



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:32 PM UTC

PDB ID : 2Z95 / pdb_00002z95
Title : Crystal Structure of GDP-D-Mannose Dehydratase from Aquifex aeolicus VF5
Authors : Niwa, H.; Kuramitsu, S.; Yokoyama, S.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2007-09-18
Resolution : 2.60 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

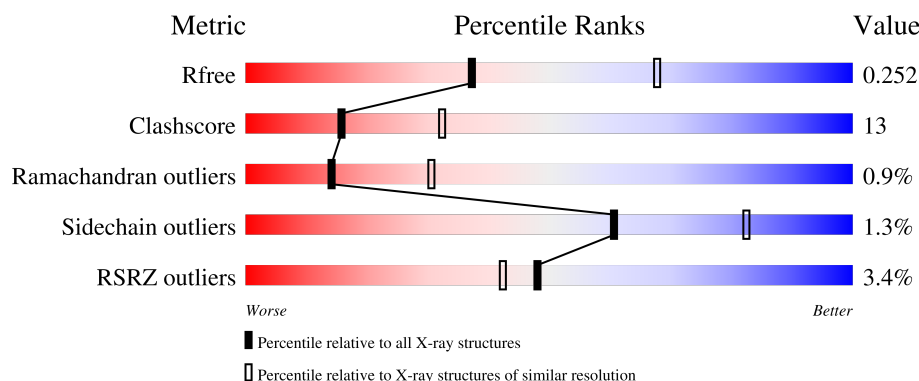
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	345	10% 68% 27% ..
1	B	345	65% 28% ..
1	C	345	68% 29% ..
1	D	345	10% 65% 30% ..

2 Entry composition [i](#)

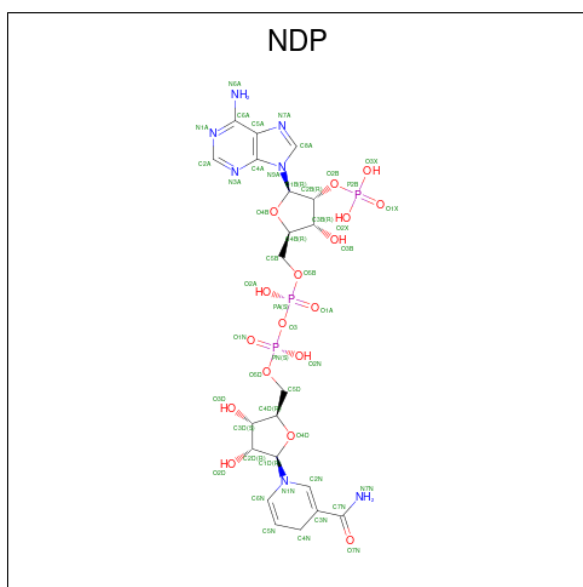
There are 4 unique types of molecules in this entry. The entry contains 10913 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 GDP-D-mannose dehydratase.

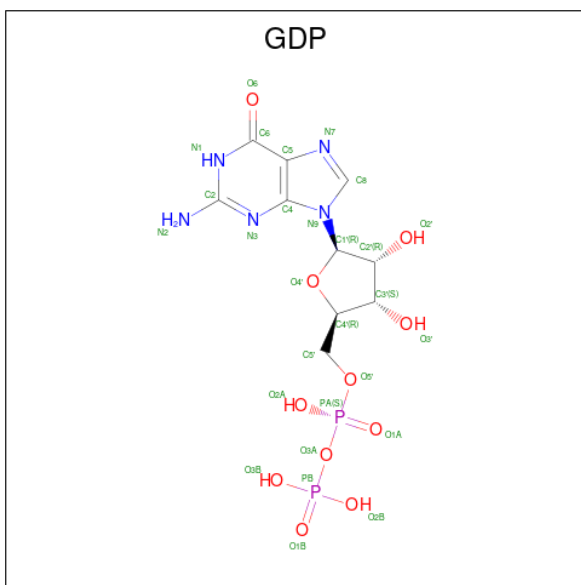
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	333	Total	C	N	O	S	0	0	0
			2662	1713	447	492	10			
1	B	332	Total	C	N	O	S	0	0	0
			2656	1711	444	491	10			
1	C	338	Total	C	N	O	S	0	0	0
			2675	1723	444	498	10			
1	D	332	Total	C	N	O	S	0	0	0
			2573	1652	429	482	10			

- Molecule 2 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is GUANOSINE-5'-DIPHOSPHATE (CCD ID: GDP) (formula: $C_{10}H_{15}N_5O_{11}P_2$).



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

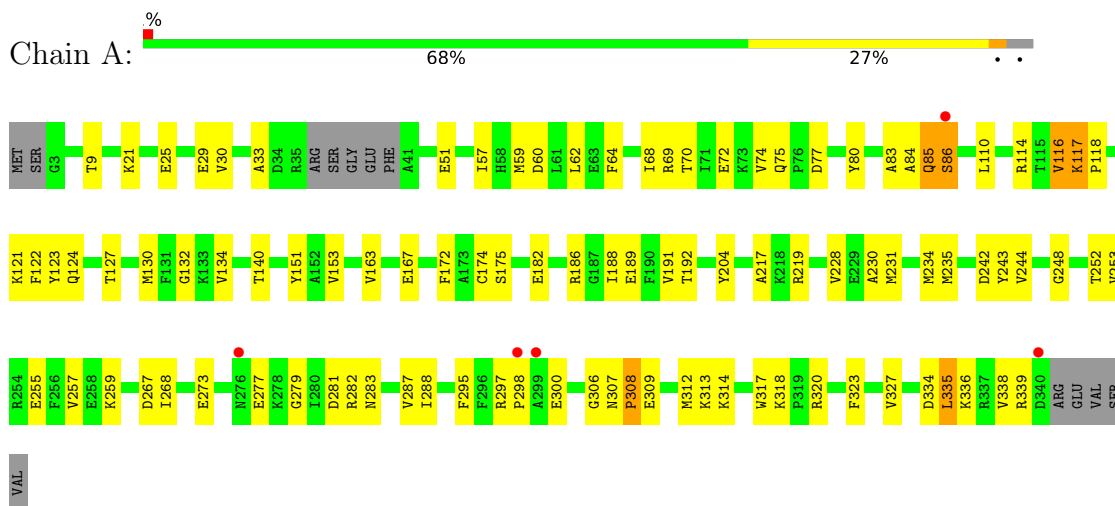
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	73	Total O 73 73	0	0
4	B	93	Total O 93 93	0	0
4	C	64	Total O 64 64	0	0
4	D	41	Total O 41 41	0	0

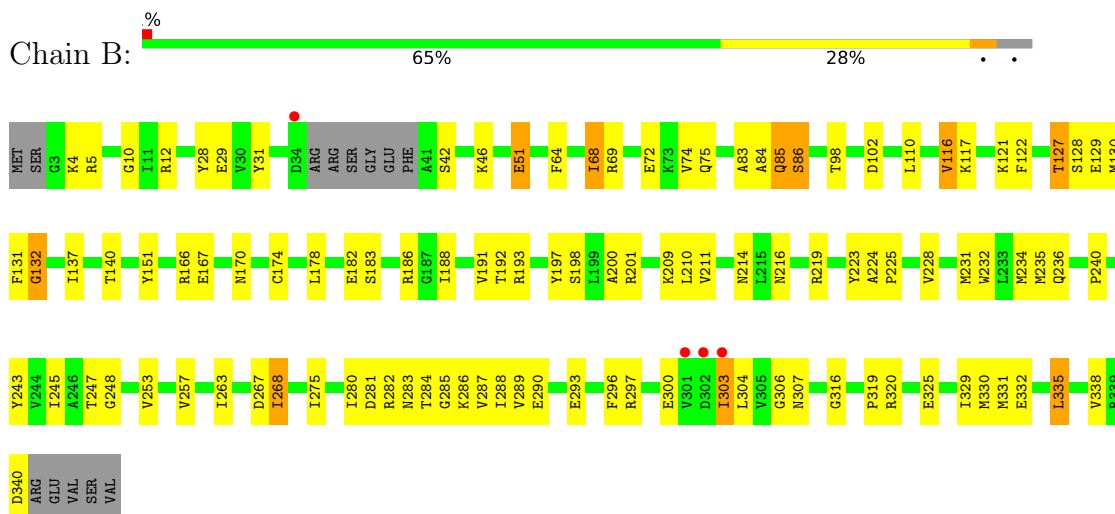
3 Residue-property plots

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: GDP-D-mannose dehydratase



• Molecule 1: GDP-D-mannose dehydratase

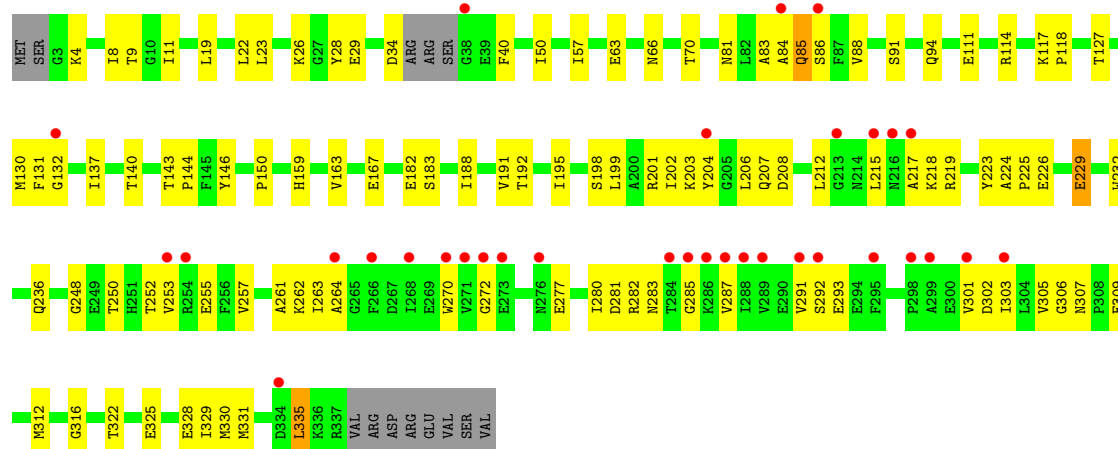


• Molecule 1: GDP-D-mannose dehydratase





● Molecule 1: GDP-D-mannose dehydratase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	80.48Å 101.18Å 186.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.16 – 2.60 49.16 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.2 (49.16-2.60) 99.3 (49.16-2.60)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.87 (at 2.61Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.200 , 0.253 0.200 , 0.252	Depositor DCC
R_{free} test set	2378 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	40.0	Xtriage
Anisotropy	0.558	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 59.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10913	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 23.09 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.9918e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: NDP, GDP

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.42	0/2719	0.94	13/3678 (0.4%)
1	B	0.40	0/2713	0.90	10/3669 (0.3%)
1	C	0.40	0/2734	0.89	5/3703 (0.1%)
1	D	0.39	0/2627	0.91	8/3559 (0.2%)
All	All	0.40	0/10793	0.91	36/14609 (0.2%)

There are no bond length outliers.

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	85	GLN	N-CA-C	-8.49	96.88	109.62
1	A	64	PHE	N-CA-C	7.68	120.78	111.40
1	B	64	PHE	N-CA-C	7.30	120.31	111.40
1	A	29	GLU	N-CA-C	-7.28	97.38	109.24
1	C	30	VAL	N-CA-C	7.25	118.32	107.80
1	B	29	GLU	N-CA-C	-6.94	97.42	108.73
1	A	314	LYS	N-CA-C	6.84	118.74	111.28
1	D	117	LYS	CA-C-N	6.63	126.88	119.32
1	D	117	LYS	C-N-CA	6.63	126.88	119.32
1	C	335	LEU	N-CA-C	-6.59	104.18	111.36
1	B	116	VAL	N-CA-C	6.37	119.33	112.96
1	A	117	LYS	CA-C-N	6.23	127.63	119.84
1	A	117	LYS	C-N-CA	6.23	127.63	119.84
1	C	29	GLU	N-CA-C	-6.10	98.79	108.73
1	A	85	GLN	N-CA-C	-6.09	95.11	107.41
1	D	29	GLU	N-CA-C	-6.01	99.40	108.96
1	A	116	VAL	N-CA-C	5.91	118.87	112.96
1	A	191	VAL	N-CA-C	5.74	116.47	110.62
1	B	117	LYS	CA-C-N	5.69	126.96	119.84
1	B	117	LYS	C-N-CA	5.69	126.96	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	335	LEU	N-CA-C	-5.62	105.06	111.07
1	A	335	LEU	N-CA-C	-5.60	105.32	111.82
1	A	30	VAL	N-CA-C	5.54	115.84	107.80
1	C	77	ASP	N-CA-C	-5.53	106.67	113.41
1	A	86	SER	N-CA-C	5.51	120.00	113.16
1	B	335	LEU	N-CA-C	-5.49	104.93	111.69
1	C	85	GLN	N-CA-C	-5.42	103.14	110.68
1	D	86	SER	N-CA-C	5.33	117.62	109.62
1	B	86	SER	N-CA-C	5.29	119.92	112.72
1	A	244	VAL	N-CA-C	-5.28	100.85	108.87
1	B	85	GLN	N-CA-C	-5.14	99.62	108.56
1	B	268	ILE	N-CA-C	5.10	116.81	109.51
1	B	316	GLY	N-CA-C	-5.05	107.31	115.08
1	D	50	ILE	N-CA-C	-5.04	108.36	113.20
1	A	77	ASP	N-CA-C	-5.02	106.32	112.90
1	D	229	GLU	N-CA-C	-5.02	105.89	111.36

There are no chirality outliers.

There are no planarity outliers.

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	2662	0	2654	64	0
1	B	2656	0	2654	75	0
1	C	2675	0	2634	68	0
1	D	2573	0	2494	81	0
2	A	48	0	26	1	0
3	D	28	0	12	0	0
4	A	73	0	0	0	0
4	B	93	0	0	4	0
4	C	64	0	0	2	0
4	D	41	0	0	0	0
All	All	10913	0	10474	280	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 13.

All (280) 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:A:281:ASP:HB2	1:A:288:ILE:HD11	1.52	0.91
1:D:182:GLU:HG3	1:D:192:THR:HG21	1.60	0.83
1:C:40:PHE:HD2	1:D:40:PHE:HB3	1.46	0.81
1:C:182:GLU:HG3	1:C:192:THR:HG21	1.64	0.79
1:D:195:ILE:HD13	1:D:257:VAL:HG22	1.62	0.79
1:A:72:GLU:HG2	1:A:116:VAL:HG11	1.65	0.79
1:B:209:LYS:HE3	1:B:287:VAL:O	1.85	0.77
1:B:42:SER:HB2	4:B:348:HOH:O	1.85	0.74
1:B:186:ARG:O	1:B:193:ARG:HG3	1.86	0.74
1:A:297:ARG:HB3	1:A:298:PRO:HD2	1.70	0.74
1:D:312:MET:HE3	1:D:312:MET:O	1.90	0.72
1:B:137:ILE:HD13	1:B:303:ILE:HD12	1.73	0.71
1:C:272:GLY:HA3	1:C:277:GLU:HA	1.73	0.70
1:D:57:ILE:HD12	1:D:70:THR:HG23	1.74	0.70
1:C:59:MET:HA	1:C:66:ASN:OD1	1.93	0.69
1:D:325:GLU:O	1:D:329:ILE:HG13	1.92	0.69
1:D:88:VAL:O	1:D:91:SER:HB3	1.92	0.68
1:D:307:ASN:OD1	1:D:309:GLU:HB2	1.94	0.67
1:B:166:ARG:HD2	1:B:240:PRO:HG2	1.77	0.67
1:D:137:ILE:HD13	1:D:303:ILE:HD13	1.78	0.66
1:A:234:MET:HE3	1:A:243:TYR:HD2	1.60	0.66
1:B:234:MET:HE3	1:B:243:TYR:HD2	1.61	0.65
1:D:199:LEU:HD13	1:D:261:ALA:HA	1.77	0.65
1:C:21:LYS:O	1:C:25:GLU:HG3	1.96	0.64
1:D:252:THR:OG1	1:D:255:GLU:HG3	1.97	0.64
1:D:114:ARG:O	1:D:118:PRO:HG3	1.98	0.64
1:B:183:SER:O	1:B:330:MET:HE3	1.97	0.64
1:B:248:GLY:HA2	1:B:306:GLY:O	1.98	0.63
1:B:281:ASP:HB2	1:B:288:ILE:HD11	1.81	0.63
1:B:127:THR:OG1	1:B:129:GLU:HG2	1.97	0.63
1:D:199:LEU:C	1:D:201:ARG:H	2.06	0.63
1:B:284:THR:O	1:B:286:LYS:HG3	1.99	0.63
1:C:325:GLU:O	1:C:329:ILE:HG13	1.98	0.63
1:A:163:VAL:O	1:A:167:GLU:HG3	1.99	0.62
1:B:86:SER:HA	1:B:151:TYR:CE1	2.34	0.62
1:D:232:TRP:O	1:D:236:GLN:HG2	1.98	0.62
1:D:291:VAL:HG12	1:D:292:SER:N	2.14	0.62
1:D:4:LYS:HD3	1:D:28:TYR:CZ	2.34	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:203:LYS:O	1:D:204:TYR:HD2	1.83	0.61
1:A:174:CYS:HB3	1:A:234:MET:HE2	1.84	0.60
1:C:247:THR:HA	1:C:319:PRO:HG2	1.82	0.60
1:C:248:GLY:HA2	1:C:306:GLY:O	2.02	0.60
1:C:224:ALA:HB3	1:C:225:PRO:HD3	1.84	0.60
1:C:46:LYS:HG2	1:C:51:GLU:OE1	2.02	0.60
1:B:72:GLU:HG3	1:B:116:VAL:HG11	1.84	0.59
1:D:263:ILE:HG22	1:D:263:ILE:O	2.03	0.59
1:B:186:ARG:HB3	1:B:193:ARG:HG2	1.85	0.58
1:D:331:MET:HE3	1:D:331:MET:O	2.04	0.58
1:D:212:LEU:HB2	1:D:215:LEU:HD21	1.86	0.57
1:C:232:TRP:O	1:C:236:GLN:HG2	2.05	0.57
1:D:248:GLY:HA2	1:D:306:GLY:O	2.04	0.57
1:D:291:VAL:HG12	1:D:292:SER:H	1.69	0.57
1:D:9:THR:OG1	1:D:81:ASN:HA	2.05	0.57
1:A:308:PRO:HG2	1:A:312:MET:HE2	1.86	0.56
1:B:281:ASP:HB2	1:B:288:ILE:CD1	2.36	0.56
1:C:234:MET:HG2	1:C:243:TYR:CD2	2.40	0.56
1:A:273:GLU:HA	1:A:277:GLU:HG3	1.87	0.55
1:B:219:ARG:HD3	4:B:391:HOH:O	2.05	0.55
1:C:254:ARG:HH21	1:C:277:GLU:CD	2.14	0.55
1:D:163:VAL:O	1:D:167:GLU:HG3	2.06	0.55
1:C:202:ILE:HG21	1:C:288:ILE:HD12	1.87	0.55
1:D:206:LEU:O	1:D:207:GLN:HB2	2.07	0.55
1:A:127:THR:O	1:A:130:MET:HG3	2.07	0.55
1:D:8:ILE:HD13	1:D:19:LEU:HD23	1.88	0.55
1:C:4:LYS:HE3	1:C:28:TYR:CZ	2.42	0.55
1:D:218:LYS:HG2	1:D:252:THR:HG22	1.89	0.54
1:A:287:VAL:HG23	1:A:287:VAL:O	2.05	0.54
1:D:203:LYS:HD2	1:D:264:ALA:HB1	1.89	0.54
1:A:80:TYR:CE1	1:A:235:MET:HG3	2.42	0.54
1:D:281:ASP:C	1:D:283:ASN:H	2.16	0.54
1:C:9:THR:OG1	1:C:81:ASN:HA	2.07	0.54
1:D:293:GLU:H	1:D:293:GLU:CD	2.15	0.54
1:B:182:GLU:HG3	1:B:192:THR:HG21	1.89	0.54
1:A:121:LYS:HG2	1:A:172:PHE:CD2	2.43	0.54
1:B:51:GLU:CD	1:B:51:GLU:H	2.15	0.54
1:B:232:TRP:O	1:B:236:GLN:HG2	2.08	0.54
1:D:226:GLU:O	1:D:229:GLU:HB3	2.07	0.54
1:B:268:ILE:HA	1:B:280:ILE:O	2.07	0.54
1:D:83:ALA:O	1:D:84:ALA:HB2	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:263:ILE:HB	1:D:331:MET:HG2	1.89	0.54
1:A:234:MET:HG2	1:A:243:TYR:CD2	2.43	0.54
1:B:46:LYS:HA	1:B:51:GLU:OE1	2.08	0.54
1:B:296:PHE:O	1:B:297:ARG:HD3	2.08	0.53
1:A:182:GLU:HG3	1:A:192:THR:HG21	1.88	0.53
1:A:9:THR:HA	1:A:33:ALA:HB3	1.91	0.53
1:B:247:THR:HA	1:B:319:PRO:HG2	1.89	0.53
1:D:91:SER:OG	1:D:150:PRO:HB2	2.08	0.53
1:A:281:ASP:C	1:A:283:ASN:H	2.16	0.53
1:B:110:LEU:HD21	1:B:122:PHE:CD2	2.43	0.53
1:B:166:ARG:O	1:B:170:ASN:HA	2.08	0.53
1:D:322:THR:OG1	1:D:325:GLU:HG3	2.09	0.52
1:D:272:GLY:O	1:D:277:GLU:HA	2.10	0.52
1:D:22:LEU:HD11	1:D:26:LYS:HE2	1.92	0.52
1:D:224:ALA:HB3	1:D:225:PRO:HD3	1.91	0.52
1:A:281:ASP:HB2	1:A:288:ILE:CD1	2.34	0.52
1:C:130:MET:HE3	1:C:145:PHE:CE1	2.44	0.52
1:C:203:LYS:HD2	1:C:264:ALA:O	2.10	0.52
1:C:209:LYS:HE3	1:C:287:VAL:HB	1.92	0.52
1:C:300:GLU:HB3	4:C:403:HOH:O	2.09	0.52
1:A:318:LYS:O	1:A:320:ARG:NH1	2.44	0.51
1:D:280:ILE:HA	1:D:287:VAL:HA	1.91	0.51
1:D:331:MET:O	1:D:335:LEU:HG	2.11	0.51
1:C:57:ILE:HD12	1:C:70:THR:HG23	1.92	0.51
1:B:331:MET:HE2	1:B:335:LEU:HG	1.94	0.50
1:A:267:ASP:HB2	1:A:283:ASN:ND2	2.27	0.50
1:C:172:PHE:CD1	1:C:240:PRO:HD3	2.46	0.50
1:C:312:MET:HE3	1:C:312:MET:O	2.11	0.50
1:B:281:ASP:O	1:B:285:GLY:N	2.44	0.50
1:B:12:ARG:HG2	1:B:42:SER:HB3	1.94	0.50
1:D:203:LYS:HD2	1:D:264:ALA:CB	2.42	0.50
1:B:191:VAL:HG23	4:B:430:HOH:O	2.11	0.49
1:B:178:LEU:HG	1:B:245:ILE:HB	1.93	0.49
1:D:11:ILE:HG22	1:D:34:ASP:HB2	1.93	0.49
1:A:336:LYS:HA	1:A:339:ARG:NH1	2.26	0.49
1:A:110:LEU:HD21	1:A:122:PHE:CD2	2.48	0.49
1:C:40:PHE:CD2	1:D:40:PHE:HB3	2.36	0.49
1:B:253:VAL:O	1:B:257:VAL:HG23	2.13	0.49
1:A:186:ARG:NH1	2:A:1001:NDP:H72N	2.11	0.49
1:A:323:PHE:O	1:A:327:VAL:HG23	2.13	0.49
1:B:83:ALA:O	1:B:84:ALA:HB2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:198:SER:O	1:D:201:ARG:HB2	2.12	0.48
1:B:129:GLU:HG3	1:B:151:TYR:HE2	1.79	0.48
1:A:83:ALA:O	1:A:84:ALA:HB2	2.13	0.48
1:B:178:LEU:HD13	4:B:378:HOH:O	2.12	0.48
1:B:228:VAL:O	1:B:231:MET:HB2	2.12	0.48
1:C:312:MET:O	1:C:316:GLY:HA2	2.13	0.48
1:B:325:GLU:O	1:B:329:ILE:HG13	2.13	0.48
1:C:97:LEU:O	1:C:101:VAL:HG23	2.14	0.48
1:D:192:THR:HB	1:D:330:MET:HE1	1.95	0.48
1:A:123:TYR:CZ	1:A:231:MET:HG2	2.49	0.48
1:A:255:GLU:O	1:A:259:LYS:HG3	2.14	0.48
1:B:140:THR:HG22	1:B:307:ASN:HB3	1.96	0.48
1:D:199:LEU:C	1:D:201:ARG:N	2.70	0.48
1:A:189:GLU:H	1:A:189:GLU:CD	2.22	0.47
1:C:86:SER:HA	1:C:151:TYR:CE1	2.49	0.47
1:C:128:SER:HB3	1:C:179:PHE:CD2	2.49	0.47
1:A:234:MET:HE3	1:A:243:TYR:CD2	2.45	0.47
1:C:131:PHE:O	1:C:132:GLY:C	2.57	0.47
1:D:283:ASN:C	1:D:285:GLY:H	2.22	0.47
1:D:331:MET:HE2	1:D:335:LEU:HG	1.95	0.47
1:C:186:ARG:O	1:C:193:ARG:HG3	2.14	0.47
1:A:140:THR:HG22	1:A:307:ASN:HB3	1.95	0.47
1:A:69:ARG:NH2	1:D:111:GLU:HG3	2.30	0.47
1:B:263:ILE:HB	1:B:331:MET:HG2	1.95	0.47
1:C:281:ASP:OD2	1:C:283:ASN:HB3	2.14	0.47
1:D:132:GLY:O	1:D:301:VAL:HG13	2.14	0.47
1:D:253:VAL:O	1:D:257:VAL:HG23	2.15	0.47
1:D:312:MET:HE3	1:D:316:GLY:HA2	1.96	0.47
1:A:248:GLY:HA2	1:A:306:GLY:O	2.15	0.47
1:C:123:TYR:CZ	1:C:231:MET:HG2	2.49	0.47
1:A:153:VAL:HG12	1:C:157:PHE:CD1	2.50	0.47
1:A:217:ALA:O	1:A:252:THR:HA	2.15	0.47
1:A:334:ASP:O	1:A:338:VAL:HG23	2.15	0.47
1:D:140:THR:HG22	1:D:307:ASN:HB3	1.96	0.47
1:D:223:TYR:CD2	1:D:225:PRO:HD2	2.50	0.47
1:A:252:THR:OG1	1:A:255:GLU:HG3	2.15	0.47
1:B:128:SER:HB2	1:B:304:LEU:HD22	1.97	0.47
1:C:186:ARG:HB3	1:C:193:ARG:HG2	1.96	0.47
1:C:23:LEU:O	1:C:28:TYR:HB2	2.15	0.46
1:B:84:ALA:HB1	1:B:102:ASP:OD1	2.14	0.46
1:C:203:LYS:HD3	1:C:204:TYR:CE2	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:200:ALA:HB1	1:C:338:VAL:HG21	1.97	0.46
1:D:218:LYS:HB3	1:D:250:THR:OG1	2.15	0.46
1:A:219:ARG:NH2	1:A:300:GLU:OE1	2.49	0.46
1:C:259:LYS:O	1:C:263:ILE:HG12	2.15	0.46
1:C:221:TRP:O	1:C:247:THR:HG21	2.15	0.46
1:C:323:PHE:O	1:C:327:VAL:HG23	2.15	0.46
1:A:188:ILE:HG22	1:A:189:GLU:OE2	2.16	0.46
1:D:4:LYS:HB3	1:D:28:TYR:CD1	2.51	0.46
1:A:72:GLU:HG2	1:A:116:VAL:CG1	2.40	0.46
1:A:230:ALA:O	1:A:234:MET:HG3	2.16	0.46
1:B:200:ALA:CB	1:B:338:VAL:HG21	2.46	0.46
1:B:198:SER:CB	1:B:210:LEU:HD13	2.46	0.45
1:A:335:LEU:O	1:A:339:ARG:HG3	2.16	0.45
1:A:21:LYS:O	1:A:25:GLU:HG3	2.15	0.45
1:D:218:LYS:HG2	1:D:252:THR:CG2	2.46	0.45
1:B:131:PHE:O	1:B:132:GLY:C	2.59	0.45
1:C:39:GLU:HB3	1:D:188:ILE:HG12	1.99	0.45
1:A:267:ASP:HB2	1:A:283:ASN:HD21	1.82	0.45
1:C:39:GLU:O	1:C:40:PHE:HD1	1.99	0.45
1:C:203:LYS:HD3	1:C:204:TYR:HE2	1.82	0.45
1:C:338:VAL:C	1:C:340:ASP:H	2.24	0.45
1:A:312:MET:HA	1:A:317:TRP:O	2.17	0.45
1:B:174:CYS:HB3	1:B:234:MET:HE2	1.98	0.45
1:B:197:TYR:CZ	1:B:201:ARG:HD2	2.52	0.45
1:C:214:ASN:OD1	1:C:216:ASN:HB2	2.17	0.45
1:D:143:THR:O	1:D:144:PRO:C	2.59	0.45
1:A:51:GLU:OE1	1:A:51:GLU:N	2.49	0.44
1:B:188:ILE:HD12	1:B:197:TYR:CD2	2.52	0.44
1:C:193:ARG:HG3	1:C:193:ARG:HH11	1.82	0.44
1:D:191:VAL:HG13	1:D:192:THR:N	2.31	0.44
1:B:127:THR:O	1:B:130:MET:HG3	2.17	0.44
1:B:234:MET:HG2	1:B:243:TYR:CD2	2.52	0.44
1:C:263:ILE:HB	1:C:331:MET:HG2	2.00	0.44
1:B:72:GLU:CG	1:B:116:VAL:HG11	2.47	0.44
1:A:114:ARG:O	1:A:118:PRO:HG3	2.17	0.44
1:C:247:THR:HA	1:C:319:PRO:CG	2.46	0.44
1:B:98:THR:HG23	1:B:102:ASP:OD2	2.18	0.44
1:A:57:ILE:HD12	1:A:70:THR:HG23	1.98	0.44
1:C:198:SER:O	1:C:202:ILE:HG13	2.17	0.44
1:D:203:LYS:O	1:D:204:TYR:CD2	2.67	0.44
1:A:60:ASP:OD1	1:A:62:LEU:HB2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:18:TYR:CZ	1:C:185:LEU:HD12	2.53	0.44
1:C:178:LEU:HD12	1:C:178:LEU:N	2.33	0.44
1:D:131:PHE:O	1:D:132:GLY:C	2.61	0.44
1:A:86:SER:HA	1:A:151:TYR:CE1	2.53	0.43
1:B:10:GLY:N	1:B:83:ALA:HB2	2.33	0.43
1:B:51:GLU:CD	1:B:51:GLU:N	2.76	0.43
1:B:214:ASN:OD1	1:B:216:ASN:HB2	2.19	0.43
1:C:189:GLU:CD	1:C:189:GLU:H	2.26	0.43
1:D:305:VAL:HG13	1:D:305:VAL:O	2.19	0.43
1:C:275:ILE:O	1:C:275:ILE:HG12	2.18	0.43
1:A:69:ARG:HH21	1:D:111:GLU:HG3	1.84	0.43
1:B:5:ARG:CG	1:B:31:TYR:HE2	2.32	0.43
1:B:223:TYR:CD1	1:B:225:PRO:HD2	2.53	0.43
1:B:74:VAL:O	1:B:75:GLN:C	2.61	0.43
1:A:57:ILE:CD1	1:A:74:VAL:HG21	2.48	0.43
1:A:124:GLN:HA	1:A:124:GLN:OE1	2.19	0.43
1:C:334:ASP:O	1:C:338:VAL:HG23	2.18	0.43
1:D:4:LYS:HD3	1:D:28:TYR:CE1	2.54	0.43
1:C:11:ILE:HG23	1:C:12:ARG:HG3	2.00	0.43
1:C:63:GLU:OE2	1:D:94:GLN:HA	2.19	0.43
1:C:211:VAL:CG1	1:C:292:SER:HB3	2.49	0.43
1:D:219:ARG:HG2	1:D:219:ARG:HH11	1.83	0.43
1:A:175:SER:HB3	1:A:242:ASP:OD1	2.19	0.42
1:B:228:VAL:HA	1:B:231:MET:HE3	2.00	0.42
1:B:68:ILE:HG22	1:B:69:ARG:N	2.33	0.42
1:B:320:ARG:HG3	1:B:320:ARG:HH11	1.83	0.42
1:D:262:LYS:C	1:D:264:ALA:H	2.27	0.42
1:A:268:ILE:HD11	1:A:279:GLY:HA3	2.01	0.42
1:D:307:ASN:OD1	1:D:309:GLU:CB	2.65	0.42
1:C:178:LEU:HD13	4:C:347:HOH:O	2.19	0.42
1:D:183:SER:HB3	1:D:224:ALA:HB3	2.01	0.42
1:A:74:VAL:O	1:A:75:GLN:C	2.61	0.42
1:B:46:LYS:HG2	1:B:51:GLU:HB3	2.00	0.42
1:C:161:ILE:HD12	1:C:161:ILE:HA	1.91	0.42
1:D:63:GLU:HG2	1:D:66:ASN:HB2	2.02	0.42
1:D:195:ILE:CD1	1:D:257:VAL:HG22	2.41	0.42
1:D:322:THR:HG23	1:D:325:GLU:OE1	2.20	0.42
1:B:12:ARG:HD2	1:B:42:SER:HB3	2.02	0.42
1:B:267:ASP:HB2	1:B:283:ASN:ND2	2.35	0.41
1:D:263:ILE:CD1	1:D:328:GLU:HA	2.50	0.41
1:A:228:VAL:HA	1:A:231:MET:HE3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:281:ASP:C	1:A:283:ASN:N	2.78	0.41
1:C:134:VAL:HG23	1:C:134:VAL:O	2.19	0.41
1:C:281:ASP:O	1:C:285:GLY:N	2.53	0.41
1:D:127:THR:O	1:D:130:MET:HG3	2.19	0.41
1:D:202:ILE:HG23	1:D:208:ASP:O	2.20	0.41
1:B:167:GLU:OE2	1:D:146:TYR:OH	2.35	0.41
1:B:224:ALA:HB3	1:B:225:PRO:HD3	2.03	0.41
1:B:332:GLU:OE1	1:B:332:GLU:HA	2.19	0.41
1:D:212:LEU:O	1:D:291:VAL:HA	2.20	0.41
1:D:291:VAL:CG1	1:D:292:SER:N	2.82	0.41
1:A:309:GLU:O	1:A:313:LYS:HB2	2.20	0.41
1:B:211:VAL:HA	1:B:290:GLU:O	2.21	0.41
1:B:282:ARG:O	1:B:283:ASN:C	2.63	0.41
1:A:295:PHE:HA	1:A:297:ARG:HH12	1.86	0.41
1:B:121:LYS:HB3	1:B:235:MET:CE	2.51	0.41
1:B:200:ALA:HB1	1:B:338:VAL:HG21	2.02	0.41
1:B:303:ILE:H	1:B:303:ILE:HG13	1.56	0.41
1:C:140:THR:HG22	1:C:307:ASN:HB3	2.02	0.41
1:C:331:MET:HE2	1:C:335:LEU:HG	2.03	0.41
1:A:68:ILE:HG22	1:A:69:ARG:HD3	2.03	0.41
1:A:253:VAL:O	1:A:257:VAL:HG23	2.21	0.41
1:B:300:GLU:H	1:B:300:GLU:CD	2.29	0.41
1:B:257:VAL:HG11	1:B:289:VAL:HG21	2.01	0.40
1:D:23:LEU:O	1:D:28:TYR:HB2	2.21	0.40
1:B:275:ILE:HD11	1:B:293:GLU:CA	2.51	0.40
1:A:9:THR:HG22	1:A:59:MET:HE1	2.04	0.40
1:A:117:LYS:N	1:A:118:PRO:HD3	2.36	0.40
1:B:5:ARG:HD2	1:B:74:VAL:O	2.21	0.40
1:C:174:CYS:HB3	1:C:234:MET:HE2	2.03	0.40
1:B:4:LYS:HE3	1:B:28:TYR:CZ	2.57	0.40
1:C:262:LYS:O	1:C:264:ALA:N	2.54	0.40
1:D:159:HIS:O	1:D:163:VAL:HG23	2.21	0.40
1:A:204:TYR:CZ	1:A:339:ARG:HG2	2.56	0.40
1:A:339:ARG:HG3	1:A:339:ARG:HH11	1.86	0.40
1:C:172:PHE:CE1	1:C:240:PRO:HD3	2.56	0.40
1:C:178:LEU:N	1:C:178:LEU:CD1	2.85	0.40
1:C:178:LEU:HG	1:C:245:ILE:HB	2.04	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	329/345 (95%)	311 (94%)	15 (5%)	3 (1%)	14	30
1	B	328/345 (95%)	309 (94%)	18 (6%)	1 (0%)	36	58
1	C	336/345 (97%)	313 (93%)	19 (6%)	4 (1%)	10	23
1	D	328/345 (95%)	288 (88%)	36 (11%)	4 (1%)	10	23
All	All	1321/1380 (96%)	1221 (92%)	88 (7%)	12 (1%)	14	30

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	86	SER
1	D	302	ASP
1	C	83	ALA
1	A	132	GLY
1	D	217	ALA
1	D	270	TRP
1	B	132	GLY
1	C	84	ALA
1	D	282	ARG
1	A	282	ARG
1	A	134	VAL
1	C	263	ILE

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	279/294 (95%)	277 (99%)	2 (1%)	76	89
1	B	279/294 (95%)	273 (98%)	6 (2%)	45	72
1	C	277/294 (94%)	272 (98%)	5 (2%)	51	76
1	D	258/294 (88%)	257 (100%)	1 (0%)	84	93
All	All	1093/1176 (93%)	1079 (99%)	14 (1%)	61	82

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	85	GLN
1	A	308	PRO
1	B	51	GLU
1	B	68	ILE
1	B	85	GLN
1	B	127	THR
1	B	303	ILE
1	B	340	ASP
1	C	51	GLU
1	C	191	VAL
1	C	208	ASP
1	C	219	ARG
1	C	313	LYS
1	D	85	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	75	GLN
1	B	180	ASN
1	C	251	HIS
1	D	66	ASN
1	D	170	ASN
1	D	214	ASN
1	D	251	HIS

5.3.3 RNA ⓘ

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](#)

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	GDP	D	1002	-	29,30,30	0.95	1 (3%)	45,47,47	0.61	0
2	NDP	A	1001	-	51,52,52	1.59	6 (11%)	71,80,80	0.98	3 (4%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GDP	D	1002	-	-	1/16/32/32	0/3/3/3
2	NDP	A	1001	-	-	8/34/77/77	0/5/5/5

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	NDP	PN-O3	5.69	1.65	1.59
2	A	1001	NDP	PA-O3	5.18	1.65	1.59
2	A	1001	NDP	C7N-C3N	-4.59	1.38	1.48
3	D	1002	GDP	PA-O3A	3.48	1.63	1.59
2	A	1001	NDP	P2B-O2B	-3.13	1.54	1.59
2	A	1001	NDP	C4N-C3N	2.65	1.55	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	NDP	C4N-C5N	2.45	1.55	1.49

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	NDP	O3-PN-O1N	-2.68	102.65	110.70
2	A	1001	NDP	O2B-P2B-O1X	-2.62	100.01	109.33
2	A	1001	NDP	C2D-C1D-N1N	2.22	118.77	113.31

There are no chirality outliers.

All (9) torsion outliers are listed below:

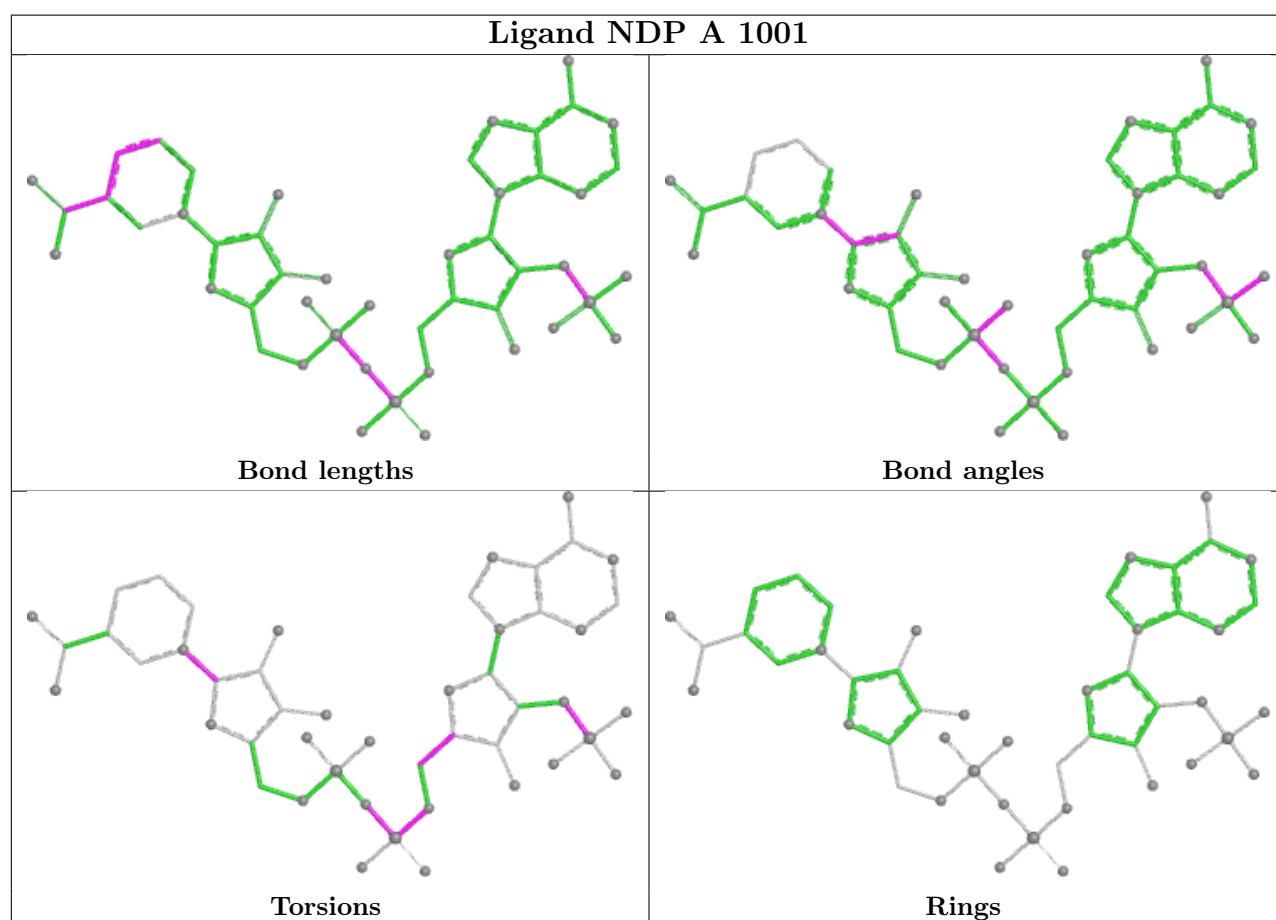
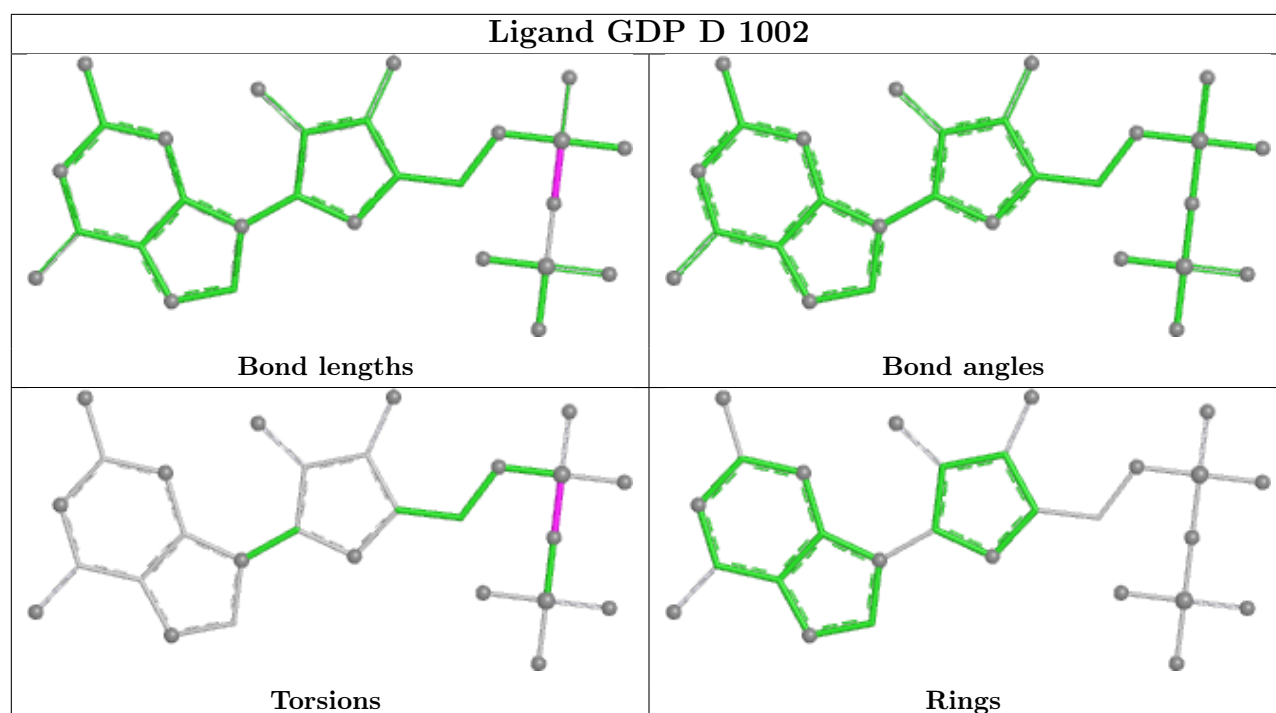
Mol	Chain	Res	Type	Atoms
2	A	1001	NDP	C5B-O5B-PA-O2A
2	A	1001	NDP	C5B-O5B-PA-O3
2	A	1001	NDP	O4B-C4B-C5B-O5B
2	A	1001	NDP	C3B-C4B-C5B-O5B
3	D	1002	GDP	PB-O3A-PA-O5'
2	A	1001	NDP	O4D-C1D-N1N-C6N
2	A	1001	NDP	C2B-O2B-P2B-O1X
2	A	1001	NDP	PN-O3-PA-O1A
2	A	1001	NDP	PN-O3-PA-O2A

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1001	NDP	1	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	333/345 (96%)	-0.17	5 (1%) 72 68	16, 37, 69, 85	0
1	B	332/345 (96%)	-0.30	4 (1%) 76 73	18, 37, 63, 76	0
1	C	338/345 (97%)	-0.15	3 (0%) 81 78	21, 39, 78, 91	0
1	D	332/345 (96%)	0.26	33 (9%) 13 10	23, 46, 97, 103	0
All	All	1335/1380 (96%)	-0.09	45 (3%) 48 42	16, 39, 82, 103	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	34	ASP	3.7
1	D	270	TRP	3.3
1	C	86	SER	3.2
1	A	298	PRO	3.1
1	B	301	VAL	3.0
1	D	288	ILE	2.9
1	D	217	ALA	2.9
1	D	213	GLY	2.8
1	D	289	VAL	2.8
1	D	273	GLU	2.6
1	A	299	ALA	2.6
1	C	85	GLN	2.6
1	D	303	ILE	2.6
1	D	204	TYR	2.6
1	D	299	ALA	2.5
1	D	284	THR	2.5
1	D	86	SER	2.5
1	C	340	ASP	2.4
1	D	287	VAL	2.4
1	D	254	ARG	2.4
1	A	340	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	38	GLY	2.4
1	B	303	ILE	2.3
1	D	268	ILE	2.3
1	D	301	VAL	2.3
1	D	295	PHE	2.3
1	A	276	ASN	2.2
1	D	253	VAL	2.2
1	A	86	SER	2.2
1	D	216	ASN	2.2
1	D	272	GLY	2.2
1	D	298	PRO	2.2
1	B	302	ASP	2.2
1	D	266	PHE	2.2
1	D	286	LYS	2.1
1	D	215	LEU	2.1
1	D	292	SER	2.1
1	D	84	ALA	2.1
1	D	264	ALA	2.1
1	D	132	GLY	2.1
1	D	285	GLY	2.1
1	D	271	VAL	2.1
1	D	291	VAL	2.1
1	D	334	ASP	2.0
1	D	276	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](#)

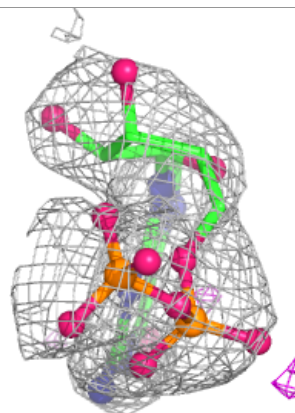
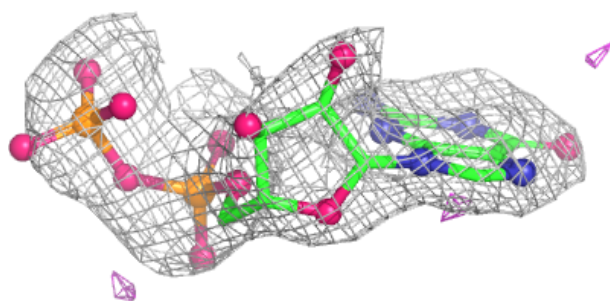
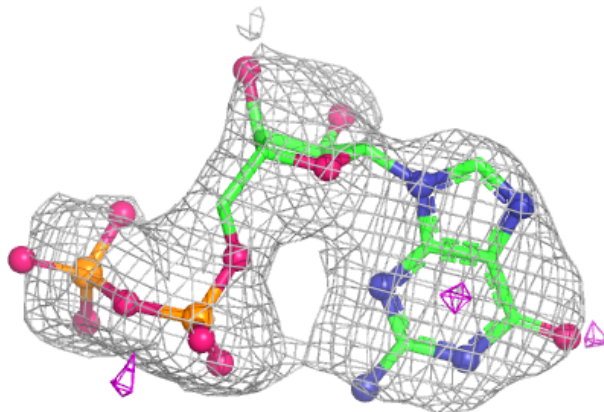
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	GDP	D	1002	28/28	0.83	0.13	85,90,103,103	0
2	NDP	A	1001	48/48	0.84	0.15	76,79,86,88	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.

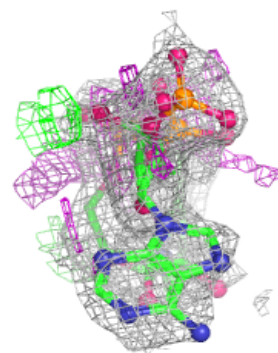
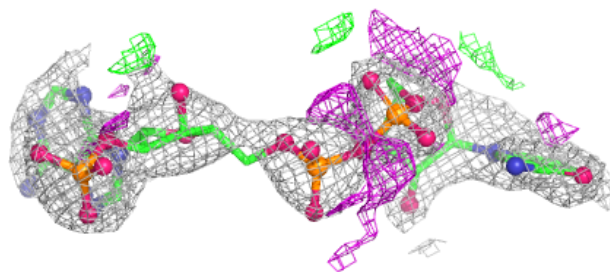
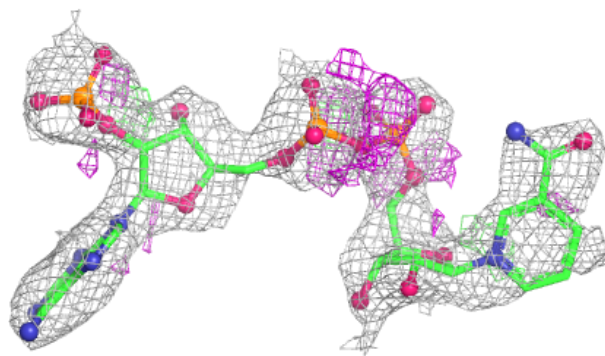
Electron density around GDP D 1002:

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 NDP A 1001:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.