



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 12:00 AM GMT

PDB ID : 2NYN
Title : Crystal structure of phenylalanine ammonia-lyase from *Anabaena variabilis*
Authors : Louie, G.V.; Moffitt, M.C.; Bowman, M.E.; Pence, J.; Noel, J.P.; Moore, B.S.
Deposited on : 2006-11-21
Resolution : 1.90 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

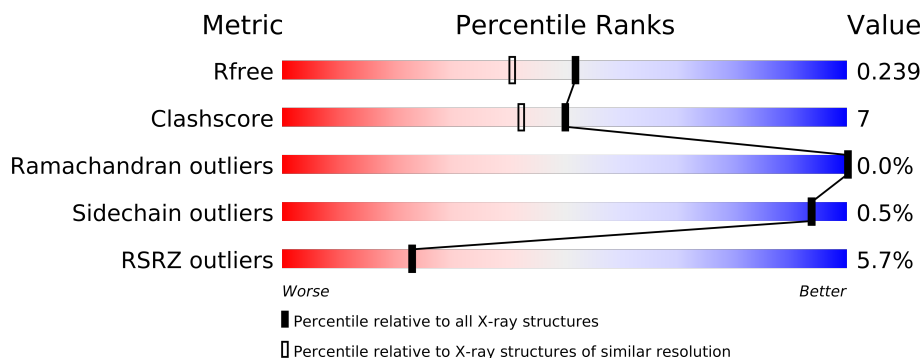
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **FAILED**
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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}	66092	3684 (1.90-1.90)
Clashscore	79885	4465 (1.90-1.90)
Ramachandran outliers	78287	4413 (1.90-1.90)
Sidechain outliers	78261	4414 (1.90-1.90)
RSRZ outliers	66119	3686 (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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	565	
1	B	565	
1	C	565	
1	D	565	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 16261 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Phenylalanine/histidineammonia-lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	512	Total	C	N	O	S	0	0	0
			3930	2477	690	744	19			
1	B	512	Total	C	N	O	S	0	0	0
			3930	2477	690	744	19			
1	C	512	Total	C	N	O	S	0	0	0
			3930	2477	690	744	19			
1	D	512	Total	C	N	O	S	0	0	0
			3930	2477	690	744	19			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	167	MDO	ALA	SEE REMARK 999	UNP Q3M5Z3
A	167	MDO	SER	SEE REMARK 999	UNP Q3M5Z3
A	167	MDO	GLY	SEE REMARK 999	UNP Q3M5Z3
B	167	MDO	ALA	SEE REMARK 999	UNP Q3M5Z3
B	167	MDO	SER	SEE REMARK 999	UNP Q3M5Z3
B	167	MDO	GLY	SEE REMARK 999	UNP Q3M5Z3
C	167	MDO	ALA	SEE REMARK 999	UNP Q3M5Z3
C	167	MDO	SER	SEE REMARK 999	UNP Q3M5Z3
C	167	MDO	GLY	SEE REMARK 999	UNP Q3M5Z3
D	167	MDO	ALA	SEE REMARK 999	UNP Q3M5Z3
D	167	MDO	SER	SEE REMARK 999	UNP Q3M5Z3
D	167	MDO	GLY	SEE REMARK 999	UNP Q3M5Z3

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	142	Total	O	0	0
			142	142		
2	B	131	Total	O	0	0
			131	131		

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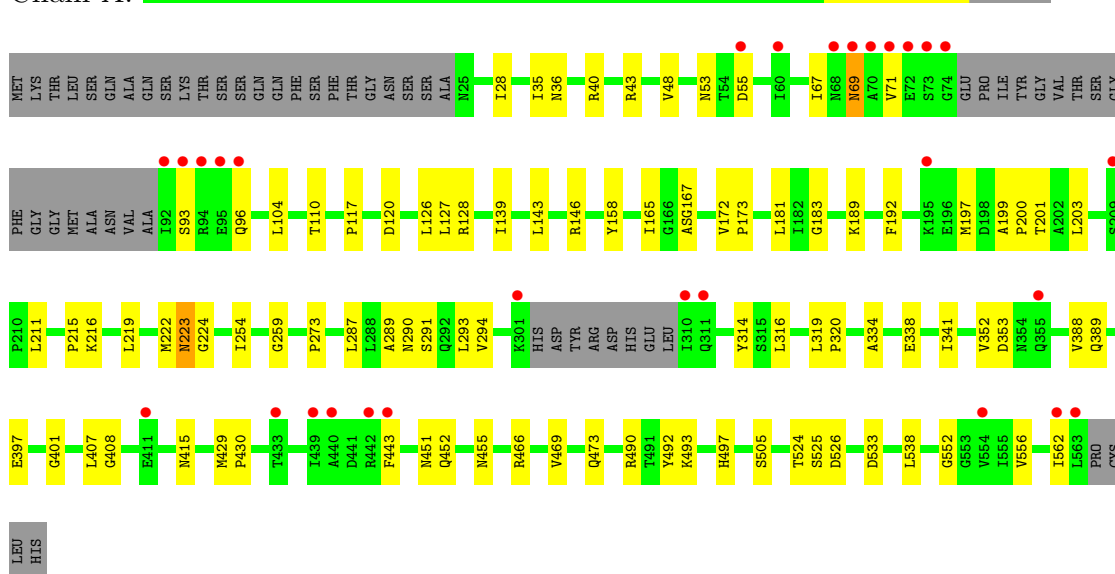
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	143	Total 143	O 143	0	0
2	D	125	Total 125	O 125	0	0

3 Residue-property plots

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.

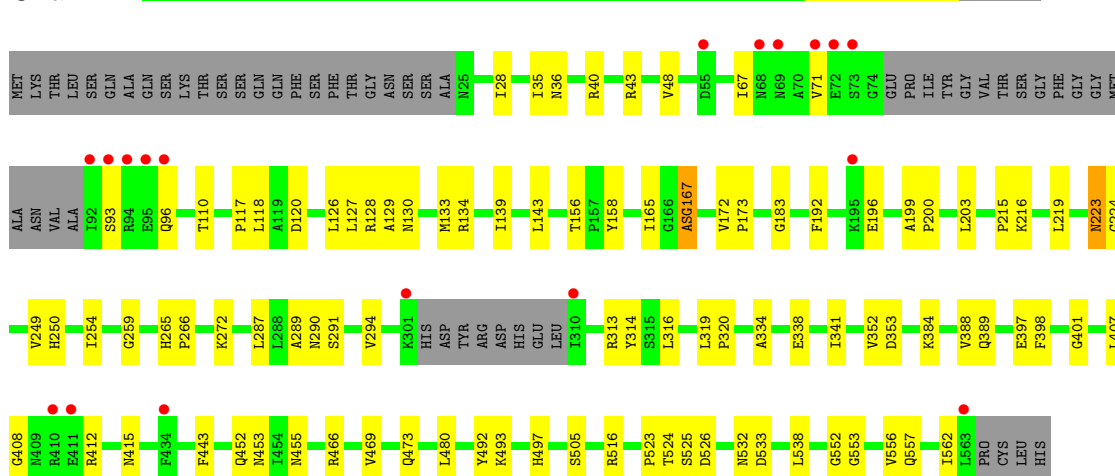
• Molecule 1: Phenylalanine/histidineammonia-lyase

Chain A:



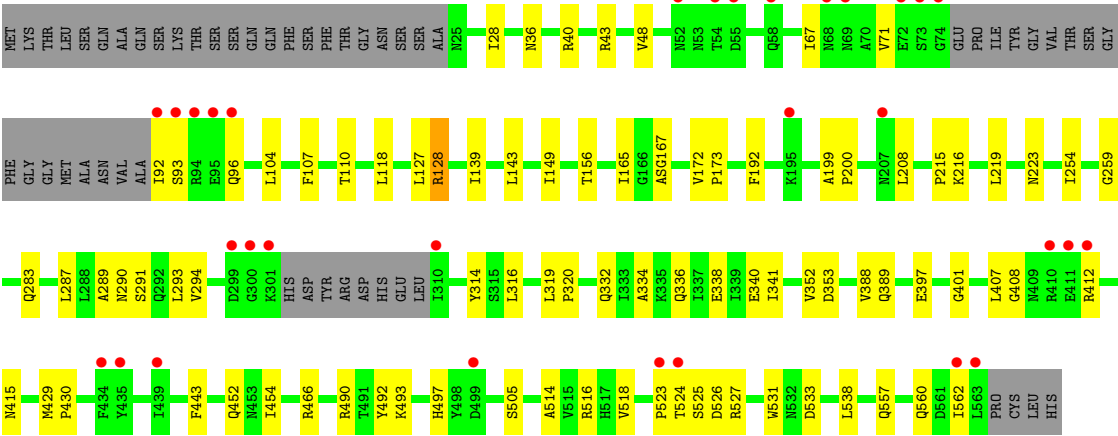
• Molecule 1: Phenylalanine/histidineammonia-lyase

Chain B:



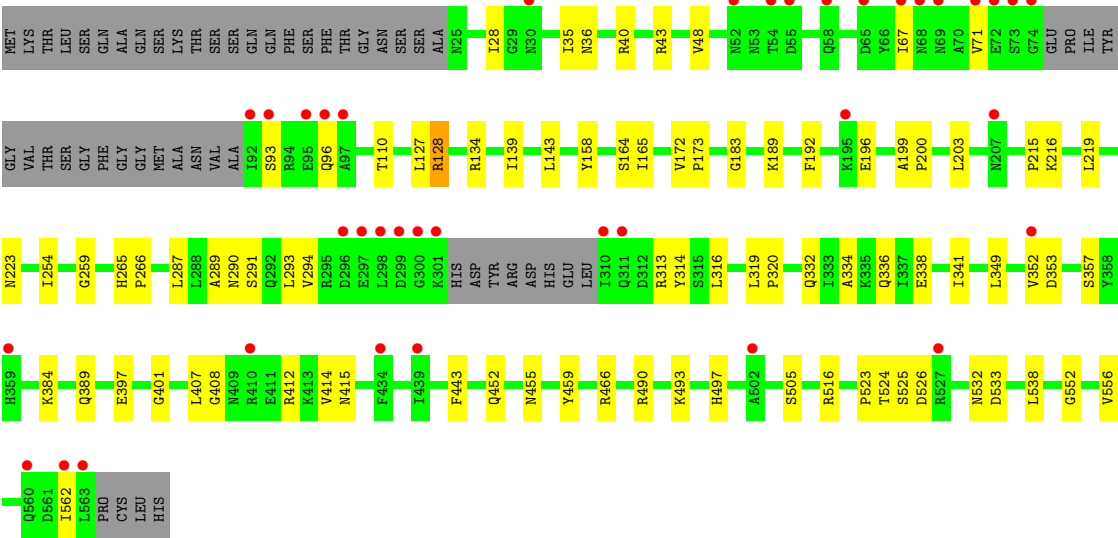
• Molecule 1: Phenylalanine/histidineammonia-lyase

Chain C:



• Molecule 1: Phenylalanine/histidineammonia-lyase

Chain D:



4 Data and refinement statistics

Xtriage (Phenix) failed to run properly - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	78.19Å 88.52Å 90.06Å 103.52° 97.82° 116.24°	Depositor
Resolution (Å)	100.00 – 1.90 48.56 – 1.74	Depositor EDS
% Data completeness (in resolution range)	(Not available) (100.00-1.90) 84.0 (48.56-1.74)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.223 , 0.242 0.220 , 0.239	Depositor DCC
R_{free} test set	7620 reflections (5.29%)	DCC
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 42.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	16261	wwPDB-VP
Average B, all atoms (Å ²)	30.0	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: MDO

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	0.33	0/3984	0.57	0/5406
1	B	0.33	0/3984	0.56	0/5406
1	C	0.33	0/3984	0.56	0/5406
1	D	0.33	0/3984	0.56	0/5406
All	All	0.33	0/15936	0.56	0/21624

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	167	MDO	Mainchain
1	B	167	MDO	Mainchain

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3930	0	3927	63	0
1	B	3930	0	3926	70	0
1	C	3930	0	3927	63	0
1	D	3930	0	3927	59	0
2	A	142	0	0	1	0
2	B	131	0	0	5	0
2	C	143	0	0	1	0
2	D	125	0	0	5	0
All	All	16261	0	15707	235	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 7.

All (235) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:287:LEU:HD21	1:D:562:ILE:HD11	1.39	1.04
1:C:287:LEU:HD21	1:C:562:ILE:HD11	1.46	0.98
1:A:287:LEU:HD21	1:A:562:ILE:HD11	1.45	0.98
1:C:110:THR:HG22	1:D:538:LEU:HB2	1.51	0.92
1:B:287:LEU:HD21	1:B:562:ILE:HD11	1.51	0.91
1:C:538:LEU:HB2	1:D:110:THR:HG22	1.56	0.88
1:B:128:ARG:HH12	1:B:224:GLY:HA3	1.41	0.84
1:D:128:ARG:HA	1:D:128:ARG:HE	1.45	0.81
1:A:538:LEU:HB2	1:B:110:THR:HG22	1.63	0.80
1:C:128:ARG:HE	1:C:128:ARG:HA	1.48	0.78
1:D:524:THR:HG22	1:D:526:ASP:H	1.49	0.78
1:C:127:LEU:HD22	1:C:341:ILE:HG23	1.65	0.77
1:A:110:THR:HG22	1:B:538:LEU:HB2	1.65	0.77
1:A:524:THR:HG22	1:A:526:ASP:H	1.50	0.77
1:C:524:THR:HG22	1:C:526:ASP:H	1.52	0.74
1:B:524:THR:HG22	1:B:526:ASP:H	1.50	0.74
1:D:127:LEU:HD22	1:D:341:ILE:HG23	1.72	0.70
1:C:28:ILE:HD11	1:C:48:VAL:HG11	1.72	0.69
1:A:128:ARG:HH12	1:A:224:GLY:HA3	1.58	0.69

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:452:GLN:HG2	1:C:314:TYR:CE2	2.29	0.67
1:D:414:VAL:HG13	2:D:1393:HOH:O	1.95	0.66
1:C:557:GLN:HA	1:C:560:GLN:HG3	1.79	0.65
1:B:128:ARG:NH1	1:B:224:GLY:HA3	2.11	0.64
1:A:452:GLN:HG2	1:D:314:TYR:CE2	2.33	0.63
1:D:316:LEU:HD23	1:D:389:GLN:HE22	1.63	0.63
1:C:316:LEU:HD23	1:C:389:GLN:HE22	1.64	0.62
1:A:69:ASN:N	1:A:69:ASN:HD22	1.96	0.62
1:C:407:LEU:HD12	1:C:408:GLY:N	2.15	0.62
1:C:407:LEU:HD12	1:C:408:GLY:H	1.65	0.61
1:D:352:VAL:HG22	2:D:1509:HOH:O	1.99	0.61
1:B:165:ILE:HD11	1:B:452:GLN:OE1	2.01	0.61
1:B:199:ALA:HB3	1:B:200:PRO:HD3	1.83	0.60
1:A:314:TYR:CE2	1:D:452:GLN:HG2	2.36	0.60
1:D:172:VAL:HB	1:D:173:PRO:HD3	1.83	0.59
1:A:127:LEU:HD22	1:A:341:ILE:HG23	1.84	0.59
1:A:291:SER:HB2	1:A:505:SER:HB3	1.85	0.59
1:A:452:GLN:HG2	1:D:314:TYR:CZ	2.38	0.59
1:D:219:LEU:O	1:D:223:ASN:HB2	2.03	0.59
1:A:199:ALA:HB3	1:A:200:PRO:HD3	1.86	0.58
1:B:172:VAL:HB	1:B:173:PRO:HD3	1.86	0.58
1:B:127:LEU:HD22	1:B:341:ILE:HG23	1.86	0.57
1:C:43:ARG:HH21	1:C:43:ARG:HG3	1.69	0.57
1:D:28:ILE:HD11	1:D:48:VAL:HG11	1.86	0.57
1:C:92:ILE:O	1:C:93:SER:HB3	2.03	0.57
1:C:497:HIS:HB2	1:C:525:SER:HB2	1.86	0.57
1:A:172:VAL:HB	1:A:173:PRO:HD3	1.87	0.57
1:B:139:ILE:HD11	1:B:143:LEU:HD23	1.86	0.57
1:A:43:ARG:HG3	1:A:43:ARG:HH21	1.69	0.57
1:C:516:ARG:NH2	1:C:523:PRO:HA	2.20	0.57
1:B:452:GLN:HG2	1:C:314:TYR:CZ	2.41	0.56
1:D:139:ILE:HD11	1:D:143:LEU:HD23	1.87	0.56
1:C:219:LEU:O	1:C:223:ASN:HB2	2.05	0.56
1:C:199:ALA:HB3	1:C:200:PRO:HD3	1.86	0.56
1:A:139:ILE:HD11	1:A:143:LEU:HD23	1.87	0.56
1:C:36:ASN:O	1:C:40:ARG:HG3	2.06	0.55
1:B:314:TYR:CE2	1:C:452:GLN:HG2	2.41	0.55
1:A:28:ILE:HD11	1:A:126:LEU:HD12	1.88	0.55
1:A:67:ILE:O	1:A:71:VAL:HG23	2.06	0.55
1:C:93:SER:HB3	1:C:96:GLN:HB3	1.89	0.55
1:B:28:ILE:HD11	1:B:48:VAL:HG11	1.87	0.55
1:A:36:ASN:O	1:A:40:ARG:HG3	2.06	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:36:ASN:O	1:D:40:ARG:HG3	2.07	0.55
1:C:67:ILE:O	1:C:71:VAL:HG23	2.07	0.55
1:C:291:SER:HB2	1:C:505:SER:HB3	1.90	0.54
1:B:407:LEU:HD12	1:B:408:GLY:N	2.22	0.54
1:C:139:ILE:HD11	1:C:143:LEU:HD23	1.89	0.54
1:D:67:ILE:O	1:D:71:VAL:HG23	2.07	0.54
1:B:314:TYR:CZ	1:C:452:GLN:HG2	2.43	0.54
1:A:314:TYR:CZ	1:D:452:GLN:HG2	2.43	0.54
1:B:407:LEU:HD12	1:B:408:GLY:H	1.73	0.54
1:A:165:ILE:HD11	1:A:452:GLN:OE1	2.07	0.54
1:D:334:ALA:O	1:D:338:GLU:HG3	2.08	0.54
1:D:254:ILE:HG12	1:D:316:LEU:HD22	1.90	0.53
1:A:289:ALA:O	1:A:290:ASN:HB2	2.09	0.53
1:D:128:ARG:NE	1:D:128:ARG:HA	2.20	0.53
1:C:514:ALA:O	1:C:518:VAL:HG23	2.09	0.53
1:C:334:ALA:O	1:C:338:GLU:HG3	2.09	0.53
1:B:36:ASN:O	1:B:40:ARG:HG3	2.08	0.53
1:A:316:LEU:HD23	1:A:389:GLN:HE22	1.74	0.53
1:C:110:THR:HG22	1:D:538:LEU:CB	2.29	0.53
1:C:289:ALA:O	1:C:290:ASN:HB2	2.08	0.53
1:B:319:LEU:HB3	1:B:320:PRO:HD3	1.90	0.52
1:D:199:ALA:HB3	1:D:200:PRO:HD3	1.90	0.52
1:B:289:ALA:O	1:B:290:ASN:HB2	2.10	0.52
1:C:172:VAL:HB	1:C:173:PRO:HD3	1.91	0.52
1:D:407:LEU:HD12	1:D:408:GLY:H	1.75	0.52
1:A:407:LEU:HD12	1:A:408:GLY:N	2.25	0.52
1:A:183:GLY:HA3	1:A:203:LEU:HD12	1.91	0.51
1:A:319:LEU:HB3	1:A:320:PRO:HD3	1.91	0.51
1:D:401:GLY:HA3	1:D:493:LYS:HE3	1.91	0.51
1:D:43:ARG:HH21	1:D:43:ARG:HG3	1.76	0.51
1:C:319:LEU:HB3	1:C:320:PRO:HD3	1.93	0.51
1:A:429:MET:HB3	1:A:430:PRO:HD3	1.93	0.51
1:D:407:LEU:HD12	1:D:408:GLY:N	2.26	0.51
1:B:43:ARG:HH21	1:B:43:ARG:HG3	1.75	0.51
1:D:165:ILE:HD11	1:D:452:GLN:OE1	2.11	0.51
1:B:67:ILE:O	1:B:71:VAL:HG23	2.10	0.51
1:A:497:HIS:HB2	1:A:525:SER:HB2	1.91	0.51
1:D:516:ARG:NH2	1:D:523:PRO:HA	2.26	0.50
1:D:291:SER:HB2	1:D:505:SER:HB3	1.93	0.50
1:A:28:ILE:HD11	1:A:48:VAL:HG11	1.92	0.50
1:C:412:ARG:HB2	1:C:415:ASN:ND2	2.27	0.50
1:D:259:GLY:O	1:D:294:VAL:HG13	2.12	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:407:LEU:HD12	1:A:408:GLY:H	1.77	0.49
1:C:562:ILE:HG22	1:C:562:ILE:O	2.13	0.49
1:A:562:ILE:O	1:A:562:ILE:HG22	2.12	0.49
1:B:291:SER:HB2	1:B:505:SER:HB3	1.94	0.49
1:C:283:GLN:NE2	2:C:1406:HOH:O	2.44	0.49
1:A:215:PRO:O	1:A:216:LYS:HB3	2.13	0.49
1:A:397:GLU:HG2	1:C:397:GLU:HB3	1.95	0.49
1:A:197:MET:HE3	1:A:201:THR:HG22	1.95	0.49
1:D:497:HIS:HB2	1:D:525:SER:HB2	1.94	0.48
1:B:334:ALA:O	1:B:338:GLU:HG3	2.13	0.48
1:B:316:LEU:HD23	1:B:389:GLN:HE22	1.78	0.48
1:D:215:PRO:O	1:D:216:LYS:HB3	2.14	0.48
1:B:516:ARG:NH2	1:B:523:PRO:HA	2.29	0.48
1:A:43:ARG:NH2	1:A:43:ARG:HG3	2.28	0.47
1:A:254:ILE:HG12	1:A:316:LEU:HD22	1.95	0.47
1:B:28:ILE:HD11	1:B:126:LEU:HD12	1.96	0.47
1:B:35:ILE:HG13	1:B:338:GLU:HG2	1.96	0.47
1:B:497:HIS:HB2	1:B:525:SER:HB2	1.95	0.47
1:D:93:SER:CB	1:D:96:GLN:HE21	2.28	0.47
1:A:104:LEU:HD23	1:A:222:MET:HE3	1.95	0.47
1:B:401:GLY:HA3	1:B:493:LYS:HE3	1.96	0.47
1:D:293:LEU:HD22	1:D:490:ARG:HG3	1.96	0.47
1:A:334:ALA:O	1:A:338:GLU:HG3	2.15	0.47
1:D:562:ILE:O	1:D:562:ILE:HG22	2.15	0.47
1:D:35:ILE:HG13	1:D:338:GLU:HG2	1.96	0.47
1:D:319:LEU:HB3	1:D:320:PRO:HD3	1.97	0.47
1:B:352:VAL:HG23	1:B:353:ASP:N	2.30	0.47
1:A:146:ARG:HG3	1:A:181:LEU:HD22	1.96	0.46
1:C:332:GLN:O	1:C:336:GLN:HG3	2.15	0.46
1:B:183:GLY:HA3	1:B:203:LEU:HD12	1.97	0.46
1:A:492:TYR:HA	1:A:497:HIS:O	2.16	0.46
1:B:249:VAL:HG22	1:B:480:LEU:HD23	1.97	0.46
1:C:127:LEU:CD2	1:C:341:ILE:HG23	2.42	0.46
1:B:412:ARG:NE	2:B:1390:HOH:O	2.47	0.46
1:B:93:SER:CB	1:B:96:GLN:HE21	2.29	0.46
1:C:43:ARG:NH2	1:C:43:ARG:HG3	2.29	0.46
1:C:429:MET:HB3	1:C:430:PRO:HD3	1.98	0.46
1:B:562:ILE:O	1:B:562:ILE:HG22	2.16	0.46
1:B:259:GLY:O	1:B:294:VAL:HG13	2.15	0.46
1:A:128:ARG:NH1	1:A:224:GLY:HA3	2.27	0.45
1:C:215:PRO:O	1:C:216:LYS:HB3	2.16	0.45
1:D:43:ARG:HG3	1:D:43:ARG:NH2	2.31	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:293:LEU:HD22	1:A:490:ARG:HG3	1.97	0.45
1:D:289:ALA:O	1:D:290:ASN:HB2	2.17	0.45
1:B:167:MDO:HB21	1:C:314:TYR:OH	2.17	0.45
1:B:388:VAL:HG13	2:B:1120:HOH:O	2.15	0.45
1:A:219:LEU:O	1:A:223:ASN:HB2	2.16	0.45
1:B:117:PRO:HD2	1:B:120:ASP:OD2	2.17	0.45
1:B:492:TYR:HA	1:B:497:HIS:O	2.16	0.45
1:D:415:ASN:HA	1:D:533:ASP:OD2	2.16	0.45
1:B:469:VAL:O	1:B:473:GLN:HG3	2.17	0.45
1:C:165:ILE:HD11	1:C:452:GLN:OE1	2.17	0.45
1:A:469:VAL:O	1:A:473:GLN:HG3	2.17	0.45
1:D:332:GLN:O	1:D:336:GLN:HG3	2.17	0.45
1:D:412:ARG:HD2	1:D:532:ASN:OD1	2.17	0.44
1:C:254:ILE:HG12	1:C:316:LEU:HD22	1.99	0.44
1:B:412:ARG:CZ	2:B:1390:HOH:O	2.65	0.44
1:C:352:VAL:HG23	1:C:353:ASP:N	2.33	0.44
1:A:259:GLY:O	1:A:294:VAL:HG13	2.16	0.44
1:D:134:ARG:HD3	2:D:1278:HOH:O	2.17	0.44
1:D:352:VAL:HG23	1:D:353:ASP:N	2.33	0.44
1:A:466:ARG:HD3	1:A:466:ARG:O	2.18	0.44
1:B:552:GLY:HA2	1:B:556:VAL:HG21	2.00	0.44
1:C:415:ASN:HA	1:C:533:ASP:OD2	2.17	0.44
1:C:407:LEU:HD13	1:C:415:ASN:HB3	1.99	0.44
1:B:412:ARG:HD2	1:B:532:ASN:OD1	2.17	0.44
1:C:466:ARG:HD3	1:C:466:ARG:O	2.18	0.43
1:D:466:ARG:HD3	1:D:466:ARG:O	2.17	0.43
1:A:415:ASN:HA	1:A:533:ASP:OD2	2.17	0.43
1:C:28:ILE:HD11	1:C:48:VAL:CG1	2.46	0.43
1:C:407:LEU:CD1	1:C:415:ASN:HB3	2.49	0.43
1:B:415:ASN:HA	1:B:533:ASP:OD2	2.17	0.43
1:A:352:VAL:HG23	1:A:353:ASP:N	2.33	0.43
1:C:92:ILE:O	1:C:93:SER:CB	2.66	0.43
1:B:412:ARG:HB2	1:B:415:ASN:ND2	2.34	0.43
1:B:466:ARG:O	1:B:466:ARG:HD3	2.19	0.43
1:B:28:ILE:CG2	1:B:130:ASN:HA	2.48	0.43
1:B:397:GLU:HB3	1:D:397:GLU:HG2	2.01	0.43
1:D:349:LEU:O	1:D:357:SER:HA	2.19	0.43
1:B:219:LEU:O	1:B:223:ASN:HB2	2.19	0.43
1:A:158:TYR:CD2	1:A:189:LYS:HB2	2.54	0.43
1:B:28:ILE:CD1	1:B:48:VAL:HG11	2.49	0.43
1:C:443:PHE:CE2	1:C:454:ILE:HA	2.54	0.43
1:B:553:GLY:O	1:B:557:GLN:HG3	2.19	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:43:ARG:NH2	1:B:43:ARG:HG3	2.33	0.43
1:C:401:GLY:HA3	1:C:493:LYS:HE3	2.00	0.43
1:A:69:ASN:N	1:A:69:ASN:ND2	2.67	0.42
1:A:117:PRO:HD2	1:A:120:ASP:OD2	2.18	0.42
1:B:215:PRO:O	1:B:216:LYS:HB3	2.18	0.42
1:B:407:LEU:HD13	1:B:415:ASN:HB3	2.01	0.42
1:B:134:ARG:NH2	2:B:1176:HOH:O	2.53	0.42
1:B:250:HIS:O	1:B:254:ILE:HG13	2.20	0.42
1:A:28:ILE:HD11	1:A:126:LEU:CD1	2.49	0.42
1:D:158:TYR:CD2	1:D:189:LYS:HB2	2.55	0.42
1:D:158:TYR:CE2	1:D:196:GLU:HG3	2.55	0.42
1:D:443:PHE:HB2	1:D:455:ASN:OD1	2.19	0.42
1:A:93:SER:CB	1:A:96:GLN:HE21	2.33	0.42
1:A:146:ARG:NH2	1:A:211:LEU:HD11	2.35	0.42
1:C:527:ARG:HD2	1:C:531:TRP:CD2	2.55	0.42
1:B:265:HIS:HA	1:B:266:PRO:HD3	1.86	0.41
1:A:388:VAL:HG13	2:D:1091:HOH:O	2.20	0.41
1:B:314:TYR:OH	1:C:167:MDO:HB21	2.20	0.41
1:A:552:GLY:HA2	1:A:556:VAL:HG21	2.02	0.41
1:B:129:ALA:O	1:B:133:MET:HG2	2.21	0.41
1:C:293:LEU:HD22	1:C:490:ARG:HG3	2.03	0.41
1:A:443:PHE:HB2	1:A:455:ASN:OD1	2.21	0.41
1:A:35:ILE:HG13	1:A:338:GLU:HG2	2.02	0.41
1:B:384:LYS:HE3	2:B:1326:HOH:O	2.21	0.41
1:D:265:HIS:HA	1:D:266:PRO:HD3	1.90	0.41
1:D:266:PRO:HG3	2:D:1210:HOH:O	2.19	0.41
1:A:451:ASN:OD1	1:D:313:ARG:NH2	2.53	0.41
1:A:28:ILE:CD1	1:A:48:VAL:HG11	2.51	0.41
1:A:319:LEU:HD23	1:A:319:LEU:C	2.41	0.41
1:A:104:LEU:HB3	1:A:222:MET:HE1	2.03	0.41
1:B:313:ARG:NH1	1:B:398:PHE:HD1	2.19	0.41
1:D:552:GLY:HA2	1:D:556:VAL:HG21	2.03	0.41
1:C:149:ILE:HG21	1:C:208:LEU:HD11	2.01	0.41
1:A:53:ASN:OD1	1:A:55:ASP:HB2	2.21	0.41
1:B:158:TYR:CE2	1:B:196:GLU:HG3	2.56	0.41
1:C:259:GLY:O	1:C:294:VAL:HG13	2.21	0.41
1:D:183:GLY:HA3	1:D:203:LEU:HD12	2.03	0.40
1:B:272:LYS:HE3	1:C:340:GLU:OE1	2.21	0.40
1:B:443:PHE:HB2	1:B:455:ASN:OD1	2.21	0.40
1:C:492:TYR:HA	1:C:497:HIS:O	2.21	0.40
1:B:407:LEU:CD1	1:B:415:ASN:HB3	2.51	0.40
1:D:164:SER:HB3	1:D:459:TYR:CG	2.57	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:104:LEU:HD12	1:C:107:PHE:CE1	2.57	0.40
1:C:118:LEU:HD22	1:C:156:THR:HB	2.03	0.40
2:A:1225:HOH:O	1:D:384:LYS:HE3	2.22	0.40
1:B:453:ASN:ND2	1:C:388:VAL:CG1	2.85	0.40
1:A:401:GLY:HA3	1:A:493:LYS:HE3	2.03	0.40
1:B:118:LEU:HD22	1:B:156:THR:HB	2.03	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	505/565 (89%)	484 (96%)	20 (4%)	1 (0%)	56	44
1	B	505/565 (89%)	486 (96%)	19 (4%)	0	100	100
1	C	505/565 (89%)	484 (96%)	21 (4%)	0	100	100
1	D	505/565 (89%)	485 (96%)	20 (4%)	0	100	100
All	All	2020/2260 (89%)	1939 (96%)	80 (4%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	273	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	423/468 (90%)	420 (99%)	3 (1%)	91	90

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	423/468 (90%)	421 (100%)	2 (0%)	94	94
1	C	423/468 (90%)	421 (100%)	2 (0%)	94	94
1	D	423/468 (90%)	421 (100%)	2 (0%)	94	94
All	All	1692/1872 (90%)	1683 (100%)	9 (0%)	94	94

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	69	ASN
1	A	192	PHE
1	A	223	ASN
1	B	192	PHE
1	B	223	ASN
1	C	128	ARG
1	C	192	PHE
1	D	128	ARG
1	D	192	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (41) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	ASN
1	A	68	ASN
1	A	69	ASN
1	A	96	GLN
1	A	205	GLN
1	A	207	ASN
1	A	223	ASN
1	A	240	GLN
1	A	258	ASN
1	A	283	GLN
1	A	292	GLN
1	A	474	ASN
1	B	44	ASN
1	B	68	ASN
1	B	96	GLN
1	B	205	GLN
1	B	207	ASN
1	B	223	ASN
1	B	240	GLN

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Mol	Chain	Res	Type
1	B	258	ASN
1	B	283	GLN
1	B	292	GLN
1	B	474	ASN
1	C	44	ASN
1	C	68	ASN
1	C	207	ASN
1	C	240	GLN
1	C	258	ASN
1	C	283	GLN
1	C	292	GLN
1	C	474	ASN
1	D	44	ASN
1	D	68	ASN
1	D	96	GLN
1	D	205	GLN
1	D	207	ASN
1	D	240	GLN
1	D	258	ASN
1	D	283	GLN
1	D	292	GLN
1	D	474	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

4 non-standard protein/DNA/RNA residues 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$
1	MDO	A	167	1	13,13,14	6.03	8 (61%)	16,18,20	2.30	6 (37%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	MDO	B	167	1	13,13,14	5.98	8 (61%)	16,18,20	2.34	6 (37%)
1	MDO	C	167	1	13,13,14	5.87	9 (69%)	16,18,20	2.31	6 (37%)
1	MDO	D	167	1	13,13,14	6.08	8 (61%)	16,18,20	2.31	6 (37%)

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
1	MDO	A	167	1	-	0/4/23/24	0/1/1/1
1	MDO	B	167	1	-	0/4/23/24	0/1/1/1
1	MDO	C	167	1	-	0/4/23/24	0/1/1/1
1	MDO	D	167	1	-	0/4/23/24	0/1/1/1

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	167	MDO	O-C	18.17	1.23	1.11
1	A	167	MDO	O-C	17.93	1.23	1.11
1	B	167	MDO	O-C	17.64	1.23	1.11
1	C	167	MDO	O-C	17.40	1.23	1.11
1	A	167	MDO	CB2-CA2	7.88	1.52	1.34
1	D	167	MDO	CB2-CA2	7.87	1.52	1.34
1	B	167	MDO	CB2-CA2	7.83	1.52	1.34
1	C	167	MDO	CB2-CA2	7.80	1.52	1.34
1	D	167	MDO	O2-C2	4.92	1.34	1.23
1	B	167	MDO	O2-C2	4.90	1.34	1.23
1	C	167	MDO	O2-C2	4.81	1.33	1.23
1	A	167	MDO	O2-C2	4.81	1.33	1.23
1	B	167	MDO	CA2-N2	-4.26	1.34	1.38
1	B	167	MDO	CA3-C	4.23	1.53	1.48
1	C	167	MDO	CA2-N2	-4.02	1.34	1.38
1	D	167	MDO	CA2-N2	-4.01	1.34	1.38
1	A	167	MDO	CA2-N2	-3.94	1.35	1.38
1	A	167	MDO	CA3-C	3.88	1.53	1.48
1	D	167	MDO	C2-N3	-3.84	1.31	1.39
1	A	167	MDO	C2-N3	-3.82	1.31	1.39
1	C	167	MDO	C2-N3	-3.74	1.31	1.39
1	D	167	MDO	CA3-C	3.66	1.53	1.48
1	B	167	MDO	C2-N3	-3.57	1.32	1.39
1	C	167	MDO	CA3-C	3.46	1.52	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	167	MDO	CA3-N3	-3.23	1.44	1.47
1	B	167	MDO	CA3-N3	-2.90	1.44	1.47
1	D	167	MDO	CA3-N3	-2.90	1.44	1.47
1	C	167	MDO	CA3-N3	-2.58	1.45	1.47
1	D	167	MDO	CA2-C2	-2.45	1.37	1.43
1	B	167	MDO	CA2-C2	-2.42	1.37	1.43
1	A	167	MDO	CA2-C2	-2.29	1.38	1.43
1	C	167	MDO	CA2-C2	-2.24	1.38	1.43
1	C	167	MDO	C1-N2	2.00	1.35	1.32

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	167	MDO	O2-C2-CA2	-4.80	128.18	130.96
1	D	167	MDO	O2-C2-CA2	-4.77	128.20	130.96
1	B	167	MDO	O2-C2-CA2	-4.49	128.36	130.96
1	A	167	MDO	CB2-CA2-C2	4.30	129.26	122.29
1	B	167	MDO	CB2-CA2-C2	4.30	129.26	122.29
1	A	167	MDO	O2-C2-CA2	-4.25	128.50	130.96
1	C	167	MDO	CB2-CA2-C2	3.99	128.75	122.29
1	D	167	MDO	CB2-CA2-C2	3.95	128.70	122.29
1	A	167	MDO	C2-CA2-N2	-3.60	106.17	108.96
1	D	167	MDO	C2-CA2-N2	-3.52	106.22	108.96
1	B	167	MDO	C2-CA2-N2	-3.52	106.23	108.96
1	C	167	MDO	C2-CA2-N2	-3.38	106.34	108.96
1	D	167	MDO	N3-C1-N2	-3.14	109.21	111.51
1	A	167	MDO	N3-C1-N2	-3.08	109.25	111.51
1	B	167	MDO	N3-C1-N2	-3.07	109.26	111.51
1	C	167	MDO	N3-C1-N2	-2.98	109.32	111.51
1	B	167	MDO	CA2-N2-C1	2.96	107.94	105.36
1	A	167	MDO	CA2-N2-C1	2.85	107.85	105.36
1	D	167	MDO	CA2-N2-C1	2.68	107.70	105.36
1	C	167	MDO	CA2-C2-N3	2.57	109.39	104.40
1	D	167	MDO	CA2-C2-N3	2.54	109.33	104.40
1	A	167	MDO	CA2-C2-N3	2.50	109.25	104.40
1	B	167	MDO	CA2-C2-N3	2.50	109.24	104.40
1	C	167	MDO	CA2-N2-C1	2.42	107.47	105.36

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	512/565 (90%)	0.46	29 (5%)	23	23	13, 28, 53, 88	0
1	B	512/565 (90%)	0.43	18 (3%)	42	43	13, 28, 53, 87	0
1	C	512/565 (90%)	0.40	31 (6%)	21	20	14, 28, 51, 88	0
1	D	512/565 (90%)	0.46	38 (7%)	14	14	14, 27, 53, 88	0
All	All	2048/2260 (90%)	0.44	116 (5%)	23	23	13, 28, 53, 88	0

All (116) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	310	ILE	11.0
1	C	310	ILE	9.8
1	D	310	ILE	9.7
1	A	92	ILE	9.1
1	C	92	ILE	9.0
1	B	310	ILE	8.4
1	B	92	ILE	7.2
1	D	92	ILE	6.9
1	B	71	VAL	6.8
1	D	301	LYS	6.2
1	C	301	LYS	5.8
1	C	72	GLU	5.4
1	B	95	GLU	5.3
1	B	301	LYS	5.3
1	A	73	SER	5.1
1	A	301	LYS	4.9
1	D	68	ASN	4.9
1	B	73	SER	4.8
1	D	93	SER	4.4
1	A	195	LYS	4.3
1	B	93	SER	4.1

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Mol	Chain	Res	Type	RSRZ
1	D	73	SER	4.0
1	B	410	ARG	4.0
1	D	72	GLU	3.9
1	C	68	ASN	3.9
1	C	94	ARG	3.8
1	C	300	GLY	3.8
1	C	95	GLU	3.7
1	D	96	GLN	3.6
1	B	72	GLU	3.6
1	D	300	GLY	3.6
1	B	195	LYS	3.5
1	C	299	ASP	3.5
1	B	96	GLN	3.4
1	A	68	ASN	3.4
1	A	94	ARG	3.4
1	D	69	ASN	3.3
1	C	73	SER	3.3
1	A	95	GLU	3.3
1	D	71	VAL	3.2
1	A	93	SER	3.2
1	C	96	GLN	3.2
1	D	95	GLU	3.2
1	A	72	GLU	3.1
1	D	299	ASP	3.0
1	D	439	ILE	3.0
1	A	74	GLY	3.0
1	C	52	ASN	2.9
1	C	410	ARG	2.9
1	A	55	ASP	2.9
1	D	65	ASP	2.9
1	C	93	SER	2.9
1	C	69	ASN	2.9
1	D	52	ASN	2.9
1	C	563	LEU	2.9
1	C	439	ILE	2.8
1	D	560	GLN	2.8
1	D	97	ALA	2.8
1	D	207	ASN	2.8
1	D	195	LYS	2.8
1	A	311	GLN	2.8
1	B	55	ASP	2.8
1	D	55	ASP	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	68	ASN	2.7
1	A	563	LEU	2.7
1	C	412	ARG	2.6
1	A	71	VAL	2.6
1	D	410	ARG	2.6
1	D	311	GLN	2.6
1	D	359	HIS	2.6
1	A	355	GLN	2.6
1	A	439	ILE	2.5
1	C	562	ILE	2.5
1	A	96	GLN	2.5
1	C	195	LYS	2.5
1	C	411	GLU	2.5
1	B	69	ASN	2.5
1	D	563	LEU	2.5
1	D	562	ILE	2.5
1	B	94	ARG	2.5
1	D	58	GLN	2.5
1	D	30	ASN	2.4
1	C	207	ASN	2.4
1	C	499	ASP	2.4
1	C	74	GLY	2.3
1	C	524	THR	2.3
1	A	440	ALA	2.3
1	D	352	VAL	2.3
1	A	69	ASN	2.3
1	A	562	ILE	2.3
1	D	67	ILE	2.3
1	D	54	THR	2.3
1	A	433	THR	2.2
1	D	298	LEU	2.2
1	C	434	PHE	2.2
1	D	434	PHE	2.2
1	D	502	ALA	2.2
1	C	435	TYR	2.2
1	A	209	SER	2.1
1	B	411	GLU	2.1
1	A	443	PHE	2.1
1	D	74	GLY	2.1
1	C	55	ASP	2.1
1	A	70	ALA	2.1
1	A	554	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	296	ASP	2.1
1	A	60	ILE	2.1
1	A	411	GLU	2.1
1	B	563	LEU	2.1
1	A	442	ARG	2.0
1	B	434	PHE	2.0
1	C	54	THR	2.0
1	D	297	GLU	2.0
1	D	527	ARG	2.0
1	C	58	GLN	2.0
1	C	523	PRO	2.0

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

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	MDO	A	167	13/14	0.12	-0.20	22,27,30,31	0
1	MDO	B	167	13/14	0.11	-0.32	23,26,27,29	0
1	MDO	C	167	13/14	0.11	-0.34	25,27,30,32	0
1	MDO	D	167	13/14	0.12	-0.42	25,28,29,30	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.