



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 26, 2023 – 04:27 AM EDT

PDB ID : 1AS8
Title : STRUCTURE OF NITRITE BOUND TO REDUCED ALCALIGENES FAECALIS NITRITE REDUCTASE AT CRYO TEMPERATURE
Authors : Murphy, M.E.P.; Adman, E.T.; Turley, S.
Deposited on : 1997-08-13
Resolution : 1.85 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.34
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.34

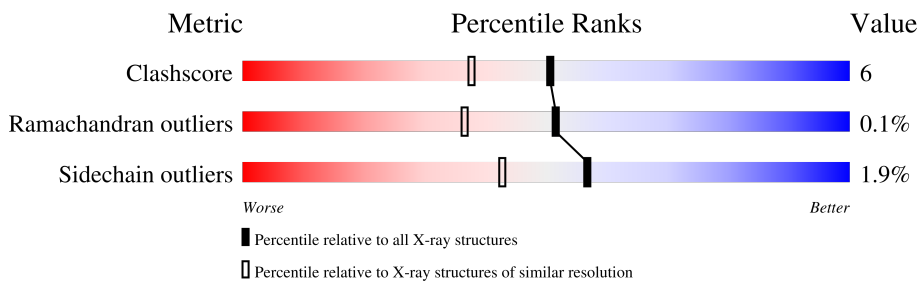
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)

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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	343	
1	B	343	
1	C	343	

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
3	NO2	B	503	-	-	X	-
3	NO2	C	503	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11412 atoms, of which 2994 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.

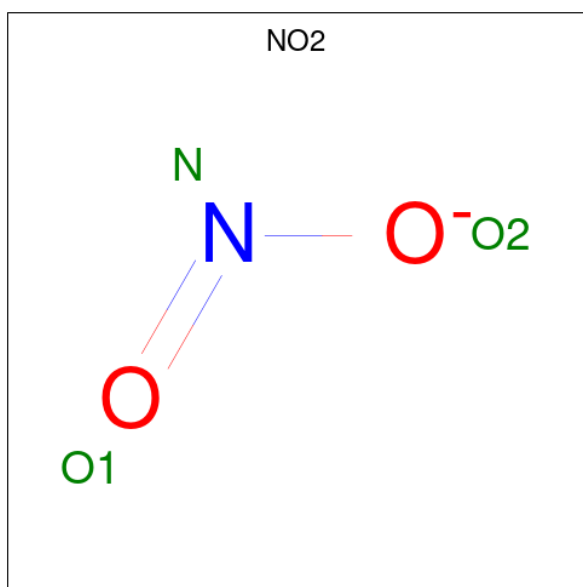
- Molecule 1 is a protein called NITRITE REDUCTASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	336	3076	1639	516	430	480	11	0	0	0
1	B	336	3076	1639	516	430	480	11	0	0	0
1	C	336	3076	1639	516	430	480	11	0	0	0

- Molecule 2 is COPPER (II) ION (three-letter code: 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 NITRITE ION (three-letter code: NO2) (formula: NO₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	N	O	0	0
			3	1	2		
3	B	1	Total	N	O	0	0
			3	1	2		
3	C	1	Total	N	O	0	0
			3	1	2		


- Molecule 4 is water.

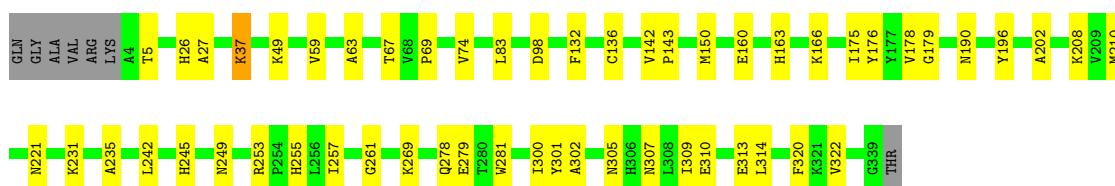
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	258	Total	H	O	0	0
			774	516	258		
4	B	231	Total	H	O	2	0
			693	462	231		
4	C	234	Total	H	O	0	0
			702	468	234		

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. 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. 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.

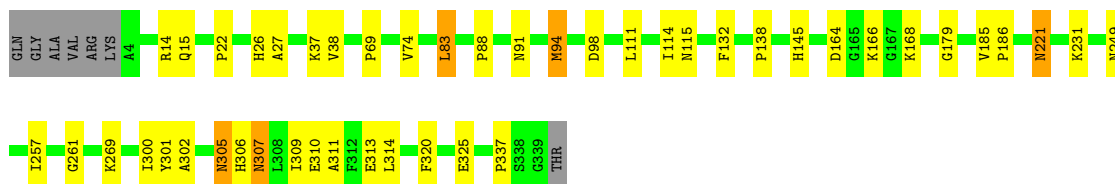
- Molecule 1: NITRITE REDUCTASE

Chain A:  82% 15%




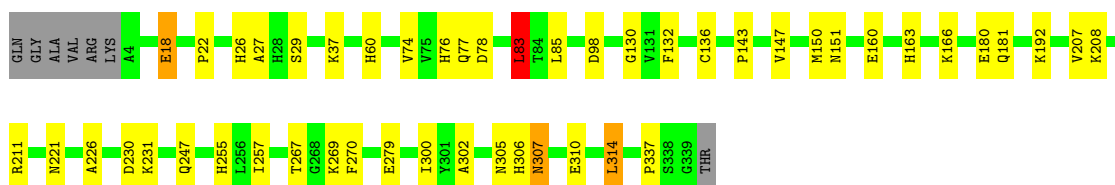
- Molecule 1: NITRITE REDUCTASE

Chain B:  85% 12%



- Molecule 1: NITRITE REDUCTASE

Chain C:  84% 13%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	61.10Å 102.80Å 146.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 1.85 19.88 – 1.82	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-1.85) 71.2 (19.88-1.82)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.37 (at 1.82Å)	Xtrriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.194 , (Not available) 0.268 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	14.9	Xtrriage
Anisotropy	0.425	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 45.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	11412	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.05% 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: NO2, 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	0.52	0/2631	0.78	4/3588 (0.1%)
1	B	0.51	0/2631	0.77	5/3588 (0.1%)
1	C	0.51	0/2631	0.78	4/3588 (0.1%)
All	All	0.51	0/7893	0.78	13/10764 (0.1%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	307	ASN	N-CA-C	-7.75	90.06	111.00
1	A	307	ASN	N-CA-C	-7.19	91.58	111.00
1	B	307	ASN	N-CA-C	-7.17	91.63	111.00
1	B	305	ASN	N-CA-C	-6.13	94.44	111.00
1	A	261	GLY	N-CA-C	-6.04	98.00	113.10
1	A	305	ASN	N-CA-C	-5.96	94.90	111.00
1	C	98	ASP	CB-CG-OD2	5.81	123.53	118.30
1	B	83	LEU	CA-CB-CG	5.62	128.24	115.30
1	C	305	ASN	N-CA-C	-5.59	95.90	111.00
1	C	83	LEU	CA-CB-CG	5.51	127.97	115.30
1	B	98	ASP	CB-CG-OD2	5.40	123.16	118.30
1	A	98	ASP	CB-CG-OD2	5.28	123.06	118.30
1	B	261	GLY	N-CA-C	-5.20	100.10	113.10

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	2560	516	2485	29	8
1	B	2560	516	2485	27	5
1	C	2560	516	2485	34	2
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
3	A	3	0	0	1	0
3	B	3	0	0	2	0
3	C	3	0	0	2	0
4	A	258	516	0	1	13
4	B	231	462	0	1	6
4	C	234	468	0	2	9
All	All	8418	2994	7455	86	22

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 6.

All (86) 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:C:151:ASN:HD21	1:C:181:GLN:HE22	1.24	0.85
1:A:257:ILE:HG12	3:C:503:NO2:N	1.91	0.85
1:A:26:HIS:HE1	1:A:74:VAL:H	1.26	0.83
1:B:26:HIS:HE1	1:B:74:VAL:H	1.27	0.82
1:C:26:HIS:HE1	1:C:74:VAL:H	1.30	0.79
1:C:257:ILE:HD12	1:C:302:ALA:HB3	1.66	0.76
1:B:257:ILE:HD12	1:B:302:ALA:HB3	1.67	0.76
1:C:26:HIS:HD2	1:C:27:ALA:O	1.68	0.75
1:C:26:HIS:CE1	1:C:74:VAL:H	2.10	0.70
1:C:160:GLU:HG2	1:C:163:HIS:CE1	2.27	0.70
1:B:14:ARG:HG2	1:B:38:VAL:HB	1.74	0.67
3:A:503:NO2:N	1:B:257:ILE:HG12	2.09	0.67
1:B:26:HIS:CE1	1:B:74:VAL:H	2.10	0.67
1:A:136:CYS:HB2	1:A:150:MET:HG2	1.76	0.66
1:C:160:GLU:HG2	1:C:163:HIS:HE1	1.61	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:257:ILE:HD12	1:A:302:ALA:HB3	1.77	0.65
1:B:22:PRO:HB2	1:B:221:ASN:HD21	1.61	0.65
1:A:208:LYS:HD2	4:A:695:HOH:O	1.97	0.64
3:B:503:NO2:N	1:C:257:ILE:HG12	2.12	0.64
1:B:301:TYR:HB2	1:B:320:PHE:HB2	1.79	0.63
1:A:26:HIS:CE1	1:A:74:VAL:H	2.12	0.63
1:C:22:PRO:HB2	1:C:221:ASN:HD21	1.63	0.62
1:B:26:HIS:HD2	1:B:27:ALA:O	1.85	0.60
3:B:503:NO2:O2	1:C:257:ILE:HG12	2.02	0.60
1:B:132:PHE:CE1	1:B:269:LYS:HE3	2.38	0.58
1:B:310:GLU:HA	1:B:314:LEU:HB2	1.87	0.57
1:C:226:ALA:O	1:C:231:LYS:HB2	2.05	0.57
1:C:60:HIS:HD2	4:C:647:HOH:O	1.87	0.56
1:C:300:ILE:HD12	1:C:300:ILE:O	2.06	0.55
1:A:235:ALA:O	1:A:322:VAL:HA	2.08	0.53
1:C:255:HIS:ND1	1:C:279:GLU:O	2.42	0.53
1:A:26:HIS:HD2	1:A:27:ALA:O	1.92	0.53
1:C:22:PRO:HB2	1:C:221:ASN:ND2	2.24	0.52
1:C:310:GLU:HA	1:C:314:LEU:HB2	1.90	0.52
1:C:143:PRO:O	1:C:147:VAL:HG22	2.09	0.52
1:C:26:HIS:CD2	1:C:27:ALA:O	2.57	0.52
1:C:130:GLY:HA2	1:C:270:PHE:CD1	2.46	0.51
1:C:151:ASN:HD21	1:C:181:GLN:NE2	2.02	0.51
1:C:29:SER:O	1:C:76:HIS:CE1	2.64	0.50
1:A:37:LYS:N	1:A:37:LYS:HD3	2.26	0.50
1:A:249:ASN:O	1:B:307:ASN:HA	2.12	0.50
1:C:207:VAL:O	1:C:211:ARG:HG3	2.13	0.49
1:A:257:ILE:HG12	3:C:503:NO2:O2	2.13	0.48
1:A:310:GLU:HA	1:A:314:LEU:HB2	1.95	0.47
1:B:111:LEU:C	4:B:731:HOH:O	2.51	0.47
1:A:132:PHE:CE1	1:A:269:LYS:HE3	2.50	0.47
1:A:196:TYR:CD2	1:A:202:ALA:HB2	2.50	0.47
1:C:18:GLU:H	1:C:18:GLU:CD	2.17	0.47
1:B:88:PRO:HD2	1:B:91:ASN:ND2	2.30	0.46
1:A:309:ILE:O	1:A:313:GLU:HB2	2.15	0.46
1:B:69:PRO:HG3	1:B:179:GLY:HA3	1.97	0.46
1:A:142:VAL:HB	1:A:143:PRO:HD3	1.97	0.46
1:A:253:ARG:HA	1:A:281:TRP:O	2.16	0.46
1:A:301:TYR:HB2	1:A:320:PHE:HB2	1.98	0.46
1:C:136:CYS:HB2	1:C:150:MET:HG2	1.98	0.46
1:B:132:PHE:CD1	1:B:269:LYS:HE3	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:255:HIS:ND1	1:A:279:GLU:O	2.49	0.45
1:B:300:ILE:O	1:B:300:ILE:HG13	2.16	0.45
1:A:160:GLU:HG2	1:A:163:HIS:HE1	1.82	0.45
1:A:63:ALA:HB1	1:A:67:THR:O	2.17	0.45
1:A:166:LYS:O	1:A:166:LYS:HG2	2.15	0.45
1:A:278:GLN:OE1	1:C:267:THR:HB	2.16	0.44
1:C:130:GLY:HA2	1:C:270:PHE:CE1	2.52	0.44
1:B:15:GLN:O	1:B:15:GLN:HG3	2.17	0.44
1:C:77:GLN:O	1:C:78:ASP:HB2	2.17	0.44
1:B:185:VAL:HA	1:B:186:PRO:HD3	1.89	0.44
1:B:94:MET:HB2	1:B:115:ASN:ND2	2.33	0.43
1:C:208:LYS:HA	1:C:208:LYS:HD3	1.70	0.43
1:B:114:ILE:HG22	1:C:337:PRO:HB3	2.00	0.43
1:B:138:PRO:HG2	1:B:145:HIS:CE1	2.54	0.43
1:B:305:ASN:O	1:B:311:ALA:HB2	2.18	0.43
1:C:132:PHE:CE1	1:C:269:LYS:HE3	2.54	0.43
1:B:22:PRO:HB2	1:B:221:ASN:ND2	2.31	0.42
1:C:180:GLU:HB3	1:C:247:GLN:HG2	2.01	0.42
1:A:69:PRO:HG3	1:A:179:GLY:HA3	2.02	0.42
1:A:143:PRO:HB2	1:A:210:MET:HE1	2.02	0.42
1:B:249:ASN:O	1:C:307:ASN:HA	2.19	0.41
1:A:178:VAL:O	1:A:245:HIS:HA	2.21	0.41
1:A:49:LYS:HA	1:A:59:VAL:O	2.20	0.41
1:A:175:ILE:HA	1:A:242:LEU:O	2.21	0.41
1:B:309:ILE:O	1:B:313:GLU:HB2	2.20	0.41
1:C:83:LEU:HD13	1:C:85:LEU:HB2	2.03	0.41
1:C:143:PRO:HG2	4:C:560:HOH:O	2.21	0.41
1:A:160:GLU:HG2	1:A:163:HIS:CE1	2.57	0.40
1:B:164:ASP:OD2	1:B:168:LYS:HB3	2.22	0.40
1:B:306:HIS:O	1:B:306:HIS:ND1	2.54	0.40

All (22) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:555:HOH:O	4:B:568:HOH:H1[4_556]	1.19	0.41
4:A:651:HOH:O	4:C:530:HOH:O[3_645]	1.86	0.34
4:A:651:HOH:O	4:C:530:HOH:H1[3_645]	1.27	0.33
1:A:27:ALA:H	4:C:700:HOH:H2[3_645]	1.31	0.29
4:B:650:HOH:H1	4:B:719:HOH:O[4_556]	1.32	0.28

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:5:THR:OG1	1:B:231:LYS:HZ3[4_456]	1.41	0.19
1:B:325:GLU:H	4:A:679:HOH:O[4_456]	1.41	0.19
1:C:230:ASP:H	4:A:701:HOH:O[3_655]	1.43	0.17
1:A:176:TYR:OH	4:C:645:HOH:O[3_645]	2.08	0.12
4:A:606:HOH:O	4:C:599:HOH:H1[3_645]	1.51	0.09
1:A:5:THR:OG1	1:B:231:LYS:NZ[4_456]	2.12	0.08
4:A:555:HOH:O	4:B:568:HOH:O[4_556]	2.13	0.07
1:A:176:TYR:OH	4:C:645:HOH:H2[3_645]	1.54	0.06
4:A:563:HOH:O	4:B:601:HOH:O[4_556]	2.14	0.06
1:A:190:ASN:OD1	1:B:166:LYS:HZ3[1_655]	1.55	0.05
1:A:231:LYS:HZ3	1:C:226:ALA:O[3_645]	1.55	0.05
1:A:190:ASN:OD1	1:B:166:LYS:NZ[1_655]	2.16	0.04
4:A:543:HOH:H2	4:B:603:HOH:O[4_556]	1.56	0.04
4:A:744:HOH:O	4:C:528:HOH:H2[3_645]	1.56	0.04
4:A:518:HOH:O	4:C:537:HOH:O[3_645]	2.17	0.03
4:A:538:HOH:O	4:C:541:HOH:H1[3_645]	1.59	0.01
4:A:752:HOH:O	4:B:716:HOH:H1[4_556]	1.59	0.01

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	334/343 (97%)	326 (98%)	8 (2%)	0	100	100
1	B	334/343 (97%)	326 (98%)	8 (2%)	0	100	100
1	C	334/343 (97%)	327 (98%)	6 (2%)	1 (0%)	41	26
All	All	1002/1029 (97%)	979 (98%)	22 (2%)	1 (0%)	51	36

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	306	HIS

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	264/269 (98%)	260 (98%)	4 (2%)	65	53
1	B	264/269 (98%)	259 (98%)	5 (2%)	57	43
1	C	264/269 (98%)	258 (98%)	6 (2%)	50	34
All	All	792/807 (98%)	777 (98%)	15 (2%)	57	43

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	LYS
1	A	83	LEU
1	A	221	ASN
1	A	300	ILE
1	B	37	LYS
1	B	83	LEU
1	B	94	MET
1	B	221	ASN
1	B	337	PRO
1	C	18	GLU
1	C	37	LYS
1	C	83	LEU
1	C	166	LYS
1	C	192	LYS
1	C	314	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	HIS
1	A	77	GLN
1	A	163	HIS
1	A	221	ASN
1	A	296	GLN
1	B	15	GLN

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Mol	Chain	Res	Type
1	B	26	HIS
1	B	115	ASN
1	B	163	HIS
1	B	221	ASN
1	C	26	HIS
1	C	60	HIS
1	C	77	GLN
1	C	115	ASN
1	C	163	HIS
1	C	181	GLN
1	C	221	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 6 are monoatomic - leaving 3 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	$\# Z > 2$
3	NO2	C	503	2	1,2,2	4.04	1 (100%)	0,1,1	-	-
3	NO2	A	503	2	1,2,2	4.12	1 (100%)	0,1,1	-	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NO2	B	503	2	1,2,2	4.24	1 (100%)	0,1,1	-	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	503	NO2	O1-N	4.24	1.43	1.22
3	A	503	NO2	O1-N	4.12	1.43	1.22
3	C	503	NO2	O1-N	4.04	1.42	1.22

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	503	NO2	2	0
3	A	503	NO2	1	0
3	B	503	NO2	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

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.