



# Full wwPDB X-ray Structure Validation Report ⓘ

Dec 8, 2016 – 01:21 PM EST

PDB ID : 2XDR  
Title : CRYSTALLOGRAPHIC STRUCTURE OF BETAIN ALDEHYDE DEHYDROGENASE MUTANT E252A FROM PSEUDOMONAS AERUGINOSA  
Authors : Diaz-Sanchez, A.G.; Gonzalez-Segura, L.; Rudino-Pinera, E.; Lira-Rocha, A.; Munoz-Clares, R.A.  
Deposited on : 2010-05-06  
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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20028442  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20028442

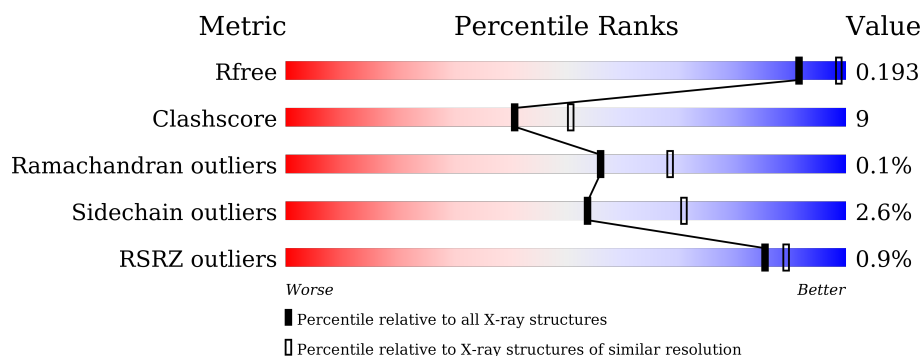
# 1 Overall quality at a glance

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))
$R_{free}$	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	489	<div> <div>85%</div> <div>13%</div> <div>.</div> </div>
1	B	489	<div> <div>%</div> <div>85%</div> <div>14%</div> <div>.</div> </div>
1	C	489	<div> <div>2%</div> <div>83%</div> <div>16%</div> <div>.</div> </div>
1	D	489	<div> <div>%</div> <div>86%</div> <div>13%</div> <div>.</div> </div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GOL	A	1491	-	-	-	X
2	GOL	B	1491	-	-	-	X
2	GOL	B	1492	-	-	-	X
2	GOL	C	1492	-	-	-	X
2	GOL	C	1493	-	-	-	X
2	GOL	C	1494	-	-	-	X
2	GOL	C	1495	-	-	-	X
2	GOL	D	1492	-	-	-	X
2	GOL	D	1493	-	-	-	X
3	TOE	A	1492	-	-	X	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 17057 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 BETAINE ALDEHYDE DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	489	Total	C	N	O	S	0	6	0
			3782	2373	665	729	15			
1	B	489	Total	C	N	O	S	0	4	0
			3775	2369	667	724	15			
1	C	489	Total	C	N	O	S	0	5	0
			3792	2379	669	728	16			
1	D	489	Total	C	N	O	S	0	6	0
			3815	2390	675	735	15			

There are 4 discrepancies between the modelled and reference sequences:

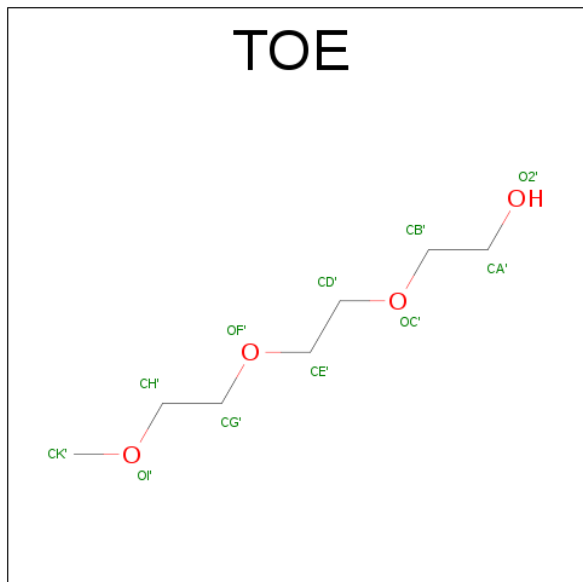
Chain	Residue	Modelled	Actual	Comment	Reference
A	252	ALA	GLU	ENGINEERED MUTATION	UNP Q9HTJ1
B	252	ALA	GLU	ENGINEERED MUTATION	UNP Q9HTJ1
C	252	ALA	GLU	ENGINEERED MUTATION	UNP Q9HTJ1
D	252	ALA	GLU	ENGINEERED MUTATION	UNP Q9HTJ1

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

- Molecule 3 is 2-[2-(2-METHOXY-ETHOXY)-ETHOXY]-ETHOXYL (three-letter code: TOE) (formula:  $C_7H_{16}O_4$ ).

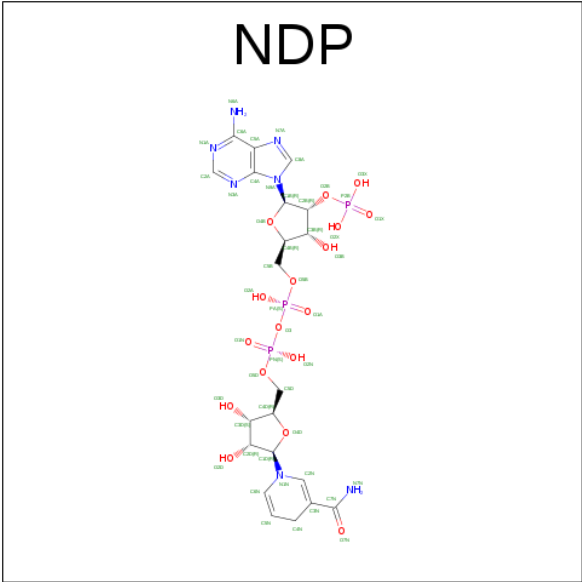


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			11	7	4		

- Molecule 4 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	2	Total	K	0	0
			2	2		
4	A	2	Total	K	0	0
			2	2		
4	D	2	Total	K	0	0
			2	2		
4	C	2	Total	K	0	0
			2	2		

- Molecule 5 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NDP) (formula:  $C_{21}H_{30}N_7O_{17}P_3$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
5	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
5	C	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
5	D	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

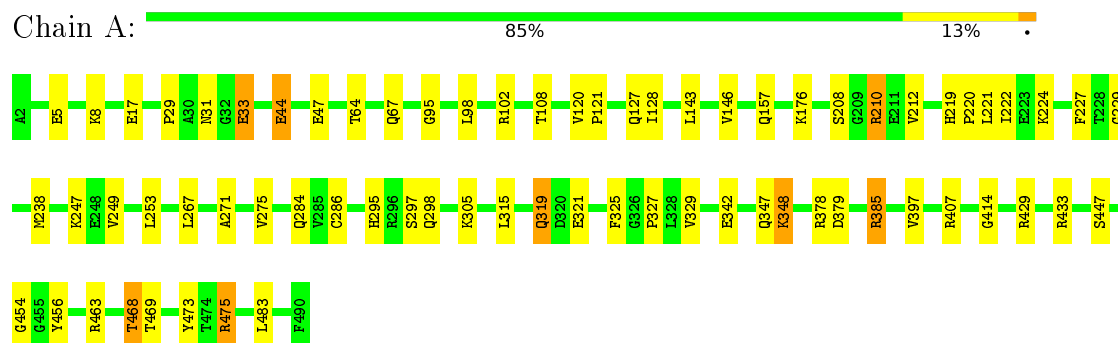
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	545	Total	O	0	0
			545	545		
6	B	390	Total	O	0	0
			390	390		
6	C	337	Total	O	0	0
			337	337		
6	D	410	Total	O	0	0
			410	410		

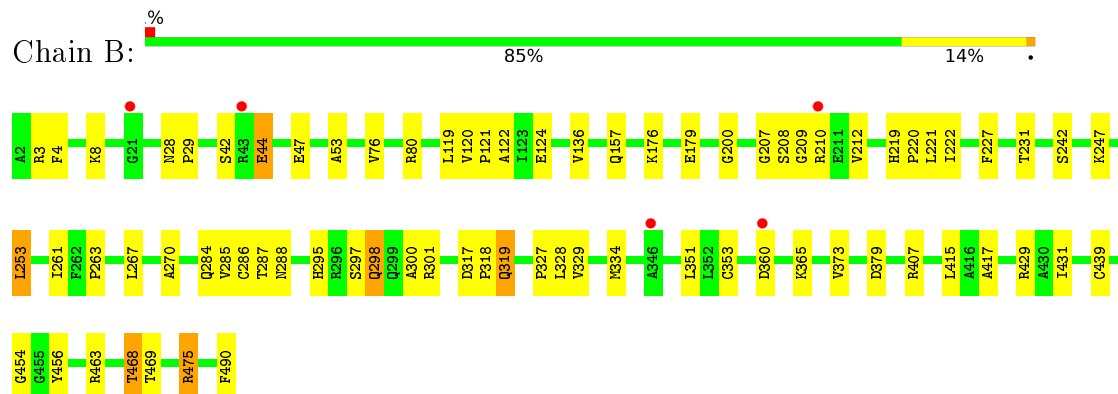
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

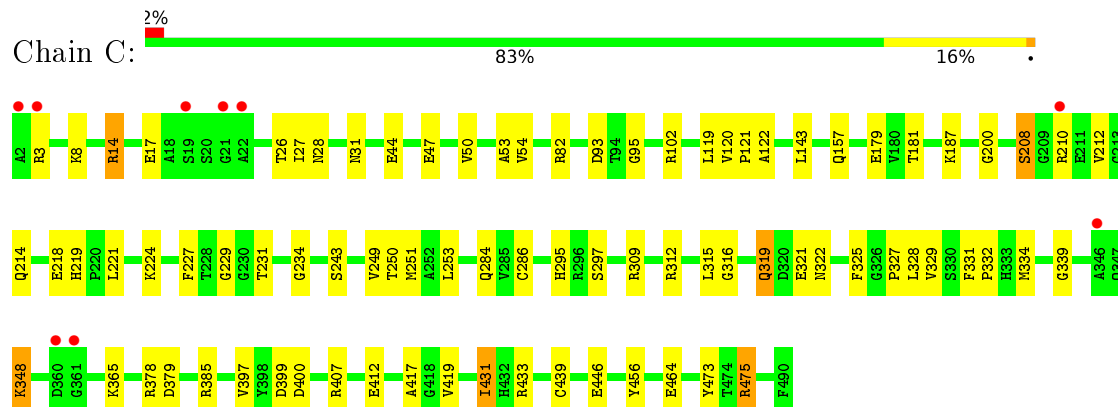
#### • Molecule 1: BETAINE ALDEHYDE DEHYDROGENASE



#### • Molecule 1: BETAINE ALDEHYDE DEHYDROGENASE

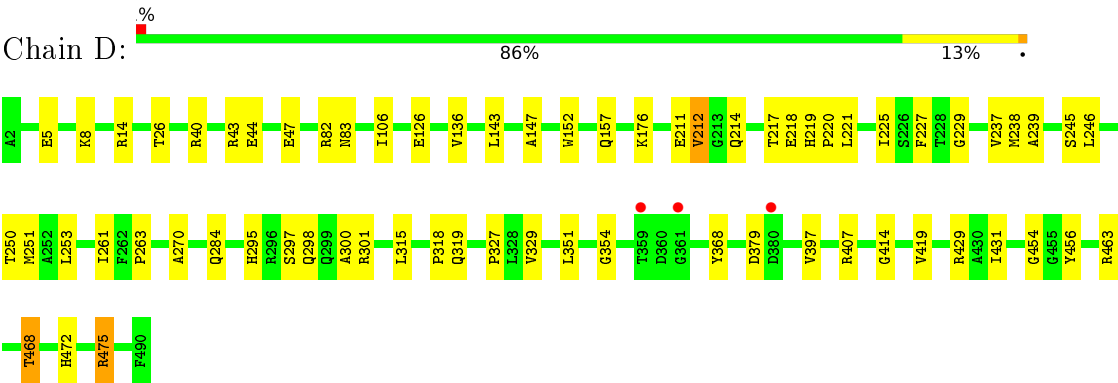


#### • Molecule 1: BETAINE ALDEHYDE DEHYDROGENASE





● Molecule 1: BETAINE ALDEHYDE DEHYDROGENASE



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	151.56Å 151.56Å 242.41Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.88 – 2.30 29.88 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.0 (29.88-2.30) 97.7 (29.88-2.30)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.35 (at 2.31Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.160 , 0.200 0.150 , 0.193	Depositor DCC
$R_{free}$ test set	7013 reflections (5.03%)	DCC
Wilson B-factor (Å <sup>2</sup> )	31.8	Xtriage
Anisotropy	0.092	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 58.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.005 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	17057	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.41% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: GOL, K, NDP, TOE

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	0.34	0/3853	0.51	0/5221
1	B	0.30	0/3843	0.48	0/5207
1	C	0.29	0/3860	0.48	0/5229
1	D	0.31	0/3883	0.48	0/5259
All	All	0.31	0/15439	0.49	0/20916

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 [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3782	0	3753	68	1
1	B	3775	0	3751	66	0
1	C	3792	0	3765	74	0
1	D	3815	0	3779	54	0
2	A	6	0	8	1	0
2	B	24	0	32	4	0
2	C	30	0	40	6	0
2	D	24	0	32	3	0
3	A	11	0	16	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	2	0	0	0	0
4	B	2	0	0	0	0
4	C	2	0	0	0	0
4	D	2	0	0	0	0
5	A	27	0	11	3	0
5	B	27	0	11	2	0
5	C	27	0	11	3	0
5	D	27	0	11	2	0
6	A	545	0	0	22	0
6	B	390	0	0	15	0
6	C	337	0	0	15	0
6	D	410	0	0	13	0
All	All	17057	0	15220	262	1

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 9.

All (262) 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:301[B]:ARG:HH21	1:B:301[B]:ARG:HG3	1.07	1.15
1:A:98:LEU:HD22	6:A:2374:HOH:O	1.55	1.04
1:A:219:HIS:HD2	1:A:221:LEU:H	1.14	0.94
1:C:27:ILE:HG22	1:C:28:ASN:H	1.30	0.93
1:B:301[B]:ARG:NH2	1:B:301[B]:ARG:HG3	1.85	0.85
1:D:284:GLN:HE22	1:D:329:VAL:H	1.26	0.84
1:D:219:HIS:HD2	1:D:221:LEU:H	1.26	0.83
1:C:143:LEU:HD23	1:C:475:ARG:HG2	1.60	0.82
6:C:2115:HOH:O	1:D:468:THR:HG22	1.78	0.81
1:B:298:GLN:HA	1:B:301[B]:ARG:NH2	1.96	0.81
1:B:261:ILE:HD13	1:B:270:ALA:HB1	1.63	0.80
1:B:176:LYS:HD2	1:B:212:VAL:HG13	1.63	0.79
1:C:400:ASP:HB3	6:C:2251:HOH:O	1.83	0.78
1:C:219:HIS:HD2	1:C:221:LEU:H	1.31	0.77
1:C:27:ILE:HG22	1:C:28:ASN:N	1.98	0.77
1:B:298:GLN:HA	1:B:301[B]:ARG:HH22	1.49	0.77
1:A:468:THR:HG22	6:B:2182:HOH:O	1.85	0.77
1:A:468:THR:HG23	6:A:2513:HOH:O	1.85	0.76
6:A:2224:HOH:O	1:B:468:THR:HG22	1.85	0.75
1:A:286[A]:CYS:SG	6:A:2310:HOH:O	2.43	0.74
1:A:157:GLN:OE1	3:A:1492:TOE:H4	1.87	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:286[A]:CYS:SG	6:B:2251:HOH:O	2.46	0.74
1:A:176:LYS:HD2	1:A:212:VAL:HG13	1.70	0.74
1:B:219:HIS:HD2	1:B:221:LEU:H	1.34	0.73
1:B:379:ASP:OD2	1:B:407:ARG:HD2	1.88	0.73
1:A:284:GLN:HE22	1:A:329:VAL:H	1.33	0.73
1:B:284:GLN:HE22	1:B:329:VAL:H	1.37	0.72
1:A:219:HIS:CD2	1:A:221:LEU:H	2.05	0.72
1:C:187:LYS:NZ	2:C:1491:GOL:H12	2.06	0.71
1:C:284:GLN:HE22	1:C:329:VAL:H	1.38	0.71
1:B:267:LEU:HD11	1:B:298:GLN:HG2	1.73	0.71
1:D:157:GLN:OE1	2:D:1492:GOL:H12	1.91	0.70
1:C:219:HIS:CD2	1:C:221:LEU:H	2.10	0.70
1:C:27:ILE:CG2	1:C:28:ASN:H	2.05	0.69
1:D:219:HIS:CD2	1:D:221:LEU:H	2.10	0.69
1:A:447:SER:H	3:A:1492:TOE:H16	1.57	0.68
1:D:468:THR:HG23	6:D:2373:HOH:O	1.93	0.67
2:B:1494:GOL:H31	6:B:2384:HOH:O	1.93	0.67
1:D:379:ASP:OD2	1:D:407:ARG:HD2	1.93	0.67
1:B:219:HIS:CD2	1:B:221:LEU:H	2.12	0.67
1:A:456:TYR:CZ	1:B:475:ARG:HD3	2.31	0.66
1:A:295:HIS:CD2	1:A:297:SER:H	2.15	0.65
1:A:29:PRO:HG3	6:A:2388:HOH:O	1.97	0.65
1:A:475:ARG:HD3	1:B:456:TYR:CZ	2.32	0.65
2:D:1491:GOL:H31	6:D:2399:HOH:O	1.97	0.64
1:B:212:VAL:HB	6:B:2074:HOH:O	1.97	0.64
1:C:446:GLU:OE1	6:C:2289:HOH:O	2.15	0.64
1:D:14[B]:ARG:CG	1:D:14[B]:ARG:HH11	2.10	0.64
1:A:447:SER:H	3:A:1492:TOE:CK'	2.11	0.64
1:A:143:LEU:HD23	1:A:475:ARG:HG2	1.79	0.63
1:D:295:HIS:CD2	1:D:297:SER:H	2.17	0.63
2:D:1491:GOL:H11	6:D:2399:HOH:O	1.99	0.63
1:B:47:GLU:OE1	1:B:219:HIS:HE1	1.82	0.62
1:A:212:VAL:HB	6:A:2090:HOH:O	1.99	0.62
2:A:1491:GOL:H12	3:A:1492:TOE:H5	1.82	0.62
1:C:157:GLN:OE1	2:C:1492:GOL:H12	2.00	0.62
1:C:295:HIS:HD2	1:C:297:SER:OG	1.83	0.61
1:C:31:ASN:O	1:C:365:LYS:HG2	2.00	0.61
1:C:456:TYR:CZ	1:D:475:ARG:HD3	2.35	0.61
1:D:284:GLN:HE22	1:D:329:VAL:N	1.97	0.61
1:D:5:GLU:HG2	6:D:2398:HOH:O	2.01	0.60
1:A:483:LEU:O	1:D:429[C]:ARG:NH2	2.35	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:284:GLN:HE22	1:A:329:VAL:N	2.00	0.59
1:D:261:ILE:HD12	1:D:270:ALA:HB1	1.84	0.59
1:A:229:GLY:O	1:A:253:LEU:HA	2.02	0.59
1:D:14[B]:ARG:HG2	1:D:14[B]:ARG:HH11	1.67	0.59
1:B:468:THR:HG23	6:B:2359:HOH:O	2.02	0.58
1:C:231:THR:HA	1:C:253:LEU:HD13	1.85	0.58
1:C:50:VAL:O	1:C:54:VAL:HG23	2.04	0.58
1:B:208:SER:HB2	6:B:2242:HOH:O	2.04	0.57
1:C:328:LEU:HD12	1:C:334:MET:HA	1.86	0.57
1:A:475:ARG:NH1	6:A:2520:HOH:O	2.04	0.57
1:B:231:THR:HA	1:B:253:LEU:HG	1.86	0.57
2:B:1493:GOL:H12	6:B:2007:HOH:O	2.04	0.57
1:C:187:LYS:HZ3	2:C:1491:GOL:H12	1.68	0.57
1:C:295:HIS:CD2	1:C:297:SER:H	2.24	0.56
1:B:463:ARG:HB2	6:B:2359:HOH:O	2.06	0.56
1:C:284:GLN:HE22	1:C:329:VAL:N	2.04	0.56
1:D:47:GLU:OE1	1:D:219:HIS:HE1	1.89	0.56
1:B:328:LEU:HD12	1:B:334:MET:HA	1.88	0.56
1:D:44[A]:GLU:HG3	6:D:2091:HOH:O	2.05	0.56
1:D:82:ARG:NH2	6:D:2155:HOH:O	2.32	0.56
1:A:102:ARG:NH2	1:A:321:GLU:HG2	2.20	0.55
1:A:379:ASP:OD2	1:A:407:ARG:HD2	2.07	0.55
1:B:157:GLN:OE1	2:B:1492:GOL:H12	2.05	0.55
1:C:249[A]:VAL:HG22	6:C:2182:HOH:O	2.06	0.55
1:B:176:LYS:HD2	1:B:212:VAL:CG1	2.34	0.55
1:B:179:GLU:HG3	1:B:207:GLY:O	2.06	0.55
1:D:14[B]:ARG:NH1	1:D:14[B]:ARG:HG2	2.21	0.54
1:A:397:VAL:O	1:A:407:ARG:NH2	2.39	0.54
1:A:227:PHE:HE1	5:A:5502:NDP:O4B	1.89	0.54
1:B:210:ARG:HG3	6:B:2242:HOH:O	2.06	0.54
1:D:229:GLY:O	1:D:253:LEU:HA	2.08	0.54
1:B:219:HIS:CD2	1:B:220:PRO:HD2	2.43	0.54
1:B:122:ALA:HB2	1:C:119:LEU:HD21	1.90	0.53
1:C:397:VAL:O	1:C:407:ARG:NH2	2.41	0.53
1:B:301[B]:ARG:HH21	1:B:301[B]:ARG:CG	1.96	0.53
1:B:227:PHE:HE1	5:B:5502:NDP:O4B	1.91	0.53
1:C:412:GLU:HG2	6:C:2132:HOH:O	2.09	0.52
1:D:318:PRO:HG3	1:D:327:PRO:HD2	1.91	0.52
1:C:227:PHE:HE1	5:C:5502:NDP:O4B	1.93	0.52
1:D:397:VAL:O	1:D:407:ARG:NH2	2.43	0.52
1:B:4:PHE:CD1	2:B:1493:GOL:H2	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:250:THR:HG21	6:D:2371:HOH:O	2.10	0.52
1:D:43:ARG:NH2	1:D:218[B]:GLU:OE1	2.43	0.51
1:A:429:ARG:HD3	6:A:2480:HOH:O	2.10	0.51
1:C:218:GLU:HG2	6:C:2178:HOH:O	2.10	0.51
1:A:47:GLU:OE1	1:A:219:HIS:HE1	1.93	0.51
1:A:319:GLN:HA	6:A:2374:HOH:O	2.10	0.51
1:B:42:SER:OG	1:B:44:GLU:HG2	2.10	0.51
1:C:417:ALA:O	1:C:439:CYS:HA	2.11	0.51
1:D:143:LEU:HD23	1:D:475:ARG:HG2	1.93	0.50
1:A:379:ASP:OD2	1:A:407:ARG:CD	2.59	0.50
1:C:378:ARG:NH1	6:C:2235:HOH:O	2.43	0.50
1:B:379:ASP:OD1	1:B:407:ARG:NH1	2.45	0.50
1:C:102:ARG:NH2	1:C:321:GLU:HG2	2.27	0.50
5:A:5502:NDP:H52A	6:A:2543:HOH:O	2.11	0.50
1:B:469:THR:HA	6:B:2361:HOH:O	2.12	0.50
1:C:102:ARG:HH21	1:C:321:GLU:HG2	1.76	0.50
1:C:219:HIS:CD2	1:C:221:LEU:HB2	2.46	0.50
1:A:210:ARG:HD3	6:A:2544:HOH:O	2.12	0.49
1:C:319:GLN:HE21	1:C:319:GLN:HA	1.78	0.49
1:C:187:LYS:HZ1	2:C:1491:GOL:H12	1.76	0.49
1:D:219:HIS:CD2	1:D:220:PRO:HD2	2.47	0.49
1:B:295:HIS:HD2	1:B:297:SER:OG	1.96	0.49
1:D:14[B]:ARG:NH1	1:D:14[B]:ARG:CG	2.74	0.49
1:C:31:ASN:HB2	1:C:365:LYS:HD3	1.95	0.48
1:C:315:LEU:HA	1:C:325:PHE:O	2.13	0.48
1:D:14[B]:ARG:HH11	1:D:14[B]:ARG:CB	2.25	0.48
1:A:219:HIS:CD2	1:A:220:PRO:HD2	2.49	0.48
1:C:95:GLY:O	1:C:327:PRO:HD2	2.13	0.48
1:A:102:ARG:HH22	1:A:321:GLU:HG2	1.78	0.48
1:A:8:LYS:HE2	1:A:17:GLU:OE2	2.13	0.48
1:C:210:ARG:HG3	6:C:2334:HOH:O	2.14	0.48
5:C:5502:NDP:O3	6:C:2336:HOH:O	2.20	0.48
5:D:5502:NDP:H52A	6:D:2223:HOH:O	2.13	0.48
1:C:419:VAL:HG22	1:C:431:ILE:HD13	1.96	0.48
1:B:119:LEU:HD21	1:C:122:ALA:HB2	1.95	0.48
6:B:2194:HOH:O	1:C:433:ARG:HD2	2.13	0.48
1:C:475:ARG:HD3	1:D:456:TYR:CZ	2.49	0.48
1:D:152:TRP:CE3	1:D:329:VAL:HG11	2.49	0.47
1:D:301:ARG:HD3	6:D:2286:HOH:O	2.14	0.47
1:D:472:HIS:CE1	6:D:2374:HOH:O	2.67	0.47
6:C:2182:HOH:O	1:D:238:MET:HE3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:321:GLU:HG3	6:A:2381:HOH:O	2.15	0.47
1:C:224:LYS:HE3	1:C:473:TYR:CZ	2.50	0.47
1:B:120:VAL:HB	1:B:121:PRO:HD3	1.95	0.47
1:B:353:CYS:HB2	1:B:373:VAL:H	1.80	0.47
1:C:120:VAL:HB	1:C:121:PRO:HD3	1.96	0.47
1:D:472:HIS:HE1	6:D:2374:HOH:O	1.98	0.47
1:A:146:VAL:HG12	1:A:224:LYS:HB3	1.97	0.46
1:B:286[B]:CYS:SG	1:B:287:THR:HG23	2.55	0.46
1:A:44[A]:GLU:HG3	6:A:2097:HOH:O	2.15	0.46
1:A:5:GLU:HG2	6:A:2002:HOH:O	2.15	0.46
1:B:318:PRO:HG3	1:B:327:PRO:HD2	1.98	0.46
1:C:8:LYS:HE2	1:C:17:GLU:OE2	2.15	0.46
1:A:271:ALA:O	1:A:275[A]:VAL:HG22	2.15	0.46
1:C:348:LYS:NZ	6:C:2225:HOH:O	2.48	0.46
1:A:210:ARG:NH2	6:A:2294:HOH:O	2.48	0.46
1:B:76:VAL:O	1:B:80:ARG:HG2	2.16	0.46
1:A:329:VAL:HG22	6:A:2388:HOH:O	2.15	0.46
1:A:95:GLY:O	1:A:327:PRO:HD2	2.16	0.46
1:D:176:LYS:HD2	1:D:212:VAL:HG22	1.97	0.46
1:D:219:HIS:HA	1:D:220:PRO:HD3	1.75	0.46
1:C:47:GLU:OE1	1:C:219:HIS:HE1	1.98	0.46
1:D:14[B]:ARG:HH11	1:D:14[B]:ARG:HB3	1.80	0.46
1:B:285:VAL:HB	1:B:288:ASN:OD1	2.15	0.46
1:C:251[B]:MET:HE3	1:C:253:LEU:HD21	1.96	0.46
1:B:319:GLN:HA	1:B:319:GLN:HE21	1.81	0.45
1:C:339:GLY:HA3	2:C:1495:GOL:H12	1.99	0.45
1:D:217:THR:HG21	1:D:237:VAL:HG13	1.99	0.45
1:D:327:PRO:HD3	1:D:368:TYR:CE1	2.51	0.45
1:B:301[B]:ARG:NH2	6:B:2279:HOH:O	2.49	0.45
1:B:208:SER:O	1:B:212:VAL:HG12	2.17	0.45
1:A:224:LYS:HE3	1:A:473:TYR:CE1	2.52	0.45
1:C:14:ARG:HE	1:C:14:ARG:HB3	1.60	0.45
1:A:222:ILE:O	1:A:247:LYS:HE3	2.17	0.45
1:D:14[B]:ARG:NH2	6:D:2023:HOH:O	2.50	0.45
1:C:53:ALA:HA	1:C:200:GLY:O	2.16	0.45
1:C:295:HIS:HE1	1:C:399:ASP:O	2.00	0.44
1:A:267:LEU:HB3	1:A:305:LYS:HD2	1.99	0.44
1:A:342:GLU:HG2	6:A:2400:HOH:O	2.16	0.44
1:C:229:GLY:O	1:C:253:LEU:HA	2.17	0.44
1:B:53:ALA:HA	1:B:200:GLY:O	2.17	0.44
1:A:454:GLY:HA3	1:A:463:ARG:HD3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:44[A]:GLU:HG3	6:C:2033:HOH:O	2.16	0.44
1:C:315:LEU:C	1:C:315:LEU:HD23	2.38	0.44
1:C:214:GLN:O	1:C:218:GLU:HG3	2.17	0.44
5:C:5502:NDP:H52A	6:C:2335:HOH:O	2.17	0.44
1:B:365:LYS:HE3	1:B:365:LYS:HB3	1.73	0.44
1:A:219:HIS:HA	1:A:220:PRO:HD3	1.73	0.44
1:C:331:PHE:N	1:C:332:PRO:CD	2.81	0.44
1:A:120:VAL:HB	1:A:121:PRO:HD3	2.00	0.43
1:B:379:ASP:CG	1:B:407:ARG:HH11	2.20	0.43
1:D:284:GLN:NE2	1:D:329:VAL:H	2.06	0.43
1:B:417:ALA:O	1:B:439:CYS:HA	2.19	0.43
1:C:243:SER:HB3	1:D:239:ALA:HB2	2.00	0.43
3:A:1492:TOE:H2	6:A:2342:HOH:O	2.19	0.43
1:D:147:ALA:HB3	1:D:225:ILE:HD13	2.01	0.43
3:A:1492:TOE:H15	6:A:2086:HOH:O	2.18	0.43
1:A:31:ASN:OD1	1:A:33:GLU:HB3	2.18	0.43
1:B:263:PRO:HD2	6:B:2331:HOH:O	2.19	0.43
1:C:295:HIS:CD2	1:C:297:SER:OG	2.68	0.43
1:B:284:GLN:HE22	1:B:329:VAL:N	2.12	0.43
1:C:464:GLU:OE1	2:C:1493:GOL:H31	2.18	0.43
1:C:179:GLU:OE1	1:C:208:SER:HB2	2.17	0.43
1:A:102:ARG:HH22	1:A:321:GLU:CG	2.32	0.42
1:A:385:ARG:HH11	1:A:385:ARG:HB3	1.84	0.42
1:C:93:ASP:O	1:C:181:THR:HG22	2.19	0.42
1:C:82:ARG:HD3	6:C:2077:HOH:O	2.18	0.42
1:A:108:THR:HG21	3:A:1492:TOE:H12	2.01	0.42
1:A:347:GLN:O	1:A:348:LYS:HB2	2.18	0.42
1:B:454:GLY:HA3	1:B:463:ARG:HD3	2.01	0.42
1:C:379:ASP:OD2	1:C:407:ARG:CD	2.67	0.42
1:C:379:ASP:OD2	1:C:407:ARG:HD2	2.18	0.42
1:C:475:ARG:H	1:C:475:ARG:HG3	1.58	0.42
1:A:275[B]:VAL:HG12	1:B:490:PHE:CD2	2.53	0.42
1:D:263:PRO:HA	1:D:298:GLN:OE1	2.19	0.42
1:A:295:HIS:HD2	1:A:297:SER:OG	2.02	0.42
1:B:351:LEU:HD11	1:B:353:CYS:O	2.19	0.42
1:B:300:ALA:HB3	6:B:2279:HOH:O	2.20	0.42
1:C:286[A]:CYS:SG	6:C:2175:HOH:O	2.62	0.42
1:C:295:HIS:HD2	1:C:297:SER:H	1.67	0.42
1:B:222:ILE:O	1:B:247:LYS:HE3	2.20	0.42
1:B:28:ASN:HA	1:B:29:PRO:HD2	1.96	0.42
1:D:245:SER:O	1:D:246:LEU:HB2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:348:LYS:HA	1:A:348:LYS:HD2	1.84	0.42
1:A:64:THR:OG1	1:A:67:GLN:HG3	2.20	0.42
1:A:315:LEU:HA	1:A:325:PHE:O	2.20	0.41
1:D:315:LEU:C	1:D:315:LEU:HD23	2.41	0.41
1:A:238:MET:HB3	1:B:242:SER:HB3	2.01	0.41
1:B:286[B]:CYS:SG	1:B:415:LEU:HD21	2.60	0.41
1:D:83:ASN:ND2	1:D:106:ILE:HB	2.35	0.41
1:D:419:VAL:HG22	1:D:431:ILE:HD13	2.02	0.41
1:A:208[A]:SER:O	1:A:212:VAL:HG12	2.20	0.41
1:B:317:ASP:HA	1:B:318:PRO:HD3	1.96	0.41
1:C:315:LEU:HD23	1:C:316:GLY:N	2.35	0.41
1:C:431:ILE:HG13	1:C:439:CYS:HB3	2.03	0.41
1:B:124:GLU:OE1	1:D:126:GLU:OE2	2.37	0.41
1:D:300:ALA:HB3	6:D:2288:HOH:O	2.19	0.41
1:C:309:ARG:O	1:C:312:ARG:HG2	2.21	0.41
1:C:31:ASN:O	1:C:365:LYS:O	2.39	0.41
1:D:295:HIS:HD2	1:D:297:SER:H	1.65	0.41
1:C:250:THR:C	1:C:251[B]:MET:HG2	2.41	0.41
1:A:284:GLN:NE2	1:A:329:VAL:H	2.09	0.41
1:B:44:GLU:H	1:B:44:GLU:CD	2.24	0.41
1:C:234:GLY:C	1:C:251[B]:MET:HE1	2.42	0.41
1:C:3[C]:ARG:HG2	1:C:3[C]:ARG:H	1.57	0.41
1:A:433:ARG:HD2	6:A:2484:HOH:O	2.20	0.41
1:B:431:ILE:HD12	1:B:431:ILE:HA	1.88	0.41
1:A:127:GLN:C	1:A:128:ILE:HG13	2.42	0.40
1:A:238:MET:HG3	1:A:249:VAL:HG21	2.04	0.40
1:A:298:GLN:NE2	6:A:2327:HOH:O	2.40	0.40
1:A:227:PHE:CE1	5:A:5502:NDP:O4B	2.72	0.40
1:B:209:GLY:HA3	5:B:5502:NDP:C4A	2.51	0.40
1:B:295:HIS:CD2	1:B:297:SER:H	2.39	0.40
1:D:227:PHE:CE1	5:D:5502:NDP:O4B	2.74	0.40
1:A:98:LEU:HD21	1:A:102:ARG:NH1	2.36	0.40
1:D:351:LEU:HD21	1:D:354:GLY:O	2.20	0.40
1:D:454:GLY:HA3	1:D:463:ARG:HD3	2.02	0.40
1:A:469:THR:HA	6:A:2516:HOH:O	2.21	0.40
1:B:468:THR:CG2	6:B:2359:HOH:O	2.64	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:378:ARG:NH1	1:A:378:ARG:NH1[6_555]	2.15	0.05

## 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	493/489 (101%)	476 (97%)	16 (3%)	1 (0%)	52	64
1	B	491/489 (100%)	472 (96%)	19 (4%)	0	100	100
1	C	493/489 (101%)	474 (96%)	19 (4%)	0	100	100
1	D	495/489 (101%)	478 (97%)	16 (3%)	1 (0%)	52	64
All	All	1972/1956 (101%)	1900 (96%)	70 (4%)	2 (0%)	56	68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	414	GLY
1	D	414	GLY

### 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	395/389 (102%)	386 (98%)	9 (2%)	58	75
1	B	393/389 (101%)	380 (97%)	13 (3%)	45	61
1	C	395/389 (102%)	385 (98%)	10 (2%)	55	73

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	397/389 (102%)	385 (97%)	12 (3%)	48	65
All	All	1580/1556 (102%)	1536 (97%)	44 (3%)	54	68

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	GLU
1	A	44[A]	GLU
1	A	44[B]	GLU
1	A	210	ARG
1	A	319	GLN
1	A	348	LYS
1	A	385	ARG
1	A	468	THR
1	A	475	ARG
1	B	3	ARG
1	B	8	LYS
1	B	44	GLU
1	B	136[A]	VAL
1	B	136[B]	VAL
1	B	253	LEU
1	B	298	GLN
1	B	319	GLN
1	B	360	ASP
1	B	429[A]	ARG
1	B	429[B]	ARG
1	B	468	THR
1	B	475	ARG
1	C	14	ARG
1	C	26	THR
1	C	208	SER
1	C	212	VAL
1	C	319	GLN
1	C	322	ASN
1	C	348	LYS
1	C	385	ARG
1	C	431	ILE
1	C	475	ARG
1	D	8	LYS
1	D	26	THR
1	D	40	ARG
1	D	136	VAL

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Mol	Chain	Res	Type
1	D	211	GLU
1	D	212	VAL
1	D	214[A]	GLN
1	D	214[B]	GLN
1	D	251	MET
1	D	319	GLN
1	D	468	THR
1	D	475	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	GLN
1	A	83	ASN
1	A	219	HIS
1	A	284	GLN
1	A	295	HIS
1	A	319	GLN
1	B	7	GLN
1	B	83	ASN
1	B	219	HIS
1	B	284	GLN
1	B	295	HIS
1	B	319	GLN
1	C	7	GLN
1	C	83	ASN
1	C	219	HIS
1	C	284	GLN
1	C	295	HIS
1	C	319	GLN
1	D	7	GLN
1	D	83	ASN
1	D	219	HIS
1	D	284	GLN
1	D	295	HIS
1	D	319	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 ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 27 ligands modelled in this entry, 8 are monoatomic - leaving 19 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	$\# Z  > 2$
2	GOL	A	1491	-	5,5,5	0.50	0	5,5,5	1.41	1 (20%)
3	TOE	A	1492	-	10,10,10	0.53	0	9,9,9	1.39	0
5	NDP	A	5502	-	25,29,52	2.70	11 (44%)	31,45,80	2.52	6 (19%)
2	GOL	B	1491	-	5,5,5	0.46	0	5,5,5	1.40	1 (20%)
2	GOL	B	1492	-	5,5,5	0.51	0	5,5,5	1.27	0
2	GOL	B	1493	-	5,5,5	0.50	0	5,5,5	1.26	0
2	GOL	B	1494	-	5,5,5	0.50	0	5,5,5	1.27	0
5	NDP	B	5502	-	25,29,52	2.73	11 (44%)	31,45,80	2.52	9 (29%)
2	GOL	C	1491	-	5,5,5	0.52	0	5,5,5	1.22	0
2	GOL	C	1492	-	5,5,5	0.49	0	5,5,5	1.27	0
2	GOL	C	1493	-	5,5,5	0.47	0	5,5,5	1.31	0
2	GOL	C	1494	-	5,5,5	0.46	0	5,5,5	1.24	0
2	GOL	C	1495	-	5,5,5	0.53	0	5,5,5	1.25	0
5	NDP	C	5502	-	25,29,52	2.71	11 (44%)	31,45,80	2.58	8 (25%)
2	GOL	D	1491	-	5,5,5	0.51	0	5,5,5	1.29	0
2	GOL	D	1492	-	5,5,5	0.50	0	5,5,5	1.21	0
2	GOL	D	1493	-	5,5,5	0.54	0	5,5,5	1.37	0
2	GOL	D	1494	-	5,5,5	0.47	0	5,5,5	1.07	0
5	NDP	D	5502	-	25,29,52	2.70	12 (48%)	31,45,80	2.47	7 (22%)

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	1491	-	-	0/4/4/4	0/0/0/0
3	TOE	A	1492	-	-	0/8/8/8	0/0/0/0
5	NDP	A	5502	-	-	0/11/31/77	0/3/3/5
2	GOL	B	1491	-	-	0/4/4/4	0/0/0/0
2	GOL	B	1492	-	-	0/4/4/4	0/0/0/0
2	GOL	B	1493	-	-	0/4/4/4	0/0/0/0
2	GOL	B	1494	-	-	0/4/4/4	0/0/0/0
5	NDP	B	5502	-	-	0/11/31/77	0/3/3/5
2	GOL	C	1491	-	-	0/4/4/4	0/0/0/0
2	GOL	C	1492	-	-	0/4/4/4	0/0/0/0
2	GOL	C	1493	-	-	0/4/4/4	0/0/0/0
2	GOL	C	1494	-	-	0/4/4/4	0/0/0/0
2	GOL	C	1495	-	-	0/4/4/4	0/0/0/0
5	NDP	C	5502	-	-	0/11/31/77	0/3/3/5
2	GOL	D	1491	-	-	0/4/4/4	0/0/0/0
2	GOL	D	1492	-	-	0/4/4/4	0/0/0/0
2	GOL	D	1493	-	-	0/4/4/4	0/0/0/0
2	GOL	D	1494	-	-	0/4/4/4	0/0/0/0
5	NDP	D	5502	-	-	0/11/31/77	0/3/3/5

All (45) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	5502	NDP	P2B-O3X	-2.84	1.44	1.54
5	D	5502	NDP	P2B-O3X	-2.83	1.45	1.54
5	B	5502	NDP	P2B-O3X	-2.81	1.45	1.54
5	A	5502	NDP	P2B-O3X	-2.80	1.45	1.54
5	C	5502	NDP	PA-O3	-2.76	1.45	1.54
5	B	5502	NDP	PA-O3	-2.62	1.45	1.54
5	A	5502	NDP	PA-O3	-2.56	1.45	1.54
5	D	5502	NDP	PA-O3	-2.54	1.46	1.54
5	D	5502	NDP	C2B-C1B	-2.08	1.47	1.53
5	D	5502	NDP	PA-O2A	2.08	1.62	1.54
5	C	5502	NDP	PA-O2A	2.12	1.62	1.54
5	B	5502	NDP	PA-O2A	2.12	1.62	1.54
5	A	5502	NDP	PA-O2A	2.14	1.62	1.54
5	D	5502	NDP	O4B-C4B	2.55	1.50	1.45
5	A	5502	NDP	O4B-C4B	2.68	1.51	1.45
5	C	5502	NDP	O4B-C4B	2.73	1.51	1.45
5	B	5502	NDP	O4B-C4B	2.73	1.51	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	5502	NDP	C6A-N6A	3.08	1.46	1.34
5	C	5502	NDP	C6A-N6A	3.18	1.47	1.34
5	A	5502	NDP	C6A-N6A	3.20	1.47	1.34
5	B	5502	NDP	C6A-N6A	3.20	1.47	1.34
5	D	5502	NDP	P2B-O2X	3.52	1.66	1.54
5	C	5502	NDP	P2B-O2X	3.60	1.67	1.54
5	A	5502	NDP	C3B-C4B	3.61	1.62	1.53
5	C	5502	NDP	C3B-C4B	3.63	1.62	1.53
5	A	5502	NDP	P2B-O2X	3.63	1.67	1.54
5	D	5502	NDP	C3B-C4B	3.65	1.62	1.53
5	B	5502	NDP	P2B-O2X	3.68	1.67	1.54
5	B	5502	NDP	C3B-C4B	3.82	1.63	1.53
5	A	5502	NDP	P2B-O1X	4.11	1.63	1.50
5	C	5502	NDP	PA-O1A	4.27	1.64	1.50
5	B	5502	NDP	P2B-O1X	4.32	1.64	1.50
5	A	5502	NDP	PA-O1A	4.38	1.64	1.50
5	C	5502	NDP	P2B-O1X	4.41	1.64	1.50
5	D	5502	NDP	P2B-O1X	4.42	1.64	1.50
5	B	5502	NDP	PA-O1A	4.52	1.65	1.50
5	D	5502	NDP	PA-O1A	4.59	1.65	1.50
5	D	5502	NDP	C8A-N7A	5.39	1.45	1.34
5	A	5502	NDP	C8A-N7A	5.45	1.45	1.34
5	B	5502	NDP	C8A-N7A	5.51	1.45	1.34
5	C	5502	NDP	C8A-N7A	5.53	1.45	1.34
5	C	5502	NDP	C2A-N3A	6.11	1.43	1.32
5	B	5502	NDP	C2A-N3A	6.17	1.43	1.32
5	D	5502	NDP	C2A-N3A	6.21	1.43	1.32
5	A	5502	NDP	C2A-N3A	6.36	1.43	1.32

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	5502	NDP	N3A-C2A-N1A	-11.60	119.76	128.87
5	D	5502	NDP	N3A-C2A-N1A	-11.42	119.90	128.87
5	B	5502	NDP	N3A-C2A-N1A	-11.31	119.99	128.87
5	C	5502	NDP	N3A-C2A-N1A	-11.28	120.01	128.87
5	D	5502	NDP	O4B-C4B-C3B	-2.46	100.17	105.16
5	B	5502	NDP	O4B-C4B-C3B	-2.41	100.27	105.16
5	B	5502	NDP	C1B-N9A-C4A	-2.18	124.38	126.81
5	D	5502	NDP	C3B-C2B-C1B	-2.15	98.52	102.63
5	B	5502	NDP	C2B-C1B-N9A	-2.06	107.58	113.48
5	C	5502	NDP	O4B-C4B-C3B	-2.04	101.03	105.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	5502	NDP	O4B-C4B-C3B	-2.00	101.10	105.16
2	B	1491	GOL	O3-C3-C2	2.01	120.16	109.97
5	D	5502	NDP	O4B-C1B-N9A	2.06	111.99	108.11
5	B	5502	NDP	O2B-C2B-C1B	2.14	118.53	110.12
2	A	1491	GOL	O3-C3-C2	2.17	121.00	109.97
5	B	5502	NDP	O4B-C1B-N9A	2.29	112.43	108.11
5	C	5502	NDP	O4B-C4B-C5B	2.31	117.54	109.29
5	C	5502	NDP	O2B-C2B-C1B	2.32	119.21	110.12
5	B	5502	NDP	O3B-C3B-C2B	2.35	117.92	111.13
5	C	5502	NDP	O3-PA-O5B	2.38	113.68	106.72
5	D	5502	NDP	O3B-C3B-C2B	2.40	118.08	111.13
5	A	5502	NDP	O2B-C2B-C1B	2.49	119.90	110.12
5	D	5502	NDP	O2B-C2B-C3B	2.55	121.23	111.73
5	A	5502	NDP	O2B-C2B-C3B	2.69	121.73	111.73
5	B	5502	NDP	O2B-C2B-C3B	2.92	122.58	111.73
5	A	5502	NDP	O3B-C3B-C2B	2.97	119.71	111.13
5	C	5502	NDP	O2B-C2B-C3B	2.99	122.86	111.73
5	D	5502	NDP	O3B-C3B-C4B	3.24	120.68	111.01
5	B	5502	NDP	O3B-C3B-C4B	3.28	120.82	111.01
5	A	5502	NDP	O3B-C3B-C4B	3.30	120.88	111.01
5	C	5502	NDP	O3B-C3B-C2B	3.34	120.77	111.13
5	C	5502	NDP	O3B-C3B-C4B	3.46	121.33	111.01

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

15 monomers are involved in 30 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1491	GOL	1	0
3	A	1492	TOE	7	0
5	A	5502	NDP	3	0
2	B	1492	GOL	1	0
2	B	1493	GOL	2	0
2	B	1494	GOL	1	0
5	B	5502	NDP	2	0
2	C	1491	GOL	3	0
2	C	1492	GOL	1	0
2	C	1493	GOL	1	0
2	C	1495	GOL	1	0
5	C	5502	NDP	3	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	1491	GOL	2	0
2	D	1492	GOL	1	0
5	D	5502	NDP	2	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	489/489 (100%)	-0.60	0 100 100	14, 25, 43, 78	0
1	B	489/489 (100%)	-0.38	5 (1%) 84 88	14, 37, 63, 96	0
1	C	489/489 (100%)	-0.42	9 (1%) 71 78	21, 39, 60, 92	0
1	D	489/489 (100%)	-0.44	3 (0%) 90 93	17, 33, 58, 79	0
All	All	1956/1956 (100%)	-0.46	17 (0%) 85 89	14, 34, 59, 96	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	2	ALA	4.0
1	B	210	ARG	3.2
1	D	359	THR	3.0
1	C	346	ALA	3.0
1	C	21	GLY	2.8
1	D	361	GLY	2.6
1	C	19	SER	2.5
1	C	361	GLY	2.5
1	B	21	GLY	2.3
1	B	360	ASP	2.3
1	C	3[A]	ARG	2.3
1	D	380	ASP	2.3
1	C	210	ARG	2.2
1	B	346	ALA	2.2
1	C	22	ALA	2.1
1	C	360	ASP	2.1
1	B	43	ARG	2.0

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

### 6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	GOL	C	1493	6/6	0.86	0.26	10.71	60,79,86,90	0
2	GOL	D	1493	6/6	0.83	0.29	9.20	57,60,76,84	0
2	GOL	B	1491	6/6	0.77	0.32	8.78	61,64,66,71	0
2	GOL	C	1494	6/6	0.67	0.38	8.04	77,92,98,100	0
2	GOL	C	1492	6/6	0.92	0.30	7.90	49,58,68,73	0
2	GOL	B	1492	6/6	0.91	0.31	6.21	61,63,73,76	0
2	GOL	A	1491	6/6	0.84	0.23	4.72	36,48,55,60	0
2	GOL	C	1495	6/6	0.84	0.39	4.53	70,86,87,90	0
3	TOE	A	1492	11/11	0.83	0.27	4.28	22,61,79,88	0
2	GOL	D	1492	6/6	0.93	0.26	4.20	54,67,71,74	0
2	GOL	B	1493	6/6	0.94	0.13	1.83	62,65,67,68	0
4	K	B	5500	1/1	0.92	0.13	1.74	68,68,68,68	0
2	GOL	C	1491	6/6	0.91	0.20	1.60	70,72,75,78	0
4	K	D	1496	1/1	0.91	0.11	1.29	67,67,67,67	0
5	NDP	B	5502	27/48	0.87	0.19	0.90	39,55,72,94	20
5	NDP	A	5502	27/48	0.92	0.15	0.64	23,38,58,90	23
5	NDP	C	5502	27/48	0.92	0.16	0.60	34,52,80,95	20
4	K	C	1497	1/1	0.92	0.11	0.56	74,74,74,74	0
2	GOL	D	1494	6/6	0.91	0.14	0.50	53,56,60,62	0
4	K	A	1494	1/1	0.94	0.10	0.37	48,48,48,48	0
4	K	A	1493	1/1	0.97	0.09	0.18	30,30,30,30	0
4	K	B	5501	1/1	0.94	0.10	0.15	50,50,50,50	0
4	K	D	1495	1/1	0.98	0.10	-0.04	41,41,41,41	0
5	NDP	D	5502	27/48	0.97	0.10	-0.38	22,33,54,87	12
4	K	C	1496	1/1	0.93	0.10	-0.39	58,58,58,58	0
2	GOL	B	1494	6/6	0.92	0.18	-	46,72,79,82	0
2	GOL	D	1491	6/6	0.89	0.19	-	38,81,89,94	0

## 6.5 Other polymers

There are no such residues in this entry.