



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:57 PM UTC

PDB ID : 1RZQ / pdb_00001rzq
Title : Crystal Structure of C-Terminal Despentapeptide Nitrite Reductase from
Achromobacter Cycloclastes at pH5.0
Authors : Li, H.T.; Wang, C.; Chang, T.; Chang, W.C.; Liu, M.Y.; Le Gall, J.; Gui,
L.L.; Zhang, J.P.; An, X.M.; Chang, W.R.
Deposited on : 2003-12-26
Resolution : 2.20 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

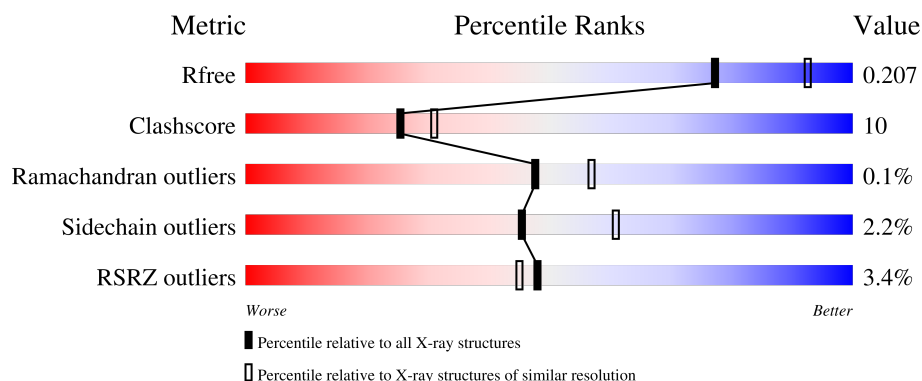
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

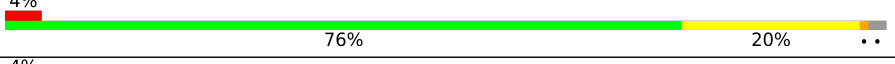
The reported resolution of this entry is 2.20 Å.

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}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	335	
1	B	335	
1	C	335	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	ACY	A	2001	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8487 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 Copper-containing nitrite reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	328	Total	C	N	O	S	0	0	0
			2540	1618	435	478	9			
1	B	328	Total	C	N	O	S	0	0	0
			2540	1618	435	478	9			
1	C	332	Total	C	N	O	S	0	0	0
			2564	1634	439	482	9			

- Molecule 2 is COPPER (II) ION (CCD ID: CU) (formula: Cu).

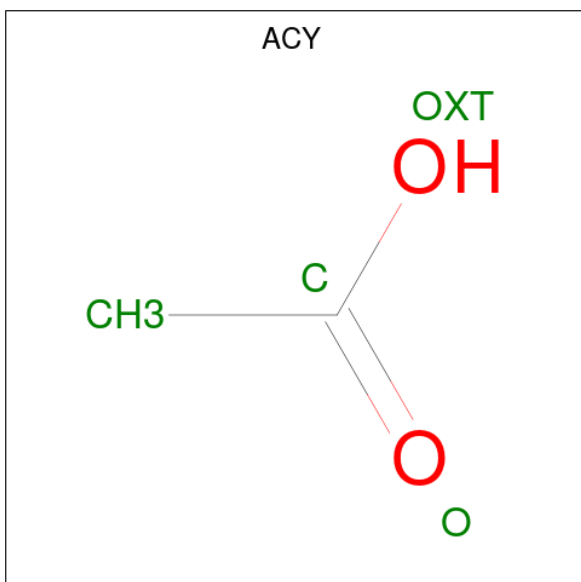
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Cu	0	0
			2	2		
2	B	2	Total	Cu	0	0
			2	2		
2	C	2	Total	Cu	0	0
			2	2		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is ACETIC ACID (CCD ID: ACY) (formula: C₂H₄O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0

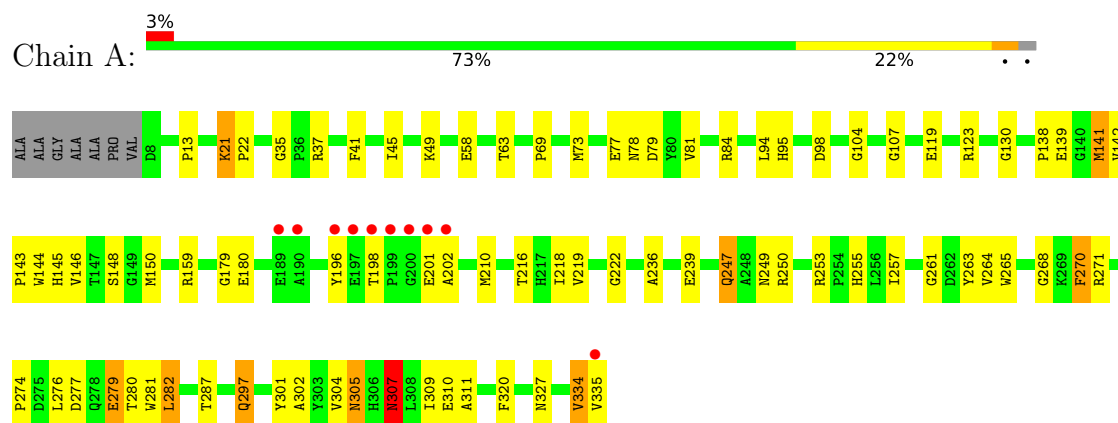
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	286	Total O 286 286	0	0
5	B	256	Total O 256 256	0	0
5	C	254	Total O 254 254	0	0

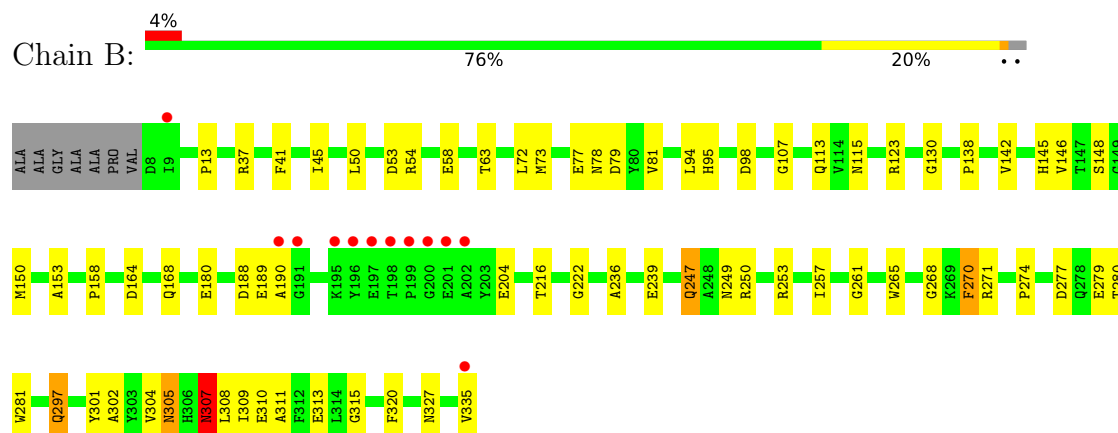
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

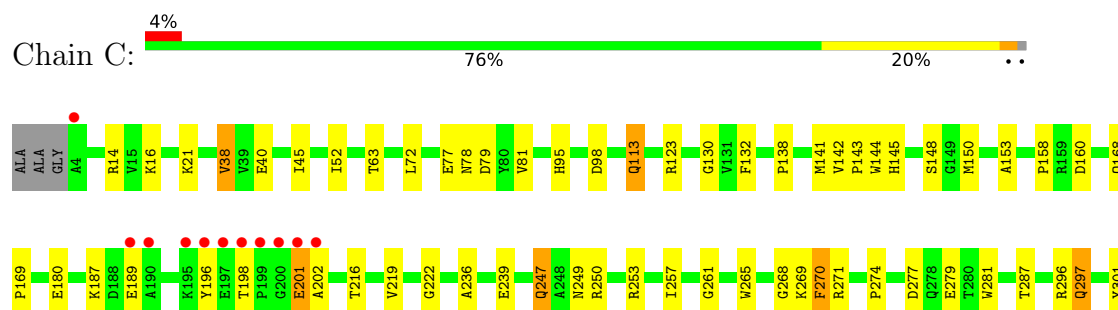
• Molecule 1: Copper-containing nitrite reductase



• Molecule 1: Copper-containing nitrite reductase



• Molecule 1: Copper-containing nitrite reductase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	99.02Å 117.44Å 122.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.20 20.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	95.2 (20.00-2.20) 95.1 (20.00-2.20)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.36 (at 2.20Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.181 , 0.208 0.181 , 0.207	Depositor DCC
R_{free} test set	2611 reflections (3.57%)	wwPDB-VP
Wilson B-factor (Å ²)	20.8	Xtriage
Anisotropy	0.292	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 47.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.013 for -h,l,k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8487	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.43% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ACY, SO4, CU

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	0.36	0/2609	1.02	16/3557 (0.4%)
1	B	0.35	0/2609	1.01	14/3557 (0.4%)
1	C	0.35	0/2634	1.01	16/3593 (0.4%)
All	All	0.35	0/7852	1.01	46/10707 (0.4%)

There are no bond length outliers.

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	307	ASN	N-CA-C	-9.97	89.21	107.75
1	B	307	ASN	N-CA-C	-9.73	89.64	107.75
1	B	268	GLY	N-CA-C	9.63	127.54	113.48
1	A	305	ASN	N-CA-C	-9.50	94.71	109.25
1	C	268	GLY	N-CA-C	9.33	127.10	113.48
1	C	307	ASN	N-CA-C	-9.14	90.74	107.75
1	A	268	GLY	N-CA-C	8.96	127.63	114.10
1	B	305	ASN	N-CA-C	-8.76	95.03	108.96
1	C	305	ASN	N-CA-C	-8.67	95.18	108.96
1	C	79	ASP	N-CA-C	-8.30	99.82	110.53
1	B	79	ASP	N-CA-C	-7.84	100.41	110.53
1	B	222	GLY	N-CA-C	7.53	124.04	115.08
1	B	270	PHE	N-CA-C	7.35	121.35	112.38
1	A	270	PHE	N-CA-C	7.03	121.31	112.87
1	C	222	GLY	N-CA-C	7.00	124.31	115.21
1	A	279	GLU	N-CA-C	-6.78	103.03	112.45
1	B	279	GLU	N-CA-C	-6.70	103.14	112.45
1	A	222	GLY	N-CA-C	6.62	122.95	115.08
1	C	279	GLU	N-CA-C	-6.55	103.35	112.45
1	A	304	VAL	N-CA-C	6.29	117.77	108.65
1	C	158	PRO	N-CA-C	-6.05	101.59	111.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	261	GLY	N-CA-C	-5.98	99.00	113.18
1	A	148	SER	N-CA-C	-5.98	105.56	112.92
1	C	270	PHE	N-CA-C	5.97	120.03	112.87
1	B	261	GLY	N-CA-C	-5.95	99.09	113.18
1	A	79	ASP	N-CA-C	-5.94	102.53	110.55
1	B	148	SER	N-CA-C	-5.93	105.69	113.17
1	B	216	THR	N-CA-C	-5.83	105.27	112.90
1	C	261	GLY	N-CA-C	-5.75	99.56	113.18
1	B	304	VAL	N-CA-C	5.74	116.97	108.65
1	C	304	VAL	N-CA-C	5.68	116.88	108.65
1	C	52	ILE	N-CA-C	5.50	118.18	112.90
1	B	63	THR	N-CA-C	5.50	118.54	109.85
1	B	315	GLY	N-CA-C	5.46	122.86	115.00
1	A	216	THR	N-CA-C	-5.39	105.84	112.90
1	C	63	THR	N-CA-C	5.38	118.35	109.85
1	C	216	THR	N-CA-C	-5.23	106.05	112.90
1	A	334	VAL	N-CA-C	5.22	115.74	110.05
1	B	158	PRO	N-CA-C	-5.18	103.14	111.38
1	C	219	VAL	N-CA-C	5.17	117.48	108.95
1	C	38	VAL	N-CA-C	-5.05	99.86	107.73
1	A	219	VAL	N-CA-C	5.04	117.27	108.95
1	A	63	THR	N-CA-C	5.04	117.82	109.85
1	A	104	GLY	N-CA-C	5.04	116.73	111.95
1	C	148	SER	N-CA-C	-5.02	106.75	112.92
1	A	264	VAL	N-CA-C	5.01	115.33	108.12

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2540	0	2464	53	0
1	B	2540	0	2464	53	0
1	C	2564	0	2490	51	0
2	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	2	0	0	0	0
2	C	2	0	0	0	0
3	A	10	0	0	0	0
3	B	10	0	0	0	0
3	C	5	0	0	0	0
4	A	12	0	9	3	0
4	C	4	0	3	0	0
5	A	286	0	0	0	0
5	B	256	0	0	6	0
5	C	254	0	0	2	0
All	All	8487	0	7430	145	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 10.

All (145) 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:A:13:PRO:HG2	1:A:37:ARG:HG2	1.46	0.97
1:B:297:GLN:HE22	1:B:327:ASN:H	1.15	0.93
1:A:297:GLN:HE22	1:A:327:ASN:H	1.16	0.90
1:C:297:GLN:HE22	1:C:327:ASN:H	1.18	0.87
4:A:2001:ACY:H1	1:B:313:GLU:OE2	1.82	0.80
1:A:196:TYR:CG	1:A:202:ALA:HB2	2.24	0.73
1:B:13:PRO:HG2	1:B:37:ARG:HD3	1.70	0.72
1:B:13:PRO:HG2	1:B:37:ARG:CD	2.22	0.69
1:C:189:GLU:H	1:C:189:GLU:CD	2.01	0.68
1:C:307:ASN:C	1:C:307:ASN:HD22	2.00	0.68
1:A:271:ARG:HD2	1:B:277:ASP:OD2	1.94	0.67
1:A:297:GLN:NE2	1:A:327:ASN:H	1.90	0.66
1:C:141:MET:HE2	1:C:145:HIS:HE2	1.60	0.65
1:A:196:TYR:CD2	1:A:202:ALA:HB2	2.32	0.65
1:C:141:MET:CE	1:C:145:HIS:HE2	2.09	0.65
1:B:335:VAL:OXT	1:B:335:VAL:HG23	1.97	0.64
1:C:257:ILE:HD12	1:C:302:ALA:HB3	1.79	0.64
1:A:277:ASP:OD2	1:C:271:ARG:HD2	1.96	0.64
1:C:297:GLN:HE21	1:C:297:GLN:HA	1.63	0.64
1:B:297:GLN:HE21	1:B:297:GLN:HA	1.62	0.63
1:A:307:ASN:HD22	1:A:307:ASN:C	2.07	0.62
1:B:271:ARG:HD2	1:C:277:ASP:OD2	2.00	0.61
1:B:307:ASN:HD22	1:B:307:ASN:C	2.09	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:198:THR:HG23	1:C:201:GLU:H	1.67	0.60
1:A:77:GLU:O	1:A:78:ASN:HB2	2.02	0.59
1:C:14:ARG:HG2	1:C:38:VAL:HB	1.82	0.59
1:A:257:ILE:HD12	1:A:302:ALA:HB3	1.83	0.59
1:B:164:ASP:OD2	1:B:168:GLN:HB2	2.03	0.58
1:C:196:TYR:CG	1:C:202:ALA:HB2	2.38	0.58
1:A:250:ARG:NH2	1:A:310:GLU:OE2	2.36	0.57
1:B:301:TYR:HB2	1:B:320:PHE:HB2	1.85	0.57
1:C:301:TYR:HB2	1:C:320:PHE:HB2	1.87	0.56
1:B:13:PRO:HG2	1:B:37:ARG:NE	2.21	0.55
1:B:250:ARG:NH1	5:B:3209:HOH:O	2.38	0.55
1:A:35:GLY:HA2	1:A:159:ARG:NH2	2.21	0.55
1:C:21:LYS:HE2	5:C:3222:HOH:O	2.06	0.55
1:A:307:ASN:HD22	1:A:309:ILE:H	1.56	0.54
1:B:189:GLU:H	1:B:189:GLU:CD	2.15	0.54
1:C:141:MET:HE2	1:C:145:HIS:NE2	2.23	0.54
1:C:297:GLN:NE2	1:C:327:ASN:H	1.98	0.54
1:A:301:TYR:HB2	1:A:320:PHE:HB2	1.89	0.54
1:A:130:GLY:HA2	1:A:270:PHE:CD2	2.44	0.53
1:A:45:ILE:HD13	1:A:95:HIS:HB2	1.90	0.53
1:A:334:VAL:HG12	1:A:335:VAL:N	2.24	0.53
1:B:45:ILE:HD13	1:B:95:HIS:HB2	1.91	0.52
1:A:297:GLN:HE21	1:A:297:GLN:HA	1.75	0.52
1:C:307:ASN:HD21	1:C:309:ILE:HB	1.74	0.52
1:B:257:ILE:HD12	1:B:302:ALA:HB3	1.92	0.52
1:B:77:GLU:O	1:B:78:ASN:HB2	2.10	0.51
1:B:249:ASN:O	1:C:307:ASN:HA	2.10	0.51
1:A:94:LEU:HD23	1:A:94:LEU:C	2.35	0.51
1:A:13:PRO:CG	1:A:37:ARG:HG2	2.30	0.51
1:A:143:PRO:HG2	4:A:2001:ACY:C	2.40	0.51
1:C:77:GLU:O	1:C:78:ASN:HB2	2.10	0.51
1:A:49:LYS:HD3	1:A:58:GLU:OE2	2.11	0.51
1:B:250:ARG:NH2	1:B:310:GLU:OE2	2.43	0.50
1:A:21:LYS:HD3	1:A:22:PRO:HD2	1.94	0.50
1:B:54:ARG:NH1	5:B:3144:HOH:O	2.44	0.50
1:C:296:ARG:HH12	1:C:330:LEU:HD21	1.76	0.50
1:C:307:ASN:C	1:C:307:ASN:ND2	2.70	0.50
1:A:142:VAL:HG12	1:B:308:LEU:HD13	1.94	0.50
1:B:45:ILE:HG21	1:B:95:HIS:CD2	2.48	0.49
1:B:189:GLU:CD	1:B:189:GLU:N	2.70	0.49
1:B:53:ASP:HB2	5:B:3041:HOH:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:297:GLN:NE2	1:B:327:ASN:H	1.96	0.49
1:A:81:VAL:O	1:A:123:ARG:HA	2.12	0.49
1:A:84:ARG:HD2	1:A:119:GLU:OE2	2.12	0.49
1:B:253:ARG:HA	1:B:281:TRP:O	2.13	0.49
1:A:141:MET:HG2	1:A:144:TRP:HE3	1.77	0.49
1:A:236:ALA:O	1:A:239:GLU:HB2	2.13	0.49
1:C:189:GLU:CD	1:C:189:GLU:N	2.69	0.48
1:B:180:GLU:HB3	1:B:247:GLN:HG2	1.96	0.48
1:B:236:ALA:O	1:B:239:GLU:HB2	2.14	0.48
1:C:236:ALA:O	1:C:239:GLU:HB2	2.14	0.48
1:A:249:ASN:O	1:B:307:ASN:HA	2.14	0.48
1:A:253:ARG:HA	1:A:281:TRP:O	2.14	0.48
1:C:98:ASP:HB2	5:C:3167:HOH:O	2.14	0.47
1:A:146:VAL:HG21	1:B:308:LEU:CD1	2.44	0.47
1:C:306:HIS:O	1:C:306:HIS:ND1	2.48	0.47
1:B:130:GLY:HA2	1:B:270:PHE:CD2	2.49	0.47
1:C:141:MET:HE3	1:C:144:TRP:CZ3	2.49	0.47
1:A:307:ASN:HA	1:C:249:ASN:O	2.14	0.47
1:B:142:VAL:HG12	1:C:308:LEU:HD13	1.97	0.47
1:A:210:MET:HE1	4:A:2001:ACY:H2	1.97	0.46
1:A:307:ASN:HD21	1:A:309:ILE:HB	1.80	0.46
1:B:138:PRO:HG2	1:B:145:HIS:CE1	2.50	0.46
1:B:94:LEU:HB2	1:B:115:ASN:ND2	2.31	0.46
1:B:113:GLN:HG3	5:B:3100:HOH:O	2.16	0.46
1:C:296:ARG:NH1	1:C:330:LEU:HD21	2.29	0.46
1:C:253:ARG:HA	1:C:281:TRP:O	2.15	0.46
1:A:180:GLU:HB3	1:A:247:GLN:HG2	1.97	0.45
1:B:81:VAL:O	1:B:123:ARG:HA	2.16	0.45
1:B:94:LEU:HD13	1:B:115:ASN:ND2	2.31	0.45
1:B:309:ILE:O	1:B:313:GLU:HB2	2.16	0.45
1:C:180:GLU:HB3	1:C:247:GLN:HG2	1.99	0.45
1:C:196:TYR:CD2	1:C:202:ALA:HB2	2.51	0.45
1:C:309:ILE:O	1:C:313:GLU:HB2	2.17	0.45
1:A:198:THR:OG1	1:A:201:GLU:HB2	2.17	0.45
1:C:81:VAL:O	1:C:123:ARG:HA	2.16	0.45
1:A:307:ASN:C	1:A:307:ASN:ND2	2.75	0.44
1:A:130:GLY:HA2	1:A:270:PHE:CE2	2.52	0.44
1:C:198:THR:CG2	1:C:201:GLU:H	2.30	0.44
1:C:265:TRP:CE2	1:C:274:PRO:HB3	2.53	0.44
1:C:72:LEU:HD12	1:C:153:ALA:O	2.18	0.44
1:A:98:ASP:OD1	1:A:107:GLY:HA3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:305:ASN:O	1:A:311:ALA:HB2	2.18	0.44
1:A:307:ASN:CG	1:A:310:GLU:HG3	2.41	0.44
1:A:307:ASN:ND2	1:A:309:ILE:H	2.16	0.43
1:B:72:LEU:HD12	1:B:153:ALA:O	2.19	0.43
1:B:146:VAL:HG21	1:C:308:LEU:CD1	2.49	0.43
1:C:307:ASN:HD22	1:C:309:ILE:H	1.65	0.43
1:A:282:LEU:C	1:A:282:LEU:HD13	2.44	0.43
1:A:218:ILE:HD12	1:A:310:GLU:HB3	2.01	0.43
1:B:307:ASN:HD22	1:B:309:ILE:H	1.66	0.43
1:B:41:PHE:CG	1:B:73:MET:HE3	2.54	0.43
1:A:265:TRP:CE2	1:A:274:PRO:HB3	2.54	0.42
1:B:98:ASP:OD1	1:B:107:GLY:HA3	2.19	0.42
1:C:113:GLN:HE21	1:C:113:GLN:HB3	1.59	0.42
1:C:168:GLN:HA	1:C:169:PRO:HD3	1.91	0.42
1:B:265:TRP:CE2	1:B:274:PRO:HB3	2.53	0.42
1:A:255:HIS:ND1	1:A:279:GLU:O	2.53	0.42
1:B:204:GLU:HB2	5:B:3052:HOH:O	2.20	0.42
1:C:138:PRO:HG2	1:C:145:HIS:CE1	2.54	0.42
1:B:204:GLU:HG3	5:B:3228:HOH:O	2.20	0.42
1:C:45:ILE:HD13	1:C:95:HIS:HB2	2.01	0.42
1:C:132:PHE:CE2	1:C:269:LYS:HE3	2.55	0.42
1:C:16:LYS:HD2	1:C:40:GLU:O	2.20	0.42
1:B:305:ASN:O	1:B:311:ALA:HB2	2.20	0.41
1:B:307:ASN:C	1:B:307:ASN:ND2	2.77	0.41
1:A:263:TYR:CD1	1:A:276:LEU:HD22	2.56	0.41
1:C:305:ASN:O	1:C:311:ALA:HB2	2.20	0.41
1:A:41:PHE:CG	1:A:73:MET:HE3	2.55	0.41
1:A:287:THR:OG1	1:B:280:THR:HB	2.21	0.41
1:B:50:LEU:O	1:B:58:GLU:HA	2.21	0.41
1:C:130:GLY:HA2	1:C:270:PHE:CD2	2.56	0.41
1:C:307:ASN:ND2	1:C:309:ILE:H	2.19	0.41
1:A:280:THR:HB	1:C:287:THR:OG1	2.21	0.41
1:A:94:LEU:HD23	1:A:95:HIS:N	2.37	0.40
1:B:188:ASP:OD1	1:B:190:ALA:HB3	2.21	0.40
1:B:307:ASN:HD21	1:B:309:ILE:HB	1.85	0.40
1:B:95:HIS:CG	1:B:150:MET:HE1	2.56	0.40
1:C:142:VAL:HB	1:C:143:PRO:CD	2.51	0.40
1:C:250:ARG:NH2	1:C:310:GLU:OE2	2.47	0.40
1:A:69:PRO:HG3	1:A:179:GLY:HA3	2.02	0.40
1:A:138:PRO:HG2	1:A:145:HIS:CE1	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	326/335 (97%)	316 (97%)	10 (3%)	0	100	100
1	B	326/335 (97%)	316 (97%)	10 (3%)	0	100	100
1	C	330/335 (98%)	319 (97%)	10 (3%)	1 (0%)	36	42
All	All	982/1005 (98%)	951 (97%)	30 (3%)	1 (0%)	48	57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	201	GLU

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	267/269 (99%)	259 (97%)	8 (3%)	36	49
1	B	267/269 (99%)	264 (99%)	3 (1%)	65	79
1	C	269/269 (100%)	262 (97%)	7 (3%)	40	55
All	All	803/807 (100%)	785 (98%)	18 (2%)	45	61

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	LYS
1	A	139	GLU

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Mol	Chain	Res	Type
1	A	141	MET
1	A	150	MET
1	A	247	GLN
1	A	282	LEU
1	A	297	GLN
1	A	307	ASN
1	B	247	GLN
1	B	297	GLN
1	B	307	ASN
1	C	113	GLN
1	C	150	MET
1	C	160	ASP
1	C	187	LYS
1	C	247	GLN
1	C	297	GLN
1	C	307	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	78	ASN
1	A	113	GLN
1	A	192	ASN
1	A	231	HIS
1	A	297	GLN
1	A	307	ASN
1	B	78	ASN
1	B	113	GLN
1	B	115	ASN
1	B	192	ASN
1	B	231	HIS
1	B	297	GLN
1	B	307	ASN
1	B	319	HIS
1	C	78	ASN
1	C	113	GLN
1	C	192	ASN
1	C	231	HIS
1	C	297	GLN
1	C	307	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 6 are monoatomic - leaving 9 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	ACY	A	2002	-	3,3,3	1.00	0	3,3,3	0.86	0
3	SO4	A	3002	-	4,4,4	0.39	0	6,6,6	0.11	0
4	ACY	A	2004	-	3,3,3	1.03	0	3,3,3	0.81	0
3	SO4	B	3001	-	4,4,4	0.36	0	6,6,6	0.11	0
3	SO4	A	3005	-	4,4,4	0.38	0	6,6,6	0.04	0
3	SO4	B	3004	-	4,4,4	0.37	0	6,6,6	0.08	0
3	SO4	C	3003	-	4,4,4	0.39	0	6,6,6	0.05	0
4	ACY	C	2003	-	3,3,3	1.01	0	3,3,3	0.80	0
4	ACY	A	2001	-	3,3,3	0.97	0	3,3,3	0.86	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	2001	ACY	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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	328/335 (97%)	-0.48	10 (3%) 52 49	11, 18, 41, 69	0
1	B	328/335 (97%)	-0.37	12 (3%) 45 42	12, 20, 48, 72	0
1	C	332/335 (99%)	-0.38	12 (3%) 46 43	11, 19, 44, 73	0
All	All	988/1005 (98%)	-0.41	34 (3%) 48 45	11, 19, 46, 73	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	199	PRO	6.2
1	B	199	PRO	5.0
1	C	198	THR	4.9
1	C	196	TYR	4.8
1	A	199	PRO	4.4
1	B	196	TYR	4.2
1	A	196	TYR	4.2
1	B	198	THR	3.8
1	A	198	THR	3.7
1	B	9	ILE	3.6
1	C	200	GLY	3.5
1	A	190	ALA	3.5
1	A	200	GLY	3.4
1	C	197	GLU	3.3
1	B	195	LYS	3.2
1	B	335	VAL	3.1
1	B	197	GLU	3.1
1	C	190	ALA	3.1
1	B	191	GLY	3.0
1	A	189	GLU	3.0
1	C	4	ALA	2.9
1	A	197	GLU	2.9
1	B	200	GLY	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	201	GLU	2.8
1	C	195	LYS	2.8
1	C	201	GLU	2.6
1	C	335	VAL	2.6
1	C	202	ALA	2.4
1	B	190	ALA	2.4
1	A	202	ALA	2.3
1	C	189	GLU	2.3
1	B	202	ALA	2.2
1	B	201	GLU	2.2
1	A	335	VAL	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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. 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	B-factors(Å ²)	Q<0.9
4	ACY	C	2003	4/4	0.79	0.16	37,38,38,39	0
4	ACY	A	2004	4/4	0.84	0.14	39,39,40,41	0
4	ACY	A	2002	4/4	0.84	0.13	29,30,32,33	0
4	ACY	A	2001	4/4	0.88	0.18	38,38,39,40	0
3	SO4	A	3005	5/5	0.89	0.13	50,50,51,51	5
3	SO4	B	3004	5/5	0.90	0.12	40,41,43,44	5
3	SO4	A	3002	5/5	0.91	0.14	49,50,51,52	0
3	SO4	B	3001	5/5	0.94	0.11	53,53,54,55	0
3	SO4	C	3003	5/5	0.96	0.09	43,45,45,45	0
2	CU	A	502	1/1	0.99	0.04	19,19,19,19	0
2	CU	B	502	1/1	0.99	0.03	21,21,21,21	0
2	CU	C	501	1/1	0.99	0.02	26,26,26,26	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CU	A	501	1/1	1.00	0.02	27,27,27,27	0
2	CU	B	501	1/1	1.00	0.01	27,27,27,27	0
2	CU	C	502	1/1	1.00	0.02	19,19,19,19	0

6.5 Other polymers [i](#)

There are no such residues in this entry.