



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 02:38 am GMT

PDB ID : 2C20
Title : CRYSTAL STRUCTURE OF UDP-GLUCOSE 4-EPIMERASE
Authors : Lebedev, A.A.; Moroz, O.V.; Blagova, E.V.; Levnikov, V.M.; Fogg, M.J.;
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Deposited on : 2005-09-22
Resolution : 2.70 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

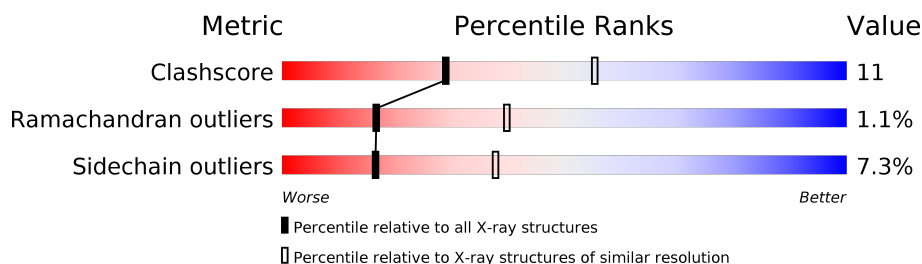
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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))
Clashscore	112137	2590 (2.70-2.70)
Ramachandran outliers	110173	2550 (2.70-2.70)
Sidechain outliers	110143	2550 (2.70-2.70)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	330	
1	B	330	
1	C	330	
1	D	330	
1	E	330	
1	F	330	

2 Entry composition [i](#)

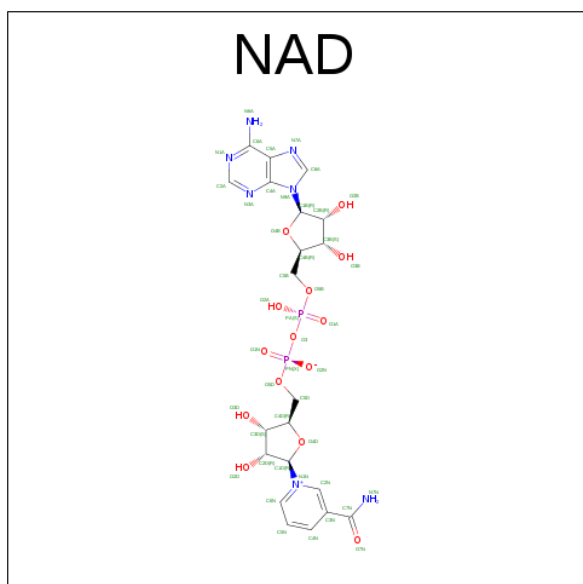
There are 4 unique types of molecules in this entry. The entry contains 16074 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 UDP-GLUCOSE 4-EPIMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	329	Total	C	N	O	S	0	1	0
			2595	1645	440	499	11			
1	B	329	Total	C	N	O	S	0	1	0
			2595	1645	440	499	11			
1	C	329	Total	C	N	O	S	0	1	0
			2595	1645	440	499	11			
1	D	329	Total	C	N	O	S	0	1	0
			2595	1645	440	499	11			
1	E	329	Total	C	N	O	S	0	1	0
			2595	1645	440	499	11			
1	F	329	Total	C	N	O	S	0	1	0
			2595	1645	440	499	11			

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	D	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	E	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	F	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	1	Total	Zn	0	0
			1	1		
3	E	1	Total	Zn	0	0
			1	1		
3	B	1	Total	Zn	0	0
			1	1		
3	C	1	Total	Zn	0	0
			1	1		
3	A	1	Total	Zn	0	0
			1	1		
3	F	1	Total	Zn	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	41	Total	O	0	0
			41	41		
4	B	41	Total	O	0	0
			41	41		
4	C	38	Total	O	0	0
			38	38		
4	D	36	Total	O	0	0
			36	36		
4	E	36	Total	O	0	0
			36	36		
4	F	42	Total	O	0	0
			42	42		

Note EDS was not executed.

Chain A:

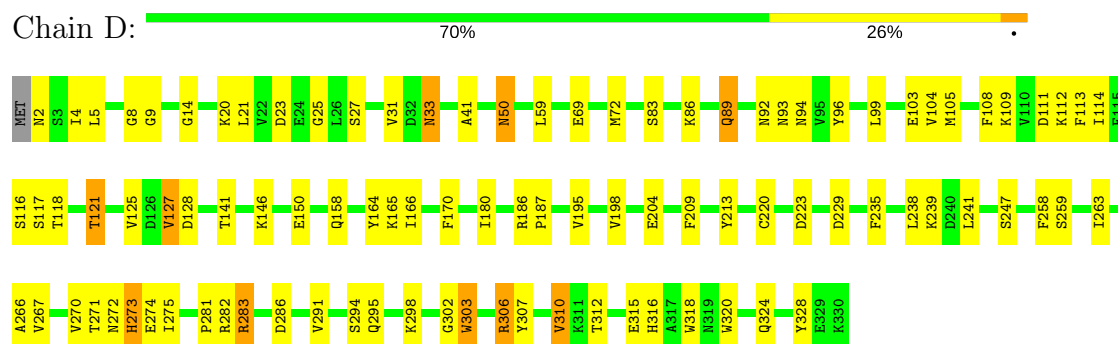
Chain B:

71% 25%

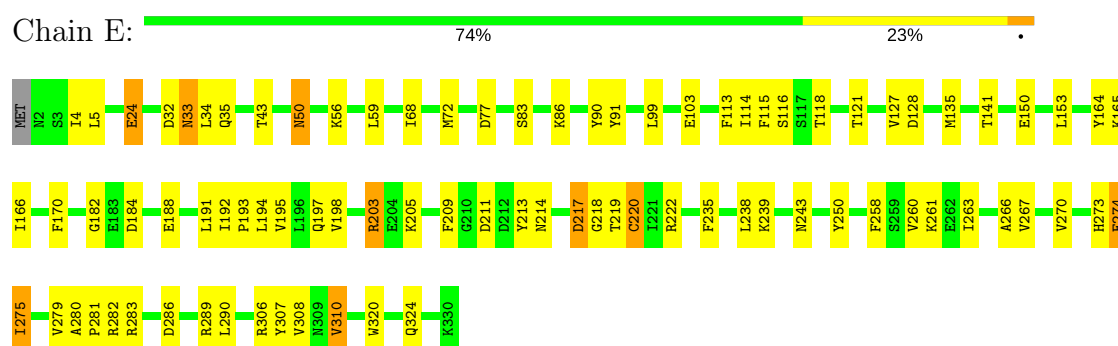
Chain B: MET A2 S3 I4 L5 I6 R20 L21 V22 D23 E24 G25 V31 D32 N33 L34 Q35 E44 N50 G51 N67 M72 L79 R80 G81 M84 S85 R86 P87 Y90 Y91 N92 Y96 I114 F115 S116 S117 T118 T121 E124 V125 D126 V127 E133 T136 T141 K146 Y164 K165 F167 F170 M171 Y172 A173 T176 P177 M178 G179 T180 E188 T189 H190 L191 P192 P193 L194 V195 V198 R203 M208 F209 T215 P216 C220 D223 H226 H234 F235 L238 K239 D240 L241 Q242 N243 S247 D248 V256

[illegible]

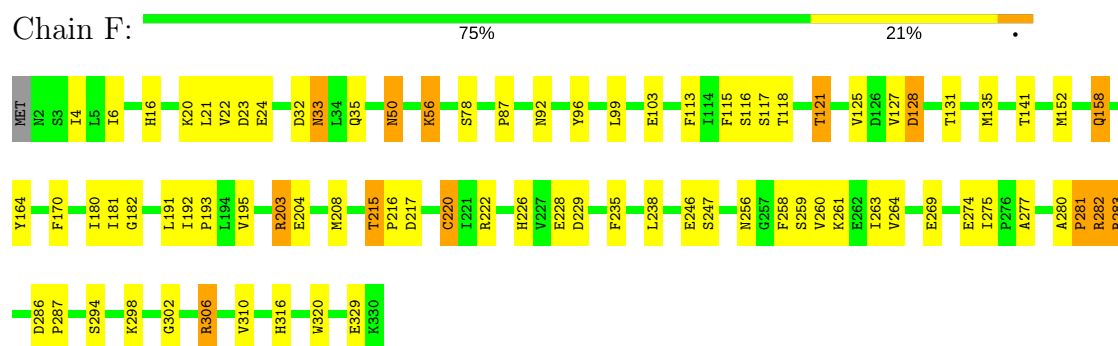
• Molecule 1: UDP-GLUCOSE 4-EPIMERASE



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• Molecule 1: UDP-GLUCOSE 4-EPIMERASE



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 32 1 2	Depositor
Cell constants a, b, c, α , β , γ	136.16 Å 136.16 Å 248.65 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.02 – 2.70	Depositor
% Data completeness (in resolution range)	73.1 (25.02-2.70)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.205 , 0.276	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	16074	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	0.67	1/2656 (0.0%)	0.72	0/3600
1	B	0.71	1/2656 (0.0%)	0.73	0/3600
1	C	0.69	1/2656 (0.0%)	0.72	0/3600
1	D	0.63	0/2656	0.71	0/3600
1	E	1.15	2/2656 (0.1%)	0.73	2/3600 (0.1%)
1	F	0.68	1/2656 (0.0%)	0.71	0/3600
All	All	0.78	6/15936 (0.0%)	0.72	2/21600 (0.0%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	274	GLU	CD-OE2	36.08	1.65	1.25
1	E	274	GLU	CD-OE1	30.32	1.59	1.25
1	B	44	GLU	CD-OE2	6.86	1.33	1.25
1	F	220	CYS	CB-SG	-6.49	1.71	1.82
1	C	220	CYS	CB-SG	-6.08	1.72	1.82
1	A	44	GLU	CD-OE2	5.16	1.31	1.25

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	274	GLU	OE1-CD-OE2	9.22	134.37	123.30
1	E	184	ASP	CB-CG-OD1	5.31	123.08	118.30

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	2595	0	2513	62	0
1	B	2595	0	2513	57	0
1	C	2595	0	2513	52	0
1	D	2595	0	2513	59	0
1	E	2595	0	2513	48	0
1	F	2595	0	2513	57	0
2	A	44	0	26	4	0
2	B	44	0	26	1	0
2	C	44	0	26	2	0
2	D	44	0	26	1	0
2	E	44	0	26	2	0
2	F	44	0	26	3	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
4	A	41	0	0	0	0
4	B	41	0	0	0	0
4	C	38	0	0	1	0
4	D	36	0	0	0	0
4	E	36	0	0	1	0
4	F	42	0	0	1	0
All	All	16074	0	15234	337	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 11.

All (337) 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:E:274:GLU:CD	1:E:274:GLU:OE2	1.65	1.35
1:C:212:ASP:C	1:C:282:ARG:HH22	1.59	1.03
1:F:158:GLN:HA	1:F:158:GLN:HE21	1.35	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:280:ALA:HB1	1:E:281:PRO:HD2	1.53	0.91
1:F:158:GLN:CA	1:F:158:GLN:HE21	1.84	0.90
1:E:33:ASN:ND2	1:E:35:GLN:H	1.73	0.87
1:F:191:LEU:O	1:F:195:VAL:HG23	1.76	0.85
1:A:215:THR:HB	1:A:216:PRO:HD2	1.58	0.85
1:A:33:ASN:HD21	1:A:35:GLN:HG2	1.43	0.83
1:F:4:ILE:HD13	1:F:238:LEU:HD21	1.64	0.80
1:F:215:THR:HB	1:F:216:PRO:CD	2.13	0.79
1:A:33:ASN:ND2	1:A:35:GLN:HG2	1.98	0.79
1:A:311:LYS:O	1:A:315:GLU:HG3	1.83	0.79
1:B:239:LYS:O	1:B:243:ASN:ND2	2.15	0.79
1:F:158:GLN:HA	1:F:158:GLN:NE2	1.96	0.79
1:F:217:ASP:OD2	1:F:261:LYS:HB3	1.84	0.77
1:F:215:THR:HB	1:F:216:PRO:HD2	1.65	0.76
1:B:215:THR:HB	1:B:216:PRO:CD	2.14	0.76
1:B:170:PHE:O	2:B:401:NAD:H4N	1.86	0.76
1:F:170:PHE:O	2:F:401:NAD:H4N	1.87	0.75
1:C:212:ASP:C	1:C:282:ARG:NH2	2.40	0.74
1:A:118:THR:O	1:A:121:THR:HB	1.88	0.73
1:C:212:ASP:OD1	1:C:282:ARG:NH2	2.21	0.73
1:D:118:THR:O	1:D:121:THR:HB	1.88	0.73
1:F:118:THR:O	1:F:121:THR:HB	1.89	0.73
1:B:209:PHE:CD1	1:B:283:ARG:NH1	2.58	0.72
1:C:125:VAL:HG12	1:C:126:ASP:N	2.05	0.71
1:A:191:LEU:HD21	1:A:260:VAL:HG13	1.73	0.71
1:C:33:ASN:ND2	1:C:35:GLN:H	1.88	0.70
1:F:275:ILE:HG22	1:F:275:ILE:O	1.91	0.70
1:F:135[B]:MET:HG3	4:F:2024:HOH:O	1.91	0.70
1:B:275:ILE:HG22	1:B:275:ILE:O	1.90	0.70
1:C:217:ASP:HB2	1:C:261:LYS:HE3	1.71	0.70
1:F:260:VAL:O	1:F:264:VAL:HG23	1.92	0.69
1:E:99:LEU:O	1:E:103:GLU:HG3	1.92	0.69
1:F:226:HIS:HD1	1:F:228:GLU:H	1.42	0.68
1:E:263:ILE:O	1:E:267:VAL:HG23	1.93	0.68
1:B:118:THR:O	1:B:121:THR:HB	1.94	0.68
1:F:220:CYS:O	1:F:259:SER:HA	1.94	0.67
1:A:127:VAL:HG22	1:A:128:ASP:N	2.10	0.67
1:B:81:GLY:O	1:B:84:MET:HB2	1.96	0.66
1:C:4:ILE:HD13	1:C:238:LEU:HD21	1.78	0.65
1:F:4:ILE:CD1	1:F:238:LEU:HD21	2.26	0.65
1:E:165:LYS:HD2	1:E:250:TYR:HE2	1.62	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:263:ILE:O	1:A:267:VAL:HG23	1.97	0.64
1:D:263:ILE:O	1:D:267:VAL:HG23	1.97	0.64
1:C:125:VAL:HG12	1:C:126:ASP:H	1.63	0.64
1:A:127:VAL:HG22	1:A:128:ASP:H	1.64	0.63
1:C:188:GLU:O	1:C:194:LEU:HD21	1.99	0.62
1:C:5:LEU:HD11	1:C:31:VAL:HG23	1.82	0.62
1:B:33:ASN:ND2	1:B:35:GLN:H	1.98	0.61
1:A:163:ARG:NH1	1:A:245:GLY:O	2.33	0.61
1:B:220:CYS:O	1:B:259:SER:HA	2.01	0.60
1:E:4:ILE:HD13	1:E:238:LEU:HD21	1.82	0.60
1:E:72:MET:HG2	1:E:114:ILE:HB	1.83	0.60
2:C:401:NAD:N7N	2:C:401:NAD:O1N	2.34	0.60
1:A:125:VAL:HG12	1:A:127:VAL:H	1.66	0.60
1:F:158:GLN:N	1:F:158:GLN:HE21	1.97	0.60
1:C:99:LEU:O	1:C:103:GLU:HG3	2.01	0.60
1:D:117:SER:HA	1:D:146:LYS:HE3	1.84	0.60
1:E:191:LEU:O	1:E:195:VAL:HG23	2.02	0.60
1:A:215:THR:HB	1:A:216:PRO:CD	2.28	0.60
1:B:235:PHE:CE1	1:B:239:LYS:HD2	2.37	0.59
1:B:33:ASN:HD22	1:B:33:ASN:C	2.04	0.59
1:C:275:ILE:O	1:C:275:ILE:HG22	2.01	0.59
1:F:4:ILE:HD13	1:F:238:LEU:CD2	2.30	0.59
1:F:24:GLU:OE1	1:F:235:PHE:CZ	2.55	0.59
1:C:284:ALA:O	1:C:286:ASP:N	2.33	0.58
1:B:259:SER:O	1:B:263:ILE:HD13	2.03	0.58
1:D:72:MET:HG2	1:D:114:ILE:HB	1.84	0.58
1:D:99:LEU:O	1:D:103:GLU:HG3	2.03	0.58
1:C:311:LYS:O	1:C:315:GLU:HG3	2.04	0.58
1:A:5:LEU:HB2	1:A:68:ILE:HD12	1.86	0.58
1:D:21:LEU:CD2	1:D:235:PHE:HD1	2.16	0.57
1:E:222:ARG:HD3	1:E:290:LEU:HB2	1.87	0.57
1:D:320:TRP:CZ3	1:D:324:GLN:HG3	2.40	0.57
1:A:146:LYS:NZ	2:A:401:NAD:HO2N	2.02	0.57
1:A:56:LYS:HE3	1:A:103:GLU:OE1	2.05	0.57
1:D:4:ILE:HD13	1:D:238:LEU:HD21	1.87	0.56
1:F:33:ASN:ND2	1:F:35:GLN:H	2.02	0.56
1:B:263:ILE:O	1:B:267:VAL:HG23	2.04	0.56
1:E:280:ALA:HB1	1:E:281:PRO:CD	2.32	0.56
1:B:215:THR:HB	1:B:216:PRO:HD3	1.86	0.56
1:B:24:GLU:OE2	1:B:235:PHE:HE2	1.87	0.55
1:E:118:THR:O	1:E:121:THR:HB	2.05	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:79:LEU:HB2	1:B:90:TYR:OH	2.05	0.55
1:F:282:ARG:HG2	1:F:283:ARG:H	1.72	0.55
1:B:6:ILE:HD11	1:B:21:LEU:HD12	1.89	0.55
1:E:320:TRP:CZ3	1:E:324:GLN:HG3	2.42	0.55
1:F:6:ILE:HD11	1:F:21:LEU:CD1	2.36	0.55
1:E:77:ASP:HB3	1:E:90:TYR:CZ	2.42	0.55
1:D:21:LEU:CD2	1:D:235:PHE:CD1	2.89	0.55
1:B:20:LYS:HB2	1:B:20:LYS:HZ2	1.71	0.55
1:D:83:SER:HB2	1:D:141:THR:HB	1.88	0.54
1:B:121:THR:HG23	1:B:136:THR:HG22	1.89	0.54
1:D:271:THR:HG22	1:D:318:TRP:CE2	2.43	0.54
1:D:9:GLY:HA2	1:D:14:GLY:HA3	1.89	0.54
1:C:125:VAL:CG1	1:C:126:ASP:N	2.71	0.54
1:B:188:GLU:HG2	1:B:193:PRO:HB2	1.90	0.53
1:E:209:PHE:CD1	1:E:283:ARG:NH2	2.76	0.53
2:A:401:NAD:N7N	2:A:401:NAD:O1N	2.41	0.53
1:A:170:PHE:O	2:A:401:NAD:H4N	2.09	0.53
1:A:41:ALA:HB2	1:A:328:TYR:CE1	2.44	0.53
1:F:246:GLU:HG3	1:F:247:SER:O	2.07	0.53
1:A:197:GLN:OE1	1:A:203:ARG:HD3	2.09	0.53
1:D:306:ARG:HG3	1:D:307:TYR:CE1	2.43	0.53
1:B:215:THR:CB	1:B:216:PRO:CD	2.86	0.52
1:A:259:SER:OG	1:A:262:GLU:HG3	2.09	0.52
1:B:171:ASN:HB2	1:B:223:ASP:O	2.09	0.52
1:A:210:GLY:HA3	1:A:282:ARG:HG2	1.89	0.52
1:D:33:ASN:HD22	1:D:33:ASN:C	2.12	0.52
1:D:50:ASN:HD22	1:D:50:ASN:C	2.13	0.52
1:B:178:ASN:OD1	1:B:180:ILE:HB	2.08	0.52
1:D:164:TYR:O	1:D:247:SER:HA	2.10	0.52
1:B:280:ALA:HB1	1:B:281:PRO:HD2	1.92	0.52
1:A:6:ILE:HD11	1:A:21:LEU:HD12	1.92	0.51
1:D:108:PHE:O	1:D:109:LYS:HB2	2.10	0.51
1:D:266:ALA:O	1:D:270:VAL:HG23	2.10	0.51
1:C:176:THR:HG21	1:C:180:ILE:HG22	1.91	0.51
1:C:258:PHE:CE2	1:C:310:VAL:HG22	2.45	0.51
1:D:21:LEU:HD22	1:D:235:PHE:CD1	2.46	0.51
1:B:191:LEU:O	1:B:195:VAL:HG23	2.11	0.51
1:A:191:LEU:O	1:A:195:VAL:HG23	2.10	0.51
1:D:302:GLY:O	1:D:303:TRP:C	2.48	0.51
1:B:173:ALA:O	1:B:226:HIS:HA	2.10	0.51
1:E:33:ASN:HD22	1:E:34:LEU:N	2.08	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:23:ASP:C	1:B:25:GLY:H	2.14	0.51
1:C:125:VAL:CG1	1:C:126:ASP:H	2.23	0.51
1:C:220:CYS:O	1:C:259:SER:HA	2.11	0.51
1:D:229:ASP:OD1	1:D:306:ARG:HG2	2.11	0.51
1:D:294:SER:O	1:D:298:LYS:HG3	2.11	0.50
1:D:271:THR:OG1	1:D:273:HIS:HB2	2.10	0.50
1:F:259:SER:O	1:F:263:ILE:HD13	2.11	0.50
1:F:258:PHE:CE1	1:F:310:VAL:HG13	2.45	0.50
1:D:21:LEU:HD22	1:D:235:PHE:CE1	2.46	0.50
1:E:150:GLU:HG2	1:E:166:ILE:HD13	1.92	0.50
1:B:33:ASN:HD22	1:B:34:LEU:N	2.10	0.50
1:B:294:SER:HB3	1:B:298:LYS:HE3	1.93	0.50
1:A:197:GLN:HB3	1:A:203:ARG:HG3	1.93	0.50
1:F:215:THR:CB	1:F:216:PRO:CD	2.85	0.50
1:D:213:TYR:CZ	1:D:282:ARG:HG3	2.46	0.50
1:C:191:LEU:O	1:C:195:VAL:HG23	2.11	0.50
1:F:92:ASN:O	1:F:96:TYR:HB3	2.12	0.50
1:A:125:VAL:HG11	1:A:130:ILE:HG12	1.94	0.49
1:C:212:ASP:H	1:C:282:ARG:HH21	1.60	0.49
1:C:4:ILE:CD1	1:C:238:LEU:HD21	2.41	0.49
1:B:247:SER:O	1:B:248:ASP:HB2	2.11	0.49
1:C:263:ILE:O	1:C:267:VAL:HG23	2.12	0.49
1:D:23:ASP:C	1:D:25:GLY:H	2.15	0.49
1:E:170:PHE:O	2:E:401:NAD:H4N	2.12	0.49
1:D:170:PHE:O	2:D:401:NAD:H4N	2.12	0.49
1:F:16:HIS:CE1	1:F:181:ILE:HD12	2.48	0.49
1:A:165:LYS:HD2	1:A:250:TYR:HE2	1.78	0.49
1:E:50:ASN:C	1:E:50:ASN:HD22	2.16	0.49
1:B:50:ASN:HD22	1:B:51:GLY:N	2.11	0.49
1:E:211:ASP:OD2	1:E:279:VAL:HG11	2.12	0.49
1:C:256:ASN:HA	1:C:291:VAL:HG11	1.93	0.48
1:C:33:ASN:HD22	1:C:34:LEU:N	2.11	0.48
1:B:215:THR:HB	1:B:216:PRO:HD2	1.92	0.48
1:D:213:TYR:HE1	1:D:220:CYS:SG	2.35	0.48
1:E:239:LYS:O	1:E:243:ASN:ND2	2.45	0.48
1:B:20:LYS:HB2	1:B:20:LYS:NZ	2.29	0.48
1:C:35:GLN:HG3	1:C:36:THR:HG23	1.95	0.48
1:A:202:GLN:O	1:A:203:ARG:HG2	2.14	0.48
1:F:99:LEU:O	1:F:103:GLU:HG3	2.13	0.48
1:B:5:LEU:HD11	1:B:31:VAL:HG23	1.95	0.48
1:D:2:ASN:HD22	1:D:69:GLU:HB2	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:229:ASP:OD1	1:F:306:ARG:HG2	2.14	0.48
1:F:117:SER:O	2:F:401:NAD:H6N	2.14	0.47
1:A:220:CYS:H	1:A:260:VAL:HG23	1.78	0.47
1:D:111:ASP:OD1	1:D:112:LYS:HG3	2.14	0.47
1:A:127:VAL:CG2	1:A:128:ASP:H	2.25	0.47
1:C:33:ASN:HD22	1:C:33:ASN:C	2.18	0.47
2:E:401:NAD:O1N	2:E:401:NAD:N7N	2.48	0.47
1:E:213:TYR:HB2	1:E:218:GLY:O	2.14	0.47
1:B:275:ILE:CG2	1:B:275:ILE:O	2.60	0.47
1:E:113:PHE:HE2	1:E:115:PHE:HB2	1.79	0.47
1:F:32:ASP:OD1	2:F:401:NAD:H1B	2.14	0.47
1:A:77:ASP:HB3	1:A:90:TYR:CE2	2.49	0.47
1:B:4:ILE:HD13	1:B:238:LEU:HD21	1.97	0.47
1:D:258:PHE:CD1	1:D:310:VAL:HG13	2.49	0.47
1:A:213:TYR:CZ	1:A:282:ARG:NE	2.83	0.47
1:C:297:ALA:O	1:C:301:LEU:HB2	2.15	0.47
1:F:208:MET:CE	1:F:277:ALA:HB1	2.44	0.47
1:E:113:PHE:O	1:E:164:TYR:HA	2.15	0.47
1:A:127:VAL:CG2	1:A:128:ASP:N	2.76	0.46
1:B:274:GLU:HG3	1:B:275:ILE:H	1.80	0.46
1:C:264:VAL:HG12	1:C:268:ARG:NH1	2.29	0.46
1:F:164:TYR:O	1:F:247:SER:HA	2.14	0.46
1:C:180:ILE:HG12	4:C:2030:HOH:O	2.15	0.46
1:F:128:ASP:OD2	1:F:128:ASP:N	2.48	0.46
1:D:312:THR:O	1:D:315:GLU:HB2	2.15	0.46
1:A:240:ASP:OD1	1:A:245:GLY:HA3	2.16	0.46
1:C:258:PHE:CD2	1:C:310:VAL:HG22	2.51	0.46
1:F:6:ILE:HD11	1:F:21:LEU:HD12	1.96	0.46
1:D:220:CYS:O	1:D:259:SER:HA	2.16	0.46
1:E:113:PHE:CE2	1:E:115:PHE:HB2	2.51	0.46
1:E:266:ALA:O	1:E:270:VAL:HG23	2.15	0.46
1:F:20:LYS:HZ3	1:F:235:PHE:HD2	1.64	0.46
1:C:125:VAL:HG12	1:C:127:VAL:H	1.81	0.46
1:E:258:PHE:CD1	1:E:310:VAL:HG13	2.51	0.46
1:B:92:ASN:O	1:B:96:TYR:HB3	2.16	0.46
1:D:312:THR:HG22	1:D:316:HIS:CE1	2.51	0.46
1:F:275:ILE:CG2	1:F:275:ILE:O	2.63	0.46
1:B:33:ASN:C	1:B:33:ASN:ND2	2.69	0.45
1:A:72:MET:HG3	1:A:114:ILE:HB	1.97	0.45
1:B:50:ASN:C	1:B:50:ASN:HD22	2.19	0.45
1:E:219:THR:HB	1:E:260:VAL:HB	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:GLU:O	1:A:26:LEU:N	2.49	0.45
1:E:192:ILE:N	1:E:193:PRO:HD2	2.31	0.45
1:C:50:ASN:HD22	1:C:50:ASN:C	2.19	0.45
1:D:209:PHE:HB3	1:D:283:ARG:HD3	1.98	0.45
1:F:298:LYS:O	1:F:302:GLY:HA2	2.17	0.45
1:D:5:LEU:HD11	1:D:31:VAL:HG23	1.99	0.45
1:E:127:VAL:HG22	1:E:128:ASP:H	1.82	0.45
1:B:320:TRP:CZ3	1:B:324:GLN:HG3	2.51	0.45
1:A:295:GLN:O	1:A:299:GLU:HG3	2.17	0.45
1:D:165:LYS:HE2	1:D:241:LEU:CD2	2.46	0.45
1:F:113:PHE:O	1:F:164:TYR:HA	2.17	0.45
1:A:198:VAL:HA	1:A:203:ARG:O	2.17	0.45
1:D:195:VAL:O	1:D:198:VAL:HB	2.17	0.44
1:A:222:ARG:HD3	1:A:290:LEU:HB2	1.98	0.44
1:B:165:LYS:HE2	1:B:241:LEU:CD2	2.47	0.44
1:C:171:ASN:HB2	1:C:223:ASP:O	2.18	0.44
1:B:4:ILE:HD13	1:B:238:LEU:CD2	2.48	0.44
1:C:215:THR:HG21	1:C:221:ILE:HG12	1.99	0.44
1:D:235:PHE:HE2	1:D:239:LYS:CE	2.30	0.44
1:A:310:VAL:HA	1:A:313:ILE:HD12	2.00	0.44
1:B:198:VAL:HG22	1:B:203:ARG:HG2	1.99	0.44
1:A:258:PHE:CD1	1:A:310:VAL:HG13	2.52	0.44
1:B:258:PHE:HB3	1:B:263:ILE:HD11	1.98	0.44
1:B:87:PRO:HA	1:B:141:THR:HG21	1.99	0.44
1:C:92:ASN:O	1:C:96:TYR:HB3	2.17	0.44
1:F:208:MET:HE3	1:F:277:ALA:HB1	1.99	0.44
1:C:163:ARG:NH1	1:C:245:GLY:O	2.50	0.44
1:C:170:PHE:O	2:C:401:NAD:H4N	2.17	0.44
1:F:113:PHE:HE2	1:F:115:PHE:HB2	1.82	0.44
1:F:33:ASN:HD22	1:F:33:ASN:C	2.19	0.44
1:A:20:LYS:HE2	1:A:235:PHE:CD2	2.53	0.44
1:A:4:ILE:HD13	1:A:238:LEU:CD2	2.48	0.44
1:B:263:ILE:HD12	1:B:263:ILE:N	2.33	0.44
1:D:223:ASP:HB2	1:D:291:VAL:HA	1.99	0.44
1:F:87:PRO:HA	1:F:141:THR:HG21	2.00	0.44
1:C:215:THR:HB	1:C:216:PRO:HD2	1.99	0.44
1:D:209:PHE:CD2	1:D:283:ARG:HD2	2.53	0.44
1:F:222:ARG:NH2	1:F:287:PRO:HD2	2.33	0.44
1:F:50:ASN:C	1:F:50:ASN:HD22	2.21	0.43
1:F:6:ILE:HD11	1:F:21:LEU:HD11	1.99	0.43
1:A:146:LYS:HA	1:A:146:LYS:HD2	1.64	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:TYR:CE2	1:A:282:ARG:NH2	2.86	0.43
1:A:150:GLU:HG2	1:A:166:ILE:HD13	2.00	0.43
1:D:209:PHE:CG	1:D:283:ARG:HD2	2.53	0.43
1:D:93:ASN:ND2	1:D:94:ASN:OD1	2.51	0.43
1:F:192:ILE:N	1:F:193:PRO:HD2	2.34	0.43
1:A:33:ASN:HD21	1:A:35:GLN:CG	2.24	0.43
1:D:125:VAL:HG12	1:D:127:VAL:H	1.83	0.43
1:A:302:GLY:O	1:A:303:TRP:C	2.56	0.43
1:B:117:SER:HA	1:B:146:LYS:HE2	2.00	0.43
1:C:153:LEU:HA	1:C:153:LEU:HD23	1.75	0.43
1:E:273:HIS:O	1:E:275:ILE:HD12	2.18	0.43
1:F:280:ALA:HB1	1:F:281:PRO:HD2	2.01	0.43
1:B:259:SER:O	1:B:260:VAL:C	2.57	0.43
1:E:217:ASP:OD2	1:E:261:LYS:HB3	2.19	0.43
1:E:33:ASN:HD21	1:E:35:GLN:H	1.62	0.43
1:A:92:ASN:O	1:A:96:TYR:HB3	2.18	0.43
1:B:164:TYR:O	1:B:247:SER:HA	2.19	0.43
1:C:212:ASP:CA	1:C:282:ARG:NH2	2.82	0.43
1:D:165:LYS:HE2	1:D:241:LEU:HD21	2.00	0.43
1:C:286:ASP:HA	1:C:287:PRO:HD3	1.83	0.43
1:A:178:ASN:OD1	1:A:178:ASN:C	2.58	0.43
1:B:258:PHE:CD1	1:B:310:VAL:HG13	2.54	0.43
1:D:4:ILE:HD13	1:D:238:LEU:CD2	2.48	0.43
1:B:72:MET:HG2	1:B:114:ILE:HB	2.00	0.42
1:F:56:LYS:HE3	1:F:103:GLU:OE1	2.19	0.42
1:C:199:ALA:HB2	1:C:275:ILE:HD11	2.02	0.42
1:C:56:LYS:HE2	1:C:103:GLU:OE1	2.19	0.42
1:D:213:TYR:HE1	1:D:220:CYS:HG	1.57	0.42
1:A:11:GLY:O	1:A:15:SER:HB3	2.18	0.42
1:C:150:GLU:O	1:C:153:LEU:HB2	2.19	0.42
1:A:83:SER:HB2	1:A:141:THR:HB	2.01	0.42
1:A:298:LYS:HG2	1:A:303:TRP:O	2.20	0.42
1:C:111:ASP:C	1:C:111:ASP:OD1	2.56	0.42
1:D:213:TYR:OH	1:D:282:ARG:HG3	2.19	0.42
1:F:20:LYS:NZ	1:F:235:PHE:HD2	2.17	0.42
1:F:226:HIS:HD1	1:F:228:GLU:N	2.14	0.42
1:A:320:TRP:CZ3	1:A:324:GLN:HG3	2.54	0.42
1:A:33:ASN:CG	1:A:35:GLN:HG2	2.39	0.42
1:D:41:ALA:HB2	1:D:328:TYR:CE1	2.54	0.42
1:D:92:ASN:O	1:D:96:TYR:HB3	2.19	0.42
1:E:258:PHE:CE1	1:E:310:VAL:HG13	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:274:GLU:HG3	1:B:275:ILE:N	2.34	0.42
1:C:164:TYR:C	1:C:164:TYR:CD2	2.93	0.42
1:C:170:PHE:HB2	1:C:223:ASP:HB3	2.01	0.42
1:C:51:GLY:HA3	1:C:58:PHE:CZ	2.55	0.42
1:B:265:ASP:OD1	1:B:268:ARG:NH1	2.53	0.41
1:C:118:THR:HG22	1:C:146:LYS:HG3	2.02	0.41
1:D:150:GLU:HG2	1:D:166:ILE:HD13	2.01	0.41
1:E:153:LEU:HD23	1:E:153:LEU:HA	1.86	0.41
1:A:169:TYR:CE2	2:A:401:NAD:C5N	3.03	0.41
1:B:223:ASP:HB2	1:B:291:VAL:HA	2.01	0.41
1:B:167:PHE:CE1	1:B:234:HIS:HA	2.55	0.41
1:D:158:GLN:HE21	1:D:158:GLN:N	2.18	0.41
1:D:271:THR:C	1:D:273:HIS:N	2.73	0.41
1:E:56:LYS:CE	1:E:103:GLU:OE1	2.68	0.41
1:A:8:GLY:O	1:A:14:GLY:HA3	2.21	0.41
1:C:118:THR:O	1:C:121:THR:HB	2.19	0.41
1:E:182:GLY:HA3	1:E:320:TRP:CG	2.55	0.41
1:A:73:HIS:HD2	1:A:115:PHE:CE1	2.38	0.41
1:A:213:TYR:CE2	1:A:282:ARG:CZ	3.04	0.41
1:D:275:ILE:HG22	1:D:275:ILE:O	2.19	0.41
1:E:195:VAL:O	1:E:198:VAL:HB	2.20	0.41
1:D:113:PHE:O	1:D:164:TYR:HA	2.20	0.41
1:D:86:LYS:HB3	1:D:89:GLN:HG2	2.03	0.41
1:E:5:LEU:HB2	1:E:68:ILE:HD12	2.01	0.41
1:E:91:TYR:CD1	1:F:152:MET:HE2	2.55	0.41
1:F:203:ARG:HH11	1:F:203:ARG:CG	2.33	0.41
1:F:286:ASP:HA	1:F:287:PRO:HD3	1.71	0.41
1:E:135[A]:MET:HG2	4:E:2018:HOH:O	2.20	0.41
1:F:203:ARG:HH11	1:F:203:ARG:HG3	1.86	0.41
1:E:306:ARG:HG3	1:E:307:TYR:CE1	2.55	0.41
1:A:41:ALA:HB2	1:A:328:TYR:CZ	2.56	0.41
1:C:56:LYS:CE	1:C:103:GLU:OE1	2.68	0.41
1:E:83:SER:HB2	1:E:141:THR:HB	2.02	0.41
1:E:188:GLU:O	1:E:194:LEU:HD21	2.21	0.41
1:E:56:LYS:HE3	1:E:103:GLU:OE1	2.21	0.41
1:D:8:GLY:O	1:D:14:GLY:HA3	2.21	0.40
1:A:207:MET:HE3	1:A:280:ALA:HB2	2.04	0.40
1:A:50:ASN:HD22	1:A:50:ASN:C	2.24	0.40
1:B:170:PHE:O	1:B:171:ASN:ND2	2.54	0.40
1:F:182:GLY:HA3	1:F:320:TRP:CD2	2.56	0.40
1:A:33:ASN:HD22	1:A:33:ASN:H	1.69	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:186:ARG:HA	1:D:187:PRO:HA	1.95	0.40
1:E:219:THR:HB	1:E:220:CYS:H	1.73	0.40
1:A:133:GLU:HG2	1:A:133:GLU:H	1.73	0.40
1:D:104:VAL:O	1:D:105:MET:C	2.60	0.40
1:E:197:GLN:HB3	1:E:203:ARG:HG3	2.04	0.40
1:E:114:ILE:HD11	1:E:238:LEU:HA	2.03	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	328/330 (99%)	297 (90%)	27 (8%)	4 (1%)	15	37
1	B	328/330 (99%)	307 (94%)	19 (6%)	2 (1%)	28	56
1	C	328/330 (99%)	309 (94%)	17 (5%)	2 (1%)	28	56
1	D	328/330 (99%)	302 (92%)	22 (7%)	4 (1%)	15	37
1	E	328/330 (99%)	291 (89%)	33 (10%)	4 (1%)	15	37
1	F	328/330 (99%)	301 (92%)	22 (7%)	5 (2%)	12	30
All	All	1968/1980 (99%)	1807 (92%)	140 (7%)	21 (1%)	17	40

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	286	ASP
1	A	25	GLY
1	D	272	ASN
1	F	281	PRO
1	F	282	ARG
1	A	285	GLY

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Mol	Chain	Res	Type
1	B	286	ASP
1	C	285	GLY
1	D	303	TRP
1	E	24	GLU
1	F	23	ASP
1	E	214	ASN
1	E	282	ARG
1	A	218	GLY
1	F	215	THR
1	B	215	THR
1	F	22	VAL
1	D	281	PRO
1	E	275	ILE
1	A	275	ILE
1	D	286	ASP

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	276/277 (100%)	258 (94%)	18 (6%)	20	44
1	B	276/277 (100%)	250 (91%)	26 (9%)	10	23
1	C	276/277 (100%)	256 (93%)	20 (7%)	17	39
1	D	276/277 (100%)	258 (94%)	18 (6%)	20	44
1	E	276/277 (100%)	259 (94%)	17 (6%)	21	46
1	F	276/277 (100%)	254 (92%)	22 (8%)	14	32
All	All	1656/1662 (100%)	1535 (93%)	121 (7%)	16	38

All (121) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ASN
1	A	19	LYS
1	A	26	LEU

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Mol	Chain	Res	Type
1	A	32	ASP
1	A	33	ASN
1	A	50	ASN
1	A	59	LEU
1	A	69	GLU
1	A	72	MET
1	A	78	SER
1	A	116	SER
1	A	121	THR
1	A	133	GLU
1	A	158	GLN
1	A	217	ASP
1	A	281	PRO
1	A	306	ARG
1	A	310	VAL
1	B	3	SER
1	B	20	LYS
1	B	33	ASN
1	B	50	ASN
1	B	67	ASN
1	B	86	LYS
1	B	116	SER
1	B	121	THR
1	B	124	GLU
1	B	125	VAL
1	B	126	ASP
1	B	127	VAL
1	B	133	GLU
1	B	176	THR
1	B	180	ILE
1	B	189	THR
1	B	203	ARG
1	B	208	MET
1	B	256	ASN
1	B	279	VAL
1	B	283	ARG
1	B	298	LYS
1	B	306	ARG
1	B	309	ASN
1	B	310	VAL
1	B	319	ASN
1	C	3	SER

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Mol	Chain	Res	Type
1	C	33	ASN
1	C	50	ASN
1	C	69	GLU
1	C	86	LYS
1	C	116	SER
1	C	121	THR
1	C	133	GLU
1	C	158	GLN
1	C	176	THR
1	C	205	LYS
1	C	215	THR
1	C	220	CYS
1	C	235	PHE
1	C	279	VAL
1	C	286	ASP
1	C	289	ARG
1	C	308	VAL
1	C	310	VAL
1	C	330	LYS
1	D	20	LYS
1	D	27	SER
1	D	33	ASN
1	D	50	ASN
1	D	59	LEU
1	D	89	GLN
1	D	116	SER
1	D	121	THR
1	D	127	VAL
1	D	128	ASP
1	D	180	ILE
1	D	204	GLU
1	D	273	HIS
1	D	274	GLU
1	D	283	ARG
1	D	295	GLN
1	D	306	ARG
1	D	310	VAL
1	E	24	GLU
1	E	32	ASP
1	E	33	ASN
1	E	43	THR
1	E	50	ASN

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Mol	Chain	Res	Type
1	E	59	LEU
1	E	86	LYS
1	E	116	SER
1	E	203	ARG
1	E	205	LYS
1	E	217	ASP
1	E	220	CYS
1	E	235	PHE
1	E	286	ASP
1	E	289	ARG
1	E	308	VAL
1	E	310	VAL
1	F	33	ASN
1	F	50	ASN
1	F	56	LYS
1	F	78	SER
1	F	116	SER
1	F	121	THR
1	F	125	VAL
1	F	127	VAL
1	F	128	ASP
1	F	131	THR
1	F	158	GLN
1	F	180	ILE
1	F	203	ARG
1	F	204	GLU
1	F	256	ASN
1	F	269	GLU
1	F	274	GLU
1	F	283	ARG
1	F	294	SER
1	F	306	ARG
1	F	316	HIS
1	F	329	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	ASN
1	A	50	ASN
1	A	73	HIS
1	A	158	GLN

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Mol	Chain	Res	Type
1	A	161	ASN
1	B	33	ASN
1	B	50	ASN
1	B	73	HIS
1	B	137	ASN
1	B	171	ASN
1	B	316	HIS
1	B	319	ASN
1	C	33	ASN
1	C	50	ASN
1	C	65	GLN
1	C	158	GLN
1	C	161	ASN
1	D	33	ASN
1	D	50	ASN
1	D	65	GLN
1	D	89	GLN
1	D	158	GLN
1	E	33	ASN
1	E	50	ASN
1	E	73	HIS
1	F	33	ASN
1	F	50	ASN
1	F	73	HIS
1	F	89	GLN
1	F	158	GLN
1	F	171	ASN
1	F	316	HIS

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 12 ligands modelled in this entry, 6 are monoatomic - leaving 6 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	NAD	A	401	1	41,48,48	1.79	3 (7%)	43,73,73	1.76	3 (6%)
2	NAD	B	401	-	41,48,48	1.74	3 (7%)	43,73,73	1.90	2 (4%)
2	NAD	C	401	-	41,48,48	1.74	3 (7%)	43,73,73	1.78	4 (9%)
2	NAD	D	401	-	41,48,48	1.80	3 (7%)	43,73,73	2.02	1 (2%)
2	NAD	E	401	-	41,48,48	1.74	3 (7%)	43,73,73	1.90	4 (9%)
2	NAD	F	401	-	41,48,48	1.68	3 (7%)	43,73,73	1.76	3 (6%)

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	NAD	A	401	1	-	0/22/62/62	0/5/5/5
2	NAD	B	401	-	-	0/22/62/62	0/5/5/5
2	NAD	C	401	-	-	0/22/62/62	0/5/5/5
2	NAD	D	401	-	-	0/22/62/62	0/5/5/5
2	NAD	E	401	-	-	0/22/62/62	0/5/5/5
2	NAD	F	401	-	-	0/22/62/62	0/5/5/5

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	NAD	C2A-N1A	2.18	1.38	1.33
2	C	401	NAD	C2A-N1A	2.22	1.38	1.33
2	A	401	NAD	C2A-N1A	2.29	1.38	1.33
2	F	401	NAD	C2A-N1A	2.34	1.38	1.33
2	D	401	NAD	C2A-N1A	2.36	1.38	1.33
2	E	401	NAD	C2A-N1A	2.57	1.38	1.33
2	F	401	NAD	C2A-N3A	3.57	1.38	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	401	NAD	C2A-N3A	3.89	1.38	1.32
2	C	401	NAD	C2A-N3A	4.04	1.38	1.32
2	B	401	NAD	C2A-N3A	4.17	1.39	1.32
2	D	401	NAD	C2A-N3A	4.18	1.39	1.32
2	A	401	NAD	C2A-N3A	4.22	1.39	1.32
2	E	401	NAD	O7N-C7N	8.50	1.41	1.24
2	B	401	NAD	O7N-C7N	8.60	1.41	1.24
2	C	401	NAD	O7N-C7N	8.61	1.41	1.24
2	F	401	NAD	O7N-C7N	8.61	1.41	1.24
2	D	401	NAD	O7N-C7N	8.91	1.42	1.24
2	A	401	NAD	O7N-C7N	8.96	1.42	1.24

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	401	NAD	N3A-C2A-N1A	-12.15	118.27	128.86
2	B	401	NAD	N3A-C2A-N1A	-10.81	119.44	128.86
2	E	401	NAD	N3A-C2A-N1A	-10.28	119.91	128.86
2	F	401	NAD	N3A-C2A-N1A	-10.02	120.13	128.86
2	A	401	NAD	N3A-C2A-N1A	-9.87	120.26	128.86
2	C	401	NAD	N3A-C2A-N1A	-9.06	120.97	128.86
2	C	401	NAD	O7N-C7N-C3N	-3.39	115.66	119.62
2	E	401	NAD	O7N-C7N-N7N	-2.58	118.91	122.58
2	B	401	NAD	O5B-C5B-C4B	-2.21	101.15	109.00
2	F	401	NAD	O5B-C5B-C4B	-2.07	101.66	109.00
2	A	401	NAD	O7N-C7N-N7N	-2.02	119.70	122.58
2	F	401	NAD	O2N-PN-O1N	2.00	122.66	112.28
2	C	401	NAD	O2B-C2B-C1B	2.04	117.99	111.61
2	E	401	NAD	C3N-C2N-N1N	2.25	122.69	120.43
2	A	401	NAD	C3N-C7N-N7N	2.64	120.79	117.77
2	E	401	NAD	C3N-C7N-N7N	2.97	121.17	117.77
2	C	401	NAD	C3N-C7N-N7N	3.78	122.09	117.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	NAD	4	0
2	B	401	NAD	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	401	NAD	2	0
2	D	401	NAD	1	0
2	E	401	NAD	2	0
2	F	401	NAD	3	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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.