



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2016 – 05:05 PM EDT

PDB ID : 3PDU
Title : Crystal structure of gamma-hydroxybutyrate dehydrogenase from *Geobacter sulfurreducens* in complex with NADP+
Authors : Zhang, Y.; Garavito, R.M.
Deposited on : 2010-10-25
Resolution : 1.89 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027674
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-20027674

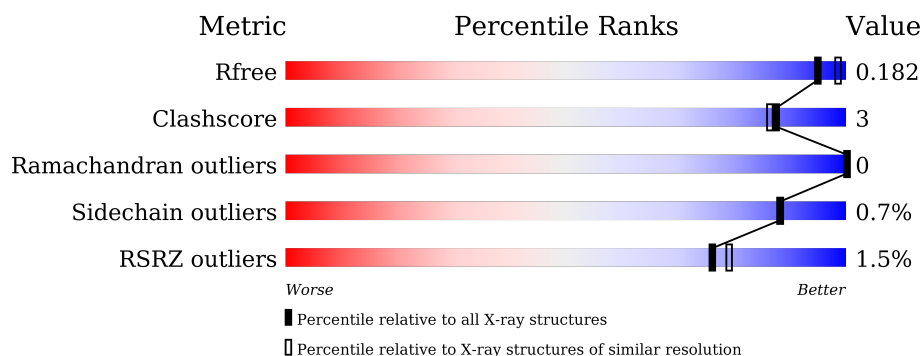
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.89 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.



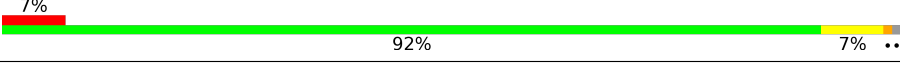
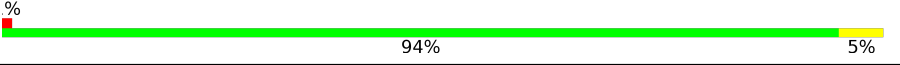
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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	287	<div> <div style="width: 95%;"></div> <div>95%</div> </div>
1	B	287	<div> <div style="width: 97%;"></div> <div>97%</div> </div>
1	C	287	<div> <div style="width: 95%;"></div> <div>95%</div> </div>
1	D	287	<div> <div style="width: 92%;"></div> <div>92%</div> </div>
1	E	287	<div> <div style="width: 95%;"></div> <div>95%</div> </div>
1	F	287	<div> <div style="width: 95%;"></div> <div>95%</div> </div>

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Mol	Chain	Length	Quality of chain
1	G	287	
1	H	287	

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	GOL	A	304	-	-	-	X
3	GOL	B	304	-	-	-	X
3	GOL	B	305	-	-	-	X
3	GOL	D	305	-	-	X	X
3	GOL	F	288	-	-	-	X
3	GOL	G	304	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 18734 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.

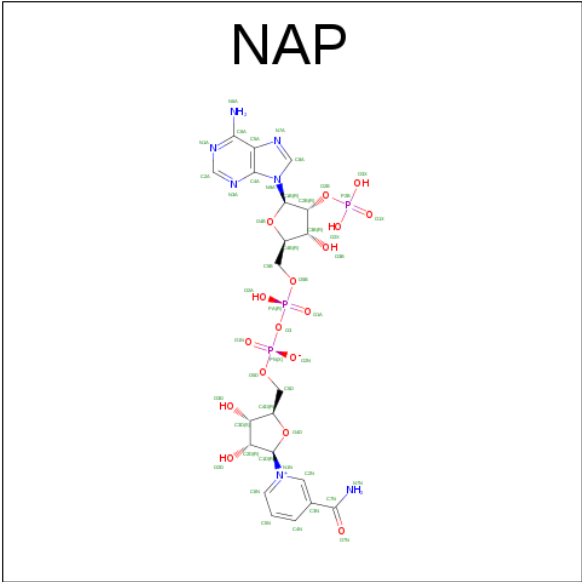
- Molecule 1 is a protein called 3-hydroxyisobutyrate dehydrogenase family protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	287	Total	C	N	O	S	0	4	0
			2082	1302	368	389	23			
1	B	286	Total	C	N	O	S	0	2	0
			2058	1287	364	385	22			
1	C	286	Total	C	N	O	S	0	0	0
			2050	1281	364	385	20			
1	D	287	Total	C	N	O	S	0	1	0
			2061	1288	365	386	22			
1	E	287	Total	C	N	O	S	0	3	0
			2070	1294	365	389	22			
1	F	286	Total	C	N	O	S	0	2	0
			2060	1288	364	386	22			
1	G	284	Total	C	N	O	S	0	1	0
			2041	1276	362	382	21			
1	H	286	Total	C	N	O	S	0	1	0
			2053	1283	364	385	21			

There are 8 discrepancies between the modelled and reference sequences:

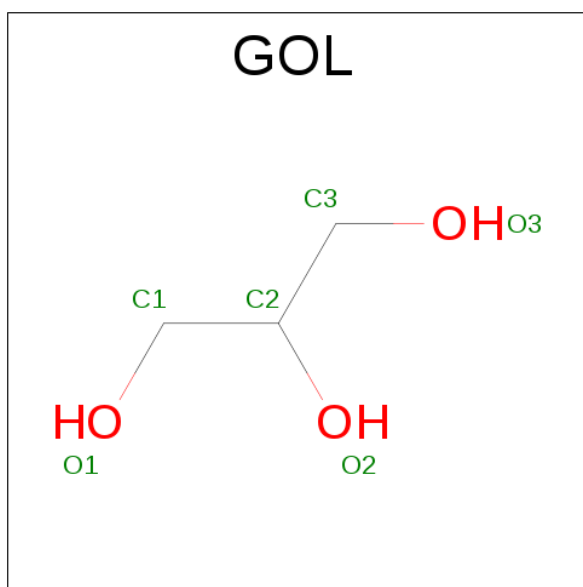
Chain	Residue	Modelled	Actual	Comment	Reference
A	287	ALA	-	EXPRESSION TAG	UNP Q74DE4
B	287	ALA	-	EXPRESSION TAG	UNP Q74DE4
C	287	ALA	-	EXPRESSION TAG	UNP Q74DE4
D	287	ALA	-	EXPRESSION TAG	UNP Q74DE4
E	287	ALA	-	EXPRESSION TAG	UNP Q74DE4
F	287	ALA	-	EXPRESSION TAG	UNP Q74DE4
G	287	ALA	-	EXPRESSION TAG	UNP Q74DE4
H	287	ALA	-	EXPRESSION TAG	UNP Q74DE4

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	C	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	D	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	E	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	F	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	G	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	H	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	E	1	Total	C	O	0	0
			6	3	3		
3	F	1	Total	C	O	0	0
			6	3	3		
3	F	1	Total	C	O	0	0
			6	3	3		
3	G	1	Total	C	O	0	0
			6	3	3		
3	H	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	288	Total	O	0	0
			288	288		

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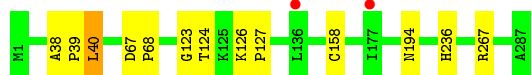
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	253	Total 253	O 253	0	0
4	C	199	Total 199	O 199	0	0
4	D	245	Total 245	O 245	0	0
4	E	222	Total 222	O 222	0	0
4	F	274	Total 274	O 274	0	0
4	G	97	Total 97	O 97	0	0
4	H	231	Total 231	O 231	0	0

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 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: 3-hydroxyisobutyrate dehydrogenase family protein



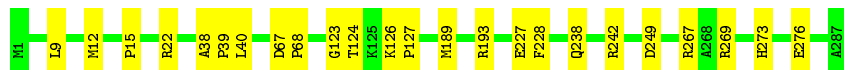
- Molecule 1: 3-hydroxyisobutyrate dehydrogenase family protein



- Molecule 1: 3-hydroxyisobutyrate dehydrogenase family protein



- Molecule 1: 3-hydroxyisobutyrate dehydrogenase family protein

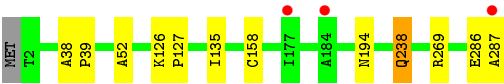


- Molecule 1: 3-hydroxyisobutyrate dehydrogenase family protein

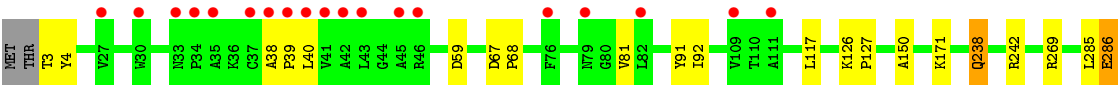
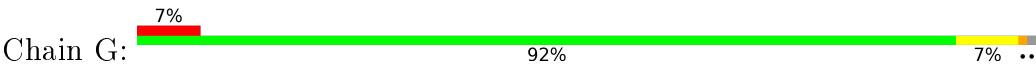


- Molecule 1: 3-hydroxyisobutyrate dehydrogenase family protein

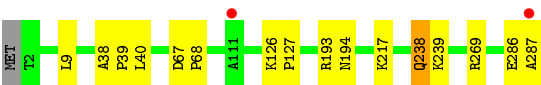




● Molecule 1: 3-hydroxyisobutyrate dehydrogenase family protein



● Molecule 1: 3-hydroxyisobutyrate dehydrogenase family protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	100.03Å 182.07Å 147.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.81 – 1.89 35.02 – 1.89	Depositor EDS
% Data completeness (in resolution range)	97.3 (28.81-1.89) 97.3 (35.02-1.89)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.23 (at 1.89Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, R_{free}	0.150 , 0.186 0.144 , 0.182	Depositor DCC
R_{free} test set	6254 reflections (3.00%)	DCC
Wilson B-factor (Å ²)	24.2	Xtriage
Anisotropy	0.082	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 52.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	18734	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.53% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.38	0/2119	0.50	1/2856 (0.0%)
1	B	0.37	0/2095	0.47	0/2825
1	C	0.36	0/2081	0.48	0/2807
1	D	0.36	0/2095	0.50	0/2825
1	E	0.35	0/2110	0.48	0/2845
1	F	0.40	0/2100	0.49	0/2833
1	G	0.27	0/2075	0.42	0/2798
1	H	0.35	0/2087	0.47	0/2815
All	All	0.36	0/16762	0.48	1/22604 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	40	LEU	CA-CB-CG	-7.26	98.61	115.30

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	2082	0	2103	10	0
1	B	2058	0	2073	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2050	0	2059	10	0
1	D	2061	0	2076	27	0
1	E	2070	0	2087	10	0
1	F	2060	0	2076	6	0
1	G	2041	0	2052	14	0
1	H	2053	0	2064	10	0
2	A	48	0	25	3	0
2	B	48	0	25	0	0
2	C	48	0	25	0	0
2	D	48	0	25	2	0
2	E	48	0	25	0	0
2	F	48	0	25	0	0
2	G	48	0	25	0	0
2	H	48	0	25	0	0
3	A	6	0	8	0	0
3	B	12	0	16	0	0
3	C	12	0	16	3	0
3	D	6	0	8	7	0
3	E	6	0	8	1	0
3	F	12	0	16	0	0
3	G	6	0	8	2	0
3	H	6	0	8	0	0
4	A	288	0	0	2	0
4	B	253	0	0	1	0
4	C	199	0	0	0	0
4	D	245	0	0	1	0
4	E	222	0	0	1	0
4	F	274	0	0	0	0
4	G	97	0	0	0	0
4	H	231	0	0	0	0
All	All	18734	0	16878	88	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 3.

All (88) 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)
3:C:288:GOL:H32	1:D:242:ARG:HD2	1.36	1.02
1:C:242:ARG:HD2	3:C:304:GOL:H32	1.48	0.94
1:C:249:ASP:OD2	1:D:242:ARG:HD3	1.73	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:242:ARG:HD3	1:D:249:ASP:OD2	1.73	0.87
3:D:305:GOL:H32	1:F:52:ALA:HB3	1.66	0.77
1:G:242:ARG:HD2	3:G:304:GOL:H32	1.71	0.70
1:D:227:GLU:HA	3:D:305:GOL:H11	1.74	0.69
1:D:238:GLN:HE21	1:D:269:ARG:HH11	1.39	0.68
1:E:3:THR:HG23	1:E:58:CYS:HA	1.79	0.64
1:D:9:LEU:HD13	1:D:40:LEU:HD12	1.81	0.62
1:G:38:ALA:HB3	1:G:39:PRO:HD3	1.82	0.62
1:G:285:LEU:O	1:G:286:GLU:HB2	2.01	0.61
1:C:126:LYS:HB3	1:C:127:PRO:HD3	1.84	0.59
1:E:163[B]:GLU:H	1:E:163[B]:GLU:CD	2.07	0.59
1:A:236:HIS:HD2	4:A:973:HOH:O	1.86	0.58
1:G:3:THR:HB	1:G:59:ASP:OD2	2.04	0.58
1:D:227:GLU:HB3	3:D:305:GOL:H31	1.85	0.57
1:G:126:LYS:HB3	1:G:127:PRO:HD3	1.86	0.57
1:F:126:LYS:HB3	1:F:127:PRO:HD3	1.88	0.56
1:F:238:GLN:HE21	1:F:269:ARG:HH11	1.53	0.56
1:G:67:ASP:HB2	1:G:68:PRO:HD2	1.88	0.56
1:H:238:GLN:HE21	1:H:269:ARG:HH11	1.52	0.56
1:E:3:THR:CG2	1:E:58:CYS:HA	2.37	0.55
1:D:238:GLN:NE2	1:D:269:ARG:HH11	2.05	0.54
1:G:238:GLN:HE21	1:G:269:ARG:HH11	1.55	0.54
1:A:236:HIS:HE1	2:A:301:NAP:O1A	1.91	0.54
1:B:126:LYS:HB3	1:B:127:PRO:HD3	1.90	0.53
1:E:67:ASP:HB2	1:E:68:PRO:HD2	1.90	0.53
1:C:249:ASP:OD2	1:D:242:ARG:CD	2.54	0.53
1:C:38:ALA:HB3	1:C:39:PRO:HD3	1.90	0.53
1:H:9:LEU:HD13	1:H:40:LEU:HD12	1.90	0.52
1:F:38:ALA:HB3	1:F:39:PRO:HD3	1.92	0.52
1:G:4:TYR:CZ	1:G:150:ALA:HB2	2.45	0.51
1:E:238:GLN:HE21	1:E:269:ARG:HH11	1.58	0.51
1:D:228:PHE:H	3:D:305:GOL:C1	2.23	0.51
1:E:1:MET:HG2	4:E:1767:HOH:O	2.11	0.51
1:H:9:LEU:HD13	1:H:40:LEU:CD1	2.40	0.51
1:D:9:LEU:HD13	1:D:40:LEU:CD1	2.41	0.51
1:D:67:ASP:HB2	1:D:68:PRO:HD2	1.92	0.51
1:E:38:ALA:HB3	1:E:39:PRO:HD3	1.93	0.51
1:C:67:ASP:HB2	1:C:68:PRO:HD2	1.92	0.50
1:D:9:LEU:HD22	1:D:40:LEU:CD1	2.42	0.49
1:A:126:LYS:HB3	1:A:127:PRO:HD3	1.94	0.49
1:A:67:ASP:HB2	1:A:68:PRO:HD2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:242:ARG:HD2	3:E:304:GOL:H31	1.94	0.48
1:H:286:GLU:O	1:H:287:ALA:HB3	2.13	0.48
1:D:227:GLU:OE1	3:D:305:GOL:H31	2.13	0.48
1:G:67:ASP:HB2	1:G:68:PRO:CD	2.43	0.48
1:G:242:ARG:HD2	3:G:304:GOL:C3	2.40	0.48
1:D:228:PHE:H	3:D:305:GOL:H2	1.77	0.47
1:A:124:THR:HA	2:A:301:NAP:O7N	2.15	0.47
1:H:38:ALA:HB3	1:H:39:PRO:HD3	1.97	0.47
1:E:67:ASP:HB2	1:E:68:PRO:CD	2.44	0.47
1:A:158[B]:CYS:SG	4:A:1035:HOH:O	2.61	0.47
1:H:126:LYS:HB3	1:H:127:PRO:HD3	1.96	0.46
1:D:126:LYS:HB3	1:D:127:PRO:HD3	1.98	0.46
1:D:67:ASP:HB2	1:D:68:PRO:CD	2.45	0.46
1:A:67:ASP:HB2	1:A:68:PRO:CD	2.46	0.45
1:H:67:ASP:HB2	1:H:68:PRO:HD2	1.98	0.45
1:A:267:ARG:NH1	1:D:267:ARG:HD3	2.32	0.45
1:G:40:LEU:N	1:G:40:LEU:HD22	2.33	0.44
1:B:67:ASP:HB2	1:B:68:PRO:HD2	1.98	0.44
1:D:189:MET:O	1:D:193:ARG:HG3	2.17	0.44
1:D:9:LEU:HD22	1:D:40:LEU:HD13	2.00	0.44
1:H:67:ASP:HB2	1:H:68:PRO:CD	2.48	0.44
1:G:81:VAL:HG21	1:G:91:TYR:CZ	2.53	0.43
1:D:22:ARG:HD2	4:D:691:HOH:O	2.17	0.43
1:D:273:HIS:ND1	1:D:276:GLU:OE1	2.49	0.43
1:A:123:GLY:O	2:A:301:NAP:H4N	2.17	0.43
1:D:123:GLY:O	2:D:301:NAP:H4N	2.18	0.43
1:H:217:LYS:HB2	1:H:217:LYS:HE3	1.85	0.43
1:F:135:ILE:HD12	1:F:158[A]:CYS:SG	2.59	0.43
1:A:38:ALA:HB3	1:A:39:PRO:HD3	2.01	0.42
1:D:228:PHE:H	3:D:305:GOL:C2	2.32	0.41
1:D:38:ALA:HB3	1:D:39:PRO:HD3	2.02	0.41
1:E:33:ASN:HA	1:E:34:PRO:HD3	1.94	0.41
1:F:286:GLU:O	1:F:287:ALA:C	2.59	0.41
1:G:92:ILE:HA	1:G:117:LEU:O	2.20	0.41
1:B:180:GLN:NE2	4:B:313:HOH:O	2.54	0.41
1:H:239:LYS:HB3	1:H:239:LYS:HE3	1.94	0.41
1:C:67:ASP:HB2	1:C:68:PRO:CD	2.50	0.41
1:D:124:THR:HA	2:D:301:NAP:O7N	2.20	0.41
1:B:67:ASP:HB2	1:B:68:PRO:CD	2.52	0.40
1:D:12:MET:O	1:D:15:PRO:HD2	2.20	0.40
1:G:171:LYS:HA	1:G:171:LYS:HD2	1.92	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:81:VAL:HG21	1:C:91:TYR:CZ	2.56	0.40
1:B:38:ALA:HB3	1:B:39:PRO:HD3	2.03	0.40
1:C:242:ARG:HD2	3:C:304:GOL:C3	2.33	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	289/287 (101%)	282 (98%)	7 (2%)	0	100	100
1	B	286/287 (100%)	280 (98%)	6 (2%)	0	100	100
1	C	284/287 (99%)	278 (98%)	6 (2%)	0	100	100
1	D	286/287 (100%)	279 (98%)	7 (2%)	0	100	100
1	E	288/287 (100%)	281 (98%)	7 (2%)	0	100	100
1	F	287/287 (100%)	281 (98%)	6 (2%)	0	100	100
1	G	283/287 (99%)	276 (98%)	7 (2%)	0	100	100
1	H	285/287 (99%)	278 (98%)	7 (2%)	0	100	100
All	All	2288/2296 (100%)	2235 (98%)	53 (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	205/201 (102%)	203 (99%)	2 (1%)	82	81
1	B	202/201 (100%)	202 (100%)	0	100	100
1	C	200/201 (100%)	198 (99%)	2 (1%)	82	81
1	D	202/201 (100%)	202 (100%)	0	100	100
1	E	204/201 (102%)	203 (100%)	1 (0%)	92	92
1	F	203/201 (101%)	201 (99%)	2 (1%)	82	81
1	G	200/201 (100%)	198 (99%)	2 (1%)	82	81
1	H	201/201 (100%)	198 (98%)	3 (2%)	72	69
All	All	1617/1608 (101%)	1605 (99%)	12 (1%)	88	88

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	LEU
1	A	194	ASN
1	C	40	LEU
1	C	286	GLU
1	E	238	GLN
1	F	194	ASN
1	F	238	GLN
1	G	238	GLN
1	G	286	GLU
1	H	193	ARG
1	H	194	ASN
1	H	238	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	180	GLN
1	A	236	HIS
1	B	180	GLN
1	C	180	GLN
1	D	180	GLN
1	D	238	GLN
1	E	238	GLN
1	F	238	GLN
1	G	238	GLN
1	H	238	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 ⓘ

19 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	NAP	A	301	-	45,52,52	1.47	3 (6%)	55,80,80	1.86	4 (7%)
3	GOL	A	304	-	5,5,5	0.22	0	5,5,5	0.42	0
2	NAP	B	301	-	45,52,52	1.51	3 (6%)	55,80,80	1.84	6 (10%)
3	GOL	B	304	-	5,5,5	0.27	0	5,5,5	0.31	0
3	GOL	B	305	-	5,5,5	0.30	0	5,5,5	0.37	0
3	GOL	C	288	-	5,5,5	0.37	0	5,5,5	0.19	0
2	NAP	C	301	-	45,52,52	1.50	3 (6%)	55,80,80	1.95	5 (9%)
3	GOL	C	304	-	5,5,5	0.33	0	5,5,5	0.23	0
2	NAP	D	301	-	45,52,52	1.51	3 (6%)	55,80,80	1.90	4 (7%)
3	GOL	D	305	-	5,5,5	0.26	0	5,5,5	0.58	0
2	NAP	E	301	-	45,52,52	1.46	3 (6%)	55,80,80	1.83	4 (7%)
3	GOL	E	304	-	5,5,5	0.33	0	5,5,5	0.74	0
3	GOL	F	288	-	5,5,5	0.36	0	5,5,5	1.01	0
2	NAP	F	301	-	45,52,52	1.45	3 (6%)	55,80,80	1.77	4 (7%)
3	GOL	F	304	-	5,5,5	0.35	0	5,5,5	0.27	0
2	NAP	G	301	-	45,52,52	1.49	3 (6%)	55,80,80	1.87	5 (9%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	G	304	-	5,5,5	0.34	0	5,5,5	0.25	0
2	NAP	H	301	-	45,52,52	1.45	3 (6%)	55,80,80	1.98	6 (10%)
3	GOL	H	305	-	5,5,5	0.25	0	5,5,5	0.35	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	NAP	A	301	-	-	0/27/67/67	0/5/5/5
3	GOL	A	304	-	-	0/4/4/4	0/0/0/0
2	NAP	B	301	-	-	0/27/67/67	0/5/5/5
3	GOL	B	304	-	-	0/4/4/4	0/0/0/0
3	GOL	B	305	-	-	0/4/4/4	0/0/0/0
3	GOL	C	288	-	-	0/4/4/4	0/0/0/0
2	NAP	C	301	-	-	0/27/67/67	0/5/5/5
3	GOL	C	304	-	-	0/4/4/4	0/0/0/0
2	NAP	D	301	-	-	0/27/67/67	0/5/5/5
3	GOL	D	305	-	-	0/4/4/4	0/0/0/0
2	NAP	E	301	-	-	0/27/67/67	0/5/5/5
3	GOL	E	304	-	-	0/4/4/4	0/0/0/0
3	GOL	F	288	-	-	0/4/4/4	0/0/0/0
2	NAP	F	301	-	-	0/27/67/67	0/5/5/5
3	GOL	F	304	-	-	0/4/4/4	0/0/0/0
2	NAP	G	301	-	-	0/27/67/67	0/5/5/5
3	GOL	G	304	-	-	0/4/4/4	0/0/0/0
2	NAP	H	301	-	-	0/27/67/67	0/5/5/5
3	GOL	H	305	-	-	0/4/4/4	0/0/0/0

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	301	NAP	C2A-N1A	2.36	1.38	1.33
2	H	301	NAP	C2A-N1A	2.42	1.38	1.33
2	G	301	NAP	C2A-N1A	2.42	1.38	1.33
2	B	301	NAP	C2A-N1A	2.43	1.38	1.33
2	D	301	NAP	C2A-N1A	2.43	1.38	1.33
2	C	301	NAP	C2A-N1A	2.43	1.38	1.33
2	A	301	NAP	C2A-N1A	2.45	1.38	1.33
2	E	301	NAP	C2A-N1A	2.49	1.38	1.33
2	H	301	NAP	C2A-N3A	3.45	1.38	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	301	NAP	C2A-N3A	3.45	1.38	1.32
2	A	301	NAP	C2A-N3A	3.47	1.38	1.32
2	E	301	NAP	C2A-N3A	3.50	1.38	1.32
2	B	301	NAP	C2A-N3A	3.55	1.38	1.32
2	G	301	NAP	C2A-N3A	3.57	1.38	1.32
2	C	301	NAP	C2A-N3A	3.61	1.38	1.32
2	D	301	NAP	C2A-N3A	3.61	1.38	1.32
2	F	301	NAP	O7N-C7N	7.27	1.39	1.24
2	E	301	NAP	O7N-C7N	7.35	1.39	1.24
2	A	301	NAP	O7N-C7N	7.36	1.39	1.24
2	H	301	NAP	O7N-C7N	7.42	1.40	1.24
2	D	301	NAP	O7N-C7N	7.52	1.40	1.24
2	C	301	NAP	O7N-C7N	7.53	1.40	1.24
2	B	301	NAP	O7N-C7N	7.63	1.40	1.24
2	G	301	NAP	O7N-C7N	7.69	1.40	1.24

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	301	NAP	N3A-C2A-N1A	-11.98	119.46	128.87
2	C	301	NAP	N3A-C2A-N1A	-11.75	119.64	128.87
2	D	301	NAP	N3A-C2A-N1A	-11.39	119.92	128.87
2	G	301	NAP	N3A-C2A-N1A	-11.34	119.96	128.87
2	B	301	NAP	N3A-C2A-N1A	-11.23	120.05	128.87
2	A	301	NAP	N3A-C2A-N1A	-11.21	120.07	128.87
2	E	301	NAP	N3A-C2A-N1A	-11.09	120.16	128.87
2	F	301	NAP	N3A-C2A-N1A	-10.99	120.24	128.87
2	H	301	NAP	C1B-N9A-C4A	-3.57	122.82	126.81
2	A	301	NAP	C1B-N9A-C4A	-3.44	122.97	126.81
2	C	301	NAP	O7N-C7N-C3N	-3.35	115.88	119.60
2	D	301	NAP	C1B-N9A-C4A	-3.31	123.11	126.81
2	H	301	NAP	O7N-C7N-C3N	-3.29	115.95	119.60
2	B	301	NAP	C1B-N9A-C4A	-3.11	123.33	126.81
2	E	301	NAP	C1B-N9A-C4A	-3.08	123.37	126.81
2	C	301	NAP	C1B-N9A-C4A	-2.72	123.77	126.81
2	F	301	NAP	C1B-N9A-C4A	-2.67	123.83	126.81
2	D	301	NAP	O7N-C7N-C3N	-2.63	116.68	119.60
2	G	301	NAP	O7N-C7N-C3N	-2.48	116.85	119.60
2	B	301	NAP	O7N-C7N-C3N	-2.20	117.16	119.60
2	E	301	NAP	O7N-C7N-C3N	-2.09	117.28	119.60
2	F	301	NAP	O2A-PA-O3	2.02	113.91	105.27
2	A	301	NAP	O2N-PN-O1N	2.04	123.19	112.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	301	NAP	O4B-C1B-N9A	2.08	112.03	108.11
2	B	301	NAP	O2A-PA-O3	2.12	114.34	105.27
2	C	301	NAP	O4B-C1B-N9A	2.16	112.18	108.11
2	B	301	NAP	O2N-PN-O1N	2.17	123.84	112.56
2	H	301	NAP	O2A-PA-O3	2.19	114.67	105.27
2	G	301	NAP	C2N-C3N-C4N	2.38	120.97	118.27
2	G	301	NAP	C3N-C7N-N7N	2.70	120.88	117.82
2	B	301	NAP	C3N-C7N-N7N	2.77	120.95	117.82
2	F	301	NAP	C3N-C7N-N7N	2.81	121.00	117.82
2	G	301	NAP	O4B-C1B-N9A	2.87	113.53	108.11
2	E	301	NAP	C3N-C7N-N7N	3.34	121.60	117.82
2	H	301	NAP	C3N-C7N-N7N	3.42	121.69	117.82
2	D	301	NAP	C3N-C7N-N7N	3.52	121.80	117.82
2	C	301	NAP	C3N-C7N-N7N	3.60	121.89	117.82
2	A	301	NAP	C3N-C7N-N7N	3.69	121.99	117.82

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	NAP	3	0
3	C	288	GOL	1	0
3	C	304	GOL	2	0
2	D	301	NAP	2	0
3	D	305	GOL	7	0
3	E	304	GOL	1	0
3	G	304	GOL	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, 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	287/287 (100%)	-0.39	2 (0%) 89 90	15, 21, 33, 57	0
1	B	286/287 (99%)	-0.25	1 (0%) 94 94	15, 23, 39, 63	0
1	C	286/287 (99%)	-0.18	4 (1%) 78 80	16, 28, 46, 69	0
1	D	287/287 (100%)	-0.38	0 100 100	15, 23, 36, 59	0
1	E	287/287 (100%)	-0.36	3 (1%) 84 86	17, 24, 40, 57	0
1	F	286/287 (99%)	-0.43	3 (1%) 84 86	13, 19, 31, 64	0
1	G	284/287 (98%)	0.21	19 (6%) 21 23	20, 41, 77, 112	0
1	H	286/287 (99%)	-0.38	2 (0%) 89 90	16, 25, 39, 67	0
All	All	2289/2296 (99%)	-0.27	34 (1%) 76 79	13, 24, 51, 112	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	43	LEU	5.1
1	G	38	ALA	4.7
1	H	287	ALA	4.3
1	G	42	ALA	4.2
1	G	41	VAL	3.6
1	G	76	PHE	3.0
1	G	46	ARG	3.0
1	C	41	VAL	2.9
1	G	39	PRO	2.7
1	G	34	PRO	2.7
1	G	40	LEU	2.7
1	B	287	ALA	2.7
1	G	35	ALA	2.6
1	H	111	ALA	2.6
1	F	287	ALA	2.6
1	G	27	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	F	177	ILE	2.5
1	G	45	ALA	2.5
1	E	177	ILE	2.4
1	G	37	CYS	2.4
1	E	183	THR	2.4
1	A	177	ILE	2.3
1	C	177	ILE	2.3
1	G	30	TRP	2.3
1	G	33	ASN	2.2
1	C	287	ALA	2.2
1	A	136	LEU	2.2
1	G	111	ALA	2.1
1	C	2	THR	2.1
1	F	184	ALA	2.1
1	G	82	LEU	2.1
1	G	109	VAL	2.1
1	E	184	ALA	2.1
1	G	79	ASN	2.1

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(Å ²)	Q<0.9
3	GOL	D	305	6/6	0.71	0.34	10.92	47,49,50,54	0
3	GOL	F	288	6/6	0.86	0.18	4.57	25,37,39,39	0
3	GOL	B	305	6/6	0.93	0.11	3.96	31,38,44,47	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	GOL	B	304	6/6	0.95	0.12	3.92	26,27,34,37	0
3	GOL	G	304	6/6	0.85	0.15	2.83	37,48,56,63	0
3	GOL	A	304	6/6	0.93	0.12	2.32	22,27,40,45	0
3	GOL	H	305	6/6	0.90	0.12	1.70	33,35,45,49	0
3	GOL	E	304	6/6	0.89	0.18	1.66	41,47,48,55	0
3	GOL	C	304	6/6	0.96	0.12	1.17	29,37,45,51	0
3	GOL	F	304	6/6	0.94	0.10	1.08	23,31,32,40	0
3	GOL	C	288	6/6	0.94	0.11	0.93	27,31,35,42	0
2	NAP	D	301	48/48	0.97	0.09	0.03	18,23,31,34	0
2	NAP	G	301	48/48	0.92	0.12	-0.01	29,45,55,61	0
2	NAP	B	301	48/48	0.96	0.09	-0.22	20,24,30,34	0
2	NAP	E	301	48/48	0.97	0.08	-0.24	21,26,31,37	0
2	NAP	F	301	48/48	0.97	0.08	-0.27	14,20,24,27	0
2	NAP	A	301	48/48	0.98	0.08	-0.45	15,21,24,27	0
2	NAP	H	301	48/48	0.97	0.08	-0.53	19,24,30,35	0
2	NAP	C	301	48/48	0.97	0.07	-0.71	23,30,36,40	0

6.5 Other polymers ⓘ

There are no such residues in this entry.