



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 02:50 PM GMT

PDB ID : 3N4L
Title : BACE-1 in complex with ELN380842
Authors : Yao, N.H.
Deposited on : 2010-05-21
Resolution : 2.70 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>

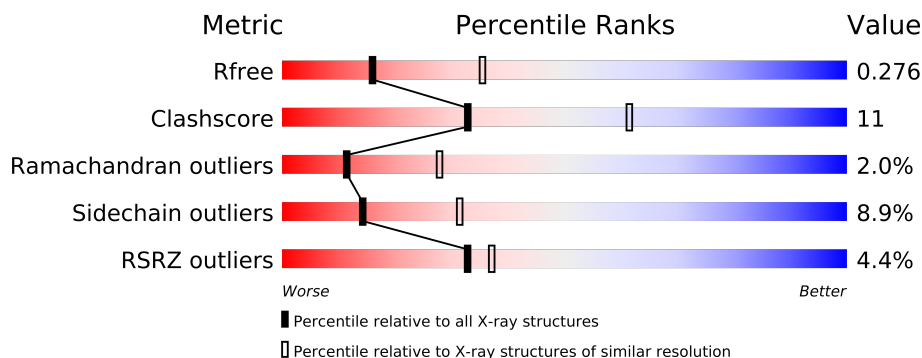
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
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 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))
R_{free}	66092	1557 (2.70-2.70)
Clashscore	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)
RSRZ outliers	66119	1559 (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. 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	406	
1	B	406	
1	C	406	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8982 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 Beta-secretase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	375	Total	C	N	O	S	0	0	0
			2953	1892	490	557	14			
1	B	375	Total	C	N	O	S	0	0	0
			2942	1887	490	551	14			
1	C	375	Total	C	N	O	S	0	0	0
			2947	1887	490	556	14			

There are 27 discrepancies between the modelled and reference sequences:

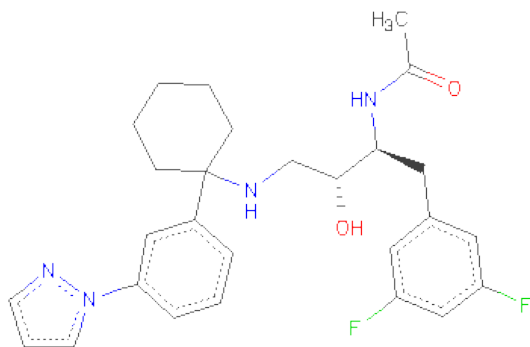
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	INITIATING METHIONINE	UNP P56817
A	399	ARG	-	EXPRESSION TAG	UNP P56817
A	400	SER	-	EXPRESSION TAG	UNP P56817
A	401	HIS	-	EXPRESSION TAG	UNP P56817
A	402	HIS	-	EXPRESSION TAG	UNP P56817
A	403	HIS	-	EXPRESSION TAG	UNP P56817
A	404	HIS	-	EXPRESSION TAG	UNP P56817
A	405	HIS	-	EXPRESSION TAG	UNP P56817
A	406	HIS	-	EXPRESSION TAG	UNP P56817
B	1	MET	-	INITIATING METHIONINE	UNP P56817
B	399	ARG	-	EXPRESSION TAG	UNP P56817
B	400	SER	-	EXPRESSION TAG	UNP P56817
B	401	HIS	-	EXPRESSION TAG	UNP P56817
B	402	HIS	-	EXPRESSION TAG	UNP P56817
B	403	HIS	-	EXPRESSION TAG	UNP P56817
B	404	HIS	-	EXPRESSION TAG	UNP P56817
B	405	HIS	-	EXPRESSION TAG	UNP P56817
B	406	HIS	-	EXPRESSION TAG	UNP P56817
C	1	MET	-	INITIATING METHIONINE	UNP P56817
C	399	ARG	-	EXPRESSION TAG	UNP P56817
C	400	SER	-	EXPRESSION TAG	UNP P56817
C	401	HIS	-	EXPRESSION TAG	UNP P56817
C	402	HIS	-	EXPRESSION TAG	UNP P56817

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Chain	Residue	Modelled	Actual	Comment	Reference
C	403	HIS	-	EXPRESSION TAG	UNP P56817
C	404	HIS	-	EXPRESSION TAG	UNP P56817
C	405	HIS	-	EXPRESSION TAG	UNP P56817
C	406	HIS	-	EXPRESSION TAG	UNP P56817

- Molecule 2 is N-[(1S,2R)-1-(3,5-DIFLUOROBENZYL)-2-HYDROXY-3-({1-[3-(1H-PYRAZOL-1-YL)PHENYL]CYCLOHEXYL}AMINO)PROPYL]ACETAMIDE (three-letter code: 842) (formula: C₂₇H₃₂F₂N₄O₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	0	0
			35	27	2	4	2		
2	B	1	Total	C	F	N	O	0	0
			35	27	2	4	2		
2	C	1	Total	C	F	N	O	0	0
			35	27	2	4	2		

- Molecule 3 is water.

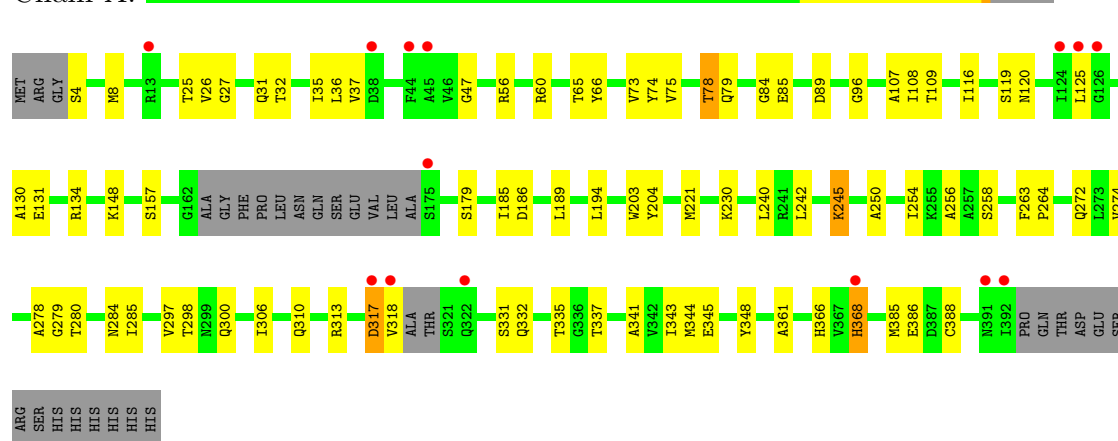
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	12	Total	O	0	0
			12	12		
3	B	14	Total	O	0	0
			14	14		
3	C	9	Total	O	0	0
			9	9		

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.

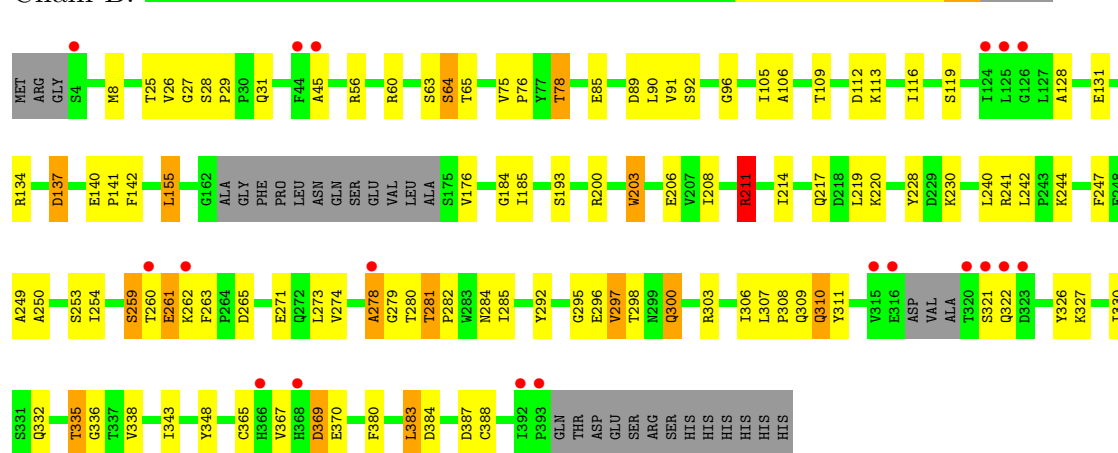
• Molecule 1: Beta-secretase 1

Chain A:



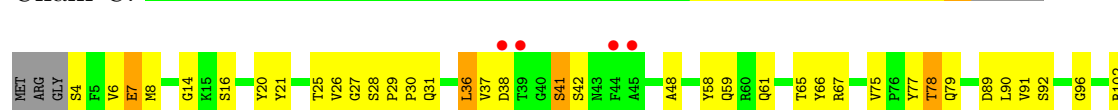
• Molecule 1: Beta-secretase 1

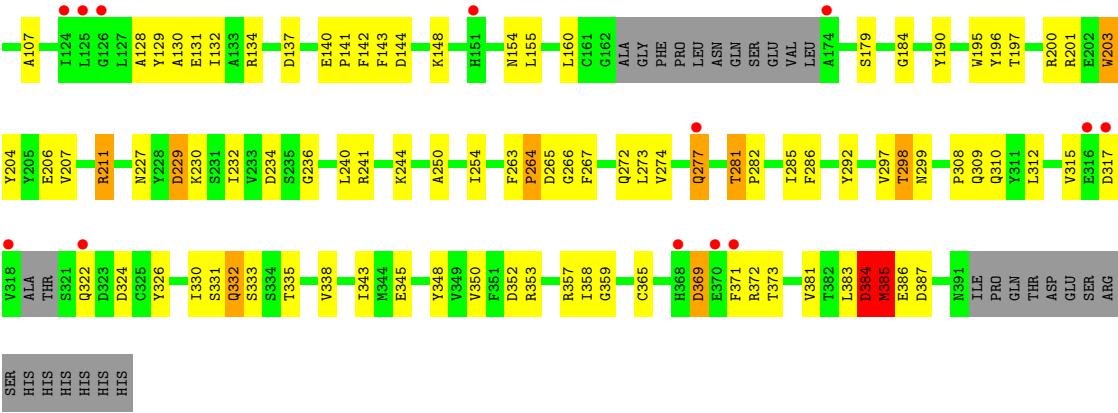
Chain B:



• Molecule 1: Beta-secretase 1

Chain C:





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.05Å 105.58Å 101.58Å 90.00° 104.63° 90.00°	Depositor
Resolution (Å)	28.00 – 2.70 28.24 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.7 (28.00-2.70) 98.8 (28.24-2.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.57 (at 2.72Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.223 , 0.283 0.216 , 0.276	Depositor DCC
R_{free} test set	2329 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	74.1	Xtriage
Anisotropy	0.044	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 35.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 46150 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8982	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.68% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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:
842

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	0.79	0/3027	0.82	1/4112 (0.0%)
1	B	0.80	1/3017 (0.0%)	0.82	1/4100 (0.0%)
1	C	0.79	0/3021	0.85	2/4105 (0.0%)
All	All	0.79	1/9065 (0.0%)	0.83	4/12317 (0.0%)

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	C	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	388	CYS	CB-SG	-5.36	1.73	1.81

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	385	MET	N-CA-C	-5.97	94.88	111.00
1	B	211	ARG	NE-CZ-NH1	5.15	122.88	120.30
1	A	344	MET	CG-SD-CE	5.12	108.39	100.20
1	C	90	LEU	CA-CB-CG	5.08	126.97	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	384	ASP	Peptide
1	C	385	MET	Peptide

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	2953	0	2862	48	0
1	B	2942	0	2850	67	0
1	C	2947	0	2849	85	0
2	A	35	0	32	3	0
2	B	35	0	32	3	0
2	C	35	0	32	3	0
3	A	12	0	0	0	0
3	B	14	0	0	0	0
3	C	9	0	0	0	0
All	All	8982	0	8657	197	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 11.

All (197) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:297:VAL:O	1:C:298:THR:HG22	1.39	1.18
1:B:211:ARG:HH11	1:B:211:ARG:HB2	1.22	1.03
1:C:384:ASP:C	1:C:386:GLU:H	1.69	0.93
1:A:317:ASP:HA	1:A:318:VAL:HB	1.50	0.93
1:B:296:GLU:H	1:B:300:GLN:HE21	1.12	0.93
1:A:32:THR:HG22	1:A:56:ARG:HH12	1.42	0.85
1:A:306:ILE:HD12	1:A:310:GLN:NE2	1.94	0.83
1:A:258:SER:HB2	1:A:285:ILE:HD12	1.61	0.82
1:C:297:VAL:O	1:C:298:THR:CG2	2.26	0.81
1:B:384:ASP:O	1:B:387:ASP:HB2	1.81	0.80
1:A:32:THR:HG22	1:A:56:ARG:NH1	1.96	0.80
1:C:384:ASP:C	1:C:386:GLU:N	2.34	0.80
1:B:327:LYS:NZ	1:B:327:LYS:HB2	1.98	0.79

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:384:ASP:CB	1:C:385:MET:HA	2.14	0.78
1:C:38:ASP:OD1	1:C:41:SER:HB3	1.84	0.77
1:A:230:LYS:NZ	1:A:335:THR:HB	1.99	0.77
1:C:155:LEU:HD12	1:C:352:ASP:HA	1.68	0.76
1:A:306:ILE:HD12	1:A:310:GLN:HE22	1.51	0.76
1:C:36:LEU:HD23	1:C:36:LEU:C	2.06	0.75
1:A:116:ILE:O	1:A:119:SER:HB2	1.88	0.73
1:B:8:MET:HG2	1:B:96:GLY:HA2	1.71	0.73
1:A:78:THR:HG22	2:A:407:842:H19	1.70	0.73
1:B:91:VAL:HG11	1:B:142:PHE:CE1	2.23	0.73
1:C:384:ASP:CB	1:C:385:MET:CA	2.67	0.72
1:C:155:LEU:HD23	1:C:184:GLY:HA2	1.71	0.71
1:B:78:THR:HG22	2:B:407:842:H19	1.71	0.71
1:A:185:ILE:HG23	1:A:348:TYR:HE2	1.55	0.71
1:C:385:MET:CE	1:C:387:ASP:HB2	2.21	0.71
1:A:8:MET:HG2	1:A:96:GLY:HA2	1.73	0.70
1:B:250:ALA:O	1:B:254:ILE:HG13	1.92	0.70
1:C:36:LEU:HD23	1:C:37:VAL:N	2.08	0.68
1:B:28:SER:O	1:B:63:SER:HA	1.94	0.68
1:B:296:GLU:N	1:B:300:GLN:HE21	1.88	0.67
1:A:306:ILE:HG13	1:A:343:ILE:HD13	1.76	0.67
1:C:196:TYR:CD2	1:C:357:ARG:HG3	2.30	0.66
1:A:341:ALA:O	1:A:345:GLU:HG3	1.95	0.66
1:B:91:VAL:HG11	1:B:142:PHE:HE1	1.59	0.66
1:C:200:ARG:HB3	1:C:206:GLU:HG2	1.79	0.65
1:A:256:ALA:HB1	1:C:371:PHE:CD1	2.32	0.65
1:C:155:LEU:CD1	1:C:352:ASP:HA	2.28	0.64
1:B:211:ARG:NH2	1:C:385:MET:HG3	2.15	0.62
1:C:263:PHE:HD2	1:C:274:VAL:HG11	1.65	0.61
1:B:230:LYS:O	1:B:336:GLY:HA3	2.01	0.61
1:C:286:PHE:CE2	1:C:312:LEU:HD11	2.35	0.61
1:C:282:PRO:HB2	1:C:285:ILE:HG12	1.83	0.60
1:C:66:TYR:O	1:C:67:ARG:HG3	2.01	0.60
1:C:281:THR:HG21	1:C:326:TYR:CE1	2.37	0.59
1:C:211:ARG:HG2	1:C:292:TYR:CD1	2.37	0.59
1:C:282:PRO:O	1:C:285:ILE:HG12	2.02	0.59
1:A:73:VAL:HG22	1:A:74:TYR:N	2.16	0.59
1:B:214:ILE:HG13	1:B:219:LEU:HD21	1.85	0.59
1:C:89:ASP:O	1:C:102:ARG:HA	2.04	0.58
1:B:308:PRO:HG2	1:B:309:GLN:OE1	2.03	0.58
1:B:327:LYS:HZ2	1:B:327:LYS:HB2	1.67	0.58
1:A:36:LEU:C	1:A:36:LEU:HD23	2.24	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:78:THR:HG23	2:A:407:842:O25	2.03	0.57
1:C:25:THR:HA	1:C:31:GLN:O	2.05	0.57
1:C:385:MET:SD	1:C:385:MET:N	2.78	0.56
1:A:230:LYS:HZ1	1:A:335:THR:HB	1.70	0.55
1:C:14:GLY:HA2	1:C:21:TYR:CE2	2.41	0.55
1:B:211:ARG:NH1	1:B:211:ARG:HB2	2.06	0.55
1:C:229:ASP:OD2	1:C:230:LYS:HG2	2.07	0.54
1:B:383:LEU:HD12	1:B:383:LEU:H	1.72	0.54
1:C:27:GLY:O	1:C:30:PRO:HA	2.07	0.54
1:B:240:LEU:HD22	1:B:343:ILE:HD11	1.89	0.54
1:A:75:VAL:HG22	1:A:134:ARG:HG3	1.89	0.54
1:C:59:GLN:HB3	1:C:61:GLN:NE2	2.23	0.54
1:A:317:ASP:HA	1:A:318:VAL:CB	2.31	0.53
1:A:368:HIS:CD2	1:A:368:HIS:H	2.25	0.53
1:A:36:LEU:HD23	1:A:37:VAL:N	2.24	0.53
1:A:73:VAL:CG2	1:A:74:TYR:N	2.70	0.53
1:B:327:LYS:HZ3	1:B:327:LYS:HB2	1.70	0.53
1:C:66:TYR:CG	1:C:67:ARG:N	2.77	0.53
1:B:282:PRO:HB2	1:B:285:ILE:HG12	1.91	0.53
1:B:281:THR:HG21	1:B:326:TYR:CE1	2.43	0.53
1:A:230:LYS:HZ3	1:A:335:THR:HB	1.74	0.52
1:B:91:VAL:CG1	1:B:142:PHE:HE1	2.23	0.52
1:B:185:ILE:HG23	1:B:348:TYR:HE2	1.74	0.52
1:B:263:PHE:HD2	1:B:274:VAL:HG11	1.73	0.52
1:A:385:MET:O	1:A:388:CYS:HB2	2.09	0.52
1:C:263:PHE:CD2	1:C:274:VAL:HG11	2.43	0.52
1:B:306:ILE:HG13	1:B:343:ILE:HD13	1.92	0.51
1:A:368:HIS:H	1:A:368:HIS:HD2	1.58	0.51
1:C:28:SER:OG	1:C:65:THR:HG21	2.11	0.51
1:C:78:THR:HG22	2:C:407:842:H19	1.93	0.51
1:B:308:PRO:HA	1:B:311:TYR:CE2	2.46	0.51
2:B:407:842:C19	2:B:407:842:H12A	2.39	0.51
1:C:38:ASP:OD2	1:C:236:GLY:HA3	2.11	0.50
1:B:309:GLN:NE2	1:B:369:ASP:HB3	2.27	0.50
1:B:297:VAL:HB	1:B:300:GLN:HG2	1.93	0.50
1:B:292:TYR:CE2	1:B:303:ARG:HB3	2.47	0.50
1:C:48:ALA:CB	1:C:107:ALA:HB1	2.42	0.50
1:C:244:LYS:HG3	1:C:332:GLN:OE1	2.12	0.49
1:A:47:GLY:HA2	1:A:108:ILE:HB	1.94	0.49
1:C:42:SER:O	1:C:141:PRO:HB3	2.12	0.49
1:B:211:ARG:HH21	1:C:385:MET:HG3	1.77	0.49
1:B:296:GLU:H	1:B:300:GLN:NE2	1.95	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:206:GLU:HG3	1:C:207:VAL:N	2.28	0.49
1:C:132:ILE:HD13	1:C:204:TYR:CE1	2.48	0.49
1:A:306:ILE:CD1	1:A:310:GLN:HE22	2.23	0.49
1:C:20:TYR:N	1:C:20:TYR:CD1	2.80	0.49
1:A:194:LEU:HD23	1:A:361:ALA:HB2	1.95	0.49
1:C:78:THR:HG23	2:C:407:842:O25	2.12	0.49
1:C:26:VAL:HG22	1:C:91:VAL:HG22	1.95	0.49
1:C:240:LEU:HD22	1:C:343:ILE:HD11	1.95	0.49
1:A:85:GLU:OE1	1:A:109:THR:HG21	2.13	0.49
1:C:365:CYS:O	1:C:365:CYS:SG	2.71	0.48
1:A:242:LEU:HD23	1:A:337:THR:HG23	1.94	0.48
1:C:241:ARG:HB3	1:C:333:SER:HB2	1.95	0.48
1:B:155:LEU:HD21	1:B:184:GLY:HA2	1.95	0.48
1:B:259:SER:C	1:B:261:GLU:H	2.16	0.48
1:B:45:ALA:HB2	1:B:106:ALA:HB3	1.94	0.48
1:B:228:TYR:O	1:B:336:GLY:HA2	2.14	0.47
1:B:292:TYR:CD2	1:B:303:ARG:HB3	2.49	0.47
1:B:89:ASP:OD1	1:B:90:LEU:N	2.41	0.47
1:B:383:LEU:HD12	1:B:383:LEU:N	2.29	0.47
1:C:42:SER:OG	1:C:128:ALA:HB3	2.15	0.47
1:A:240:LEU:HD13	1:A:343:ILE:HG13	1.97	0.47
1:C:240:LEU:HD12	1:C:338:VAL:O	2.15	0.47
1:C:240:LEU:HD23	1:C:330:ILE:HG12	1.95	0.47
1:A:186:ASP:HB3	1:A:189:LEU:HD12	1.96	0.47
1:C:250:ALA:O	1:C:254:ILE:HG13	2.15	0.47
1:C:130:ALA:O	1:C:131:GLU:C	2.53	0.47
1:C:232:ILE:HD12	1:C:234:ASP:HB2	1.97	0.47
1:A:204:TYR:CE1	2:A:407:842:H9A	2.50	0.47
1:A:27:GLY:HA2	1:A:89:ASP:OD1	2.15	0.47
1:C:200:ARG:CB	1:C:206:GLU:HG2	2.44	0.46
1:C:77:TYR:CG	2:C:407:842:H28	2.50	0.46
1:B:307:LEU:H	1:B:310:GLN:NE2	2.13	0.46
1:B:28:SER:HA	1:B:29:PRO:C	2.36	0.46
1:B:140:GLU:HA	1:B:141:PRO:HD3	1.75	0.46
1:C:154:ASN:ND2	1:C:353:ARG:HB3	2.30	0.46
1:B:203:TRP:NE1	1:B:206:GLU:OE1	2.35	0.46
1:B:116:ILE:HB	1:B:119:SER:HB3	1.97	0.46
1:A:250:ALA:O	1:A:254:ILE:HG13	2.16	0.46
1:C:308:PRO:O	1:C:310:GLN:N	2.49	0.46
1:C:201:ARG:O	1:C:203:TRP:HD1	1.99	0.46
1:A:221:MET:HE3	1:A:245:LYS:HE2	1.98	0.46
1:B:278:ALA:C	1:B:280:THR:H	2.19	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:36:LEU:CD2	1:C:36:LEU:C	2.79	0.46
1:A:185:ILE:HG23	1:A:348:TYR:CE2	2.44	0.45
1:B:112:ASP:OD2	1:B:113:LYS:HE2	2.15	0.45
1:B:25:THR:HA	1:B:31:GLN:O	2.16	0.45
1:C:385:MET:HE2	1:C:387:ASP:HB2	1.98	0.45
1:C:14:GLY:HA2	1:C:21:TYR:HE2	1.82	0.45
1:B:91:VAL:CG1	1:B:142:PHE:CE1	2.96	0.45
1:C:281:THR:CG2	1:C:326:TYR:HE1	2.29	0.45
1:B:85:GLU:OE1	1:B:109:THR:HG21	2.16	0.45
1:A:119:SER:O	1:A:120:ASN:CB	2.65	0.45
1:A:25:THR:HA	1:A:31:GLN:O	2.17	0.45
1:C:264:PRO:O	1:C:266:GLY:N	2.50	0.45
1:C:4:SER:O	1:C:7:GLU:HG3	2.18	0.44
1:A:60:ARG:CD	1:A:66:TYR:CE1	3.00	0.44
1:C:195:TRP:O	1:C:359:GLY:HA2	2.17	0.44
1:B:365:CYS:SG	1:B:365:CYS:O	2.75	0.44
1:B:295:GLY:HA3	1:B:300:GLN:NE2	2.33	0.44
2:B:407:842:H12A	2:B:407:842:H19	1.99	0.43
1:B:281:THR:CG2	1:B:326:TYR:HE1	2.31	0.43
1:B:26:VAL:HG22	1:B:91:VAL:HG22	1.99	0.43
1:A:263:PHE:HD2	1:A:274:VAL:HG11	1.82	0.43
1:C:350:VAL:O	1:C:358:ILE:HA	2.19	0.43
1:A:130:ALA:O	1:A:131:GLU:C	2.56	0.43
1:B:281:THR:HA	1:B:282:PRO:HD3	1.86	0.43
1:C:140:GLU:HA	1:C:141:PRO:HD3	1.94	0.43
1:B:200:ARG:HD2	1:B:208:ILE:HD11	2.01	0.43
1:C:8:MET:HG2	1:C:96:GLY:HA2	2.00	0.43
1:B:28:SER:HB2	1:B:64:SER:HB3	2.01	0.42
1:A:8:MET:CG	1:A:96:GLY:HA2	2.46	0.42
1:C:322:GLN:O	1:C:322:GLN:HG2	2.19	0.42
1:C:129:TYR:HD1	1:C:203:TRP:O	2.03	0.42
1:A:84:GLY:HA3	1:A:107:ALA:O	2.20	0.42
1:C:369:ASP:HB3	1:C:372:ARG:H	1.85	0.42
1:B:327:LYS:NZ	1:B:327:LYS:CB	2.76	0.42
1:B:75:VAL:HG22	1:B:134:ARG:HB2	2.02	0.42
1:B:242:LEU:HB2	1:B:247:PHE:HB2	2.00	0.42
1:C:282:PRO:HG2	1:C:285:ILE:HD11	2.02	0.42
1:B:230:LYS:NZ	1:B:335:THR:HB	2.35	0.41
1:C:281:THR:HA	1:C:282:PRO:HD3	1.91	0.41
1:C:142:PHE:O	1:C:143:PHE:C	2.57	0.41
1:C:332:GLN:HB2	1:C:332:GLN:HE21	1.56	0.41
1:B:27:GLY:HA2	1:B:89:ASP:OD1	2.21	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:190:TYR:CD1	1:C:348:TYR:CD2	3.09	0.41
1:B:263:PHE:CD2	1:B:274:VAL:HG11	2.55	0.41
1:B:380:PHE:HE2	1:C:298:THR:HG23	1.85	0.41
1:C:385:MET:SD	1:C:387:ASP:HB2	2.60	0.41
1:C:28:SER:HA	1:C:29:PRO:C	2.40	0.41
1:C:26:VAL:HG12	1:C:58:TYR:CE1	2.56	0.41
1:C:160:LEU:O	1:C:345:GLU:HA	2.20	0.41
1:B:242:LEU:HD12	1:B:330:ILE:CG2	2.51	0.41
1:C:75:VAL:HG22	1:C:134:ARG:HG3	2.03	0.41
1:B:241:ARG:HB2	1:B:338:VAL:HB	2.02	0.41
1:A:368:HIS:CD2	1:A:368:HIS:N	2.89	0.40
1:A:297:VAL:HG12	1:A:300:GLN:HB3	2.02	0.40
1:B:383:LEU:HD11	1:C:383:LEU:HD11	2.02	0.40
1:C:241:ARG:HA	1:C:331:SER:O	2.22	0.40
1:B:295:GLY:CA	1:B:300:GLN:NE2	2.84	0.40
1:A:35:ILE:CG2	1:A:125:LEU:HB2	2.52	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	369/406 (91%)	346 (94%)	19 (5%)	4 (1%)	21	49
1	B	369/406 (91%)	323 (88%)	35 (10%)	11 (3%)	7	15
1	C	369/406 (91%)	322 (87%)	40 (11%)	7 (2%)	12	29
All	All	1107/1218 (91%)	991 (90%)	94 (8%)	22 (2%)	11	28

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	370	GLU
1	A	278	ALA
1	A	279	GLY

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Mol	Chain	Res	Type
1	B	64	SER
1	B	128	ALA
1	C	265	ASP
1	C	277	GLN
1	C	309	GLN
1	A	284	ASN
1	B	260	THR
1	B	278	ALA
1	B	279	GLY
1	B	321	SER
1	C	298	THR
1	C	384	ASP
1	C	369	ASP
1	B	137	ASP
1	B	249	ALA
1	B	261	GLU
1	C	264	PRO
1	A	264	PRO
1	B	76	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	321/348 (92%)	301 (94%)	20 (6%)	26	54
1	B	318/348 (91%)	284 (89%)	34 (11%)	10	22
1	C	319/348 (92%)	288 (90%)	31 (10%)	12	27
All	All	958/1044 (92%)	873 (91%)	85 (9%)	14	31

All (85) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	SER
1	A	26	VAL
1	A	65	THR
1	A	78	THR

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Mol	Chain	Res	Type
1	A	79	GLN
1	A	148	LYS
1	A	157	SER
1	A	179	SER
1	A	203	TRP
1	A	245	LYS
1	A	272	GLN
1	A	280	THR
1	A	298	THR
1	A	313	ARG
1	A	317	ASP
1	A	331	SER
1	A	332	GLN
1	A	366	HIS
1	A	368	HIS
1	A	386	GLU
1	B	56	ARG
1	B	60	ARG
1	B	65	THR
1	B	78	THR
1	B	92	SER
1	B	105	ILE
1	B	131	GLU
1	B	137	ASP
1	B	155	LEU
1	B	176	VAL
1	B	193	SER
1	B	203	TRP
1	B	211	ARG
1	B	217	GLN
1	B	220	LYS
1	B	244	LYS
1	B	253	SER
1	B	259	SER
1	B	262	LYS
1	B	265	ASP
1	B	271	GLU
1	B	273	LEU
1	B	281	THR
1	B	284	ASN
1	B	297	VAL
1	B	298	THR

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Mol	Chain	Res	Type
1	B	300	GLN
1	B	310	GLN
1	B	322	GLN
1	B	332	GLN
1	B	335	THR
1	B	367	VAL
1	B	369	ASP
1	B	383	LEU
1	C	6	VAL
1	C	7	GLU
1	C	16	SER
1	C	36	LEU
1	C	41	SER
1	C	78	THR
1	C	79	GLN
1	C	92	SER
1	C	137	ASP
1	C	144	ASP
1	C	148	LYS
1	C	179	SER
1	C	197	THR
1	C	203	TRP
1	C	211	ARG
1	C	227	ASN
1	C	229	ASP
1	C	267	PHE
1	C	272	GLN
1	C	273	LEU
1	C	277	GLN
1	C	281	THR
1	C	299	ASN
1	C	315	VAL
1	C	317	ASP
1	C	324	ASP
1	C	332	GLN
1	C	335	THR
1	C	373	THR
1	C	381	VAL
1	C	385	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	59	GLN
1	A	95	HIS
1	A	272	GLN
1	A	277	GLN
1	A	310	GLN
1	A	368	HIS
1	B	104	ASN
1	B	217	GLN
1	B	284	ASN
1	B	300	GLN
1	C	79	GLN
1	C	104	ASN
1	C	272	GLN
1	C	299	ASN
1	C	310	GLN
1	C	332	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

3 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	842	A	407	-	38,38,38	1.31	2 (5%)	53,53,53	2.14	10 (18%)
2	842	B	407	-	38,38,38	1.26	2 (5%)	53,53,53	1.99	10 (18%)
2	842	C	407	-	38,38,38	1.28	5 (13%)	53,53,53	2.43	17 (32%)

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	842	A	407	-	-	0/28/38/38	0/4/4/4
2	842	B	407	-	-	0/28/38/38	0/4/4/4
2	842	C	407	-	-	0/28/38/38	0/4/4/4

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	407	842	C16-N20	-5.46	1.33	1.43
2	B	407	842	C16-N20	-5.35	1.33	1.43
2	C	407	842	C16-N20	-5.03	1.34	1.43
2	A	407	842	F34-C31	-2.41	1.31	1.36
2	C	407	842	C32-C31	2.24	1.41	1.37
2	C	407	842	C24-N20	-2.21	1.32	1.36
2	C	407	842	C17-C16	2.10	1.41	1.38
2	C	407	842	C28-C29	2.08	1.41	1.37
2	B	407	842	C28-C29	2.03	1.41	1.37

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	407	842	C16-N20-N21	7.37	125.35	118.87
2	C	407	842	C17-C16-C15	-7.21	115.21	121.55
2	A	407	842	C17-C16-C15	-6.88	115.50	121.55
2	A	407	842	C16-C15-C14	6.10	124.98	120.06
2	A	407	842	C16-N20-N21	5.83	124.00	118.87
2	B	407	842	C17-C16-C15	-5.71	116.53	121.55
2	B	407	842	C17-C16-N20	5.14	124.07	119.18
2	C	407	842	C27-C28-C29	5.02	123.11	118.96
2	A	407	842	C17-C16-N20	4.91	123.84	119.18
2	C	407	842	C26-C4-N3	-4.87	104.52	110.20
2	C	407	842	C17-C16-N20	4.79	123.73	119.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	407	842	C24-N20-N21	4.67	115.02	111.93
2	B	407	842	C16-C15-C14	4.51	123.70	120.06
2	C	407	842	C6-N7-C8	3.94	124.18	116.86
2	B	407	842	C4-N3-C2	3.89	128.89	122.95
2	C	407	842	C16-C15-C14	3.72	123.06	120.06
2	C	407	842	C18-C17-C16	3.57	123.48	118.70
2	C	407	842	C27-C26-C4	-3.54	106.89	113.46
2	A	407	842	C27-C26-C4	-3.45	107.06	113.46
2	B	407	842	C27-C28-C29	3.24	121.64	118.96
2	A	407	842	C24-N20-N21	2.88	113.84	111.93
2	B	407	842	C13-C12-C11	-2.88	106.84	111.28
2	A	407	842	C26-C4-N3	-2.75	107.00	110.20
2	B	407	842	C18-C17-C16	2.62	122.22	118.70
2	C	407	842	C26-C4-C5	2.61	116.47	111.37
2	A	407	842	C14-C8-N7	-2.57	104.14	110.75
2	C	407	842	C5-C6-N7	2.37	113.53	110.61
2	A	407	842	C6-N7-C8	2.32	121.17	116.86
2	C	407	842	C24-N20-N21	2.32	113.47	111.93
2	B	407	842	C30-C29-C28	-2.29	120.24	123.48
2	C	407	842	C24-N20-C16	-2.23	121.00	124.58
2	C	407	842	F33-C29-C28	2.18	121.58	118.22
2	C	407	842	F34-C31-C32	2.16	121.55	118.22
2	A	407	842	C5-C6-N7	2.14	113.25	110.61
2	C	407	842	C30-C29-C28	-2.07	120.55	123.48
2	B	407	842	C13-C8-C9	-2.06	106.93	109.36
2	C	407	842	C28-C27-C32	-2.04	115.88	118.95

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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	375/406 (92%)	0.02	14 (3%)	39 44	43, 66, 99, 111	0
1	B	375/406 (92%)	0.03	19 (5%)	27 30	44, 67, 99, 114	0
1	C	375/406 (92%)	-0.02	17 (4%)	32 36	45, 66, 99, 112	0
All	All	1125/1218 (92%)	0.01	50 (4%)	33 37	43, 66, 99, 114	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	393	PRO	5.7
1	C	174	ALA	5.0
1	C	318	VAL	4.8
1	A	392	ILE	4.4
1	A	175	SER	4.3
1	A	45	ALA	3.9
1	A	125	LEU	3.8
1	B	320	THR	3.5
1	B	316	GLU	3.4
1	B	321	SER	3.4
1	B	392	ILE	3.3
1	B	45	ALA	3.1
1	C	322	GLN	3.1
1	C	368	HIS	3.1
1	B	124	ILE	3.0
1	A	322	GLN	2.9
1	A	318	VAL	2.9
1	B	366	HIS	2.9
1	B	278	ALA	2.8
1	C	126	GLY	2.8
1	B	125	LEU	2.7
1	A	126	GLY	2.6
1	B	322	GLN	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	323	ASP	2.6
1	A	44	PHE	2.6
1	B	262	LYS	2.5
1	A	368	HIS	2.5
1	C	44	PHE	2.5
1	A	317	ASP	2.4
1	B	4	SER	2.4
1	C	125	LEU	2.4
1	B	368	HIS	2.4
1	B	126	GLY	2.4
1	C	316	GLU	2.4
1	C	370	GLU	2.3
1	A	124	ILE	2.3
1	C	151	HIS	2.2
1	C	317	ASP	2.2
1	A	13	ARG	2.2
1	C	38	ASP	2.2
1	C	371	PHE	2.2
1	C	124	ILE	2.2
1	B	44	PHE	2.2
1	A	391	ASN	2.1
1	B	260	THR	2.1
1	C	45	ALA	2.1
1	A	38	ASP	2.1
1	C	277	GLN	2.1
1	C	39	THR	2.0
1	B	315	VAL	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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
2	842	B	407	35/35	0.14	-0.73	44,50,63,65	0
2	842	C	407	35/35	0.14	-0.79	44,49,63,66	0
2	842	A	407	35/35	0.14	-0.82	44,50,61,63	0

6.5 Other polymers ⓘ

There are no such residues in this entry.