHORIC ACID-ADENYLATE ESTER (CCD ID: AGS) (formula: C₁₀H₁₆N₅O₁₂P₃S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	D	1	Total	C	N	O	P	S	0	0
			31	10	5	12	3	1		
4	D	1	Total	C	N	O	P	S	0	0
			31	10	5	12	3	1		

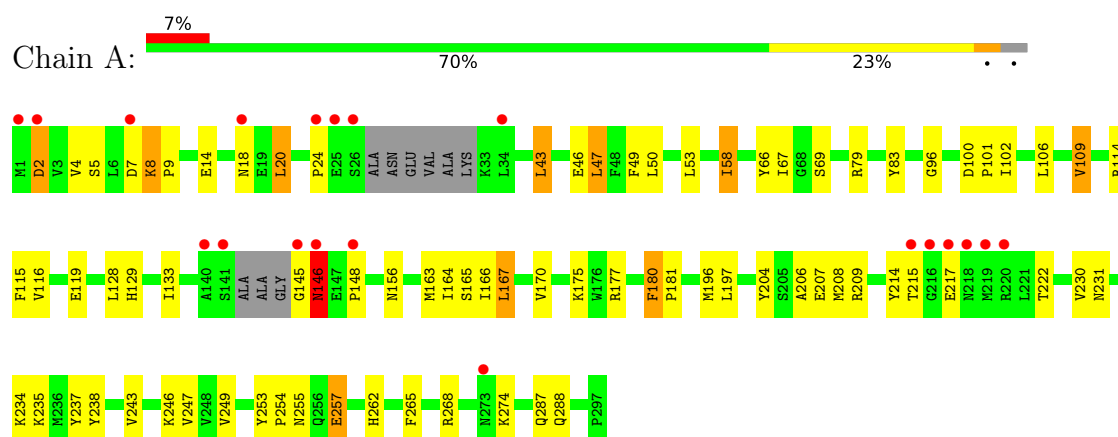
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	120	Total	O	0	0
			120	120		
5	D	140	Total	O	0	0
			140	140		

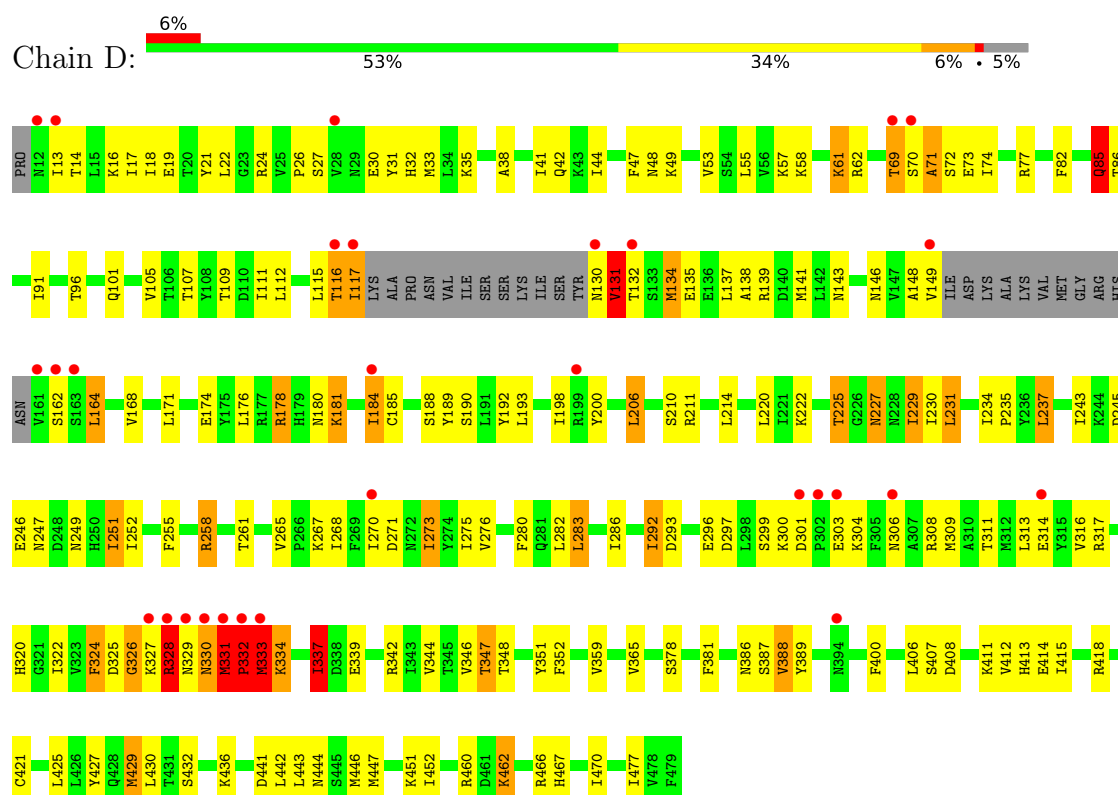
3 Residue-property plots [i](#)

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: Cap-specific mRNA (nucleoside-2'-O-)-methyltransferase



- Molecule 2: Poly(A) polymerase catalytic subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	70.05Å 91.69Å 133.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.30 50.00 – 2.31	Depositor EDS
% Data completeness (in resolution range)	93.8 (50.00-2.30) 94.6 (50.00-2.31)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.53 (at 2.31Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.222 , 0.260 0.213 , 0.254	Depositor DCC
R_{free} test set	1910 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	28.9	Xtriage
Anisotropy	0.100	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 42.7	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	6333	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.12% 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

5.1 Standard geometry

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

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	0.45	0/2444	0.96	9/3308 (0.3%)
2	D	0.48	1/3686 (0.0%)	1.02	20/4972 (0.4%)
All	All	0.47	1/6130 (0.0%)	1.00	29/8280 (0.4%)

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
1	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	333	MET	SD-CE	-6.01	1.64	1.79

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	131	VAL	N-CA-C	10.35	120.36	110.42
2	D	234	ILE	CA-C-N	9.87	129.51	119.24
2	D	234	ILE	C-N-CA	9.87	129.51	119.24
2	D	328	ARG	CA-C-N	8.59	135.38	120.87
2	D	328	ARG	C-N-CA	8.59	135.38	120.87
2	D	71	ALA	N-CA-C	-7.48	103.26	112.38
2	D	331	MET	CA-C-N	-7.47	110.50	119.84
2	D	331	MET	C-N-CA	-7.47	110.50	119.84
1	A	243	VAL	N-CA-C	7.40	118.13	110.36
1	A	180	PHE	CA-C-N	6.55	128.03	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	180	PHE	C-N-CA	6.55	128.03	119.84
1	A	133	ILE	N-CA-C	6.13	116.75	108.17
2	D	251	ILE	CB-CA-C	-5.99	105.75	111.44
1	A	249	VAL	N-CA-C	5.91	118.48	111.09
2	D	116	THR	N-CA-C	5.69	119.98	112.25
2	D	337	ILE	CB-CA-C	-5.66	103.46	111.21
1	A	247	VAL	N-CA-C	-5.37	100.65	108.17
2	D	225	THR	N-CA-C	5.35	117.19	111.36
1	A	146	ASN	CB-CA-C	-5.33	108.85	115.89
2	D	27	SER	N-CA-C	-5.32	103.86	110.41
2	D	206	LEU	N-CA-C	-5.21	100.91	109.40
2	D	328	ARG	O-C-N	5.14	129.42	122.59
2	D	85	GLN	N-CA-C	-5.13	105.77	111.36
2	D	324	PHE	N-CA-C	5.10	117.47	108.24
1	A	274	LYS	N-CA-C	-5.08	102.17	110.20
1	A	166	ILE	CB-CA-C	-5.07	105.39	112.04
2	D	339	GLU	N-CA-C	5.05	118.22	111.75
2	D	328	ARG	CA-C-O	5.04	127.71	120.51
2	D	210	SER	N-CA-C	5.02	117.15	111.02

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	204	TYR	Sidechain

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	2382	0	2389	55	0
2	D	3627	0	3689	232	0
3	D	2	0	0	0	0
4	D	62	0	24	9	0
5	A	120	0	0	11	0
5	D	140	0	0	20	0
All	All	6333	0	6102	285	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 23.

All (285) 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:13:ILE:HA	5:D:555:HOH:O	1.28	1.25
2:D:331:MET:HB3	2:D:332:PRO:HD2	1.33	1.09
2:D:325:ASP:OD2	2:D:327:LYS:HB3	1.54	1.05
2:D:283:LEU:HD22	2:D:429:MET:HE2	1.36	1.04
2:D:444:ASN:HA	2:D:447:MET:HE2	1.44	0.98
2:D:328:ARG:HG3	2:D:441:ASP:HA	1.50	0.94
1:A:181:PRO:HB3	5:A:375:HOH:O	1.69	0.93
2:D:309:MET:HE2	2:D:313:LEU:HG	1.49	0.92
2:D:331:MET:HE2	2:D:331:MET:HA	1.51	0.92
2:D:331:MET:O	2:D:332:PRO:C	2.07	0.92
2:D:304:LYS:HG3	2:D:308:ARG:HH21	1.34	0.91
2:D:413:HIS:HD2	2:D:415:ILE:H	1.17	0.90
2:D:48:ASN:N	4:D:483:AGS:H2	1.89	0.88
2:D:231:LEU:CD2	2:D:243:ILE:HG12	2.04	0.87
2:D:21:TYR:CZ	2:D:85:GLN:HG2	2.09	0.87
2:D:328:ARG:CG	2:D:441:ASP:HA	2.05	0.85
2:D:229:ILE:HD13	2:D:229:ILE:H	1.42	0.85
2:D:331:MET:HG3	2:D:442:LEU:HA	1.59	0.85
2:D:444:ASN:HD22	2:D:447:MET:HE1	1.42	0.85
2:D:328:ARG:CB	2:D:441:ASP:HA	2.05	0.84
2:D:333:MET:CE	2:D:348:THR:HG22	2.08	0.83
2:D:184:ILE:HD13	2:D:185:CYS:N	1.92	0.83
2:D:222:LYS:HE2	5:D:595:HOH:O	1.78	0.83
2:D:331:MET:HB3	2:D:332:PRO:CD	2.09	0.82
2:D:444:ASN:HD22	2:D:447:MET:CE	1.93	0.82
2:D:314:GLU:OE2	2:D:317:ARG:NH2	2.14	0.81
2:D:283:LEU:HD22	2:D:429:MET:CE	2.11	0.80
2:D:184:ILE:HD11	2:D:276:VAL:N	1.98	0.78
2:D:328:ARG:HB2	2:D:444:ASN:HB2	1.66	0.78
2:D:70:SER:HB2	2:D:73:GLU:HB2	1.67	0.77
2:D:329:ASN:O	2:D:331:MET:N	2.17	0.77
2:D:328:ARG:HG2	5:D:489:HOH:O	1.84	0.76
2:D:48:ASN:H	4:D:483:AGS:H2	1.50	0.76
2:D:328:ARG:HB3	2:D:441:ASP:HA	1.67	0.76
2:D:62:ARG:HD3	5:D:588:HOH:O	1.85	0.75
2:D:331:MET:HA	2:D:331:MET:CE	2.17	0.74
2:D:231:LEU:HD22	2:D:243:ILE:HG12	1.68	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:470:ILE:HG12	2:D:477:ILE:HD12	1.68	0.74
2:D:326:GLY:HA2	2:D:447:MET:HE3	1.69	0.74
2:D:328:ARG:HG3	2:D:441:ASP:OD1	1.86	0.74
2:D:328:ARG:CG	5:D:489:HOH:O	2.35	0.74
2:D:325:ASP:O	2:D:327:LYS:N	2.21	0.73
2:D:227:ASN:OD1	2:D:227:ASN:C	2.30	0.73
2:D:337:ILE:HD12	2:D:337:ILE:H	1.52	0.73
2:D:325:ASP:C	2:D:327:LYS:H	1.95	0.72
1:A:66:TYR:CD1	1:A:69:SER:HB3	2.24	0.72
2:D:130:ASN:HB3	2:D:314:GLU:OE2	1.90	0.72
2:D:270:ILE:O	2:D:273:ILE:HG23	1.91	0.71
2:D:328:ARG:HG3	2:D:441:ASP:CA	2.20	0.70
2:D:31:TYR:O	2:D:35:LYS:HB2	1.91	0.70
2:D:324:PHE:HA	2:D:444:ASN:HD21	1.57	0.69
2:D:378:SER:HB3	2:D:388:VAL:HG13	1.73	0.69
2:D:331:MET:HG3	2:D:442:LEU:CA	2.23	0.68
1:A:7:ASP:HB2	1:A:8:LYS:NZ	2.09	0.68
2:D:235:PRO:HG3	2:D:466:ARG:O	1.93	0.67
2:D:432:SER:HA	5:D:587:HOH:O	1.93	0.67
2:D:184:ILE:HD11	2:D:275:ILE:C	2.19	0.67
2:D:58:LYS:O	2:D:61:LYS:HG3	1.95	0.67
2:D:309:MET:HE2	2:D:313:LEU:CG	2.23	0.67
2:D:69:THR:HG22	2:D:73:GLU:OE1	1.95	0.66
2:D:184:ILE:HD12	2:D:276:VAL:HG23	1.78	0.66
1:A:255:ASN:OD1	1:A:257:GLU:HB2	1.96	0.66
2:D:19:GLU:HG3	2:D:24:ARG:O	1.96	0.65
2:D:85:GLN:H	2:D:85:GLN:NE2	1.95	0.65
2:D:143:ASN:O	2:D:451:LYS:HE2	1.96	0.65
2:D:306:ASN:HB2	5:D:622:HOH:O	1.97	0.65
2:D:333:MET:HE2	2:D:348:THR:HG22	1.77	0.64
2:D:280:PHE:HA	2:D:429:MET:CE	2.27	0.64
2:D:333:MET:HE3	2:D:348:THR:HA	1.77	0.64
2:D:231:LEU:HD21	2:D:243:ILE:HG12	1.78	0.64
2:D:317:ARG:HH11	2:D:317:ARG:HG2	1.61	0.64
2:D:44:ILE:HD13	2:D:101:GLN:HE21	1.63	0.63
2:D:282:LEU:HD11	2:D:309:MET:HE1	1.81	0.63
2:D:329:ASN:C	2:D:331:MET:H	2.06	0.63
2:D:109:THR:HG21	4:D:483:AGS:O2A	1.98	0.62
2:D:314:GLU:CD	2:D:317:ARG:NH2	2.57	0.62
2:D:280:PHE:HA	2:D:429:MET:HE1	1.82	0.62
1:A:8:LYS:HD2	1:A:8:LYS:O	1.98	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:LEU:HB3	1:A:58:ILE:HG12	1.81	0.62
2:D:258:ARG:HA	5:D:580:HOH:O	2.00	0.62
2:D:470:ILE:HG23	2:D:477:ILE:CD1	2.30	0.61
2:D:462:LYS:HD2	5:D:606:HOH:O	1.99	0.60
2:D:408:ASP:HB3	2:D:411:LYS:HD2	1.83	0.60
2:D:316:VAL:CG1	2:D:322:ILE:HD12	2.32	0.60
2:D:309:MET:CE	2:D:313:LEU:HG	2.30	0.59
2:D:297:ASP:HA	2:D:300:LYS:HE3	1.84	0.59
1:A:9:PRO:HG3	1:A:262:HIS:HA	1.84	0.58
1:A:2:ASP:OD1	1:A:246:LYS:HD3	2.04	0.58
2:D:189:TYR:O	2:D:193:LEU:HD23	2.04	0.58
2:D:229:ILE:H	2:D:229:ILE:CD1	2.15	0.58
2:D:245:ASP:OD1	2:D:249:ASN:HB2	2.03	0.58
2:D:325:ASP:C	2:D:327:LYS:N	2.62	0.58
2:D:328:ARG:HB3	2:D:441:ASP:CA	2.32	0.58
2:D:328:ARG:NH1	2:D:329:ASN:HD21	2.02	0.58
2:D:171:LEU:HD23	2:D:171:LEU:C	2.29	0.58
2:D:314:GLU:CD	2:D:317:ARG:HH21	2.12	0.57
1:A:230:VAL:CG2	5:A:375:HOH:O	2.52	0.57
2:D:14:THR:HG23	2:D:38:ALA:HB2	1.86	0.57
2:D:164:LEU:O	2:D:168:VAL:HG23	2.05	0.57
2:D:171:LEU:HD23	2:D:171:LEU:O	2.03	0.57
2:D:337:ILE:HG23	2:D:344:VAL:HG22	1.86	0.57
2:D:47:PHE:O	2:D:49:LYS:HE3	2.03	0.57
2:D:162:SER:OG	2:D:251:ILE:HD13	2.05	0.57
1:A:215:THR:HG23	2:D:365:VAL:HG22	1.85	0.57
2:D:331:MET:CB	2:D:332:PRO:CD	2.82	0.57
1:A:231:ASN:HD21	1:A:235:LYS:HE3	1.69	0.56
2:D:13:ILE:HD13	5:D:555:HOH:O	2.05	0.56
2:D:436:LYS:HB2	5:D:593:HOH:O	2.06	0.56
2:D:413:HIS:CD2	2:D:415:ILE:H	2.09	0.56
1:A:230:VAL:HG22	5:A:375:HOH:O	2.05	0.56
2:D:109:THR:HG21	4:D:483:AGS:PA	2.46	0.56
2:D:304:LYS:CG	2:D:308:ARG:HH21	2.11	0.56
2:D:132:THR:HA	2:D:135:GLU:HG3	1.87	0.56
2:D:148:ALA:O	2:D:149:VAL:HB	2.06	0.55
1:A:58:ILE:HD12	1:A:170:VAL:HG23	1.88	0.55
2:D:176:LEU:HD13	2:D:273:ILE:HD12	1.88	0.55
2:D:334:LYS:HB3	2:D:347:THR:OG1	2.06	0.55
2:D:337:ILE:HD11	2:D:412:VAL:HG23	1.89	0.55
2:D:282:LEU:CD1	2:D:309:MET:HE1	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:LYS:HD2	1:A:8:LYS:C	2.30	0.55
2:D:211:ARG:NH1	5:D:566:HOH:O	2.39	0.55
2:D:55:LEU:CD2	4:D:483:AGS:S1G	2.95	0.55
2:D:378:SER:HB3	2:D:388:VAL:CG1	2.35	0.55
1:A:96:GLY:HA3	1:A:115:PHE:CE1	2.41	0.54
2:D:348:THR:HB	2:D:352:PHE:HD1	1.72	0.54
2:D:325:ASP:O	2:D:444:ASN:ND2	2.39	0.54
1:A:7:ASP:HB2	1:A:8:LYS:HZ2	1.71	0.54
2:D:320:HIS:HB3	2:D:436:LYS:HD2	1.89	0.54
2:D:329:ASN:C	2:D:331:MET:N	2.66	0.54
2:D:359:VAL:CG1	2:D:400:PHE:HB2	2.38	0.54
2:D:18:ILE:HG22	2:D:26:PRO:HG3	1.89	0.54
2:D:303:GLU:HG2	5:D:579:HOH:O	2.08	0.54
1:A:175:LYS:HE3	1:A:207:GLU:OE2	2.08	0.54
2:D:460:ARG:HD2	5:D:512:HOH:O	2.08	0.53
2:D:85:GLN:H	2:D:85:GLN:CD	2.16	0.53
1:A:102:ILE:HG23	5:A:396:HOH:O	2.08	0.53
2:D:32:HIS:HD2	2:D:33:MET:HE2	1.73	0.53
2:D:116:THR:O	2:D:117:ILE:HG23	2.08	0.53
1:A:4:VAL:HG22	1:A:5:SER:N	2.24	0.52
1:A:14:GLU:HG2	1:A:14:GLU:O	2.08	0.52
2:D:229:ILE:HD13	2:D:229:ILE:N	2.19	0.52
1:A:146:ASN:OD1	1:A:146:ASN:N	2.42	0.52
1:A:253:TYR:OH	1:A:287:GLN:NE2	2.42	0.52
2:D:189:TYR:CE2	2:D:193:LEU:HD21	2.45	0.52
2:D:70:SER:HB2	2:D:73:GLU:H	1.75	0.52
2:D:333:MET:HE3	2:D:348:THR:HG22	1.91	0.52
1:A:43:LEU:C	1:A:43:LEU:HD12	2.34	0.52
2:D:337:ILE:HG13	2:D:412:VAL:CG2	2.39	0.52
2:D:57:LYS:HE2	2:D:71:ALA:HB1	1.92	0.52
2:D:227:ASN:ND2	2:D:246:GLU:HB2	2.25	0.52
1:A:79:ARG:HD3	1:A:79:ARG:C	2.35	0.52
2:D:131:VAL:CG1	2:D:311:THR:HA	2.39	0.52
2:D:146:ASN:HD21	2:D:299:SER:HA	1.75	0.52
2:D:184:ILE:HG22	2:D:206:LEU:HB2	1.92	0.52
2:D:225:THR:HG22	2:D:227:ASN:HB3	1.91	0.52
2:D:301:ASP:C	2:D:301:ASP:OD1	2.53	0.51
2:D:413:HIS:HE1	2:D:452:ILE:O	1.93	0.51
2:D:261:THR:O	2:D:265:VAL:HG23	2.10	0.51
2:D:327:LYS:O	2:D:328:ARG:HG2	2.11	0.51
2:D:330:ASN:C	2:D:331:MET:O	2.52	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:ASP:OD1	1:A:101:PRO:HD2	2.11	0.51
2:D:333:MET:HE1	2:D:352:PHE:HE1	1.76	0.50
2:D:134:MET:HE2	2:D:137:LEU:HD22	1.92	0.50
2:D:346:VAL:O	2:D:348:THR:HG23	2.12	0.50
2:D:134:MET:HE2	2:D:137:LEU:CD2	2.41	0.50
2:D:184:ILE:CD1	2:D:276:VAL:HG23	2.41	0.50
2:D:331:MET:HG3	2:D:442:LEU:N	2.26	0.50
2:D:30:GLU:HG2	2:D:91:ILE:HD11	1.93	0.49
1:A:7:ASP:HB2	1:A:8:LYS:HZ1	1.76	0.49
2:D:117:ILE:O	2:D:178:ARG:NH2	2.44	0.49
2:D:141:MET:HE3	2:D:446:MET:SD	2.52	0.49
2:D:193:LEU:HD12	2:D:271:ASP:OD1	2.13	0.49
2:D:49:LYS:O	2:D:53:VAL:HG23	2.12	0.49
1:A:145:GLY:C	1:A:146:ASN:OD1	2.56	0.48
1:A:215:THR:CG2	2:D:365:VAL:HG22	2.43	0.48
2:D:184:ILE:HD13	2:D:185:CYS:O	2.13	0.48
2:D:327:LYS:O	2:D:328:ARG:CG	2.62	0.48
2:D:271:ASP:O	2:D:273:ILE:HG22	2.14	0.48
2:D:206:LEU:HD23	2:D:255:PHE:HB2	1.95	0.48
2:D:381:PHE:CG	2:D:430:LEU:HD13	2.48	0.48
2:D:280:PHE:HA	2:D:429:MET:HE3	1.95	0.48
1:A:196:MET:HE2	1:A:209:ARG:HB2	1.95	0.48
2:D:188:SER:HB3	2:D:200:TYR:HB3	1.95	0.48
2:D:180:ASN:HB2	2:D:181:LYS:NZ	2.29	0.47
2:D:333:MET:O	2:D:334:LYS:C	2.57	0.47
2:D:180:ASN:HB2	2:D:181:LYS:HZ2	1.78	0.47
2:D:268:ILE:HG22	2:D:270:ILE:HG12	1.97	0.47
2:D:306:ASN:CB	5:D:622:HOH:O	2.60	0.47
2:D:22:LEU:O	2:D:86:THR:HA	2.15	0.47
2:D:55:LEU:HD22	4:D:483:AGS:S1G	2.54	0.47
2:D:138:ALA:HB3	2:D:306:ASN:OD1	2.15	0.47
2:D:184:ILE:HD12	2:D:276:VAL:CG2	2.44	0.47
1:A:156:ASN:ND2	5:A:397:HOH:O	2.45	0.46
2:D:316:VAL:HG13	2:D:322:ILE:HD12	1.97	0.46
2:D:96:THR:HG23	2:D:477:ILE:HG12	1.95	0.46
2:D:229:ILE:C	2:D:230:ILE:HD12	2.40	0.46
2:D:328:ARG:HG3	5:D:489:HOH:O	2.10	0.46
2:D:48:ASN:H	4:D:483:AGS:C2	2.26	0.46
5:A:409:HOH:O	2:D:258:ARG:HB3	2.15	0.46
2:D:16:LYS:HD2	5:D:555:HOH:O	2.15	0.46
2:D:184:ILE:CG2	2:D:206:LEU:HB2	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:328:ARG:HG3	2:D:441:ASP:CG	2.41	0.46
2:D:189:TYR:CZ	2:D:193:LEU:HD21	2.51	0.46
2:D:17:ILE:HD13	5:D:558:HOH:O	2.15	0.46
2:D:55:LEU:HD21	4:D:483:AGS:PG	2.56	0.46
2:D:70:SER:O	2:D:74:ILE:HG12	2.16	0.46
2:D:214:LEU:HD22	2:D:243:ILE:HG13	1.98	0.46
2:D:342:ARG:NH1	5:D:520:HOH:O	2.48	0.46
1:A:46:GLU:OE2	1:A:66:TYR:OH	2.25	0.45
2:D:48:ASN:O	4:D:483:AGS:N1	2.49	0.45
2:D:337:ILE:CD1	2:D:412:VAL:HG23	2.45	0.45
2:D:245:ASP:CG	2:D:249:ASN:HB2	2.42	0.45
2:D:414:GLU:O	2:D:418:ARG:HG3	2.16	0.45
2:D:190:SER:HB2	2:D:275:ILE:HB	1.98	0.45
2:D:252:ILE:HD12	5:D:547:HOH:O	2.16	0.45
2:D:53:VAL:HG12	2:D:57:LYS:HE3	1.98	0.45
2:D:82:PHE:C	2:D:85:GLN:HE22	2.24	0.45
2:D:230:ILE:HD12	2:D:230:ILE:N	2.33	0.44
2:D:444:ASN:ND2	2:D:447:MET:HE1	2.22	0.44
2:D:225:THR:CG2	2:D:227:ASN:HB3	2.47	0.44
2:D:13:ILE:O	2:D:17:ILE:HG12	2.18	0.44
2:D:237:LEU:HD12	2:D:237:LEU:HA	1.86	0.44
2:D:325:ASP:C	2:D:447:MET:HE1	2.43	0.44
2:D:443:LEU:HA	2:D:446:MET:HE3	2.00	0.44
2:D:21:TYR:CE1	2:D:85:GLN:HG2	2.50	0.44
2:D:192:TYR:HA	2:D:198:ILE:HB	1.99	0.44
2:D:328:ARG:HH11	2:D:329:ASN:HD21	1.65	0.44
2:D:444:ASN:ND2	2:D:447:MET:CE	2.73	0.44
1:A:18:ASN:HB3	1:A:238:TYR:CE2	2.53	0.43
2:D:283:LEU:HD21	2:D:425:LEU:HB3	1.99	0.43
1:A:177:ARG:HD2	5:A:352:HOH:O	2.17	0.43
2:D:317:ARG:HH11	2:D:317:ARG:CG	2.30	0.43
2:D:112:LEU:HD13	2:D:220:LEU:HD22	2.00	0.43
2:D:69:THR:HG21	2:D:77:ARG:HH22	1.84	0.43
2:D:317:ARG:HG2	2:D:317:ARG:NH1	2.30	0.43
2:D:328:ARG:HA	2:D:328:ARG:HD3	1.27	0.43
1:A:53:LEU:HD13	1:A:58:ILE:HD11	2.01	0.43
2:D:139:ARG:HG2	2:D:306:ASN:ND2	2.34	0.43
1:A:106:LEU:HB2	1:A:109:VAL:HG22	2.00	0.42
2:D:292:ILE:HG22	2:D:293:ASP:N	2.33	0.42
2:D:107:THR:O	2:D:111:ILE:HG12	2.20	0.42
1:A:164:ILE:HD12	1:A:214:TYR:OH	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:184:ILE:HD13	2:D:184:ILE:C	2.43	0.42
1:A:4:VAL:HG23	5:A:331:HOH:O	2.19	0.42
2:D:386:ASN:O	2:D:387:SER:HB2	2.20	0.42
2:D:331:MET:O	2:D:333:MET:N	2.49	0.42
2:D:330:ASN:O	2:D:331:MET:O	2.37	0.42
2:D:408:ASP:CB	2:D:411:LYS:HD2	2.48	0.42
1:A:43:LEU:O	1:A:47:LEU:HB2	2.19	0.42
1:A:235:LYS:HE3	5:A:399:HOH:O	2.20	0.42
2:D:443:LEU:HD23	2:D:446:MET:HE3	2.01	0.42
1:A:67:ILE:HD12	1:A:116:VAL:HG11	2.02	0.42
2:D:303:GLU:HA	2:D:303:GLU:OE1	2.19	0.42
1:A:253:TYR:CD2	1:A:254:PRO:HD2	2.55	0.42
2:D:21:TYR:CG	2:D:41:ILE:HG23	2.55	0.42
1:A:20:LEU:HB3	1:A:237:TYR:HD2	1.83	0.41
1:A:100:ASP:HA	1:A:101:PRO:HD3	1.93	0.41
1:A:206:ALA:O	1:A:208:MET:HE2	2.20	0.41
1:A:181:PRO:CB	5:A:375:HOH:O	2.47	0.41
2:D:292:ILE:O	2:D:296:GLU:HG3	2.19	0.41
2:D:348:THR:HB	2:D:352:PHE:CD1	2.53	0.41
1:A:43:LEU:HD12	1:A:43:LEU:O	2.20	0.41
2:D:17:ILE:HD12	2:D:42:GLN:HB2	2.03	0.41
2:D:47:PHE:CD1	2:D:105:VAL:HG13	2.56	0.41
1:A:49:PHE:CZ	1:A:53:LEU:HD11	2.55	0.41
1:A:180:PHE:HA	1:A:181:PRO:HD3	1.91	0.41
2:D:17:ILE:HD11	2:D:42:GLN:OE1	2.20	0.41
2:D:115:LEU:CD1	2:D:171:LEU:HD21	2.51	0.41
2:D:286:ILE:CD1	2:D:446:MET:HE1	2.51	0.41
1:A:230:VAL:CG1	1:A:234:LYS:HE2	2.50	0.41
1:A:265:PHE:O	1:A:268:ARG:NH1	2.54	0.41
2:D:148:ALA:O	2:D:149:VAL:CB	2.68	0.41
2:D:331:MET:O	2:D:332:PRO:O	2.37	0.41
2:D:389:TYR:CE2	2:D:427:TYR:HB2	2.56	0.41
1:A:129:HIS:HA	5:A:374:HOH:O	2.20	0.41
2:D:466:ARG:HG2	2:D:467:HIS:N	2.36	0.41
2:D:181:LYS:HE3	2:D:181:LYS:HA	2.03	0.40
1:A:163:MET:O	1:A:167:LEU:HB2	2.20	0.40
1:A:253:TYR:HE2	1:A:287:GLN:HE22	1.69	0.40
2:D:331:MET:SD	2:D:421:CYS:SG	3.07	0.40
2:D:331:MET:HE2	2:D:351:TYR:HE1	1.86	0.40
2:D:388:VAL:HG13	2:D:389:TYR:N	2.37	0.40
1:A:66:TYR:CG	1:A:69:SER:HB3	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:79:ARG:O	1:A:83:TYR:HB2	2.21	0.40
2:D:247:ASN:HB2	2:D:249:ASN:ND2	2.36	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	282/297 (95%)	270 (96%)	10 (4%)	2 (1%)	18	23
2	D	439/469 (94%)	413 (94%)	21 (5%)	5 (1%)	11	13
All	All	721/766 (94%)	683 (95%)	31 (4%)	7 (1%)	12	15

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	330	ASN
2	D	331	MET
2	D	332	PRO
2	D	326	GLY
2	D	334	LYS
1	A	2	ASP
1	A	24	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	266/270 (98%)	247 (93%)	19 (7%)	13	19
2	D	422/443 (95%)	390 (92%)	32 (8%)	12	16
All	All	688/713 (96%)	637 (93%)	51 (7%)	13	17

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	LYS
1	A	20	LEU
1	A	43	LEU
1	A	47	LEU
1	A	50	LEU
1	A	58	ILE
1	A	109	VAL
1	A	114	ARG
1	A	119	GLU
1	A	128	LEU
1	A	146	ASN
1	A	148	PRO
1	A	165	SER
1	A	167	LEU
1	A	197	LEU
1	A	217	GLU
1	A	222	THR
1	A	257	GLU
1	A	288	GLN
2	D	61	LYS
2	D	69	THR
2	D	72	SER
2	D	85	GLN
2	D	117	ILE
2	D	131	VAL
2	D	134	MET
2	D	164	LEU
2	D	174	GLU
2	D	178	ARG
2	D	181	LYS
2	D	184	ILE
2	D	227	ASN
2	D	229	ILE
2	D	231	LEU
2	D	237	LEU

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Mol	Chain	Res	Type
2	D	258	ARG
2	D	267	LYS
2	D	273	ILE
2	D	283	LEU
2	D	292	ILE
2	D	328	ARG
2	D	331	MET
2	D	332	PRO
2	D	333	MET
2	D	337	ILE
2	D	347	THR
2	D	388	VAL
2	D	406	LEU
2	D	407	SER
2	D	429	MET
2	D	462	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	GLN
1	A	84	ASN
1	A	156	ASN
1	A	160	GLN
1	A	183	GLN
1	A	192	HIS
1	A	194	ASN
1	A	231	ASN
1	A	287	GLN
1	A	288	GLN
2	D	32	HIS
2	D	59	ASN
2	D	85	GLN
2	D	87	GLN
2	D	101	GLN
2	D	167	ASN
2	D	179	HIS
2	D	197	ASN
2	D	228	ASN
2	D	239	ASN
2	D	247	ASN
2	D	329	ASN

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Mol	Chain	Res	Type
2	D	392	HIS
2	D	413	HIS
2	D	444	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 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	AGS	D	482	3	32,33,33	1.27	3 (9%)	45,52,52	2.20	14 (31%)
4	AGS	D	483	-	32,33,33	1.83	7 (21%)	45,52,52	1.95	8 (17%)

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	AGS	D	482	3	-	3/21/38/38	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	AGS	D	483	-	-	3/21/38/38	0/3/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	483	AGS	PA-O3A	4.82	1.64	1.59
4	D	483	AGS	PB-O3A	4.24	1.64	1.59
4	D	483	AGS	PB-O3B	4.19	1.64	1.59
4	D	482	AGS	PB-O3B	3.19	1.62	1.59
4	D	483	AGS	C5-C4	2.81	1.44	1.39
4	D	483	AGS	PG-O2G	2.58	1.63	1.54
4	D	482	AGS	PG-S1G	-2.55	1.85	1.90
4	D	483	AGS	C5-N7	-2.46	1.34	1.39
4	D	482	AGS	C5-C4	2.22	1.43	1.39
4	D	483	AGS	PG-S1G	-2.07	1.86	1.90

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	483	AGS	O2G-PG-O3B	-8.74	75.46	104.64
4	D	482	AGS	N3-C2-N1	-6.03	119.45	128.58
4	D	482	AGS	C5-C4-N3	-4.60	120.39	126.72
4	D	482	AGS	PB-O3B-PG	-4.45	116.88	133.17
4	D	482	AGS	N3-C4-N9	4.28	134.44	127.17
4	D	482	AGS	C2-N3-C4	4.02	121.64	111.83
4	D	482	AGS	O2B-PB-O3B	3.67	117.20	107.27
4	D	483	AGS	N3-C2-N1	-3.58	123.17	128.58
4	D	483	AGS	O2A-PA-O3A	3.42	116.51	107.27
4	D	482	AGS	O2B-PB-O3A	3.33	116.27	107.27
4	D	482	AGS	C5-N7-C8	2.96	108.10	103.45
4	D	483	AGS	C5-C4-N3	-2.80	122.86	126.72
4	D	482	AGS	O3A-PB-O1B	-2.66	102.71	110.70
4	D	482	AGS	O2G-PG-O3B	2.61	113.37	104.64
4	D	482	AGS	C2-N1-C6	2.58	122.97	118.73
4	D	483	AGS	C2-N1-C6	2.55	122.91	118.73
4	D	482	AGS	PA-O5'-C5'	-2.36	107.83	121.35
4	D	483	AGS	C2-N3-C4	2.31	117.47	111.83
4	D	483	AGS	PB-O3B-PG	-2.12	125.39	133.17
4	D	483	AGS	C4-C5-N7	-2.06	108.23	110.58
4	D	482	AGS	C4-C5-N7	-2.05	108.24	110.58
4	D	482	AGS	O4'-C1'-N9	2.01	111.96	108.09

There are no chirality outliers.

All (6) torsion outliers are listed below:

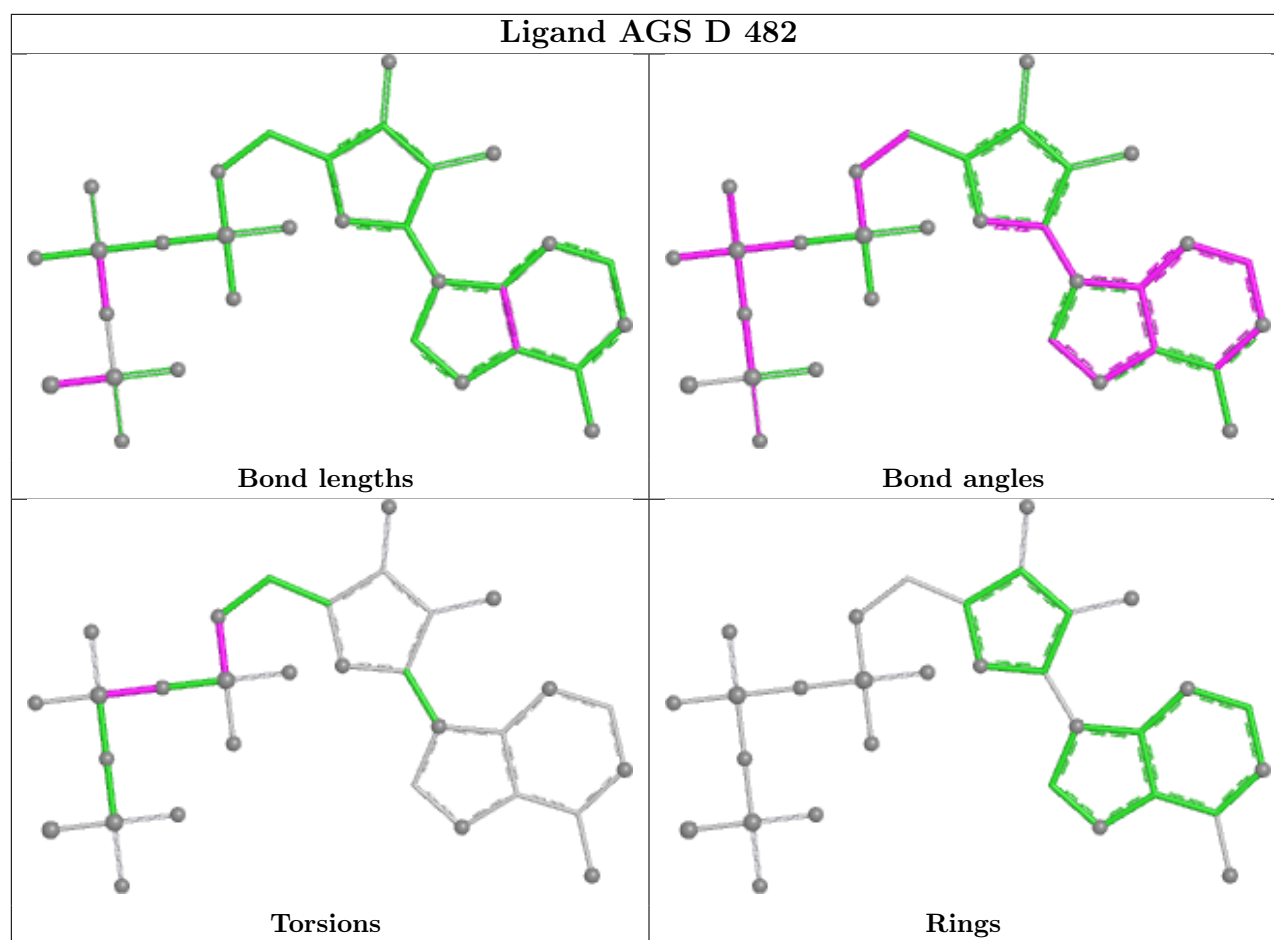
Mol	Chain	Res	Type	Atoms
4	D	482	AGS	C5'-O5'-PA-O2A
4	D	482	AGS	C5'-O5'-PA-O3A
4	D	482	AGS	PA-O3A-PB-O1B
4	D	483	AGS	O4'-C4'-C5'-O5'
4	D	483	AGS	C3'-C4'-C5'-O5'
4	D	483	AGS	PB-O3A-PA-O1A

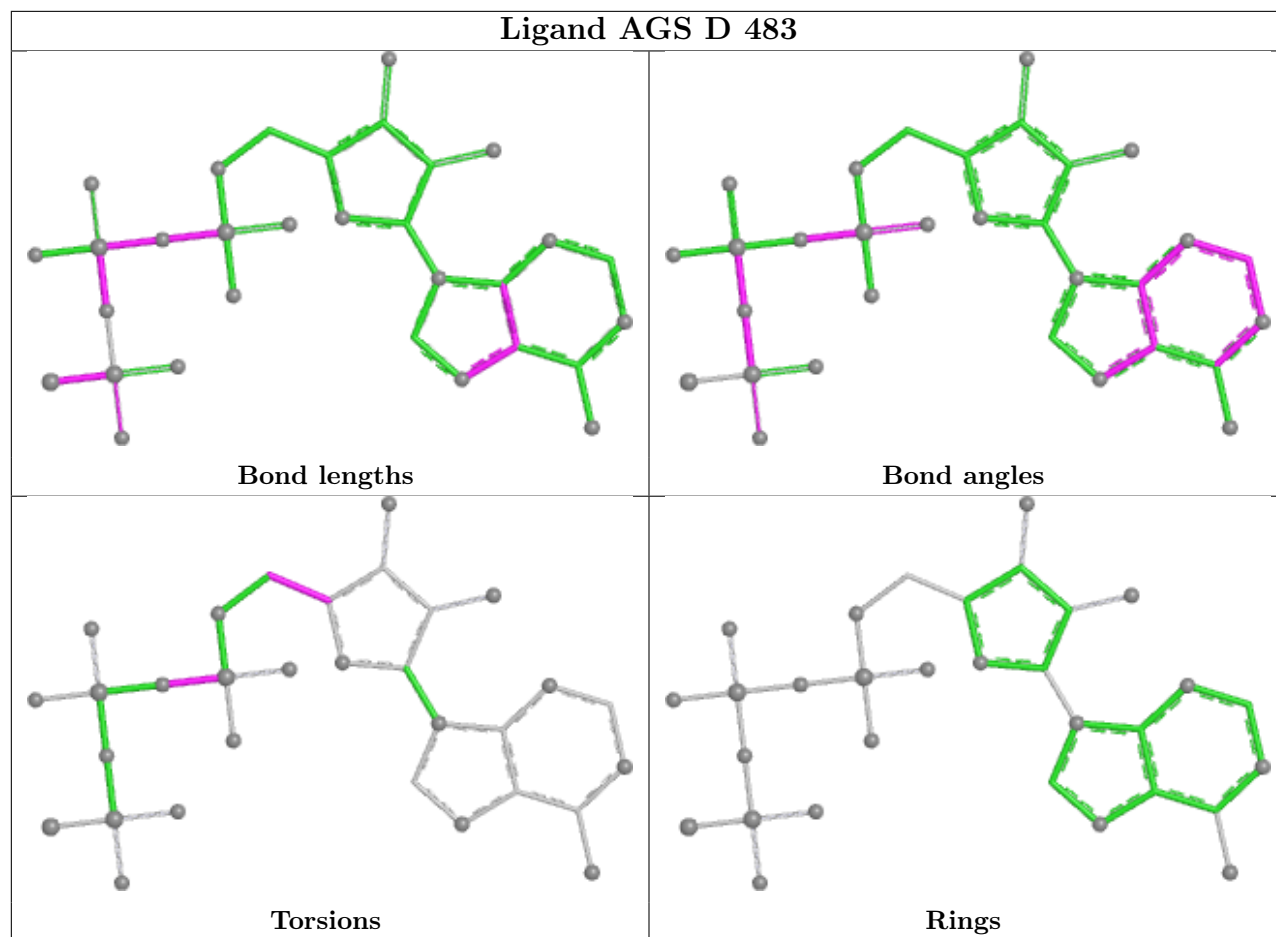
There are no ring outliers.

1 monomer is involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	483	AGS	9	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	288/297 (96%)	0.27	20 (6%) 23 25	15, 29, 56, 72	0
2	D	445/469 (94%)	0.32	29 (6%) 25 27	17, 33, 55, 66	0
All	All	733/766 (95%)	0.30	49 (6%) 24 26	15, 31, 56, 72	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	117	ILE	6.3
2	D	149	VAL	5.5
2	D	332	PRO	5.4
1	A	145	GLY	5.4
1	A	141	SER	5.2
2	D	328	ARG	5.1
1	A	217	GLU	4.6
2	D	331	MET	4.5
1	A	140	ALA	4.4
2	D	161	VAL	4.2
1	A	218	ASN	4.0
2	D	327	LYS	3.7
1	A	1	MET	3.6
1	A	24	PRO	3.5
1	A	220	ARG	3.4
2	D	303	GLU	3.4
2	D	12	ASN	3.3
2	D	329	ASN	3.3
1	A	2	ASP	2.9
2	D	28	VAL	2.9
1	A	146	ASN	2.9
1	A	148	PRO	2.8
2	D	314	GLU	2.8
1	A	219	MET	2.8

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Mol	Chain	Res	Type	RSRZ
2	D	163	SER	2.6
2	D	70	SER	2.6
2	D	330	ASN	2.5
1	A	216	GLY	2.5
1	A	34	LEU	2.5
2	D	270	ILE	2.5
2	D	394	ASN	2.4
2	D	306	ASN	2.4
2	D	301	ASP	2.3
1	A	215	THR	2.3
2	D	13	ILE	2.3
2	D	302	PRO	2.2
1	A	273	ASN	2.2
1	A	26	SER	2.2
2	D	116	THR	2.2
2	D	162	SER	2.2
1	A	18	ASN	2.1
2	D	69	THR	2.1
2	D	132	THR	2.1
2	D	184	ILE	2.1
2	D	333	MET	2.1
1	A	7	ASP	2.1
1	A	25	GLU	2.0
2	D	199	ARG	2.0
2	D	130	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](#)

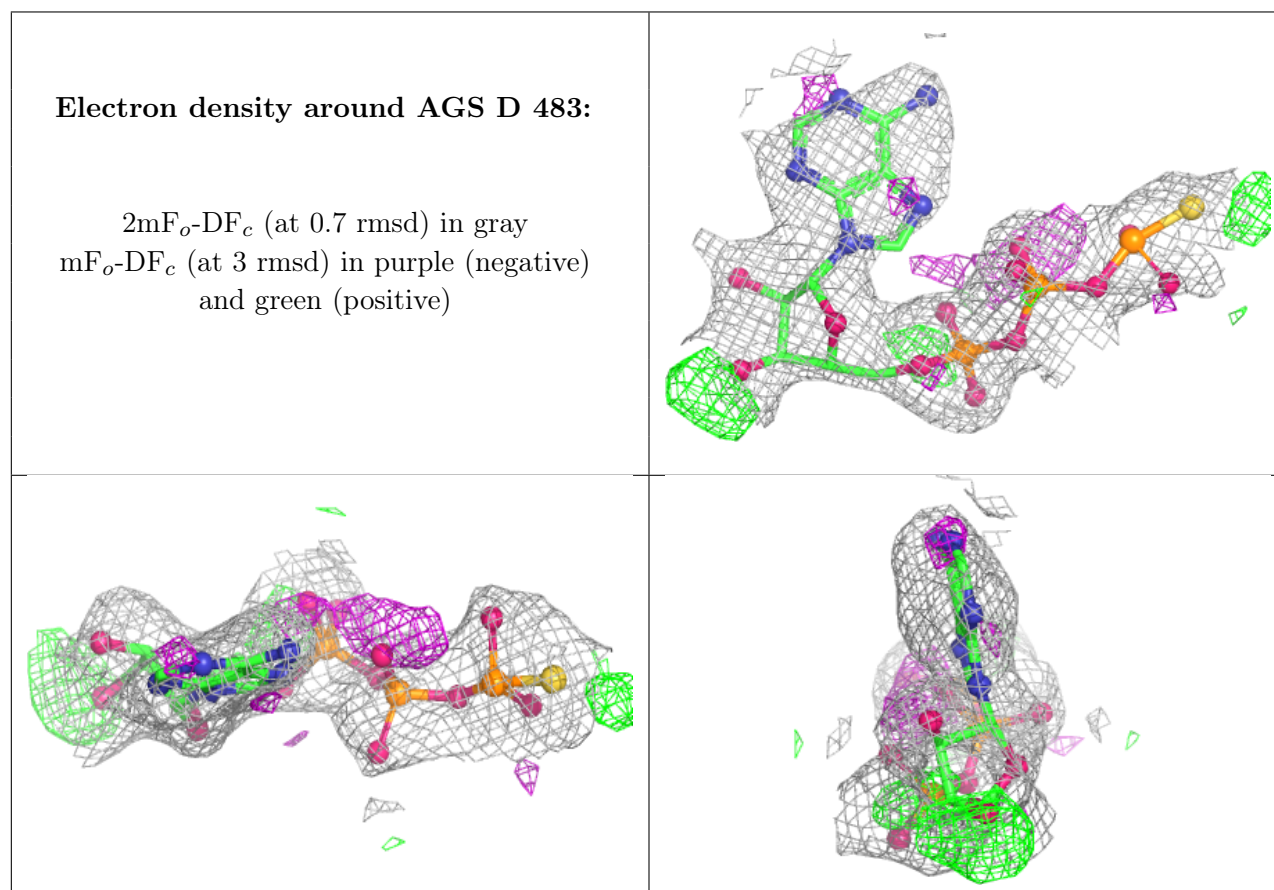
There are no oligosaccharides in this entry.

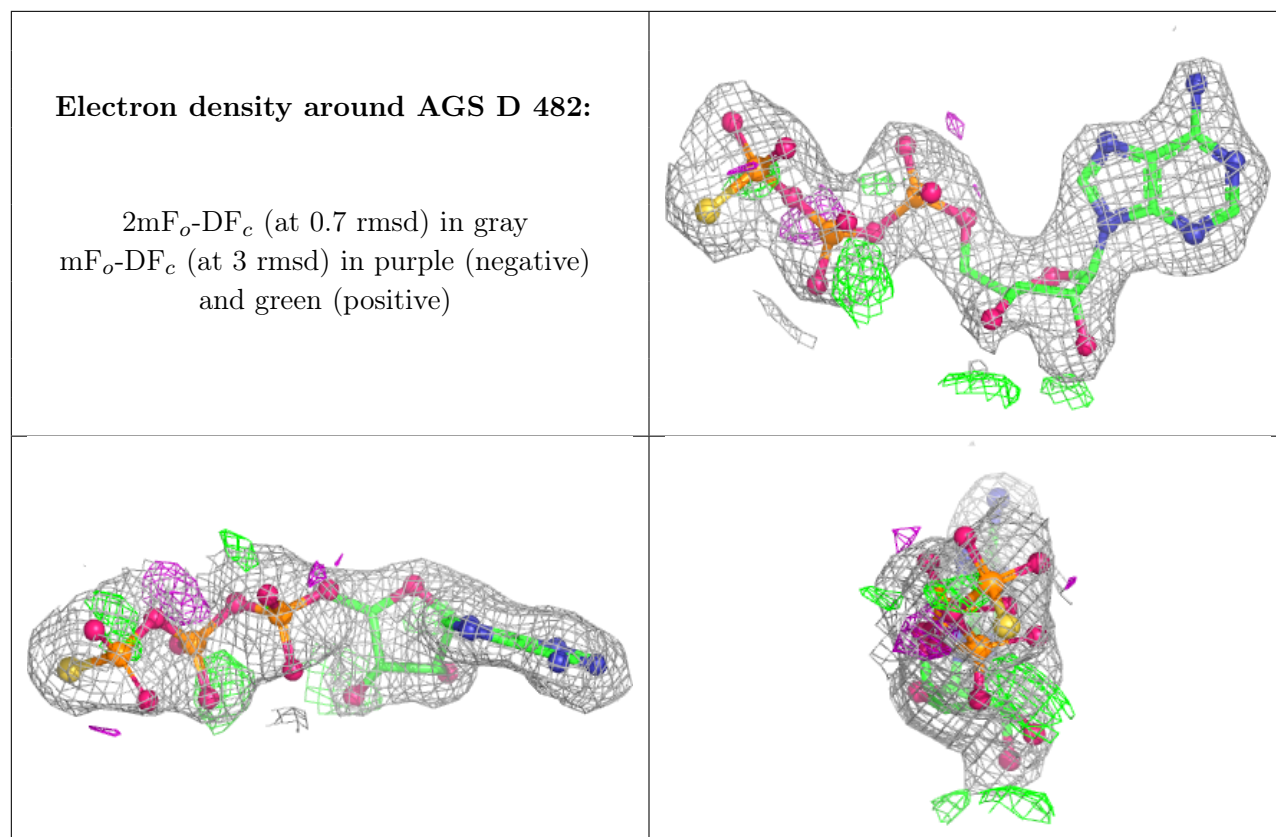
6.4 Ligands [i](#)

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	CA	D	481	1/1	0.68	0.18	71,71,71,71	0
4	AGS	D	483	31/31	0.71	0.17	55,63,83,85	0
4	AGS	D	482	31/31	0.89	0.12	38,46,57,58	0
3	CA	D	480	1/1	0.91	0.08	36,36,36,36	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.