



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 05:41 am GMT

PDB ID : 4NMK
Title : Thermostable aldehyde dehydrogenase from *Pyrobaculum* sp. crystallized in microgravity (complex with NADP+)
Authors : Petrova, T.; Boyko, K.M.; Bezsudnova, E.Y.; Mardanov, A.V.; Gumerov, V.M.; Ravin, N.V.; Popov, V.O.
Deposited on : 2013-11-15
Resolution : 1.90 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

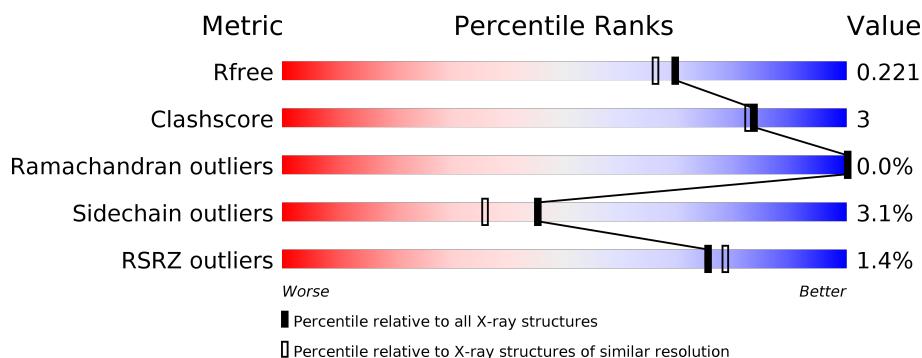
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	5047 (1.90-1.90)
Clashscore	112137	5731 (1.90-1.90)
Ramachandran outliers	110173	5669 (1.90-1.90)
Sidechain outliers	110143	5670 (1.90-1.90)
RSRZ outliers	101464	5100 (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	491	<div> <div>92%</div> <div>7%</div> <div>•</div> </div>
1	B	491	<div> <div>92%</div> <div>6%</div> <div>•</div> </div>
1	C	491	<div> <div>91%</div> <div>9%</div> <div>•</div> </div>
1	D	491	<div> <div>2%</div> <div>87%</div> <div>10%</div> <div>••</div> </div>
1	E	491	<div> <div>2%</div> <div>89%</div> <div>9%</div> <div>••</div> </div>
1	F	491	<div> <div>2%</div> <div>91%</div> <div>8%</div> <div>•</div> </div>

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Mol	Chain	Length	Quality of chain
1	G	491	<div> <div></div> <div>90%</div> <div>8%</div> <div>..</div> </div>
1	H	491	<div> <div></div> <div>92%</div> <div>7%</div> <div>..</div> </div>

2 Entry composition

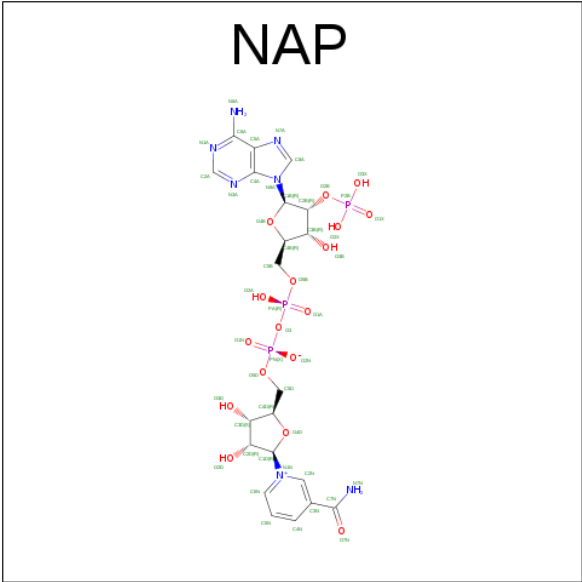
There are 3 unique types of molecules in this entry. The entry contains 33972 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	488	Total	C	N	O	S	0	2	0
			3810	2429	664	705	12			
1	B	484	Total	C	N	O	S	0	2	0
			3784	2410	661	701	12			
1	C	490	Total	C	N	O	S	0	10	0
			3866	2464	672	717	13			
1	D	485	Total	C	N	O	S	0	2	0
			3787	2412	661	702	12			
1	E	485	Total	C	N	O	S	0	2	0
			3789	2413	662	702	12			
1	F	490	Total	C	N	O	S	0	2	0
			3830	2439	669	710	12			
1	G	484	Total	C	N	O	S	0	2	0
			3780	2409	658	701	12			
1	H	488	Total	C	N	O	S	0	8	0
			3854	2459	671	711	13			

- 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	1
			48	21	7	17	3		
2	B	1	Total	C	N	O	P	0	1
			48	21	7	17	3		
2	C	1	Total	C	N	O	P	0	1
			48	21	7	17	3		
2	D	1	Total	C	N	O	P	0	1
			48	21	7	17	3		
2	E	1	Total	C	N	O	P	0	1
			48	21	7	17	3		
2	F	1	Total	C	N	O	P	0	1
			48	21	7	17	3		
2	G	1	Total	C	N	O	P	0	1
			48	21	7	17	3		
2	H	1	Total	C	N	O	P	0	1
			48	21	7	17	3		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	431	Total	O	0	7
			431	431		
3	B	385	Total	O	0	7
			385	385		
3	C	420	Total	O	0	10
			420	420		
3	D	356	Total	O	0	2
			356	356		

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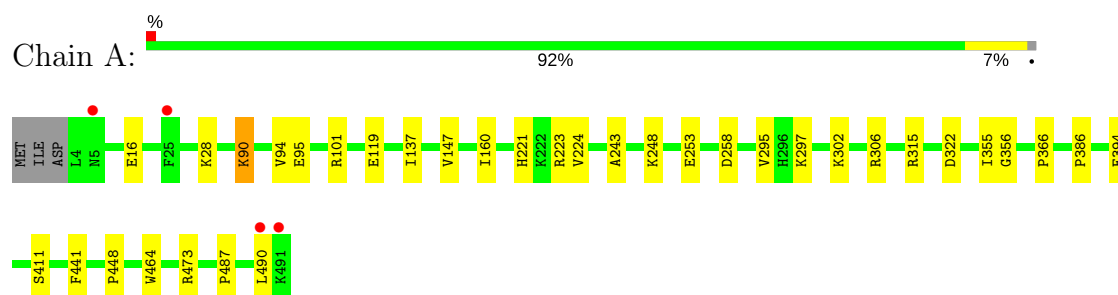
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	331	Total 331	O 331	0	7
3	F	382	Total 382	O 382	0	4
3	G	363	Total 363	O 363	0	6
3	H	420	Total 420	O 420	0	5

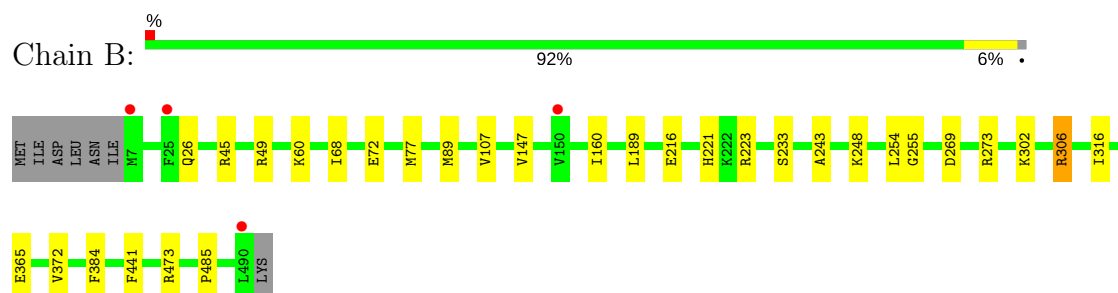
3 Residue-property plots [i](#)

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.

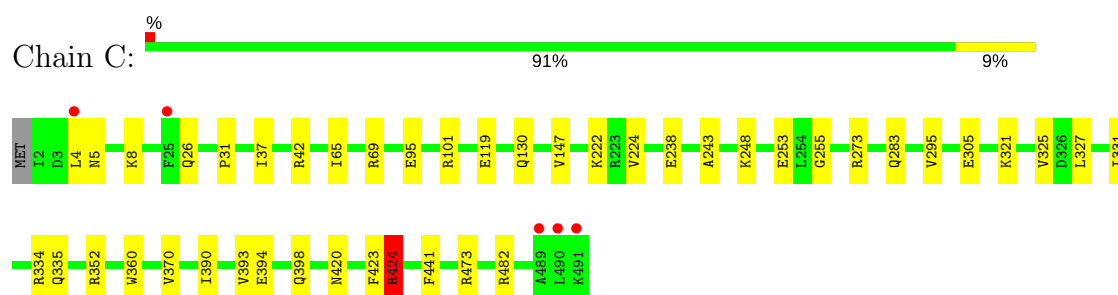
- Molecule 1: Aldehyde dehydrogenase



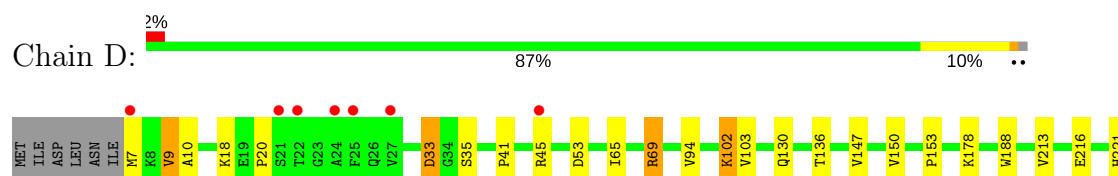
- Molecule 1: Aldehyde dehydrogenase

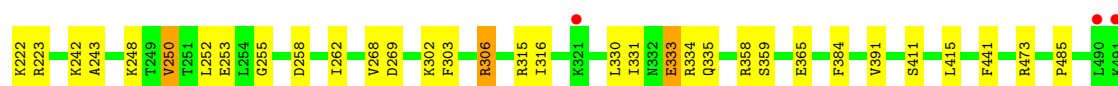


- Molecule 1: Aldehyde dehydrogenase

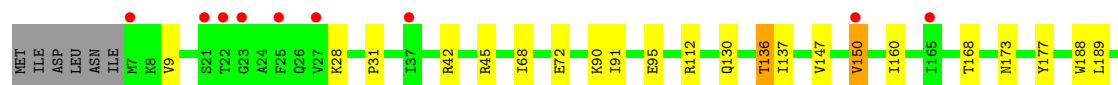
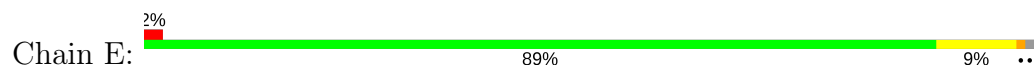


- Molecule 1: Aldehyde dehydrogenase

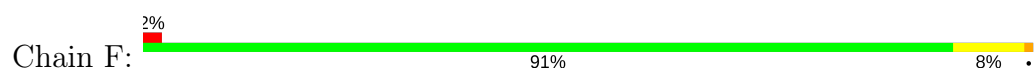




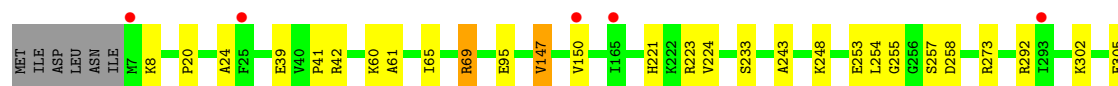
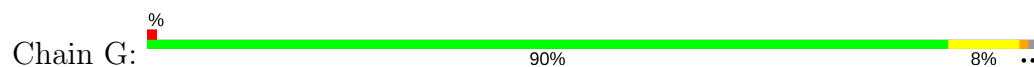
● Molecule 1: Aldehyde dehydrogenase



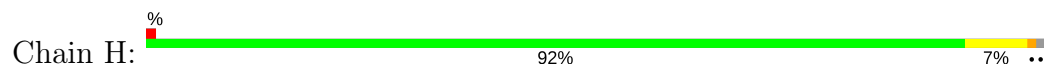
● Molecule 1: Aldehyde dehydrogenase



● Molecule 1: Aldehyde dehydrogenase



● Molecule 1: Aldehyde dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	163.81Å 185.36Å 208.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.14 – 1.90 29.63 – 1.90	Depositor EDS
% Data completeness (in resolution range)	87.6 (29.14-1.90) 87.6 (29.63-1.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 1.89Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.189 , 0.221 0.187 , 0.221	Depositor DCC
R_{free} test set	21738 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	25.0	Xtriage
Anisotropy	0.176	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	33972	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 47.96 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 9.2060e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.39	0/3889	0.53	0/5271
1	B	0.38	0/3863	0.52	0/5236
1	C	0.48	2/3947 (0.1%)	0.55	1/5351 (0.0%)
1	D	0.39	0/3866	0.52	0/5241
1	E	0.36	0/3868	0.51	0/5243
1	F	0.38	0/3909	0.53	0/5298
1	G	0.37	0/3859	0.52	0/5231
1	H	0.40	0/3937	0.53	0/5335
All	All	0.39	2/31138 (0.0%)	0.53	1/42206 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	8[A]	LYS	CD-CE	11.55	1.80	1.51
1	C	8[B]	LYS	CD-CE	11.55	1.80	1.51

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	424	ARG	NE-CZ-NH1	5.75	123.17	120.30

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	3810	0	3815	19	0
1	B	3784	0	3784	18	0
1	C	3866	0	3860	22	0
1	D	3787	0	3779	26	0
1	E	3789	0	3786	30	0
1	F	3830	0	3831	23	0
1	G	3780	0	3777	24	0
1	H	3854	0	3864	17	0
2	A	48	0	13	1	0
2	B	48	0	13	3	0
2	C	48	0	13	1	0
2	D	48	0	13	2	0
2	E	48	0	12	5	0
2	F	48	0	13	1	0
2	G	48	0	13	1	0
2	H	48	0	13	2	0
3	A	431	0	0	0	0
3	B	385	0	0	3	0
3	C	420	0	0	4	0
3	D	356	0	0	3	0
3	E	331	0	0	1	0
3	F	382	0	0	2	0
3	G	363	0	0	1	0
3	H	420	0	0	2	0
All	All	33972	0	30599	170	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 (170) 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:H:255:GLY:H	2:H:500[C]:NAP:H71N	1.27	0.83
1:E:255:GLY:H	2:E:500[C]:NAP:H71N	1.26	0.80
1:D:20:PRO:HG3	1:D:41:PRO:HB3	1.65	0.79
1:G:255:GLY:H	2:G:500[C]:NAP:H71N	1.36	0.74
1:A:315:ARG:NH2	1:A:322:ASP:OD2	2.23	0.69
1:G:20:PRO:HG3	1:G:41:PRO:HB3	1.73	0.69
1:B:255:GLY:H	2:B:500[C]:NAP:H71N	1.39	0.68
1:E:45:ARG:HH12	1:E:216:GLU:HG2	1.58	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:316:ILE:HD12	1:B:365:GLU:HG2	1.77	0.66
1:G:420:ASN:O	1:G:424:ARG:HD3	1.98	0.64
1:F:292:ARG:NH2	1:F:402:VAL:O	2.33	0.62
1:D:268:VAL:HG21	1:D:302:LYS:HE3	1.82	0.61
1:G:147:VAL:HG22	1:G:224:VAL:HA	1.81	0.61
1:E:28:LYS:HE2	1:E:95:GLU:HG3	1.83	0.61
1:G:305:GLU:HG2	1:G:306:ARG:HH12	1.66	0.60
1:H:33:ASP:HB3	1:H:35:SER:H	1.65	0.60
1:A:297:LYS:HD3	1:A:394:GLU:HG2	1.83	0.60
1:F:232:GLU:HG3	2:F:500[C]:NAP:H4D	1.84	0.60
1:B:302:LYS:NZ	3:B:832:HOH:O	2.35	0.59
1:F:221:HIS:HE1	1:F:223:ARG:HG3	1.68	0.57
1:D:33:ASP:HB3	1:D:35:SER:H	1.70	0.57
1:A:90:LYS:O	1:A:94:VAL:HG23	2.05	0.56
1:D:316:ILE:HD12	1:D:365:GLU:HG2	1.85	0.56
1:D:384:PHE:CE1	2:D:500[C]:NAP:H2D	2.41	0.56
1:C:5[B]:ASN:OD1	1:C:5[B]:ASN:N	2.38	0.56
1:E:160:ILE:HD11	2:E:500[C]:NAP:H5N	1.87	0.56
1:C:255:GLY:H	2:C:500[C]:NAP:H71N	1.54	0.56
1:H:315:ARG:NH1	3:H:1003:HOH:O	2.39	0.55
1:G:292:ARG:NH2	1:G:402:VAL:O	2.39	0.55
1:C:420:ASN:O	1:C:424:ARG:HD3	2.07	0.55
1:G:273:ARG:CZ	1:H:490:LEU:HD21	2.37	0.55
1:E:490:LEU:HD13	1:F:313:MET:HE1	1.89	0.54
1:E:393:VAL:HB	1:E:398:GLN:HB3	1.89	0.53
1:C:4[A]:LEU:HD22	1:C:37:ILE:HG23	1.91	0.53
1:B:89:MET:HG3	1:B:107:VAL:HG21	1.92	0.52
1:A:490:LEU:HD13	1:B:273:ARG:NH1	2.25	0.52
1:H:4[B]:LEU:HD22	1:H:37:ILE:HG23	1.91	0.51
1:F:18:LYS:NZ	1:F:50:GLU:OE2	2.35	0.51
1:F:243:ALA:HB1	1:F:248:LYS:HG3	1.92	0.50
1:H:69:ARG:HA	1:H:69:ARG:HE	1.76	0.50
1:G:370:VAL:HG12	1:G:390:ILE:HB	1.94	0.50
1:C:352:ARG:NH1	3:C:741:HOH:O	2.32	0.49
1:F:45:ARG:NH1	1:F:217:GLU:OE1	2.44	0.49
1:F:102:LYS:HE3	1:F:282:HIS:CE1	2.47	0.49
1:A:221:HIS:HE1	1:A:223:ARG:HG3	1.77	0.49
1:E:130:GLN:HG3	3:H:726:HOH:O	2.11	0.49
1:F:221:HIS:CE1	1:F:223:ARG:HG3	2.48	0.49
1:D:222:LYS:NZ	3:D:765:HOH:O	2.42	0.49
1:A:147:VAL:HG22	1:A:224:VAL:HA	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:487:PRO:HD2	1:A:490:LEU:HD12	1.95	0.49
1:C:273:ARG:NH2	3:C:953:HOH:O	2.46	0.49
1:B:269:ASP:OD1	1:B:306:ARG:NH1	2.46	0.48
1:C:243:ALA:HB1	1:C:248:LYS:HG3	1.95	0.48
1:A:243:ALA:HB1	1:A:248:LYS:HG3	1.95	0.48
1:H:250:VAL:HG13	1:H:252:LEU:HG	1.94	0.48
1:E:147:VAL:HG22	1:E:224:VAL:HA	1.94	0.48
1:D:358:ARG:HG3	1:D:359:SER:N	2.29	0.48
1:E:221:HIS:HE1	1:E:223:ARG:HG3	1.78	0.47
1:F:487:PRO:HD2	1:F:490:LEU:HD12	1.96	0.47
1:E:233:SER:HA	1:E:254:LEU:HB3	1.96	0.47
1:F:297:LYS:HE3	1:F:394:GLU:HG2	1.97	0.47
1:E:45:ARG:HD3	1:E:213:VAL:HG13	1.97	0.47
1:D:53:ASP:OD1	1:D:223:ARG:NH2	2.47	0.47
1:A:258:ASP:OD2	1:A:411:SER:HB3	2.15	0.47
1:D:243:ALA:HB1	1:D:248:LYS:HG3	1.97	0.47
1:E:91:ILE:HG13	1:E:189:LEU:HD11	1.97	0.47
1:G:223:ARG:HD3	3:G:788:HOH:O	2.13	0.47
1:F:306:ARG:HA	1:F:306:ARG:HD3	1.62	0.46
1:G:490:LEU:HD13	1:H:273:ARG:NH1	2.29	0.46
1:E:68:ILE:O	1:E:72:GLU:HG3	2.16	0.46
1:E:384:PHE:CE1	2:E:500[C]:NAP:H2D	2.51	0.46
1:C:65:ILE:HD12	1:C:69:ARG:HB3	1.98	0.46
1:G:65:ILE:HD12	1:G:69:ARG:HB3	1.97	0.46
1:A:101:ARG:HG2	1:B:485:PRO:HB2	1.98	0.46
1:D:69:ARG:HA	1:D:69:ARG:HE	1.79	0.46
1:D:102:LYS:HG3	1:D:103:VAL:N	2.31	0.46
1:A:119:GLU:HG2	1:C:119:GLU:HA	1.97	0.45
1:C:335:GLN:NE2	3:C:815:HOH:O	2.40	0.45
1:C:393:VAL:HB	1:C:398:GLN:HB3	1.98	0.45
1:G:8:LYS:HE3	1:G:39:GLU:OE2	2.16	0.45
1:A:137:ILE:HD12	1:C:423:PHE:CD2	2.51	0.45
1:C:147:VAL:HG22	1:C:224:VAL:HA	1.97	0.45
1:F:68:ILE:O	1:F:72:GLU:HG3	2.17	0.45
1:H:243:ALA:HB1	1:H:248:LYS:HG3	1.98	0.45
1:F:60:LYS:HB3	1:F:60:LYS:HE2	1.77	0.45
1:G:257:SER:OG	1:G:292:ARG:HD2	2.17	0.45
1:F:257:SER:OG	1:F:292:ARG:HD2	2.17	0.45
1:E:9:VAL:HG13	1:E:188:TRP:CG	2.52	0.45
1:E:112:ARG:NH1	3:E:736:HOH:O	2.50	0.45
1:E:255:GLY:HA3	2:E:500[C]:NAP:O2D	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:223:ARG:HD3	3:F:918:HOH:O	2.17	0.45
1:B:243:ALA:HB1	1:B:248:LYS:HG3	1.99	0.44
1:C:370:VAL:HG12	1:C:390:ILE:HB	1.99	0.44
1:G:306:ARG:HD3	1:G:306:ARG:HA	1.74	0.44
1:D:9:VAL:HB	1:D:188:TRP:CG	2.52	0.44
1:B:26:GLN:NE2	3:B:801:HOH:O	2.50	0.44
1:D:315:ARG:NH1	3:D:852:HOH:O	2.49	0.44
1:C:101:ARG:HG2	1:D:485:PRO:HB2	1.99	0.44
1:B:233:SER:HA	1:B:254:LEU:HB3	1.99	0.44
1:F:331:ILE:HG22	1:F:335:GLN:HG3	2.00	0.44
1:B:384:PHE:CZ	2:B:500[C]:NAP:H3D	2.52	0.44
1:F:233:SER:HA	1:F:254:LEU:HB3	1.98	0.44
1:F:262:ILE:HG12	1:F:415:LEU:HD12	2.00	0.44
1:H:258:ASP:OD2	1:H:411:SER:HB3	2.16	0.44
1:F:315:ARG:NH2	1:F:322:ASP:OD1	2.33	0.44
1:G:224:VAL:O	1:G:248:LYS:HE3	2.17	0.44
1:D:333:GLU:HG2	1:D:334:ARG:N	2.31	0.44
1:G:233:SER:HA	1:G:254:LEU:HB3	2.00	0.44
1:C:424:ARG:HG2	1:C:424:ARG:HH11	1.83	0.43
1:E:90:LYS:HE3	1:E:90:LYS:HB2	1.86	0.43
1:B:45:ARG:NH2	3:B:915:HOH:O	2.49	0.43
1:B:160:ILE:HD11	2:B:500[C]:NAP:H5N	2.00	0.43
1:D:258:ASP:OD2	1:D:411:SER:HB3	2.18	0.43
1:B:45:ARG:O	1:B:49:ARG:HG3	2.18	0.43
1:F:89:MET:HG3	1:F:107:VAL:HG21	2.00	0.43
1:D:255:GLY:H	2:D:500[C]:NAP:H71N	1.67	0.43
1:A:356:GLY:HA3	1:A:366:PRO:O	2.19	0.43
1:E:136:THR:HG21	3:F:830:HOH:O	2.18	0.43
1:B:68:ILE:O	1:B:72:GLU:HG3	2.19	0.43
1:D:331:ILE:HG22	1:D:335:GLN:HG3	2.01	0.43
1:C:26:GLN:OE1	1:C:42:ARG:HG3	2.19	0.42
1:G:24:ALA:HB1	1:G:42:ARG:NH1	2.34	0.42
1:A:160:ILE:HD11	2:A:500[C]:NAP:H5N	2.01	0.42
1:F:313:MET:HB3	1:F:313:MET:HE3	1.81	0.42
1:B:189:LEU:HD23	1:B:189:LEU:HA	1.92	0.42
1:E:31:PRO:HB3	1:E:331:ILE:O	2.19	0.42
1:H:384:PHE:CE1	2:H:500[C]:NAP:H3D	2.55	0.42
1:C:394:GLU:N	1:C:398:GLN:OE1	2.43	0.42
1:E:243:ALA:HB1	1:E:248:LYS:HG3	2.01	0.42
1:G:221:HIS:HE1	1:G:223:ARG:HG3	1.85	0.42
1:A:302:LYS:HE2	1:A:302:LYS:HB3	1.79	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:221:HIS:HE1	1:B:223:ARG:HG3	1.85	0.42
1:C:224:VAL:O	1:C:248:LYS:HE3	2.19	0.42
1:D:221:HIS:HE1	1:D:223:ARG:HG3	1.85	0.42
1:G:243:ALA:HB1	1:G:248:LYS:HG3	2.02	0.42
1:D:10:ALA:HB1	1:D:18:LYS:O	2.20	0.42
1:D:65:ILE:HD12	1:D:69:ARG:HB3	2.02	0.42
1:D:250:VAL:HG13	1:D:252:LEU:HG	2.02	0.41
1:G:258:ASP:OD2	1:G:411:SER:HB3	2.20	0.41
1:B:77:MET:HB3	1:B:77:MET:HE3	1.93	0.41
1:A:448:PRO:HB3	1:A:464:TRP:CG	2.55	0.41
1:C:283:GLN:HG3	1:C:327:LEU:HD22	2.01	0.41
1:H:331:ILE:HG22	1:H:335:GLN:HG3	2.03	0.41
1:A:355:ILE:HG13	1:A:356:GLY:N	2.36	0.41
1:D:262:ILE:HD13	1:D:303:PHE:CE1	2.55	0.41
1:E:224:VAL:O	1:E:248:LYS:HE3	2.20	0.41
1:E:137:ILE:HD12	1:G:423:PHE:CD2	2.56	0.41
1:E:222:LYS:HE2	1:E:222:LYS:HB2	1.88	0.41
1:E:297:LYS:HD3	1:E:394:GLU:HG2	2.01	0.41
1:E:160:ILE:CD1	2:E:500[C]:NAP:H5N	2.50	0.41
1:E:306:ARG:HA	1:E:306:ARG:HD3	1.78	0.41
1:D:94:VAL:HB	3:D:896:HOH:O	2.21	0.41
1:E:168:THR:HG22	1:E:173:ASN:HB2	2.03	0.41
1:E:356:GLY:HA3	1:E:366:PRO:O	2.20	0.41
1:H:7[B]:MET:HE3	1:H:38:ALA:HB2	2.03	0.41
1:C:130:GLN:HG3	3:C:919:HOH:O	2.21	0.40
1:F:222:LYS:HB3	1:F:222:LYS:HE2	1.73	0.40
1:A:306:ARG:HD3	1:A:306:ARG:HA	1.85	0.40
1:A:366:PRO:HA	1:A:386:PRO:HB2	2.03	0.40
1:G:393:VAL:HB	1:G:398:GLN:HB3	2.03	0.40
1:D:269:ASP:OD1	1:D:306:ARG:NH1	2.55	0.40
1:G:61:ALA:O	1:G:65:ILE:HG23	2.21	0.40
1:H:306:ARG:HD3	1:H:306:ARG:HA	1.72	0.40
1:C:31:PRO:HB3	1:C:331:ILE:O	2.21	0.40
1:D:45:ARG:NH1	1:D:213:VAL:HA	2.36	0.40
1:H:77:MET:HE3	1:H:77:MET:HB3	1.95	0.40
1:E:150:VAL:HG22	1:E:177:TYR:HD1	1.86	0.40
1:G:302:LYS:HB2	1:G:302:LYS:HE3	1.78	0.40
1:H:262:ILE:HD12	1:H:303:PHE:CE1	2.57	0.40
1:H:313:MET:HE3	1:H:313:MET:HB3	1.90	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	488/491 (99%)	475 (97%)	13 (3%)	0	100	100
1	B	484/491 (99%)	472 (98%)	12 (2%)	0	100	100
1	C	496/491 (101%)	483 (97%)	13 (3%)	0	100	100
1	D	485/491 (99%)	475 (98%)	9 (2%)	1 (0%)	51	41
1	E	485/491 (99%)	473 (98%)	12 (2%)	0	100	100
1	F	490/491 (100%)	478 (98%)	12 (2%)	0	100	100
1	G	484/491 (99%)	470 (97%)	14 (3%)	0	100	100
1	H	493/491 (100%)	481 (98%)	12 (2%)	0	100	100
All	All	3905/3928 (99%)	3807 (98%)	97 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	153	PRO

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	404/408 (99%)	395 (98%)	9 (2%)	57	51
1	B	401/408 (98%)	394 (98%)	7 (2%)	66	62
1	C	410/408 (100%)	395 (96%)	15 (4%)	39	28
1	D	400/408 (98%)	378 (94%)	22 (6%)	25	14

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	401/408 (98%)	389 (97%)	12 (3%)	46	37
1	F	406/408 (100%)	395 (97%)	11 (3%)	50	42
1	G	400/408 (98%)	386 (96%)	14 (4%)	41	30
1	H	409/408 (100%)	391 (96%)	18 (4%)	33	22
All	All	3231/3264 (99%)	3123 (97%)	108 (3%)	45	33

All (108) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	GLU
1	A	28	LYS
1	A	90	LYS
1	A	95	GLU
1	A	253[A]	GLU
1	A	253[B]	GLU
1	A	295	VAL
1	A	441	PHE
1	A	473	ARG
1	B	60	LYS
1	B	147	VAL
1	B	216	GLU
1	B	306	ARG
1	B	372	VAL
1	B	441	PHE
1	B	473	ARG
1	C	95	GLU
1	C	222	LYS
1	C	238	GLU
1	C	253[A]	GLU
1	C	253[B]	GLU
1	C	295	VAL
1	C	305	GLU
1	C	321	LYS
1	C	325	VAL
1	C	334	ARG
1	C	360	TRP
1	C	424	ARG
1	C	441	PHE
1	C	473	ARG
1	C	482	ARG
1	D	7	MET

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Mol	Chain	Res	Type
1	D	9	VAL
1	D	33	ASP
1	D	69	ARG
1	D	102	LYS
1	D	130	GLN
1	D	136	THR
1	D	147	VAL
1	D	150	VAL
1	D	178	LYS
1	D	216	GLU
1	D	242	LYS
1	D	250	VAL
1	D	253[A]	GLU
1	D	253[B]	GLU
1	D	306	ARG
1	D	330	LEU
1	D	333	GLU
1	D	391	VAL
1	D	415	LEU
1	D	441	PHE
1	D	473	ARG
1	E	42	ARG
1	E	136	THR
1	E	150	VAL
1	E	216	GLU
1	E	253[A]	GLU
1	E	253[B]	GLU
1	E	302	LYS
1	E	321	LYS
1	E	324	LYS
1	E	441	PHE
1	E	473	ARG
1	E	490	LEU
1	F	18	LYS
1	F	46	GLU
1	F	102	LYS
1	F	147	VAL
1	F	223	ARG
1	F	253[A]	GLU
1	F	253[B]	GLU
1	F	358	ARG
1	F	441	PHE

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Mol	Chain	Res	Type
1	F	446	HIS
1	F	473	ARG
1	G	60	LYS
1	G	69	ARG
1	G	95	GLU
1	G	147	VAL
1	G	150	VAL
1	G	253[A]	GLU
1	G	253[B]	GLU
1	G	306	ARG
1	G	325	VAL
1	G	360	TRP
1	G	424	ARG
1	G	441	PHE
1	G	457	VAL
1	G	473	ARG
1	H	5[A]	ASN
1	H	5[B]	ASN
1	H	33	ASP
1	H	42	ARG
1	H	69	ARG
1	H	86	GLU
1	H	95	GLU
1	H	102	LYS
1	H	150	VAL
1	H	250	VAL
1	H	253[A]	GLU
1	H	253[B]	GLU
1	H	330	LEU
1	H	358	ARG
1	H	372	VAL
1	H	441	PHE
1	H	446	HIS
1	H	473	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	26	GLN
1	D	26	GLN
1	E	130	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 ⓘ

8 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	500[C]	-	44,52,52	2.84	15 (34%)	51,80,80	1.95	10 (19%)
2	NAP	B	500[C]	-	44,52,52	2.82	14 (31%)	51,80,80	2.10	11 (21%)
2	NAP	C	500[C]	-	44,52,52	2.78	16 (36%)	51,80,80	1.85	6 (11%)
2	NAP	D	500[C]	-	44,52,52	2.77	15 (34%)	51,80,80	1.93	9 (17%)
2	NAP	E	500[C]	-	44,52,52	2.81	16 (36%)	51,80,80	2.13	10 (19%)
2	NAP	F	500[C]	-	44,52,52	2.71	15 (34%)	51,80,80	2.35	8 (15%)
2	NAP	G	500[C]	-	44,52,52	2.78	17 (38%)	51,80,80	1.81	10 (19%)
2	NAP	H	500[C]	-	44,52,52	2.79	15 (34%)	51,80,80	1.90	12 (23%)

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	500[C]	-	-	0/27/67/67	0/5/5/5
2	NAP	B	500[C]	-	-	0/27/67/67	0/5/5/5
2	NAP	C	500[C]	-	-	0/27/67/67	0/5/5/5
2	NAP	D	500[C]	-	-	0/27/67/67	0/5/5/5
2	NAP	E	500[C]	-	-	0/27/67/67	0/5/5/5
2	NAP	F	500[C]	-	-	0/27/67/67	0/5/5/5
2	NAP	G	500[C]	-	-	0/27/67/67	0/5/5/5
2	NAP	H	500[C]	-	-	0/27/67/67	0/5/5/5

All (123) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500[C]	NAP	C3B-C2B	-5.13	1.41	1.53
2	C	500[C]	NAP	C3B-C2B	-5.10	1.41	1.53
2	F	500[C]	NAP	C3B-C2B	-5.03	1.41	1.53
2	G	500[C]	NAP	C3B-C2B	-4.97	1.41	1.53
2	H	500[C]	NAP	C3B-C2B	-4.94	1.42	1.53
2	B	500[C]	NAP	C3B-C2B	-4.94	1.42	1.53
2	E	500[C]	NAP	C3B-C2B	-4.87	1.42	1.53
2	D	500[C]	NAP	C3B-C2B	-4.84	1.42	1.53
2	H	500[C]	NAP	O3D-C3D	-3.66	1.34	1.43
2	C	500[C]	NAP	O3D-C3D	-3.59	1.34	1.43
2	A	500[C]	NAP	O3D-C3D	-3.53	1.34	1.43
2	B	500[C]	NAP	O3D-C3D	-3.51	1.34	1.43
2	E	500[C]	NAP	O3D-C3D	-3.42	1.35	1.43
2	G	500[C]	NAP	O3D-C3D	-3.41	1.35	1.43
2	F	500[C]	NAP	O3D-C3D	-3.39	1.35	1.43
2	D	500[C]	NAP	C2D-C1D	-3.25	1.48	1.53
2	D	500[C]	NAP	O3D-C3D	-3.11	1.35	1.43
2	H	500[C]	NAP	C2D-C3D	-3.04	1.45	1.53
2	E	500[C]	NAP	C2D-C3D	-2.97	1.45	1.53
2	B	500[C]	NAP	C2D-C3D	-2.93	1.45	1.53
2	E	500[C]	NAP	O2D-C2D	-2.85	1.36	1.43
2	C	500[C]	NAP	C5D-C4D	-2.82	1.42	1.51
2	D	500[C]	NAP	C2D-C3D	-2.78	1.46	1.53
2	B	500[C]	NAP	C3D-C4D	-2.78	1.45	1.53
2	A	500[C]	NAP	C3D-C4D	-2.77	1.45	1.53
2	F	500[C]	NAP	C2D-C3D	-2.77	1.46	1.53
2	G	500[C]	NAP	C5D-C4D	-2.74	1.43	1.51
2	C	500[C]	NAP	C2D-C3D	-2.71	1.46	1.53
2	E	500[C]	NAP	C2D-C1D	-2.69	1.49	1.53
2	A	500[C]	NAP	C2D-C3D	-2.68	1.46	1.53
2	G	500[C]	NAP	C2D-C3D	-2.67	1.46	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	500[C]	NAP	C3D-C4D	-2.65	1.46	1.53
2	G	500[C]	NAP	C3D-C4D	-2.65	1.46	1.53
2	E	500[C]	NAP	C5D-C4D	-2.64	1.43	1.51
2	C	500[C]	NAP	C3D-C4D	-2.60	1.46	1.53
2	H	500[C]	NAP	C5D-C4D	-2.60	1.43	1.51
2	F	500[C]	NAP	C5D-C4D	-2.59	1.43	1.51
2	B	500[C]	NAP	C5D-C4D	-2.59	1.43	1.51
2	A	500[C]	NAP	C5D-C4D	-2.53	1.43	1.51
2	F	500[C]	NAP	C2D-C1D	-2.48	1.49	1.53
2	F	500[C]	NAP	C3D-C4D	-2.41	1.46	1.53
2	E	500[C]	NAP	C3D-C4D	-2.40	1.46	1.53
2	D	500[C]	NAP	O2D-C2D	-2.38	1.37	1.43
2	C	500[C]	NAP	O2D-C2D	-2.33	1.37	1.43
2	G	500[C]	NAP	O2D-C2D	-2.29	1.37	1.43
2	C	500[C]	NAP	C2D-C1D	-2.28	1.50	1.53
2	F	500[C]	NAP	O2D-C2D	-2.24	1.37	1.43
2	D	500[C]	NAP	C5D-C4D	-2.23	1.44	1.51
2	H	500[C]	NAP	O4B-C4B	-2.23	1.40	1.45
2	C	500[C]	NAP	O4B-C4B	-2.20	1.40	1.45
2	G	500[C]	NAP	C2D-C1D	-2.20	1.50	1.53
2	H	500[C]	NAP	O2D-C2D	-2.19	1.37	1.43
2	C	500[C]	NAP	O5D-C5D	-2.16	1.36	1.44
2	E	500[C]	NAP	O4B-C4B	-2.15	1.40	1.45
2	A	500[C]	NAP	O2D-C2D	-2.13	1.38	1.43
2	B	500[C]	NAP	O2D-C2D	-2.13	1.38	1.43
2	A	500[C]	NAP	O4B-C4B	-2.10	1.40	1.45
2	D	500[C]	NAP	C3B-C4B	-2.09	1.47	1.53
2	A	500[C]	NAP	C3B-C4B	-2.09	1.47	1.53
2	B	500[C]	NAP	O4B-C4B	-2.08	1.40	1.45
2	G	500[C]	NAP	O4B-C4B	-2.08	1.40	1.45
2	G	500[C]	NAP	C3B-C4B	-2.08	1.47	1.53
2	F	500[C]	NAP	C3B-C4B	-2.05	1.47	1.53
2	D	500[C]	NAP	O4B-C4B	-2.04	1.40	1.45
2	H	500[C]	NAP	C3B-C4B	-2.03	1.47	1.53
2	E	500[C]	NAP	C3B-C4B	-2.03	1.47	1.53
2	G	500[C]	NAP	O5D-C5D	-2.02	1.36	1.44
2	C	500[C]	NAP	C6N-N1N	2.69	1.42	1.35
2	G	500[C]	NAP	C6N-N1N	2.74	1.42	1.35
2	D	500[C]	NAP	C6N-N1N	2.78	1.42	1.35
2	F	500[C]	NAP	C6N-N1N	2.84	1.42	1.35
2	A	500[C]	NAP	C6N-N1N	2.92	1.43	1.35
2	H	500[C]	NAP	C6N-N1N	2.93	1.43	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	500[C]	NAP	C6N-N1N	2.93	1.43	1.35
2	B	500[C]	NAP	C6N-N1N	2.99	1.43	1.35
2	D	500[C]	NAP	C6A-N6A	3.72	1.49	1.34
2	F	500[C]	NAP	C6A-N6A	3.76	1.49	1.34
2	H	500[C]	NAP	C6A-N6A	3.78	1.49	1.34
2	B	500[C]	NAP	C6A-N6A	3.83	1.49	1.34
2	E	500[C]	NAP	C6A-N6A	3.83	1.49	1.34
2	A	500[C]	NAP	C6A-N6A	3.84	1.49	1.34
2	C	500[C]	NAP	C6A-N6A	3.85	1.49	1.34
2	G	500[C]	NAP	C6A-N6A	3.86	1.49	1.34
2	C	500[C]	NAP	C6N-C5N	4.44	1.48	1.38
2	F	500[C]	NAP	C6N-C5N	4.45	1.48	1.38
2	A	500[C]	NAP	C6N-C5N	4.68	1.48	1.38
2	G	500[C]	NAP	C6N-C5N	4.74	1.49	1.38
2	B	500[C]	NAP	C6N-C5N	4.75	1.49	1.38
2	E	500[C]	NAP	C6N-C5N	4.77	1.49	1.38
2	H	500[C]	NAP	C6N-C5N	4.79	1.49	1.38
2	D	500[C]	NAP	C6N-C5N	4.82	1.49	1.38
2	F	500[C]	NAP	C5N-C4N	5.46	1.49	1.38
2	C	500[C]	NAP	C5N-C4N	5.77	1.49	1.38
2	E	500[C]	NAP	C5N-C4N	5.91	1.50	1.38
2	A	500[C]	NAP	C5N-C4N	5.99	1.50	1.38
2	H	500[C]	NAP	C5N-C4N	6.03	1.50	1.38
2	D	500[C]	NAP	C5N-C4N	6.06	1.50	1.38
2	G	500[C]	NAP	C5N-C4N	6.10	1.50	1.38
2	B	500[C]	NAP	C5N-C4N	6.15	1.50	1.38
2	F	500[C]	NAP	C4N-C3N	6.24	1.49	1.39
2	G	500[C]	NAP	C4N-C3N	6.28	1.49	1.39
2	D	500[C]	NAP	C4N-C3N	6.35	1.49	1.39
2	B	500[C]	NAP	C4N-C3N	6.39	1.49	1.39
2	A	500[C]	NAP	C4N-C3N	6.41	1.49	1.39
2	E	500[C]	NAP	C4N-C3N	6.43	1.49	1.39
2	C	500[C]	NAP	C4N-C3N	6.45	1.49	1.39
2	H	500[C]	NAP	C4N-C3N	6.51	1.50	1.39
2	H	500[C]	NAP	C7N-N7N	7.41	1.47	1.33
2	B	500[C]	NAP	C7N-N7N	7.43	1.47	1.33
2	F	500[C]	NAP	C2N-C3N	7.44	1.50	1.39
2	C	500[C]	NAP	C7N-N7N	7.53	1.47	1.33
2	G	500[C]	NAP	C7N-N7N	7.57	1.47	1.33
2	E	500[C]	NAP	C7N-N7N	7.58	1.47	1.33
2	A	500[C]	NAP	C7N-N7N	7.59	1.47	1.33
2	D	500[C]	NAP	C7N-N7N	7.65	1.47	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	500[C]	NAP	C7N-N7N	7.70	1.48	1.33
2	G	500[C]	NAP	C2N-C3N	7.81	1.50	1.39
2	D	500[C]	NAP	C2N-C3N	7.90	1.51	1.39
2	C	500[C]	NAP	C2N-C3N	7.98	1.51	1.39
2	E	500[C]	NAP	C2N-C3N	8.05	1.51	1.39
2	H	500[C]	NAP	C2N-C3N	8.09	1.51	1.39
2	B	500[C]	NAP	C2N-C3N	8.57	1.52	1.39
2	A	500[C]	NAP	C2N-C3N	8.61	1.52	1.39

All (76) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	500[C]	NAP	C5N-C4N-C3N	-8.24	110.66	120.35
2	B	500[C]	NAP	N3A-C2A-N1A	-7.32	122.49	128.86
2	D	500[C]	NAP	N3A-C2A-N1A	-7.20	122.58	128.86
2	B	500[C]	NAP	C1B-N9A-C4A	-7.16	114.26	126.64
2	F	500[C]	NAP	N3A-C2A-N1A	-7.10	122.67	128.86
2	E	500[C]	NAP	N3A-C2A-N1A	-6.99	122.77	128.86
2	F	500[C]	NAP	C1B-N9A-C4A	-6.66	115.13	126.64
2	H	500[C]	NAP	N3A-C2A-N1A	-6.62	123.09	128.86
2	C	500[C]	NAP	N3A-C2A-N1A	-6.55	123.15	128.86
2	G	500[C]	NAP	N3A-C2A-N1A	-6.47	123.22	128.86
2	A	500[C]	NAP	N3A-C2A-N1A	-6.37	123.31	128.86
2	C	500[C]	NAP	C1B-N9A-C4A	-6.34	115.69	126.64
2	D	500[C]	NAP	C1B-N9A-C4A	-6.32	115.71	126.64
2	E	500[C]	NAP	C5N-C4N-C3N	-6.11	113.16	120.35
2	E	500[C]	NAP	C1B-N9A-C4A	-6.01	116.25	126.64
2	A	500[C]	NAP	C5N-C4N-C3N	-5.85	113.47	120.35
2	G	500[C]	NAP	C1B-N9A-C4A	-5.83	116.57	126.64
2	D	500[C]	NAP	C5N-C4N-C3N	-5.75	113.59	120.35
2	C	500[C]	NAP	C5N-C4N-C3N	-5.42	113.98	120.35
2	H	500[C]	NAP	C1B-N9A-C4A	-5.41	117.29	126.64
2	E	500[C]	NAP	C4D-O4D-C1D	-5.30	104.13	109.77
2	H	500[C]	NAP	C5N-C4N-C3N	-5.27	114.15	120.35
2	G	500[C]	NAP	C5N-C4N-C3N	-5.17	114.26	120.35
2	B	500[C]	NAP	C5N-C4N-C3N	-5.15	114.30	120.35
2	A	500[C]	NAP	C1B-N9A-C4A	-4.89	118.18	126.64
2	F	500[C]	NAP	C3N-C2N-N1N	-4.89	115.51	120.43
2	F	500[C]	NAP	O7N-C7N-C3N	-4.44	114.44	119.62
2	F	500[C]	NAP	C4N-C3N-C7N	-3.42	111.97	121.07
2	A	500[C]	NAP	O7N-C7N-N7N	-3.27	117.93	122.58
2	H	500[C]	NAP	O2D-C2D-C3D	-3.20	101.58	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500[C]	NAP	O7N-C7N-N7N	-2.99	118.33	122.58
2	B	500[C]	NAP	C4D-O4D-C1D	-2.98	106.59	109.77
2	E	500[C]	NAP	O2D-C2D-C3D	-2.90	102.52	111.83
2	B	500[C]	NAP	O2D-C2D-C3D	-2.90	102.54	111.83
2	H	500[C]	NAP	O7N-C7N-N7N	-2.89	118.47	122.58
2	G	500[C]	NAP	O4B-C1B-C2B	-2.79	101.71	106.59
2	A	500[C]	NAP	O4B-C1B-C2B	-2.72	101.84	106.59
2	G	500[C]	NAP	C4D-O4D-C1D	-2.69	106.91	109.77
2	C	500[C]	NAP	O7N-C7N-N7N	-2.67	118.78	122.58
2	D	500[C]	NAP	O7N-C7N-N7N	-2.50	119.02	122.58
2	D	500[C]	NAP	C4D-O4D-C1D	-2.47	107.14	109.77
2	A	500[C]	NAP	O2D-C2D-C3D	-2.46	103.95	111.83
2	E	500[C]	NAP	O7N-C7N-N7N	-2.46	119.09	122.58
2	H	500[C]	NAP	C4A-C5A-N7A	-2.42	107.07	109.41
2	F	500[C]	NAP	C5N-C6N-N1N	-2.33	116.82	120.40
2	A	500[C]	NAP	C4A-C5A-N7A	-2.33	107.16	109.41
2	H	500[C]	NAP	O4B-C1B-C2B	-2.28	102.60	106.59
2	G	500[C]	NAP	O7N-C7N-N7N	-2.24	119.40	122.58
2	H	500[C]	NAP	C4D-O4D-C1D	-2.21	107.41	109.77
2	D	500[C]	NAP	C4A-C5A-N7A	-2.21	107.28	109.41
2	C	500[C]	NAP	C4A-C5A-N7A	-2.08	107.40	109.41
2	B	500[C]	NAP	O4D-C4D-C3D	-2.04	101.11	105.17
2	A	500[C]	NAP	C4D-O4D-C1D	-2.01	107.63	109.77
2	E	500[C]	NAP	C6N-C5N-C4N	2.02	122.49	119.44
2	D	500[C]	NAP	C2N-C3N-C4N	2.04	120.59	118.26
2	E	500[C]	NAP	C2N-C3N-C4N	2.06	120.61	118.26
2	G	500[C]	NAP	O5D-C5D-C4D	2.10	116.45	109.00
2	H	500[C]	NAP	C6N-C5N-C4N	2.17	122.72	119.44
2	B	500[C]	NAP	C2N-C3N-C4N	2.21	120.78	118.26
2	G	500[C]	NAP	C2N-C3N-C4N	2.24	120.81	118.26
2	B	500[C]	NAP	C6N-C5N-C4N	2.31	122.93	119.44
2	G	500[C]	NAP	C3N-C7N-N7N	2.34	120.44	117.77
2	G	500[C]	NAP	C6N-C5N-C4N	2.44	123.12	119.44
2	H	500[C]	NAP	C2N-C3N-C4N	2.47	121.08	118.26
2	E	500[C]	NAP	O5D-C5D-C4D	2.74	118.72	109.00
2	H	500[C]	NAP	O5D-C5D-C4D	2.79	118.88	109.00
2	B	500[C]	NAP	O5D-C5D-C4D	2.80	118.92	109.00
2	A	500[C]	NAP	O5D-C5D-C4D	2.96	119.50	109.00
2	D	500[C]	NAP	C3N-C7N-N7N	3.04	121.24	117.77
2	D	500[C]	NAP	O5D-C5D-C4D	3.15	120.18	109.00
2	H	500[C]	NAP	C3N-C7N-N7N	3.22	121.45	117.77
2	C	500[C]	NAP	C3N-C7N-N7N	3.69	121.99	117.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	500[C]	NAP	C3N-C7N-N7N	3.99	122.33	117.77
2	B	500[C]	NAP	C3N-C7N-N7N	4.23	122.61	117.77
2	A	500[C]	NAP	C3N-C7N-N7N	4.78	123.23	117.77
2	F	500[C]	NAP	C3N-C7N-N7N	5.17	123.68	117.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500[C]	NAP	1	0
2	B	500[C]	NAP	3	0
2	C	500[C]	NAP	1	0
2	D	500[C]	NAP	2	0
2	E	500[C]	NAP	5	0
2	F	500[C]	NAP	1	0
2	G	500[C]	NAP	1	0
2	H	500[C]	NAP	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	488/491 (99%)	-0.25	4 (0%) 86 87	9, 17, 33, 52	0
1	B	484/491 (98%)	-0.27	4 (0%) 86 87	10, 18, 36, 55	0
1	C	490/491 (99%)	-0.28	5 (1%) 82 84	10, 18, 32, 66	3 (0%)
1	D	485/491 (98%)	-0.18	10 (2%) 64 67	10, 21, 40, 64	0
1	E	485/491 (98%)	-0.10	12 (2%) 58 62	11, 23, 42, 67	0
1	F	490/491 (99%)	-0.29	8 (1%) 72 75	10, 19, 36, 59	0
1	G	484/491 (98%)	-0.25	7 (1%) 75 78	9, 21, 36, 48	0
1	H	488/491 (99%)	-0.29	6 (1%) 79 82	9, 17, 33, 65	1 (0%)
All	All	3894/3928 (99%)	-0.24	56 (1%) 75 78	9, 19, 36, 67	4 (0%)

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	491	LYS	7.4
1	D	7	MET	6.9
1	F	491	LYS	5.2
1	C	491	LYS	5.1
1	H	491	LYS	5.0
1	D	25	PHE	4.8
1	D	491	LYS	4.7
1	H	490	LEU	4.6
1	E	490	LEU	4.0
1	A	490	LEU	4.0
1	B	490	LEU	3.9
1	C	490	LEU	3.8
1	E	23	GLY	3.7
1	B	25	PHE	3.7
1	A	491	LYS	3.7
1	C	25	PHE	3.6

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Mol	Chain	Res	Type	RSRZ
1	F	25	PHE	3.6
1	D	22	THR	3.5
1	E	25	PHE	3.5
1	E	27	VAL	3.4
1	G	25	PHE	3.3
1	F	4	LEU	3.3
1	E	7	MET	3.2
1	G	490	LEU	3.2
1	D	490	LEU	3.1
1	B	7	MET	3.0
1	F	490	LEU	2.9
1	H	4[A]	LEU	2.9
1	D	21	SER	2.8
1	E	489	ALA	2.8
1	E	165	ILE	2.8
1	C	4[A]	LEU	2.7
1	C	489	ALA	2.7
1	H	25	PHE	2.7
1	B	150	VAL	2.6
1	G	150	VAL	2.6
1	A	25	PHE	2.6
1	D	27	VAL	2.6
1	G	165	ILE	2.6
1	D	24	ALA	2.5
1	A	5	ASN	2.5
1	G	293	ILE	2.4
1	F	21	SER	2.4
1	H	6[A]	ILE	2.3
1	E	21	SER	2.3
1	E	150	VAL	2.2
1	G	7	MET	2.2
1	F	5	ASN	2.2
1	E	37	ILE	2.1
1	D	45	ARG	2.1
1	F	297	LYS	2.1
1	E	22	THR	2.0
1	F	22	THR	2.0
1	H	150	VAL	2.0
1	G	489	ALA	2.0
1	D	321	LYS	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(Å ²)	Q<0.9
2	NAP	D	500[C]	48/48	0.92	0.15	1.59	20,26,31,38	21
2	NAP	B	500[C]	48/48	0.93	0.14	1.08	18,25,30,33	21
2	NAP	H	500[C]	48/48	0.95	0.12	1.02	15,22,29,30	17
2	NAP	C	500[C]	48/48	0.94	0.12	0.73	16,22,29,32	17
2	NAP	F	500[C]	48/48	0.95	0.12	0.71	15,25,31,33	17
2	NAP	E	500[C]	48/48	0.93	0.12	0.58	19,31,38,42	21
2	NAP	G	500[C]	48/48	0.95	0.11	0.49	17,25,32,36	17
2	NAP	A	500[C]	48/48	0.95	0.10	0.19	14,23,27,33	17

6.5 Other polymers [i](#)

There are no such residues in this entry.