



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 02:36 pm GMT

PDB ID : 2Y53
Title : Crystal structure of E257Q mutant of the box pathway encoded ALDH from Burkholderia xenovorans LB400
Authors : Bains, J.; Leon, R.; Temke, K.G.; Boulanger, M.J.
Deposited on : 2011-01-11
Resolution : 1.40 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

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

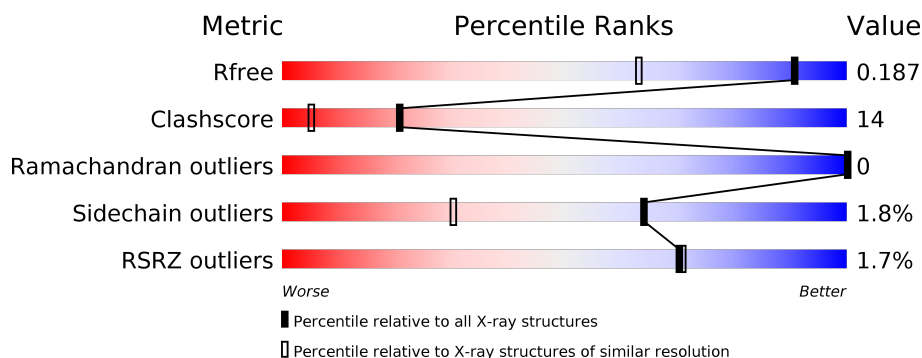
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.40 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}	100719	1307 (1.40-1.40)
Clashscore	112137	1411 (1.40-1.40)
Ramachandran outliers	110173	1373 (1.40-1.40)
Sidechain outliers	110143	1372 (1.40-1.40)
RSRZ outliers	101464	1315 (1.40-1.40)

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 ≥ 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	534	<div> <div>%</div> <div> <div></div> <div>81%</div> <div>14%</div> <div>..</div> </div> </div>
1	B	534	<div> <div>2%</div> <div> <div></div> <div>81%</div> <div>15%</div> <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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NAP	A	1534	-	-	X	X
3	NAP	B	1527	-	-	X	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9444 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.

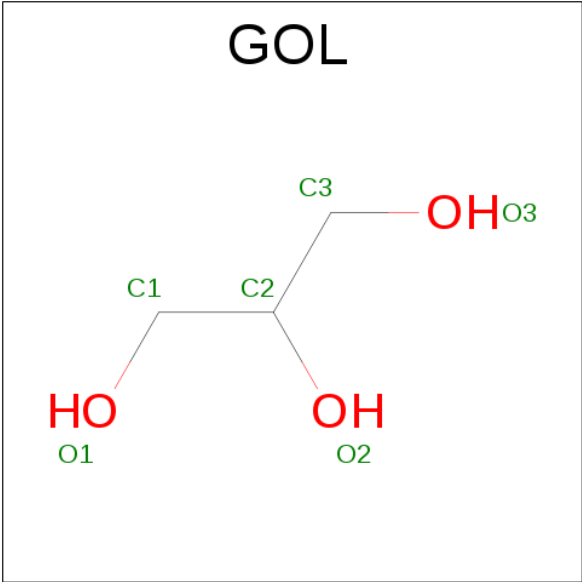
- Molecule 1 is a protein called ALDEHYDE DEHYDROGENASE (BOX PATHWAY).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	529	Total	C	N	O	S	0	1	0
			3905	2451	706	738	10			
1	B	524	Total	C	N	O	S	0	3	0
			3884	2434	703	737	10			

There are 8 discrepancies between the modelled and reference sequences:

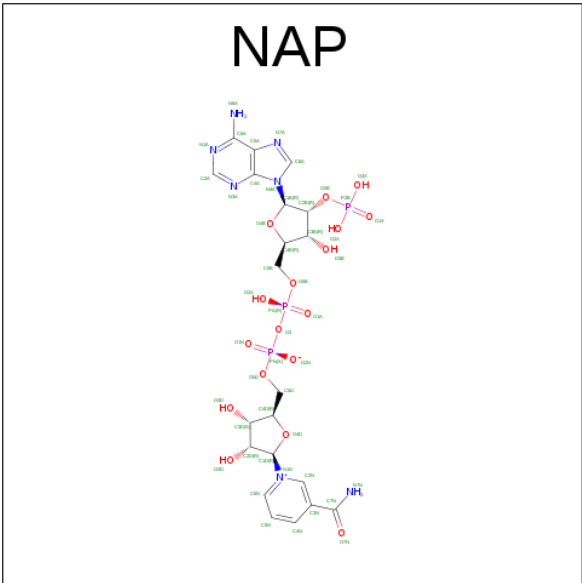
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP Q13WK4
A	-1	SER	-	EXPRESSION TAG	UNP Q13WK4
A	0	HIS	-	EXPRESSION TAG	UNP Q13WK4
A	257	GLN	GLU	ENGINEERED MUTATION	UNP Q13WK4
B	-2	GLY	-	EXPRESSION TAG	UNP Q13WK4
B	-1	SER	-	EXPRESSION TAG	UNP Q13WK4
B	0	HIS	-	EXPRESSION TAG	UNP Q13WK4
B	257	GLN	GLU	ENGINEERED MUTATION	UNP Q13WK4

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



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

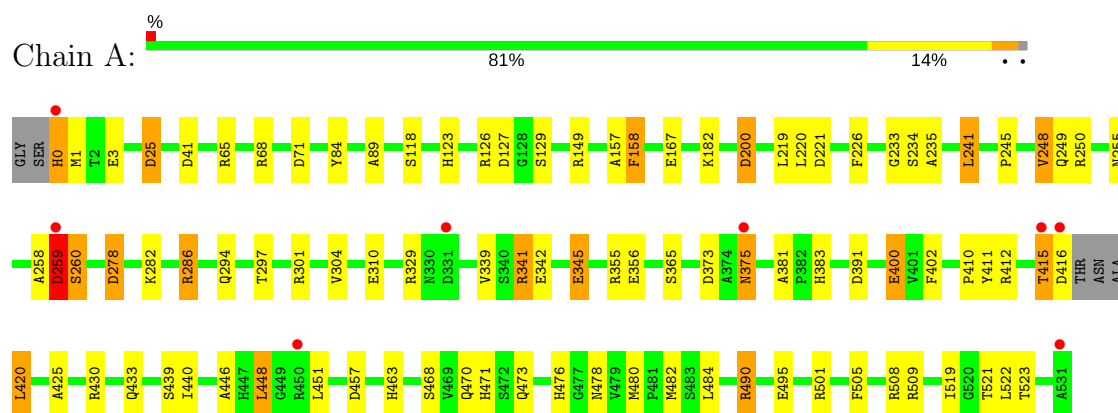
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	797	Total 797	O 797	0	0
4	B	738	Total 738	O 738	0	0

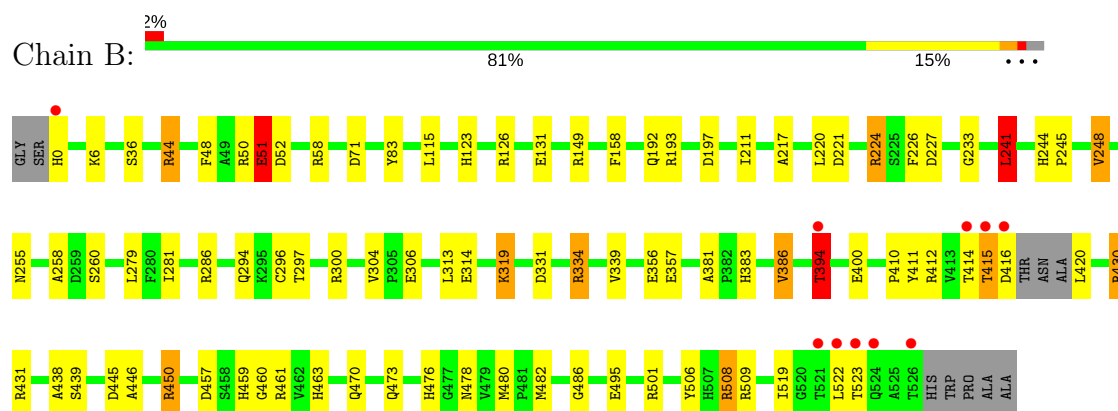
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: ALDEHYDE DEHYDROGENASE (BOX PATHWAY)



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4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	58.24Å 67.81Å 77.51Å 111.22° 90.45° 113.51°	Depositor
Resolution (Å)	23.37 – 1.40 23.37 – 1.40	Depositor EDS
% Data completeness (in resolution range)	95.6 (23.37-1.40) 83.0 (23.37-1.40)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.63 (at 1.40Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.156 , 0.185 0.157 , 0.187	Depositor DCC
R_{free} test set	9501 reflections (5.33%)	DCC
Wilson B-factor (Å ²)	8.9	Xtriage
Anisotropy	0.011	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 45.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.014 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	9444	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.93% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	1.46	28/3982 (0.7%)	1.53	49/5425 (0.9%)
1	B	1.41	15/3957 (0.4%)	1.31	32/5387 (0.6%)
All	All	1.43	43/7939 (0.5%)	1.42	81/10812 (0.7%)

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	129	SER	CB-OG	-17.04	1.20	1.42
1	A	341	ARG	CZ-NH2	11.82	1.48	1.33
1	A	412	ARG	CZ-NH2	-10.56	1.19	1.33
1	A	345	GLU	CD-OE1	-9.73	1.15	1.25
1	A	375	ASN	CB-CG	9.12	1.72	1.51
1	B	51	GLU	CB-CG	-8.80	1.35	1.52
1	A	248	VAL	CB-CG2	-8.78	1.34	1.52
1	A	226	PHE	CD1-CE1	8.35	1.55	1.39
1	A	341	ARG	CZ-NH1	8.10	1.43	1.33
1	A	341	ARG	CB-CG	8.04	1.74	1.52
1	A	260	SER	CB-OG	-8.01	1.31	1.42
1	A	342	GLU	CD-OE1	7.01	1.33	1.25
1	A	446	ALA	CA-CB	-6.96	1.37	1.52
1	B	36	SER	CB-OG	6.72	1.50	1.42
1	B	51	GLU	CG-CD	-6.71	1.41	1.51
1	A	411	TYR	CE2-CZ	6.53	1.47	1.38
1	A	508	ARG	CZ-NH2	-6.42	1.24	1.33
1	A	341	ARG	CD-NE	6.32	1.57	1.46
1	A	411	TYR	CG-CD1	6.22	1.47	1.39
1	B	131	GLU	CD-OE2	-6.10	1.19	1.25
1	B	386	VAL	CB-CG2	-6.08	1.40	1.52
1	A	118	SER	CB-OG	6.04	1.50	1.42
1	A	250	ARG	CZ-NH2	5.99	1.40	1.33
1	B	314	GLU	CB-CG	5.98	1.63	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	430	ARG	CB-CG	-5.97	1.36	1.52
1	A	167	GLU	CD-OE2	5.96	1.32	1.25
1	B	193	ARG	CZ-NH2	5.96	1.40	1.33
1	A	508	ARG	CD-NE	-5.82	1.36	1.46
1	A	341	ARG	CG-CD	5.67	1.66	1.51
1	B	306	GLU	CD-OE2	-5.66	1.19	1.25
1	B	48	PHE	CG-CD1	5.62	1.47	1.38
1	B	411	TYR	CE2-CZ	5.57	1.45	1.38
1	A	425	ALA	CA-CB	-5.53	1.40	1.52
1	B	394	THR	C-O	5.51	1.33	1.23
1	A	365	SER	CB-OG	-5.47	1.35	1.42
1	B	50	ARG	CZ-NH1	5.19	1.39	1.33
1	A	84	TYR	CB-CG	5.10	1.59	1.51
1	A	468	SER	CB-OG	5.09	1.48	1.42
1	A	400	GLU	CD-OE2	5.08	1.31	1.25
1	A	446	ALA	C-O	5.07	1.32	1.23
1	B	248	VAL	CB-CG2	-5.06	1.42	1.52
1	A	3	GLU	CD-OE2	5.05	1.31	1.25
1	B	226	PHE	CD2-CE2	5.03	1.49	1.39

All (81) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	341	ARG	NE-CZ-NH1	-31.43	104.58	120.30
1	A	412	ARG	NE-CZ-NH2	-31.08	104.76	120.30
1	A	412	ARG	NE-CZ-NH1	30.00	135.30	120.30
1	A	341	ARG	NE-CZ-NH2	12.24	126.42	120.30
1	A	508	ARG	NE-CZ-NH2	-11.25	114.67	120.30
1	B	58	ARG	NE-CZ-NH1	10.78	125.69	120.30
1	B	227	ASP	CB-CG-OD2	-10.42	108.93	118.30
1	A	241	LEU	CA-CB-CG	10.22	138.81	115.30
1	B	412	ARG	NE-CZ-NH2	-9.90	115.35	120.30
1	B	501	ARG	NE-CZ-NH2	-9.66	115.47	120.30
1	A	127	ASP	CB-CG-OD1	9.52	126.87	118.30
1	A	259	ASP	CB-CG-OD2	9.48	126.83	118.30
1	A	68	ARG	NE-CZ-NH2	-9.31	115.64	120.30
1	A	278	ASP	CB-CG-OD2	-8.83	110.35	118.30
1	B	300	ARG	NE-CZ-NH2	-8.83	115.89	120.30
1	B	412	ARG	NE-CZ-NH1	8.81	124.70	120.30
1	A	341	ARG	NH1-CZ-NH2	8.69	128.96	119.40
1	A	200	ASP	CB-CG-OD2	-8.63	110.53	118.30
1	A	149	ARG	NE-CZ-NH2	-8.45	116.08	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	260	SER	N-CA-CB	-8.38	97.93	110.50
1	A	226	PHE	CB-CG-CD1	8.30	126.61	120.80
1	A	301	ARG	NE-CZ-NH2	-8.23	116.19	120.30
1	A	490	ARG	NE-CZ-NH1	8.10	124.35	120.30
1	B	71	ASP	CB-CG-OD2	-8.09	111.02	118.30
1	A	65	ARG	NE-CZ-NH1	8.06	124.33	120.30
1	A	345	GLU	CG-CD-OE1	7.92	134.13	118.30
1	B	0	HIS	N-CA-C	-7.78	90.00	111.00
1	B	457	ASP	CB-CG-OD1	7.75	125.28	118.30
1	A	412	ARG	CD-NE-CZ	7.74	134.43	123.60
1	A	448	LEU	CB-CG-CD2	7.48	123.72	111.00
1	B	508[A]	ARG	NE-CZ-NH2	-7.47	116.56	120.30
1	B	508[B]	ARG	NE-CZ-NH2	-7.47	116.56	120.30
1	A	341	ARG	CB-CG-CD	7.20	130.32	111.60
1	A	65	ARG	NE-CZ-NH2	-6.98	116.81	120.30
1	A	457	ASP	CB-CG-OD1	6.95	124.55	118.30
1	A	329	ARG	NE-CZ-NH2	-6.86	116.87	120.30
1	A	226	PHE	CB-CG-CD2	-6.84	116.01	120.80
1	A	126	ARG	NE-CZ-NH2	-6.73	116.93	120.30
1	A	221	ASP	CB-CG-OD1	6.62	124.26	118.30
1	B	221	ASP	CB-CG-OD1	6.55	124.20	118.30
1	A	345	GLU	CG-CD-OE2	-6.47	105.35	118.30
1	B	260	SER	N-CA-CB	-6.36	100.96	110.50
1	A	259	ASP	N-CA-CB	6.33	121.99	110.60
1	A	278	ASP	CB-CG-OD1	6.26	123.94	118.30
1	B	457	ASP	CB-CG-OD2	-6.21	112.72	118.30
1	B	431	ARG	NE-CZ-NH2	-6.09	117.25	120.30
1	B	450	ARG	NE-CZ-NH1	6.06	123.33	120.30
1	B	506	TYR	CB-CG-CD2	6.04	124.62	121.00
1	B	445	ASP	CB-CG-OD1	-5.98	112.92	118.30
1	A	68	ARG	NE-CZ-NH1	5.92	123.26	120.30
1	A	373	ASP	CB-CG-OD2	-5.78	113.10	118.30
1	A	391	ASP	CB-CG-OD1	5.75	123.48	118.30
1	A	375	ASN	CB-CA-C	5.74	121.89	110.40
1	B	44	ARG	NE-CZ-NH2	-5.71	117.45	120.30
1	B	131	GLU	OE1-CD-OE2	5.70	130.14	123.30
1	A	41	ASP	CB-CG-OD2	-5.69	113.18	118.30
1	B	241	LEU	CA-CB-CG	5.66	128.31	115.30
1	B	461	ARG	NE-CZ-NH1	5.63	123.12	120.30
1	A	301	ARG	NE-CZ-NH1	5.62	123.11	120.30
1	A	375	ASN	N-CA-CB	-5.59	100.54	110.60
1	B	241	LEU	CB-CG-CD2	-5.54	101.58	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	219	LEU	CB-CG-CD2	5.42	120.22	111.00
1	A	373	ASP	CB-CG-OD1	5.39	123.15	118.30
1	A	158	PHE	CB-CG-CD2	-5.39	117.03	120.80
1	B	331	ASP	CB-CG-OD1	5.36	123.12	118.30
1	A	505	PHE	CB-CG-CD2	-5.35	117.06	120.80
1	A	355	ARG	NE-CZ-NH1	-5.28	117.66	120.30
1	B	52	ASP	CB-CG-OD1	5.24	123.02	118.30
1	A	501	ARG	NE-CZ-NH1	5.21	122.90	120.30
1	A	400	GLU	OE1-CD-OE2	5.17	129.50	123.30
1	B	197	ASP	CB-CG-OD1	5.16	122.94	118.30
1	A	25	ASP	CB-CG-OD1	5.14	122.93	118.30
1	A	286	ARG	NE-CZ-NH2	-5.13	117.73	120.30
1	A	241	LEU	CB-CG-CD2	-5.13	102.28	111.00
1	B	501	ARG	NE-CZ-NH1	5.13	122.86	120.30
1	B	83	TYR	CB-CG-CD2	-5.11	117.94	121.00
1	B	115	LEU	CB-CG-CD2	-5.10	102.33	111.00
1	A	420	LEU	CB-CG-CD2	5.09	119.66	111.00
1	B	412	ARG	CG-CD-NE	-5.09	101.11	111.80
1	B	0	HIS	C-N-CA	5.04	134.31	121.70
1	B	279	LEU	CB-CG-CD1	-5.03	102.44	111.00

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	3905	0	3908	111	0
1	B	3884	0	3892	91	0
2	A	12	0	16	0	0
2	B	12	0	16	0	0
3	A	48	0	24	24	0
3	B	48	0	24	23	0
4	A	797	0	0	51	3
4	B	738	0	0	37	4
All	All	9444	0	7880	218	4

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 14.

All (218) 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:508[A]:ARG:HG2	4:B:2709:HOH:O	1.26	1.32
1:A:402:PHE:CZ	3:A:1534:NAP:O2D	1.84	1.30
1:B:286:ARG:HD3	4:B:2462:HOH:O	1.29	1.30
1:A:416:ASP:HB3	4:A:2686:HOH:O	1.16	1.27
1:A:249:GLN:HG2	4:A:2480:HOH:O	1.28	1.26
1:A:310:GLU:HG3	4:A:2525:HOH:O	1.30	1.22
3:B:1527:NAP:H6N	4:B:2724:HOH:O	1.36	1.22
1:A:278:ASP:HB3	4:A:2503:HOH:O	1.37	1.22
1:A:241:LEU:HD12	4:A:2470:HOH:O	1.35	1.20
1:A:0:HIS:CE1	4:A:2004:HOH:O	1.99	1.15
1:A:259:ASP:HB2	3:A:1534:NAP:C3N	1.78	1.13
1:B:523:THR:HG21	4:B:2324:HOH:O	1.49	1.10
1:B:286:ARG:NH2	1:B:478[A]:ASN:HD21	1.51	1.09
3:A:1534:NAP:C5D	4:A:2795:HOH:O	2.02	1.08
1:A:241:LEU:HB2	4:A:2470:HOH:O	1.52	1.08
1:A:286:ARG:NH2	1:A:478[A]:ASN:HD21	1.52	1.05
1:B:357[B]:GLU:OE2	1:B:394:THR:OG1	1.76	1.03
1:A:430:ARG:NH1	4:A:2695:HOH:O	1.93	1.01
1:A:248:VAL:HB	4:A:2476:HOH:O	0.81	0.99
1:B:296:CYS:SG	3:B:1527:NAP:O7N	2.22	0.98
1:A:420:LEU:N	4:A:2690:HOH:O	1.95	0.98
1:B:509:ARG:NH2	4:B:2716:HOH:O	1.96	0.98
1:A:478[B]:ASN:ND2	1:A:480:MET:CE	2.27	0.98
3:B:1527:NAP:O1N	3:B:1527:NAP:C5B	2.12	0.96
3:B:1527:NAP:O3B	4:B:2730:HOH:O	1.85	0.95
1:A:484:LEU:HD21	1:B:508[A]:ARG:HH22	1.34	0.93
4:A:2467:HOH:O	1:B:248:VAL:HB	0.74	0.92
1:A:519:ILE:O	1:A:523:THR:HG23	1.68	0.92
1:A:478[B]:ASN:HD21	1:A:480:MET:HE1	1.34	0.91
1:A:478[B]:ASN:ND2	1:A:480:MET:HE1	1.85	0.91
1:A:356:GLU:HG3	4:A:2588:HOH:O	1.71	0.90
1:A:0:HIS:ND1	1:A:0:HIS:O	2.03	0.90
1:B:519:ILE:O	1:B:523:THR:HG23	1.70	0.90
1:A:375:ASN:HB3	4:A:2635:HOH:O	1.72	0.89
3:A:1534:NAP:H52N	4:A:2795:HOH:O	1.66	0.89
1:A:235:ALA:H	3:A:1534:NAP:H71N	1.21	0.89
1:B:478[B]:ASN:ND2	1:B:480:MET:HE1	1.90	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:233:GLY:O	3:B:1527:NAP:N7N	2.08	0.86
1:A:484:LEU:HD21	1:B:508[A]:ARG:NH2	1.90	0.85
3:B:1527:NAP:O1N	3:B:1527:NAP:H51A	1.77	0.84
1:B:420:LEU:N	4:B:2626:HOH:O	2.10	0.83
1:B:286:ARG:NH2	1:B:478[A]:ASN:ND2	2.27	0.83
1:B:478[B]:ASN:ND2	1:B:480:MET:CE	2.41	0.83
1:A:375:ASN:CB	4:A:2635:HOH:O	2.28	0.82
1:A:478[B]:ASN:ND2	1:A:480:MET:HE3	1.95	0.81
1:A:495:GLU:OE2	1:B:508[A]:ARG:NH1	2.12	0.81
1:B:248:VAL:HG21	4:B:2427:HOH:O	1.80	0.81
1:A:286:ARG:NH2	1:A:478[A]:ASN:ND2	2.29	0.80
1:B:149:ARG:NH1	4:B:2333:HOH:O	1.76	0.80
1:A:158:PHE:CD2	3:A:1534:NAP:H51N	2.17	0.79
1:B:450:ARG:NE	4:B:2652:HOH:O	2.10	0.79
1:A:255:ASN:HD21	1:A:495:GLU:H	1.31	0.78
1:B:6:LYS:NZ	4:B:2024:HOH:O	2.15	0.78
1:A:157:ALA:HB1	3:A:1534:NAP:H3D	1.66	0.78
1:A:356:GLU:CG	4:A:2588:HOH:O	2.29	0.77
1:A:286:ARG:HH22	1:A:478[A]:ASN:HD21	1.33	0.76
1:A:0:HIS:HE1	4:A:2004:HOH:O	1.51	0.76
1:B:383:HIS:HE1	4:B:2555:HOH:O	1.69	0.74
1:B:508[A]:ARG:CG	4:B:2709:HOH:O	1.98	0.74
1:B:255:ASN:HD21	1:B:495:GLU:H	1.32	0.74
1:A:341:ARG:HG2	4:A:2342:HOH:O	1.85	0.74
1:B:258:ALA:HA	3:B:1527:NAP:H72N	1.52	0.74
3:B:1527:NAP:O1N	3:B:1527:NAP:H52A	1.88	0.73
3:B:1527:NAP:H52N	3:B:1527:NAP:O2A	1.89	0.73
1:B:523:THR:CG2	4:B:2324:HOH:O	2.21	0.73
3:B:1527:NAP:C6N	4:B:2724:HOH:O	2.11	0.72
1:A:259:ASP:CB	3:A:1534:NAP:C3N	2.59	0.72
1:A:478[A]:ASN:ND2	4:A:2740:HOH:O	2.23	0.72
1:A:235:ALA:N	3:A:1534:NAP:H71N	1.87	0.71
1:A:356:GLU:OE2	4:A:2588:HOH:O	2.07	0.71
1:A:234:SER:HA	3:A:1534:NAP:H71N	1.55	0.71
1:B:416:ASP:C	4:B:2621:HOH:O	2.29	0.71
1:A:282:LYS:HE3	4:A:2503:HOH:O	1.91	0.70
1:A:375:ASN:CB	4:A:2634:HOH:O	2.39	0.70
1:B:244:HIS:O	1:B:248:VAL:HG22	1.91	0.70
1:A:375:ASN:HB3	4:A:2634:HOH:O	1.92	0.70
1:B:478[A]:ASN:ND2	4:B:2681:HOH:O	2.24	0.70
1:A:259:ASP:HB2	3:A:1534:NAP:C2N	2.21	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:383:HIS:HE1	4:A:2609:HOH:O	1.75	0.69
1:B:508[A]:ARG:NH1	4:B:2711:HOH:O	2.24	0.69
1:A:286:ARG:HH22	1:A:478[A]:ASN:ND2	1.88	0.68
1:A:490:ARG:HD3	4:A:2749:HOH:O	1.92	0.68
3:B:1527:NAP:O2D	3:B:1527:NAP:C2N	2.36	0.68
1:B:245:PRO:HA	1:B:248:VAL:CG2	2.23	0.68
1:A:523:THR:HG21	4:A:2359:HOH:O	1.94	0.67
1:B:450:ARG:NH2	4:B:2653:HOH:O	2.26	0.67
1:A:375:ASN:CG	4:A:2635:HOH:O	2.32	0.67
1:A:158:PHE:CE2	3:A:1534:NAP:H4D	2.31	0.66
1:A:439:SER:OG	1:A:476:HIS:HD2	1.80	0.65
1:B:51:GLU:OE2	4:B:2135:HOH:O	2.13	0.65
1:A:402:PHE:CE1	3:A:1534:NAP:O2D	2.45	0.64
1:A:259:ASP:OD2	4:A:2491:HOH:O	2.14	0.64
1:A:341:ARG:CG	4:A:2571:HOH:O	2.45	0.64
1:B:258:ALA:HA	3:B:1527:NAP:N7N	2.14	0.63
1:A:286:ARG:HH21	1:A:478[A]:ASN:HD21	1.43	0.63
1:A:375:ASN:ND2	4:A:2638:HOH:O	2.21	0.63
3:B:1527:NAP:O2A	3:B:1527:NAP:C5D	2.47	0.62
1:B:478[B]:ASN:ND2	1:B:480:MET:HE3	2.14	0.62
1:A:402:PHE:CE2	3:A:1534:NAP:O2D	2.46	0.62
1:B:439:SER:OG	1:B:476:HIS:HD2	1.83	0.61
1:B:245:PRO:HA	1:B:248:VAL:HG23	1.82	0.61
1:B:478[B]:ASN:HD21	1:B:480:MET:HE1	1.62	0.61
3:B:1527:NAP:H52N	4:B:2727:HOH:O	2.00	0.60
1:A:341:ARG:HG2	4:A:2571:HOH:O	2.02	0.59
1:A:0:HIS:O	1:A:0:HIS:CG	2.55	0.59
1:B:123:HIS:HE1	4:B:2273:HOH:O	1.85	0.59
1:A:123:HIS:HE1	4:A:2311:HOH:O	1.85	0.59
1:A:294:GLN:HE22	1:A:339:VAL:H	1.48	0.58
1:B:224:ARG:NH1	4:B:2401:HOH:O	2.36	0.58
1:A:440:ILE:HG13	1:A:448:LEU:HD22	1.84	0.58
1:B:220:LEU:HD11	1:B:241:LEU:HG	1.84	0.58
1:A:381:ALA:O	1:A:383:HIS:HD2	1.87	0.58
1:A:248:VAL:CG2	4:A:2476:HOH:O	2.13	0.57
1:A:470:GLN:HG2	4:A:2734:HOH:O	2.04	0.57
1:B:381:ALA:O	1:B:383:HIS:HD2	1.88	0.56
1:A:234:SER:CA	3:A:1534:NAP:H71N	2.18	0.56
1:A:415:THR:C	1:A:416:ASP:OD1	2.44	0.56
1:A:0:HIS:CG	4:A:2003:HOH:O	2.58	0.56
1:A:241:LEU:CB	4:A:2470:HOH:O	2.25	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:ASP:OD1	4:A:2205:HOH:O	2.18	0.56
1:A:258:ALA:HA	3:A:1534:NAP:C5N	2.36	0.56
1:A:258:ALA:HA	3:A:1534:NAP:H5N	1.89	0.55
1:B:281:ILE:HG22	1:B:319:LYS:HE3	1.88	0.55
1:A:255:ASN:HD21	1:A:495:GLU:N	2.03	0.54
1:A:463:HIS:HE1	1:A:473:GLN:OE1	1.90	0.54
1:B:415:THR:HG23	4:B:2620:HOH:O	2.07	0.53
1:B:294:GLN:HE22	1:B:339:VAL:H	1.56	0.53
1:A:345:GLU:OE2	4:A:2577:HOH:O	0.60	0.53
1:B:313:LEU:HD21	1:B:386:VAL:HG22	1.90	0.53
1:A:200:ASP:OD1	4:A:2419:HOH:O	2.19	0.53
1:A:258:ALA:CA	3:A:1534:NAP:H5N	2.34	0.52
1:A:304:VAL:O	1:A:410:PRO:HA	2.09	0.52
1:B:463:HIS:HE1	1:B:473:GLN:OE1	1.94	0.51
1:A:0:HIS:CD2	4:A:2003:HOH:O	2.62	0.51
1:A:241:LEU:CD1	4:A:2470:HOH:O	2.05	0.51
1:B:400:GLU:HA	4:B:2603:HOH:O	2.11	0.51
1:A:158:PHE:CD1	1:A:339:VAL:HG21	2.47	0.50
1:B:415:THR:O	1:B:416:ASP:CB	2.58	0.50
1:B:508[A]:ARG:HG3	1:B:509:ARG:N	2.25	0.50
1:A:249:GLN:NE2	4:A:2478:HOH:O	2.44	0.50
1:A:375:ASN:HB3	4:A:2637:HOH:O	2.10	0.50
1:B:415:THR:O	1:B:416:ASP:HB2	2.12	0.50
3:A:1534:NAP:H51N	4:A:2795:HOH:O	1.90	0.50
1:A:182:LYS:NZ	3:A:1534:NAP:O1X	2.45	0.49
1:A:471:HIS:CD2	4:A:2733:HOH:O	2.66	0.49
3:A:1534:NAP:O5B	3:A:1534:NAP:O5D	2.30	0.49
1:B:217:ALA:HB2	4:B:2383:HOH:O	2.12	0.49
1:B:470:GLN:HG2	4:B:2665:HOH:O	2.13	0.49
1:B:224:ARG:CG	1:B:224:ARG:HH11	2.25	0.49
1:B:334:ARG:NH1	1:B:334:ARG:HG3	2.27	0.49
1:B:334:ARG:NH2	4:B:2506:HOH:O	2.46	0.49
1:A:415:THR:O	1:A:416:ASP:OD1	2.30	0.49
1:B:319:LYS:HE2	1:B:319:LYS:HB2	1.66	0.48
1:B:356:GLU:OE1	1:B:356:GLU:HA	2.13	0.48
1:A:522:LEU:C	1:A:522:LEU:HD12	2.33	0.48
1:A:521:THR:HG21	1:B:446:ALA:HB1	1.94	0.48
1:A:341:ARG:HG3	4:A:2571:HOH:O	2.09	0.47
1:B:439:SER:OG	1:B:476:HIS:CD2	2.66	0.47
1:A:420:LEU:HD11	1:A:451:LEU:HA	1.96	0.47
1:A:310:GLU:CG	4:A:2525:HOH:O	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:286:ARG:CD	4:B:2462:HOH:O	2.14	0.47
1:B:233:GLY:HA2	3:B:1527:NAP:H2N	1.97	0.46
1:B:509:ARG:CZ	4:B:2716:HOH:O	2.54	0.46
1:A:521:THR:HG21	1:B:446:ALA:CB	2.45	0.46
1:A:220:LEU:HD11	1:A:241:LEU:HG	1.97	0.46
3:B:1527:NAP:H2N	3:B:1527:NAP:H2D	1.35	0.45
1:B:297:THR:O	1:B:476:HIS:CE1	2.69	0.45
1:B:286:ARG:HH21	1:B:478[A]:ASN:HD21	1.55	0.45
3:B:1527:NAP:O5D	3:B:1527:NAP:O2A	2.35	0.45
1:B:334:ARG:NH1	1:B:334:ARG:CG	2.79	0.45
1:B:297:THR:O	1:B:476:HIS:HE1	1.99	0.45
1:A:233:GLY:O	1:A:258:ALA:HA	2.17	0.45
1:B:258:ALA:CA	3:B:1527:NAP:H72N	2.23	0.45
1:B:255:ASN:HD21	1:B:495:GLU:N	2.09	0.45
1:A:433:GLN:NE2	4:A:2701:HOH:O	2.50	0.45
1:A:297:THR:O	1:A:476:HIS:CE1	2.70	0.45
3:B:1527:NAP:O1X	4:B:2728:HOH:O	2.21	0.45
1:B:439:SER:CB	1:B:476:HIS:HD2	2.29	0.45
1:A:297:THR:O	1:A:476:HIS:HE1	1.99	0.45
3:A:1534:NAP:H6N	3:A:1534:NAP:H2D	1.58	0.44
1:A:439:SER:CB	1:A:476:HIS:HD2	2.30	0.44
1:A:158:PHE:CE2	3:A:1534:NAP:H51N	2.52	0.44
3:B:1527:NAP:O1N	3:B:1527:NAP:H3D	2.18	0.44
1:A:0:HIS:HA	1:A:1:MET:HA	1.60	0.44
1:A:286:ARG:NH1	4:A:2511:HOH:O	2.22	0.44
1:B:334:ARG:CG	1:B:334:ARG:HH11	2.30	0.44
1:B:192:GLN:HA	1:B:211:ILE:HD13	2.00	0.44
1:A:381:ALA:O	1:A:383:HIS:CD2	2.70	0.43
3:B:1527:NAP:H3D	4:B:2734:HOH:O	2.16	0.43
1:A:245:PRO:HA	1:A:248:VAL:HG22	2.00	0.43
1:B:430:ARG:HG3	4:B:2662:HOH:O	2.17	0.43
1:B:123:HIS:CE1	4:B:2273:HOH:O	2.66	0.43
1:A:509:ARG:NH1	1:B:495:GLU:OE1	2.52	0.43
1:B:158:PHE:CD1	1:B:339:VAL:HG21	2.55	0.42
1:B:304:VAL:O	1:B:410:PRO:HA	2.19	0.42
1:A:259:ASP:CG	1:A:260:SER:H	2.23	0.42
1:A:439:SER:OG	1:A:476:HIS:CD2	2.66	0.42
1:A:234:SER:HA	3:A:1534:NAP:N7N	2.30	0.42
1:A:25:ASP:HA	1:A:89:ALA:O	2.20	0.42
1:B:126:ARG:HD3	1:B:523:THR:HG22	2.02	0.42
1:B:522:LEU:HD23	1:B:522:LEU:HA	1.82	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:415:THR:CG2	4:A:2675:HOH:O	2.69	0.41
1:A:341:ARG:HH11	1:A:341:ARG:HD2	1.44	0.41
1:A:439:SER:HA	1:A:463:HIS:O	2.21	0.41
1:B:286:ARG:NH1	4:B:2460:HOH:O	2.49	0.41
1:B:313:LEU:HD21	1:B:386:VAL:CG2	2.49	0.41
1:B:438:ALA:HB2	1:B:459:HIS:CD2	2.56	0.41
1:B:414:THR:HG22	1:B:416:ASP:H	1.85	0.41
3:B:1527:NAP:C5B	3:B:1527:NAP:PN	3.09	0.40
3:B:1527:NAP:C5D	4:B:2727:HOH:O	2.64	0.40
1:B:245:PRO:HA	1:B:248:VAL:HG22	2.00	0.40
1:B:460:GLY:HA3	1:B:486:GLY:O	2.21	0.40
1:B:233:GLY:O	1:B:258:ALA:HA	2.21	0.40
1:B:394:THR:OG1	4:B:2594:HOH:O	2.21	0.40
1:A:420:LEU:HD13	1:A:451:LEU:HD13	2.04	0.40
1:B:281:ILE:CG2	1:B:319:LYS:HE3	2.49	0.40
1:B:381:ALA:O	1:B:383:HIS:CD2	2.72	0.40

All (4) 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)
4:A:2238:HOH:O	4:B:2192:HOH:O[1_444]	2.05	0.15
4:A:2109:HOH:O	4:B:2003:HOH:O[1_444]	2.11	0.09
4:B:2431:HOH:O	4:B:2665:HOH:O[1_455]	2.17	0.03
4:A:2239:HOH:O	4:B:2189:HOH:O[1_444]	2.18	0.02

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	526/534 (98%)	518 (98%)	8 (2%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	523/534 (98%)	514 (98%)	9 (2%)	0	100	100
All	All	1049/1068 (98%)	1032 (98%)	17 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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	397/399 (100%)	392 (99%)	5 (1%)	73	46
1	B	396/399 (99%)	387 (98%)	9 (2%)	56	20
All	All	793/798 (99%)	779 (98%)	14 (2%)	64	30

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

Mol	Chain	Res	Type
1	A	0	HIS
1	A	259	ASP
1	A	400	GLU
1	A	415	THR
1	A	482	MET
1	B	44	ARG
1	B	51	GLU
1	B	224	ARG
1	B	241	LEU
1	B	319	LYS
1	B	334	ARG
1	B	394	THR
1	B	415	THR
1	B	482	MET

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	0	HIS
1	A	90	ASN
1	A	123	HIS
1	A	249	GLN
1	A	255	ASN
1	A	294	GLN
1	A	346	ASN
1	A	383	HIS
1	A	433	GLN
1	A	463	HIS
1	A	476	HIS
1	A	513	GLN
1	B	90	ASN
1	B	123	HIS
1	B	255	ASN
1	B	294	GLN
1	B	343	GLN
1	B	346	ASN
1	B	383	HIS
1	B	392	ASN
1	B	433	GLN
1	B	463	HIS
1	B	476	HIS
1	B	513	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

6 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	$\# Z > 2$
2	GOL	A	1532	-	5,5,5	0.93	0	5,5,5	0.57	0
2	GOL	A	1533	-	5,5,5	0.77	0	5,5,5	0.79	0
3	NAP	A	1534	-	44,52,52	1.21	3 (6%)	51,80,80	1.81	4 (7%)
3	NAP	B	1527	-	44,52,52	3.09	14 (31%)	51,80,80	3.35	16 (31%)
2	GOL	B	1528	-	5,5,5	0.82	0	5,5,5	0.88	0
2	GOL	B	1529	-	5,5,5	0.87	0	5,5,5	0.57	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	1532	-	-	0/4/4/4	0/0/0/0
2	GOL	A	1533	-	-	0/4/4/4	0/0/0/0
3	NAP	A	1534	-	-	0/27/67/67	0/5/5/5
3	NAP	B	1527	-	-	0/27/67/67	0/5/5/5
2	GOL	B	1528	-	-	0/4/4/4	0/0/0/0
2	GOL	B	1529	-	-	0/4/4/4	0/0/0/0

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1527	NAP	O3B-C3B	-9.09	1.22	1.43
3	B	1527	NAP	O4B-C1B	-8.39	1.29	1.41
3	B	1527	NAP	C4A-N3A	-6.26	1.26	1.35
3	B	1527	NAP	O4D-C4D	-5.55	1.32	1.45
3	B	1527	NAP	C5A-C4A	-4.02	1.31	1.40
3	B	1527	NAP	O4B-C4B	-3.03	1.38	1.45
3	B	1527	NAP	C5A-N7A	-2.99	1.29	1.39
3	B	1527	NAP	P2B-O1X	2.89	1.60	1.50
3	B	1527	NAP	PA-O1A	2.90	1.61	1.50
3	B	1527	NAP	C8A-N7A	2.96	1.40	1.34
3	A	1534	NAP	PN-O1N	3.04	1.62	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1534	NAP	PA-O1A	3.06	1.62	1.50
3	A	1534	NAP	P2B-O1X	3.36	1.62	1.50
3	B	1527	NAP	C3D-C4D	3.49	1.62	1.53
3	B	1527	NAP	P2B-O2B	3.98	1.66	1.59
3	B	1527	NAP	O4D-C1D	4.58	1.47	1.41
3	B	1527	NAP	O3D-C3D	7.51	1.60	1.43

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1527	NAP	C4D-O4D-C1D	-13.50	95.40	109.77
3	B	1527	NAP	N3A-C2A-N1A	-9.03	121.00	128.86
3	A	1534	NAP	N3A-C2A-N1A	-8.19	121.73	128.86
3	B	1527	NAP	C5B-C4B-C3B	-7.55	86.50	115.29
3	B	1527	NAP	O4D-C4D-C5D	-6.38	87.85	109.40
3	B	1527	NAP	O4D-C4D-C3D	-5.56	94.12	105.17
3	A	1534	NAP	C4D-O4D-C1D	-5.36	104.06	109.77
3	A	1534	NAP	C4B-O4B-C1B	-5.34	104.08	109.77
3	B	1527	NAP	O4B-C1B-C2B	-3.35	100.72	106.59
3	B	1527	NAP	C1B-N9A-C4A	-3.26	121.00	126.64
3	B	1527	NAP	O2D-C2D-C1D	-2.67	103.27	111.61
3	B	1527	NAP	C5N-C6N-N1N	-2.51	116.54	120.40
3	A	1534	NAP	C3N-C7N-N7N	-2.45	114.97	117.77
3	B	1527	NAP	O2B-C2B-C3B	2.02	119.12	111.63
3	B	1527	NAP	O3D-C3D-C4D	2.94	119.67	111.09
3	B	1527	NAP	C6N-C5N-C4N	3.21	124.28	119.44
3	B	1527	NAP	O2N-PN-O5D	3.63	125.30	108.14
3	B	1527	NAP	C5D-C4D-C3D	4.20	131.31	115.29
3	B	1527	NAP	O2D-C2D-C3D	4.85	127.37	111.83
3	B	1527	NAP	O3D-C3D-C2D	4.87	127.42	111.83

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 47 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1534	NAP	24	0
3	B	1527	NAP	23	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	529/534 (99%)	-0.28	8 (1%) 74 74	4, 8, 18, 43	0
1	B	524/534 (98%)	-0.18	10 (1%) 67 68	4, 9, 22, 50	0
All	All	1053/1068 (98%)	-0.23	18 (1%) 70 71	4, 9, 21, 50	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	523	THR	5.0
1	B	524	GLN	4.9
1	A	0	HIS	4.8
1	B	394	THR	4.4
1	A	415	THR	4.2
1	B	415	THR	4.1
1	B	521	THR	4.1
1	B	0	HIS	3.9
1	A	375	ASN	3.8
1	B	526	THR	3.3
1	B	416	ASP	3.2
1	B	522	LEU	3.1
1	A	416	ASP	2.8
1	A	450	ARG	2.6
1	B	414	THR	2.6
1	A	531	ALA	2.2
1	A	259	ASP	2.0
1	A	331	ASP	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 carbohydrates 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. 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, 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	LLDF	B-factors(\AA^2)	Q<0.9
3	NAP	A	1534	48/48	0.72	0.34	8.22	12,19,28,33	18
3	NAP	B	1527	48/48	0.86	0.23	4.71	5,22,33,38	17
2	GOL	A	1532	6/6	0.97	0.07	0.77	5,9,12,14	0
2	GOL	B	1528	6/6	0.98	0.08	0.29	10,11,13,13	0
2	GOL	B	1529	6/6	0.97	0.07	-0.01	6,10,12,15	0
2	GOL	A	1533	6/6	0.96	0.07	-0.40	8,11,11,12	0

6.5 Other polymers [i](#)

There are no such residues in this entry.