



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 10:48 PM UTC

PDB ID : 6QNZ / pdb_00006qnz
Title : Crystal structure of the site-specific DNA nickase N.BspD6I E418A Mutant
Authors : Artyukh, R.I.; Kachalova, G.S.; Yunusova, A.K.; Gabdulkhakov, A.G.; Fatkhullin, B.F.; Atanasov, B.P.; Perevyazova, T.A.; Popov, A.N.; Zhelez-naya, L.A.
Deposited on : 2019-02-12
Resolution : 2.45 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
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

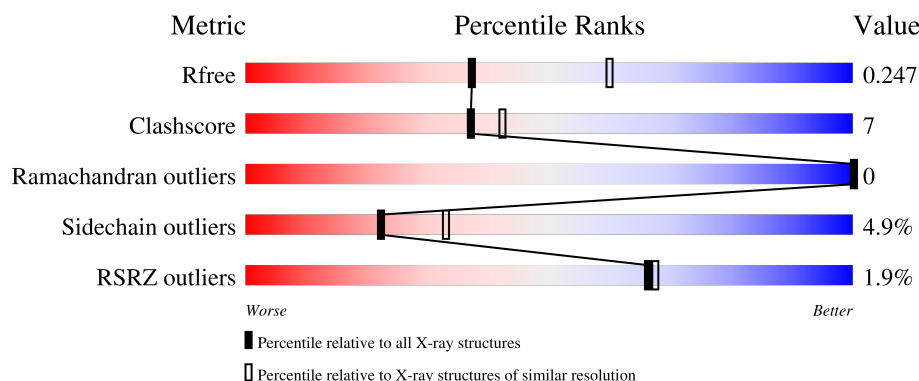
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	604	
1	B	604	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	PO4	A	703	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9515 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 Heterodimeric restriction endonuclease R.BspD6I large subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	575	Total	C	N	O	S	0	3	0
			4774	3065	806	888	15			
1	B	567	Total	C	N	O	S	0	0	0
			4695	3016	792	872	15			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	418	ALA	GLU	engineered mutation	UNP A3FEV7
B	418	ALA	GLU	engineered mutation	UNP A3FEV7

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	A	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		

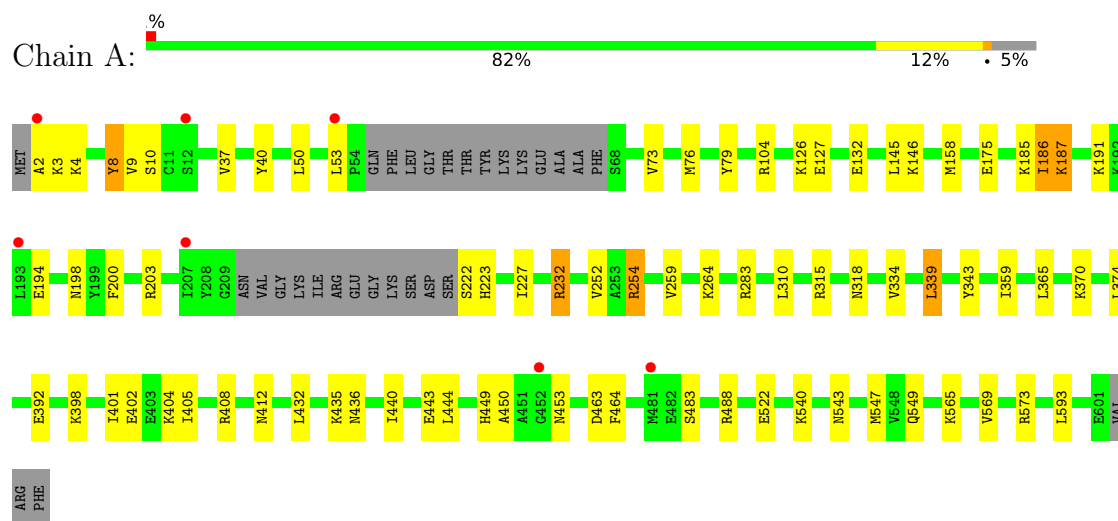
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	6	Total	O	0	0
			6	6		
4	B	3	Total	O	0	0
			3	3		

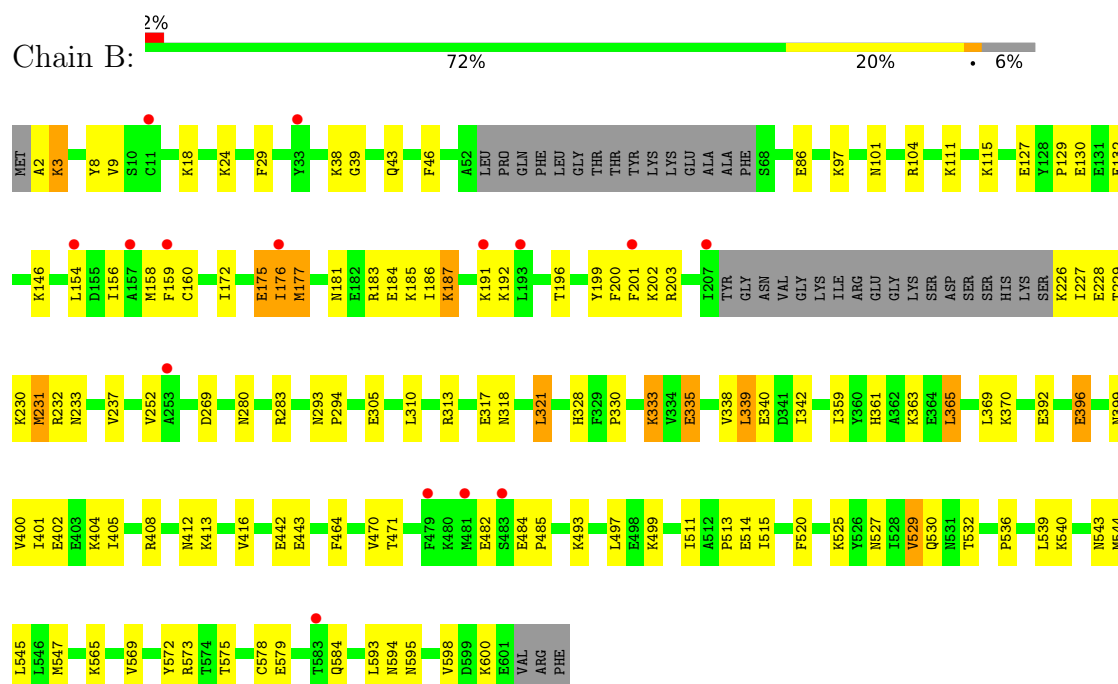
3 Residue-property plots [i](#)

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: Heterodimeric restriction endonuclease R.BspD6I large subunit



- Molecule 1: Heterodimeric restriction endonuclease R.BspD6I large subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	76.43Å 92.25Å 113.80Å 90.00° 105.61° 90.00°	Depositor
Resolution (Å)	47.11 – 2.45 47.11 – 2.45	Depositor EDS
% Data completeness (in resolution range)	87.8 (47.11-2.45) 87.9 (47.11-2.45)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 2.45Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.231 , 0.243 0.234 , 0.247	Depositor DCC
R_{free} test set	2468 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	58.8	Xtriage
Anisotropy	0.514	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9515	wwPDB-VP
Average B, all atoms (Å ²)	66.0	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 67.82 % 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 4.8327e-06. 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: GOL, PO4

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	0.22	0/4868	0.39	0/6553
1	B	0.21	0/4786	0.38	0/6442
All	All	0.22	0/9654	0.39	0/12995

There are no bond length outliers.

There are no bond angle outliers.

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	4774	0	4813	55	0
1	B	4695	0	4744	82	0
2	A	12	0	16	5	0
3	A	10	0	0	5	0
3	B	15	0	0	1	0
4	A	6	0	0	2	0
4	B	3	0	0	1	0
All	All	9515	0	9573	136	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 7.

All (136) 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:B:228:GLU:HA	1:B:231:MET:SD	1.88	1.14
1:B:228:GLU:HA	1:B:231:MET:CG	1.79	1.12
1:B:228:GLU:HA	1:B:231:MET:HG2	1.33	1.10
1:B:228:GLU:CA	1:B:231:MET:HG2	1.89	1.02
1:A:187:LYS:N	1:A:187:LYS:HE3	1.78	0.98
1:B:228:GLU:O	1:B:231:MET:HG2	1.65	0.96
1:B:228:GLU:C	1:B:231:MET:HG2	1.93	0.93
1:A:264:LYS:HE2	2:A:702:GOL:H11	1.48	0.93
1:A:412[A]:ASN:HD21	3:A:703:PO4:P	1.93	0.91
1:A:187:LYS:HE3	1:A:187:LYS:H	1.33	0.91
1:A:412[A]:ASN:ND2	3:A:703:PO4:P	2.47	0.86
1:A:3:LYS:HD3	1:A:132:GLU:OE1	1.81	0.80
1:B:396:GLU:HG2	1:B:400:VAL:HG11	1.67	0.77
1:A:264:LYS:HE2	2:A:702:GOL:C1	2.17	0.74
1:A:436:ASN:HD21	2:A:701:GOL:H12	1.54	0.73
1:B:146:LYS:NZ	1:B:269:ASP:OD1	2.21	0.73
1:A:186:ILE:N	1:A:186:ILE:HD13	2.03	0.71
1:B:181:ASN:O	1:B:185:LYS:HG2	1.92	0.69
1:B:199:TYR:HA	1:B:202:LYS:HG2	1.73	0.69
1:B:228:GLU:CA	1:B:231:MET:SD	2.77	0.67
1:A:412[A]:ASN:ND2	3:A:703:PO4:O2	2.26	0.67
1:A:404:LYS:O	1:A:408:ARG:NH1	2.27	0.67
1:B:175:GLU:OE2	1:B:203:ARG:NH1	2.28	0.65
1:B:228:GLU:O	1:B:232:ARG:N	2.25	0.65
1:A:435:LYS:NZ	4:A:802:HOH:O	2.30	0.64
1:B:232:ARG:HG3	1:B:233:ASN:N	2.13	0.64
1:B:318:ASN:HD22	1:B:339:LEU:HD21	1.61	0.64
1:A:254:ARG:HD2	1:A:259:VAL:HG21	1.80	0.62
1:B:186:ILE:O	1:B:192:LYS:NZ	2.33	0.61
1:A:412[A]:ASN:ND2	3:A:703:PO4:O1	2.33	0.61
1:A:50:LEU:HD12	1:A:53:LEU:HD12	1.82	0.61
1:B:412:ASN:OD1	1:B:413:LYS:N	2.34	0.60
1:B:529:VAL:HG12	1:B:530:GLN:HG2	1.84	0.59
1:B:201:PHE:CE1	1:B:227:ILE:HG12	2.39	0.58
1:B:228:GLU:O	1:B:231:MET:CG	2.46	0.58
1:A:175[A]:GLU:CD	1:A:203:ARG:HH12	2.10	0.58
1:A:252:VAL:HG11	1:A:443:GLU:HG3	1.85	0.58
1:B:470:VAL:HG12	1:B:511:ILE:HB	1.84	0.58
1:B:579:GLU:OE2	1:B:584:GLN:HG3	2.03	0.58
1:A:222:SER:OG	1:A:223:HIS:N	2.39	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:330:PRO:O	1:B:333:LYS:HG3	2.07	0.54
1:A:392:GLU:OE1	1:A:543:ASN:ND2	2.41	0.54
1:B:252:VAL:HG21	1:B:443:GLU:HG3	1.88	0.54
1:A:194:GLU:OE2	1:B:283:ARG:NH1	2.41	0.54
1:B:192:LYS:O	1:B:196:THR:HG23	2.08	0.53
1:B:342:ILE:HD12	1:B:365:LEU:HD22	1.90	0.53
1:B:412:ASN:O	1:B:416:VAL:HG23	2.09	0.52
1:A:2:ALA:O	1:A:127:GLU:HG2	2.09	0.52
1:B:104:ARG:NH2	3:B:701:PO4:O3	2.29	0.52
1:A:440:ILE:H	2:A:701:GOL:H32	1.74	0.52
1:A:449:HIS:HD2	1:A:450:ALA:O	1.93	0.51
1:B:229:THR:HA	1:B:232:ARG:HG2	1.93	0.51
1:A:549:GLN:NE2	4:A:801:HOH:O	2.28	0.51
1:B:172:ILE:O	1:B:176:ILE:HG23	2.09	0.51
1:B:115:LYS:NZ	1:B:293:ASN:O	2.43	0.51
1:B:159:PHE:CD2	1:B:176:ILE:HG22	2.46	0.50
1:B:176:ILE:HG13	1:B:177:MET:N	2.26	0.50
1:A:334:VAL:HG21	1:A:432:LEU:HB3	1.93	0.50
1:B:569:VAL:O	1:B:573:ARG:HG2	2.12	0.49
1:A:73:VAL:HA	1:A:76:MET:HE3	1.95	0.49
1:B:342:ILE:HD13	1:B:361:HIS:HB3	1.94	0.49
1:A:483:SER:OG	1:A:522:GLU:HG2	2.13	0.49
1:B:339:LEU:HG	1:B:365:LEU:HD23	1.95	0.49
1:A:440:ILE:H	2:A:701:GOL:C3	2.26	0.49
1:B:310:LEU:HD23	1:B:359:ILE:HD13	1.95	0.48
1:A:158:MET:HG2	1:A:200:PHE:CD1	2.49	0.48
1:B:97:LYS:HE3	1:B:101:ASN:HD21	1.79	0.48
1:B:184:GLU:OE1	1:B:187:LYS:NZ	2.32	0.48
1:B:513:PRO:HA	1:B:539:LEU:HB3	1.96	0.48
1:B:227:ILE:HG13	1:B:231:MET:HE3	1.96	0.47
1:A:310:LEU:HD23	1:A:359:ILE:HD13	1.97	0.47
1:A:318:ASN:HB3	1:A:339:LEU:HD21	1.96	0.47
1:B:280:ASN:O	1:B:280:ASN:ND2	2.47	0.47
1:A:232:ARG:HG3	1:A:232:ARG:HH11	1.80	0.47
1:B:399:ASN:OD1	1:B:400:VAL:N	2.48	0.47
1:B:2:ALA:O	1:B:127:GLU:HG2	2.15	0.47
1:A:569:VAL:O	1:A:573:ARG:HG2	2.15	0.47
1:A:401:ILE:O	1:A:405:ILE:HG12	2.15	0.46
1:A:543:ASN:O	1:A:547:MET:HG3	2.16	0.46
1:B:401:ILE:O	1:B:405:ILE:HG12	2.16	0.46
1:B:442:GLU:HG2	4:B:802:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:594:ASN:O	1:B:598:VAL:HG23	2.16	0.46
1:B:183:ARG:HE	1:B:196:THR:HG21	1.80	0.46
1:A:186:ILE:N	1:A:186:ILE:CD1	2.73	0.46
1:B:335:GLU:HG3	1:B:369:LEU:HD23	1.97	0.46
1:B:471:THR:HB	1:B:482:GLU:OE2	2.15	0.46
1:B:525:LYS:O	1:B:529:VAL:HB	2.16	0.46
1:A:175[A]:GLU:OE1	1:A:203:ARG:NH2	2.42	0.45
1:A:453:ASN:HA	1:A:488:ARG:HH12	1.81	0.45
1:B:365:LEU:HD12	1:B:365:LEU:HA	1.73	0.45
1:B:600:LYS:HA	1:B:600:LYS:HD3	1.84	0.45
1:B:156:ILE:HA	1:B:160:CYS:HB2	1.98	0.45
1:B:572:TYR:O	1:B:575:THR:OG1	2.32	0.45
1:A:398:LYS:O	1:A:402:GLU:HG3	2.17	0.45
1:B:493:LYS:O	1:B:497:LEU:HG	2.17	0.44
1:A:223:HIS:O	1:A:227:ILE:HG22	2.17	0.44
1:B:200:PHE:CE1	1:B:231:MET:HB2	2.53	0.44
1:B:540:LYS:HG2	1:B:544:MET:HE3	1.98	0.44
1:B:86:GLU:CD	1:B:283:ARG:HH21	2.25	0.44
1:B:97:LYS:HA	1:B:97:LYS:HD2	1.67	0.44
1:B:202:LYS:HG3	1:B:203:ARG:N	2.33	0.44
1:B:399:ASN:OD1	1:B:400:VAL:HG23	2.18	0.44
1:B:545:LEU:HD21	1:B:593:LEU:HD11	2.00	0.43
1:B:29:PHE:CZ	1:B:46:PHE:HA	2.54	0.43
1:A:8:TYR:CZ	1:A:10:SER:HA	2.53	0.43
1:A:315:ARG:HH11	1:A:339:LEU:HD23	1.83	0.43
1:A:2:ALA:O	1:A:127:GLU:CG	2.66	0.43
1:B:228:GLU:CA	1:B:231:MET:CG	2.61	0.43
1:B:9:VAL:H	1:B:237:VAL:HG13	1.84	0.43
1:B:313:ARG:O	1:B:317:GLU:HG3	2.19	0.43
1:B:305:GLU:OE1	1:B:305:GLU:N	2.36	0.42
1:A:318:ASN:CB	1:A:339:LEU:HD21	2.48	0.42
1:B:39:GLY:O	1:B:43:GLN:HG3	2.20	0.42
1:B:111:LYS:HD2	1:B:294:PRO:O	2.19	0.42
1:B:464:PHE:CE1	1:B:565:LYS:HB2	2.54	0.42
1:B:484:GLU:HB3	1:B:485:PRO:HD3	2.01	0.42
1:A:9:VAL:HG13	1:A:79:TYR:OH	2.20	0.42
1:A:463:ASP:OD1	1:A:463:ASP:N	2.51	0.42
1:A:175[A]:GLU:OE2	1:A:203:ARG:NH1	2.38	0.42
1:A:186:ILE:HG22	1:A:191:LYS:HB3	2.02	0.42
1:A:315:ARG:HD3	1:A:343:TYR:CE1	2.55	0.42
1:A:464:PHE:CE1	1:A:565:LYS:HB2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:520:PHE:CE2	1:B:536:PRO:HG2	2.55	0.41
1:A:365:LEU:HA	1:A:365:LEU:HD12	1.76	0.41
1:B:527:ASN:OD1	1:B:532:THR:OG1	2.36	0.41
1:A:565:LYS:O	1:A:569:VAL:HG23	2.20	0.41
1:A:104:ARG:NH2	3:A:704:PO4:O4	2.48	0.41
1:B:187:LYS:O	1:B:191:LYS:HD2	2.21	0.41
1:B:363:LYS:HA	1:B:363:LYS:HD2	1.60	0.41
1:B:97:LYS:HE3	1:B:101:ASN:ND2	2.36	0.41
1:A:37:VAL:HA	1:A:40:TYR:CE2	2.56	0.41
1:A:374:LEU:HD11	1:A:444:LEU:HD23	2.02	0.41
1:B:338:VAL:HB	1:B:365:LEU:HD21	2.03	0.41
1:B:321:LEU:HD23	1:B:321:LEU:HA	1.83	0.40
1:B:3:LYS:NZ	1:B:129:PRO:HG2	2.36	0.40
1:B:543:ASN:O	1:B:547:MET:HG3	2.21	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	571/604 (94%)	560 (98%)	11 (2%)	0	100	100
1	B	561/604 (93%)	550 (98%)	11 (2%)	0	100	100
All	All	1132/1208 (94%)	1110 (98%)	22 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	525/547 (96%)	509 (97%)	16 (3%)	36	51
1	B	516/547 (94%)	481 (93%)	35 (7%)	14	20
All	All	1041/1094 (95%)	990 (95%)	51 (5%)	22	33

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

Mol	Chain	Res	Type
1	A	4	LYS
1	A	8	TYR
1	A	126	LYS
1	A	145	LEU
1	A	146	LYS
1	A	185	LYS
1	A	186	ILE
1	A	187	LYS
1	A	198	ASN
1	A	232	ARG
1	A	254	ARG
1	A	283	ARG
1	A	339	LEU
1	A	370	LYS
1	A	540	LYS
1	A	593	LEU
1	B	3	LYS
1	B	8	TYR
1	B	18	LYS
1	B	24	LYS
1	B	38	LYS
1	B	130	GLU
1	B	132	GLU
1	B	154	LEU
1	B	158	MET
1	B	175	GLU
1	B	176	ILE
1	B	177	MET
1	B	187	LYS
1	B	226	LYS
1	B	230	LYS
1	B	231	MET
1	B	321	LEU

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Mol	Chain	Res	Type
1	B	328	HIS
1	B	333	LYS
1	B	335	GLU
1	B	339	LEU
1	B	340	GLU
1	B	365	LEU
1	B	370	LYS
1	B	392	GLU
1	B	396	GLU
1	B	402	GLU
1	B	404	LYS
1	B	408	ARG
1	B	499	LYS
1	B	514	GLU
1	B	515	ILE
1	B	529	VAL
1	B	578	CYS
1	B	595	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	101	ASN
1	A	280	ASN
1	A	516	ASN
1	B	312	HIS
1	B	445	GLN
1	B	516	ASN
1	B	530	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

7 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	PO4	B	702	-	4,4,4	1.16	0	6,6,6	0.47	0
3	PO4	B	703	-	4,4,4	1.54	1 (25%)	6,6,6	0.95	0
3	PO4	B	701	-	4,4,4	1.22	0	6,6,6	0.53	0
3	PO4	A	703	-	4,4,4	1.42	0	6,6,6	0.60	0
2	GOL	A	702	-	5,5,5	0.20	0	5,5,5	0.70	0
2	GOL	A	701	-	5,5,5	0.18	0	5,5,5	0.47	0
3	PO4	A	704	-	4,4,4	1.26	0	6,6,6	0.77	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	A	702	-	-	4/4/4/4	-
2	GOL	A	701	-	-	4/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	703	PO4	P-O3	-2.04	1.48	1.54

There are no bond angle outliers.

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701	GOL	O1-C1-C2-O2
2	A	701	GOL	O1-C1-C2-C3
2	A	701	GOL	C1-C2-C3-O3
2	A	702	GOL	O1-C1-C2-C3
2	A	701	GOL	O2-C2-C3-O3
2	A	702	GOL	O2-C2-C3-O3
2	A	702	GOL	C1-C2-C3-O3
2	A	702	GOL	O1-C1-C2-O2

There are no ring outliers.

5 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	701	PO4	1	0
3	A	703	PO4	4	0
2	A	702	GOL	2	0
2	A	701	GOL	3	0
3	A	704	PO4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	575/604 (95%)	0.18	7 (1%) 76 78	24, 58, 92, 104	6 (1%)
1	B	567/604 (93%)	0.47	15 (2%) 57 58	44, 68, 93, 107	9 (1%)
All	All	1142/1208 (94%)	0.32	22 (1%) 66 67	24, 63, 93, 107	15 (1%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2	ALA	4.0
1	B	11	CYS	3.6
1	A	207	ILE	3.2
1	A	53	LEU	3.1
1	B	207	ILE	3.1
1	B	176	ILE	3.0
1	B	157	ALA	2.8
1	B	193	LEU	2.6
1	B	33	TYR	2.6
1	B	154	LEU	2.5
1	B	481	MET	2.4
1	A	193	LEU	2.4
1	A	452	GLY	2.3
1	B	201	PHE	2.3
1	A	12	SER	2.3
1	B	479	PHE	2.2
1	B	253	ALA	2.1
1	B	191	LYS	2.1
1	A	481	MET	2.1
1	B	483	SER	2.1
1	B	159	PHE	2.0
1	B	583	THR	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	PO4	B	703	5/5	0.79	0.13	61,61,69,78	5
3	PO4	B	702	5/5	0.88	0.11	55,60,63,76	5
2	GOL	A	701	6/6	0.89	0.13	38,51,59,60	6
2	GOL	A	702	6/6	0.89	0.12	80,84,88,90	0
3	PO4	A	703	5/5	0.91	0.12	52,53,55,60	5
3	PO4	A	704	5/5	0.92	0.12	57,59,64,72	5
3	PO4	B	701	5/5	0.94	0.14	54,55,64,65	5

6.5 Other polymers [i](#)

There are no such residues in this entry.