Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 04:37 AM GMT

PDB ID	:	2NAP
Title	:	DISSIMILATORY NITRATE REDUCTASE (NAP) FROM DESULFOVIBRIO DESULFURICANS
Authors	:	Dias, J.M.; Than, M.; Humm, A.; Huber, R.; Bourenkov, G.; Bartunik, H.; Bursakov, S.; Calvete, J.; Caldeira, J.; Carneiro, C.; Moura, J.; Moura, I.; Romao, M.J.
Deposited on	:	1998-09-18
Resolution	:	1.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

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

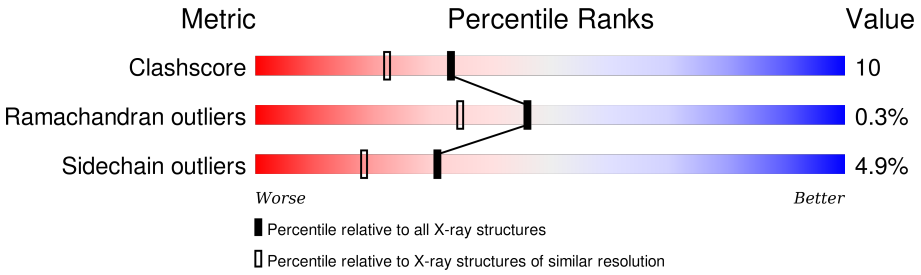
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	723	80% 18% ··

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6301 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.

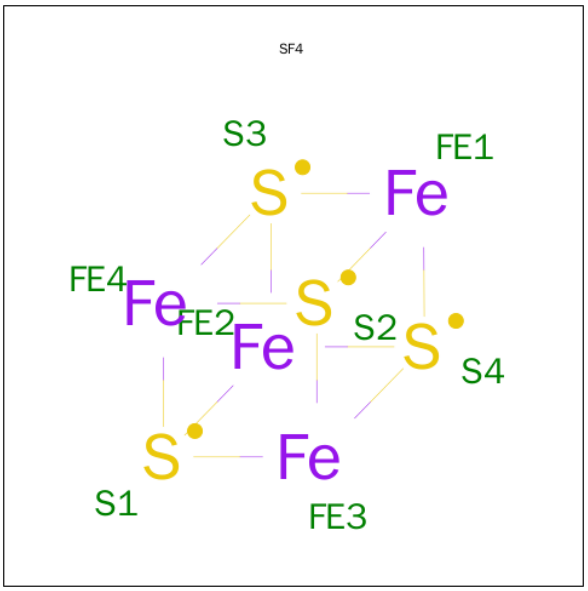
- Molecule 1 is a protein called PROTEIN (PERIPLASMIC NITRATE REDUCTASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	720	Total	C	N	O	S	33	0	0
			5591	3534	1004	1013	40			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	37	ASN	ASP	CONFLICT	GB 38209021

- Molecule 2 is IRON/SULFUR CLUSTER (three-letter code: SF4) (formula: Fe_4S_4).

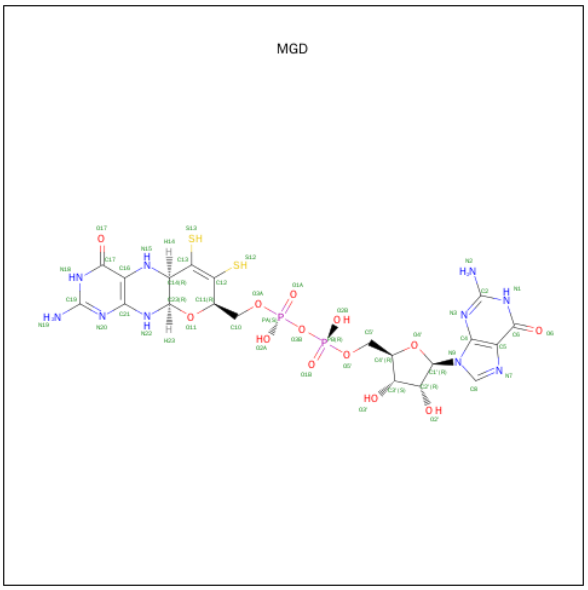


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 3 is MOLYBDENUM ATOM (three-letter code: MO) (formula: Mo).

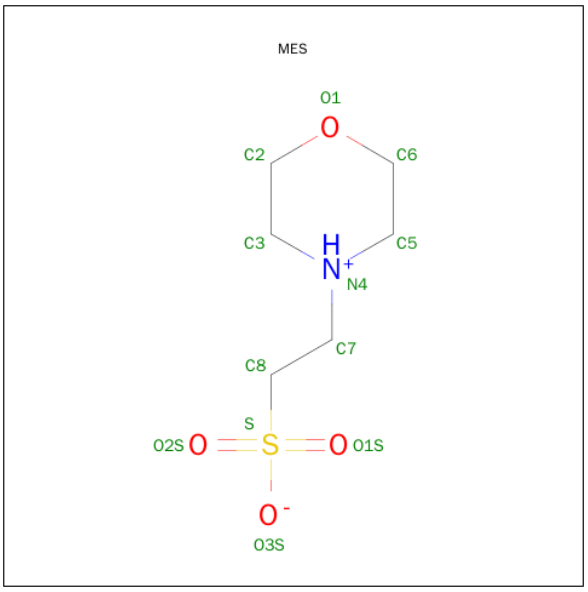
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Mo 1 1	0	0

- Molecule 4 is 2-AMINO-5,6-DIMERCAPTO-7-METHYL-3,7,8A,9-TETRAHYDRO-8-OXA-1,3,9,10-TETRAAZA-ANTHRACEN-4-ONE GUANOSINE DINUCLEOTIDE (three-letter code: MGD) (formula: $C_{20}H_{26}N_{10}O_{13}P_2S_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C N O P S 47 20 10 13 2 2	0	0
4	A	1	Total C N O P S 47 20 10 13 2 2	0	0

- Molecule 5 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	595	Total	O	0	0
			595	595		

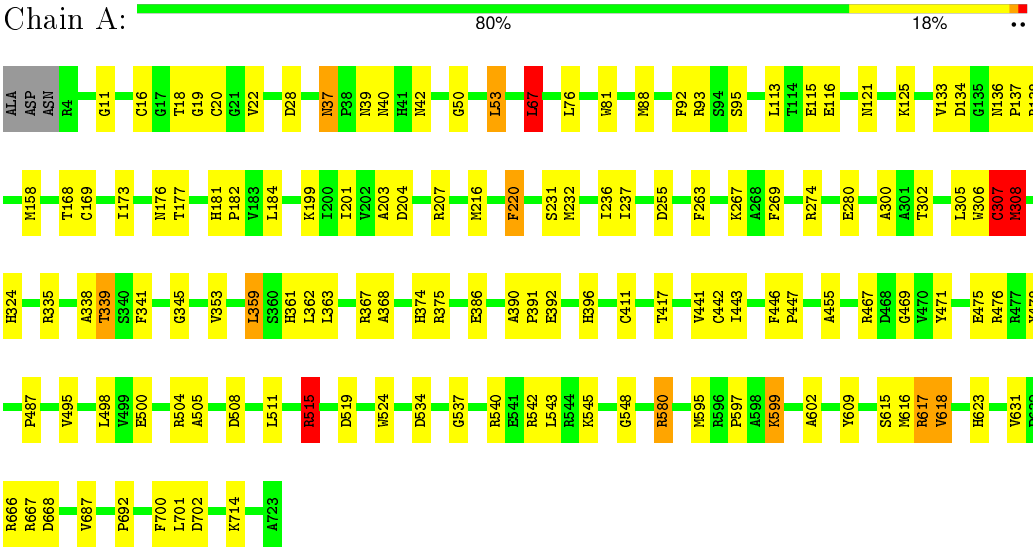
3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

Note EDS was not executed.

- Molecule 1: PROTEIN (PERIPLASMIC NITRATE REDUCTASE)

Chain A:



4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	106.26Å 106.26Å 134.91Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 1.90	Depositor
% Data completeness (in resolution range)	96.1 (20.00-1.90)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.218 , 0.267	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6301	wwPDB-VP
Average B, all atoms (Å^2)	36.0	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: MO, SF4, MES, MGD

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	0.61	2/5737 (0.0%)	0.81	6/7791 (0.1%)

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	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	308	MET	C-N	11.51	1.53	1.33
1	A	386	GLU	CB-CG	-5.29	1.42	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	308	MET	O-C-N	-15.40	97.01	123.20
1	A	307	CYS	O-C-N	-9.39	107.68	122.70
1	A	515	ARG	NE-CZ-NH2	7.31	123.96	120.30
1	A	67	LEU	CA-CB-CG	6.52	130.30	115.30
1	A	308	MET	CG-SD-CE	5.97	109.75	100.20
1	A	204	ASP	CB-CG-OD1	5.79	123.51	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	307	CYS	Mainchain
1	A	308	MET	Mainchain

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5591	0	5470	114	0
2	A	8	0	0	0	0
3	A	1	0	0	0	0
4	A	94	0	42	16	0
5	A	12	0	13	2	0
6	A	595	0	0	18	0
All	All	6301	0	5525	115	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 (115) 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:617:ARG:HH22	4:A:811:MGD:H15	1.14	0.94
1:A:339:THR:HG23	6:A:1256:HOH:O	1.76	0.85
1:A:714:LYS:HE3	4:A:812:MGD:H5'1	1.60	0.84
1:A:368:ALA:H	1:A:374:HIS:HD2	1.32	0.77
1:A:597:PRO:HG2	1:A:599:LYS:NZ	2.01	0.74
1:A:168:THR:HG21	6:A:1224:HOH:O	1.86	0.74
1:A:597:PRO:HG2	1:A:599:LYS:HZ2	1.53	0.73
1:A:140:CYS:SG	1:A:308:MET:HE1	2.31	0.69
1:A:307:CYS:SG	4:A:812:MGD:H101	2.32	0.69
1:A:392:GLU:HB3	6:A:1234:HOH:O	1.94	0.67
1:A:140:CYS:SG	1:A:308:MET:CE	2.83	0.67
1:A:368:ALA:H	1:A:374:HIS:CD2	2.13	0.67
1:A:623:HIS:CE1	4:A:812:MGD:S13	2.90	0.65
1:A:345:GLY:HA3	4:A:812:MGD:C12	2.27	0.64
1:A:134:ASP:OD1	1:A:138:ARG:HD2	1.98	0.64
1:A:617:ARG:HH11	4:A:812:MGD:H15	1.46	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:ARG:HD2	1:A:505:ALA:O	2.02	0.60
1:A:11:GLY:HA2	1:A:481:THR:HA	1.85	0.59
1:A:609:TYR:CE1	1:A:666:ARG:HG3	2.37	0.59
1:A:269:PHE:CG	1:A:595:MET:HE3	2.38	0.58
5:A:2300:MES:H32	6:A:1347:HOH:O	2.02	0.58
1:A:140:CYS:HB2	4:A:811:MGD:S13	2.43	0.58
1:A:18:THR:HG23	1:A:184:LEU:HG	1.85	0.57
1:A:617:ARG:NH2	4:A:811:MGD:H15	1.95	0.57
1:A:442:CYS:SG	1:A:455:ALA:HB2	2.45	0.57
1:A:19:GLY:HA2	6:A:1001:HOH:O	2.05	0.56
1:A:149:VAL:HG11	5:A:2300:MES:H31	1.88	0.56
1:A:345:GLY:HA3	4:A:812:MGD:S12	2.47	0.55
1:A:203:ALA:HB1	1:A:220:PHE:CE2	2.42	0.55
1:A:269:PHE:HB2	1:A:595:MET:CE	2.36	0.55
1:A:88:MET:HG3	1:A:441:VAL:HG21	1.88	0.54
1:A:615:SER:HA	1:A:687:VAL:O	2.07	0.54
1:A:207:ARG:HG3	1:A:207:ARG:HH11	1.71	0.54
1:A:201:ILE:HD12	1:A:216:MET:HE3	1.91	0.53
1:A:138:ARG:HH22	1:A:361:HIS:CE1	2.28	0.52
1:A:207:ARG:NH1	6:A:1564:HOH:O	2.43	0.52
1:A:255:ASP:HA	1:A:595:MET:HE2	1.92	0.52
1:A:269:PHE:HB2	1:A:595:MET:HE1	1.90	0.52
1:A:269:PHE:CB	1:A:595:MET:HE3	2.39	0.52
1:A:617:ARG:HD3	4:A:812:MGD:N15	2.25	0.51
1:A:308:MET:O	1:A:714:LYS:HG3	2.10	0.51
1:A:314:VAL:HG11	1:A:602:ALA:HB2	1.93	0.51
1:A:300:ALA:O	1:A:335:ARG:NH2	2.43	0.51
1:A:136:ASN:HD22	1:A:417:THR:HG23	1.76	0.51
1:A:618:VAL:HG13	6:A:1441:HOH:O	2.10	0.50
1:A:617:ARG:HD3	4:A:812:MGD:C16	2.42	0.50
1:A:121:ASN:HD21	1:A:125:LYS:NZ	2.10	0.50
1:A:714:LYS:CE	4:A:812:MGD:H5'1	2.36	0.50
1:A:500:GLU:OE1	1:A:504:ARG:HD2	2.11	0.50
1:A:140:CYS:O	1:A:308:MET:HE1	2.13	0.49
1:A:475:GLU:O	1:A:476:ARG:HB2	2.13	0.49
1:A:443:ILE:HG22	4:A:811:MGD:HN21	1.77	0.48
1:A:633:GLU:HB3	6:A:1487:HOH:O	2.13	0.48
1:A:37:ASN:ND2	6:A:1070:HOH:O	2.39	0.48
1:A:263:PHE:CE2	1:A:267:LYS:HE2	2.49	0.48
1:A:617:ARG:NH1	1:A:623:HIS:HE1	2.12	0.48
1:A:137:PRO:HG2	1:A:353:VAL:HG21	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:121:ASN:HD22	1:A:133:VAL:HG11	1.79	0.47
1:A:67:LEU:HB3	1:A:76:LEU:HG	1.97	0.47
1:A:478:TYR:O	1:A:548:GLY:HA2	2.14	0.47
1:A:362:LEU:HD22	6:A:1238:HOH:O	2.15	0.47
1:A:623:HIS:HE1	4:A:812:MGD:S13	2.36	0.47
1:A:115:GLU:OE2	1:A:540:ARG:NH2	2.46	0.46
1:A:537:GLY:O	1:A:542:ARG:HD3	2.15	0.46
1:A:138:ARG:HD3	1:A:396:HIS:HB2	1.97	0.46
1:A:476:ARG:HA	6:A:1039:HOH:O	2.15	0.46
1:A:302:THR:O	1:A:338:ALA:HA	2.15	0.46
1:A:168:THR:HG22	6:A:1473:HOH:O	2.16	0.46
1:A:20:CYS:HB3	1:A:42:ASN:ND2	2.30	0.46
1:A:37:ASN:ND2	1:A:39:ASN:H	2.14	0.46
1:A:181:HIS:N	1:A:182:PRO:HD3	2.31	0.45
1:A:482:GLU:HG3	6:A:1265:HOH:O	2.15	0.45
1:A:542:ARG:HA	1:A:545:LYS:HE3	1.98	0.45
1:A:113:LEU:HD22	1:A:469:GLY:HA3	1.98	0.45
1:A:232:MET:O	1:A:236:ILE:HG13	2.17	0.45
1:A:515:ARG:HE	1:A:515:ARG:HB2	1.48	0.45
1:A:237:ILE:HG21	1:A:267:LYS:HG2	1.98	0.45
1:A:390:ALA:HA	1:A:391:PRO:HD3	1.86	0.44
1:A:113:LEU:HB2	1:A:116:GLU:HG3	1.99	0.44
1:A:269:PHE:CB	1:A:595:MET:CE	2.95	0.44
1:A:168:THR:O	1:A:168:THR:HG22	2.19	0.43
1:A:169:CYS:HA	1:A:199:LYS:O	2.18	0.43
1:A:483:LYS:NZ	1:A:487:PRO:HD3	2.33	0.43
1:A:692:PRO:HG3	6:A:1054:HOH:O	2.18	0.43
1:A:136:ASN:ND2	1:A:417:THR:HG23	2.33	0.43
1:A:345:GLY:CA	4:A:812:MGD:C12	2.96	0.42
1:A:40:ASN:HB3	1:A:42:ASN:OD1	2.20	0.42
1:A:500:GLU:HG3	1:A:504:ARG:HD2	2.00	0.42
1:A:515:ARG:NE	1:A:519:ASP:OD2	2.52	0.42
1:A:597:PRO:HG2	1:A:599:LYS:HZ1	1.81	0.42
1:A:631:VAL:HG12	1:A:633:GLU:HG2	2.02	0.42
1:A:540:ARG:HA	1:A:540:ARG:HD3	1.73	0.42
1:A:173:ILE:O	1:A:306:TRP:HA	2.19	0.42
1:A:353:VAL:CG1	1:A:359:LEU:HD13	2.49	0.42
1:A:701:LEU:HD12	1:A:702:ASP:H	1.84	0.42
1:A:615:SER:HB3	1:A:700:PHE:CZ	2.54	0.42
1:A:116:GLU:HG2	1:A:495:VAL:HG23	2.02	0.42
1:A:508:ASP:O	1:A:511:LEU:HB2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:375:ARG:HD2	6:A:1270:HOH:O	2.19	0.42
1:A:140:CYS:SG	1:A:308:MET:HE3	2.60	0.41
1:A:363:LEU:HB2	1:A:367:ARG:HB2	2.02	0.41
1:A:305:LEU:HA	1:A:341:PHE:O	2.21	0.41
1:A:446:PHE:HA	1:A:447:PRO:HD2	1.77	0.41
1:A:580:ARG:HD3	6:A:1258:HOH:O	2.20	0.41
1:A:136:ASN:OD1	1:A:137:PRO:HD3	2.19	0.41
1:A:417:THR:CG2	6:A:1042:HOH:O	2.69	0.41
1:A:443:ILE:C	4:A:811:MGD:HN22	2.21	0.41
1:A:92:PHE:CE2	1:A:411:CYS:HB3	2.55	0.41
1:A:320:ASN:O	1:A:324:HIS:CD2	2.73	0.41
1:A:274:ARG:HD3	6:A:1361:HOH:O	2.20	0.41
1:A:203:ALA:HB1	1:A:220:PHE:HE2	1.84	0.41
1:A:367:ARG:HA	1:A:374:HIS:CD2	2.55	0.40
1:A:50:GLY:HA2	1:A:53:LEU:HD22	2.03	0.40
1:A:136:ASN:N	1:A:137:PRO:CD	2.85	0.40
1:A:667:ARG:O	1:A:668:ASP:HB2	2.22	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	718/723 (99%)	680 (95%)	36 (5%)	2 (0%)	46 35

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	28	ASP
1	A	534	ASP

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	586/596 (98%)	557 (95%)	29 (5%)	31 18

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	CYS
1	A	22	VAL
1	A	37	ASN
1	A	53	LEU
1	A	67	LEU
1	A	81	TRP
1	A	95	SER
1	A	158	MET
1	A	176	ASN
1	A	177	THR
1	A	220	PHE
1	A	231	SER
1	A	280	GLU
1	A	307	CYS
1	A	339	THR
1	A	359	LEU
1	A	467	ARG
1	A	471	TYR
1	A	498	LEU
1	A	515	ARG
1	A	524	TRP
1	A	543	LEU
1	A	580	ARG
1	A	599	LYS
1	A	616	MET
1	A	617	ARG
1	A	618	VAL
1	A	633	GLU
1	A	660	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	ASN
1	A	121	ASN
1	A	132	ASN
1	A	176	ASN
1	A	238	ASN
1	A	361	HIS
1	A	374	HIS
1	A	496	ASN
1	A	623	HIS
1	A	646	ASN
1	A	657	HIS

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 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	$\#\|Z\| > 2$
5	MES	A	2300	-	11,12,12	3.23	3 (27%)	14,16,16	2.43	4 (28%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
2	SF4	A	800	1	0,12,12	0.00	-	0,24,24	0.00	-
4	MGD	A	811	3	38,52,52	2.30	8 (21%)	43,81,81	3.73	16 (37%)
4	MGD	A	812	3	38,52,52	2.11	8 (21%)	43,81,81	3.10	21 (48%)

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
5	MES	A	2300	-	-	0/6/14/14	0/1/1/1
2	SF4	A	800	1	-	0/0/48/48	2/6/5/5
4	MGD	A	811	3	-	0/18/66/66	0/6/6/6
4	MGD	A	812	3	-	0/18/66/66	0/6/6/6

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	811	MGD	O11-C23	-9.85	1.29	1.43
4	A	812	MGD	O11-C23	-8.55	1.31	1.43
4	A	811	MGD	O17-C17	-3.05	1.17	1.24
4	A	811	MGD	C2-N2	-2.80	1.28	1.34
4	A	812	MGD	C19-N20	-2.69	1.30	1.35
4	A	812	MGD	O17-C17	-2.51	1.18	1.24
4	A	811	MGD	C17-N18	2.11	1.37	1.33
4	A	812	MGD	C16-C21	2.20	1.46	1.41
4	A	812	MGD	C6-N1	2.27	1.37	1.33
4	A	812	MGD	C2-N1	2.31	1.39	1.35
4	A	811	MGD	C6-C5	2.48	1.46	1.41
4	A	811	MGD	O4'-C1'	3.19	1.45	1.41
4	A	812	MGD	C17-N18	3.32	1.39	1.33
4	A	811	MGD	C23-C14	3.74	1.56	1.53
5	A	2300	MES	O2S-S	3.89	1.57	1.45
5	A	2300	MES	O1S-S	4.01	1.57	1.45
4	A	812	MGD	C21-N22	4.27	1.43	1.35
4	A	811	MGD	C21-N22	4.74	1.43	1.35
5	A	2300	MES	O3S-S	8.30	1.67	1.46

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	812	MGD	C4'-O4'-C1'	-9.03	99.79	109.72
4	A	811	MGD	C4'-O4'-C1'	-7.47	101.51	109.72
4	A	811	MGD	C5-C6-N1	-6.10	115.25	123.59
4	A	812	MGD	C5-C6-N1	-5.45	116.14	123.59
4	A	811	MGD	O4'-C1'-N9	-4.41	98.86	108.10
4	A	811	MGD	N18-C19-N20	-4.15	118.73	125.53
4	A	812	MGD	O3B-PB-O5'	-3.74	93.00	102.94
4	A	812	MGD	N18-C19-N20	-3.69	119.48	125.53
4	A	811	MGD	C6-C5-C4	-3.41	116.83	120.90
4	A	812	MGD	N3-C2-N1	-2.72	123.30	127.44
4	A	812	MGD	C6-C5-C4	-2.60	117.79	120.90
5	A	2300	MES	O3S-S-O2S	-2.54	105.71	111.61
4	A	812	MGD	C21-N22-C23	-2.45	118.86	123.67
4	A	812	MGD	O4'-C4'-C5'	-2.42	100.67	109.32
4	A	812	MGD	C2'-C3'-C4'	-2.40	97.68	102.61
4	A	811	MGD	O3B-PA-O3A	-2.20	97.09	102.94
4	A	811	MGD	O4'-C4'-C5'	-2.12	101.75	109.32
4	A	812	MGD	O4'-C1'-N9	-2.08	103.75	108.10
4	A	812	MGD	C16-C21-N22	2.39	120.84	118.34
5	A	2300	MES	C5-N4-C3	2.40	114.10	108.90
4	A	812	MGD	C19-N20-C21	2.59	120.36	114.54
4	A	812	MGD	O2'-C2'-C3'	2.63	120.39	111.83
4	A	811	MGD	C19-N20-C21	2.64	120.48	114.54
4	A	812	MGD	N19-C19-N20	2.71	121.69	117.20
4	A	812	MGD	C4-C5-N7	2.74	112.00	109.48
4	A	811	MGD	O2'-C2'-C3'	2.89	121.24	111.83
4	A	811	MGD	O4'-C4'-C3'	3.00	111.20	105.15
4	A	812	MGD	C17-C16-C21	3.04	117.32	114.56
4	A	812	MGD	O4'-C4'-C3'	3.43	112.06	105.15
4	A	811	MGD	C17-C16-C21	3.86	118.05	114.56
4	A	811	MGD	N19-C19-N20	3.86	123.59	117.20
4	A	812	MGD	O11-C23-C14	4.34	111.93	108.96
5	A	2300	MES	O1S-S-C8	4.44	110.69	106.91
4	A	812	MGD	C6-N1-C2	5.06	122.96	115.94
4	A	811	MGD	C6-N1-C2	5.26	123.24	115.94
4	A	811	MGD	C17-N18-C19	5.54	123.63	115.94
4	A	812	MGD	C17-N18-C19	6.27	124.64	115.94
5	A	2300	MES	O2S-S-C8	6.29	112.27	106.91
4	A	812	MGD	C2'-C1'-N9	8.80	127.74	114.29
4	A	811	MGD	C2'-C1'-N9	11.88	132.44	114.29
4	A	811	MGD	O11-C23-C14	13.04	117.88	108.96

There are no chirality outliers.

There are no torsion outliers.

All (2) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	800	SF4	FE1-FE2-S3-S4
2	A	800	SF4	FE3-FE4-S1-S2

3 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	2300	MES	2	0
4	A	811	MGD	5	0
4	A	812	MGD	11	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

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.