



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 02:45 pm GMT

PDB ID : 4DMM
Title : 3-oxoacyl-[acyl-carrier-protein] reductase from *Synechococcus elongatus* PCC 7942 in complex with NADP
Authors : Chen, C.; Zhuang, N.N.; Lee, K.H.
Deposited on : 2012-02-08
Resolution : 2.38 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 ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

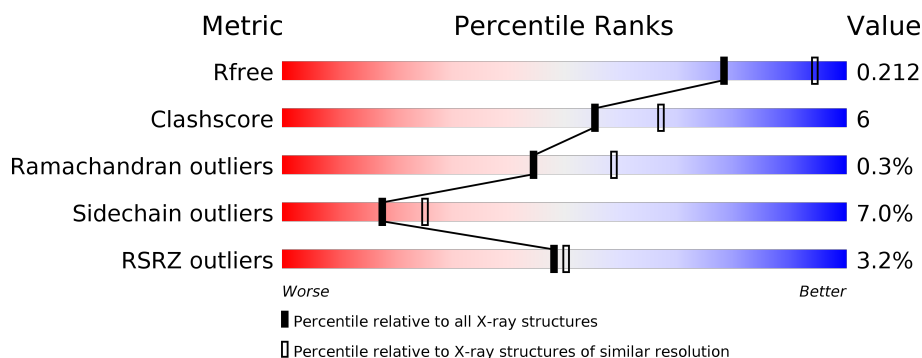
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.38 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}	100719	4388 (2.40-2.36)
Clashscore	112137	4984 (2.40-2.36)
Ramachandran outliers	110173	4907 (2.40-2.36)
Sidechain outliers	110143	4909 (2.40-2.36)
RSRZ outliers	101464	4423 (2.40-2.36)

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 $\leq 5\%$. 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	269	<div> <div>3%</div> <div> <div></div> <div>75%</div> <div>15%</div> <div>•</div> <div>9%</div> </div> </div>
1	B	269	<div> <div>2%</div> <div> <div></div> <div>78%</div> <div>10%</div> <div>•</div> <div>11%</div> </div> </div>
1	C	269	<div> <div>2%</div> <div> <div></div> <div>76%</div> <div>12%</div> <div>•</div> <div>11%</div> </div> </div>
1	D	269	<div> <div>4%</div> <div> <div></div> <div>71%</div> <div>13%</div> <div>•</div> <div>12%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7259 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 3-oxoacyl-[acyl-carrier-protein] reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	246	Total	C	N	O	S	0	0	0
			1768	1113	313	336	6			
1	B	240	Total	C	N	O	S	0	0	0
			1723	1085	306	326	6			
1	C	240	Total	C	N	O	S	0	0	0
			1728	1089	307	326	6			
1	D	237	Total	C	N	O	S	0	0	0
			1703	1075	304	319	5			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP Q31QF3
A	-18	GLY	-	EXPRESSION TAG	UNP Q31QF3
A	-17	SER	-	EXPRESSION TAG	UNP Q31QF3
A	-16	SER	-	EXPRESSION TAG	UNP Q31QF3
A	-15	HIS	-	EXPRESSION TAG	UNP Q31QF3
A	-14	HIS	-	EXPRESSION TAG	UNP Q31QF3
A	-13	HIS	-	EXPRESSION TAG	UNP Q31QF3
A	-12	HIS	-	EXPRESSION TAG	UNP Q31QF3
A	-11	HIS	-	EXPRESSION TAG	UNP Q31QF3
A	-10	HIS	-	EXPRESSION TAG	UNP Q31QF3
A	-9	SER	-	EXPRESSION TAG	UNP Q31QF3
A	-8	SER	-	EXPRESSION TAG	UNP Q31QF3
A	-7	GLY	-	EXPRESSION TAG	UNP Q31QF3
A	-6	LEU	-	EXPRESSION TAG	UNP Q31QF3
A	-5	VAL	-	EXPRESSION TAG	UNP Q31QF3
A	-4	PRO	-	EXPRESSION TAG	UNP Q31QF3
A	-3	ARG	-	EXPRESSION TAG	UNP Q31QF3
A	-2	GLY	-	EXPRESSION TAG	UNP Q31QF3
A	-1	SER	-	EXPRESSION TAG	UNP Q31QF3
A	0	HIS	-	EXPRESSION TAG	UNP Q31QF3
B	-19	MET	-	EXPRESSION TAG	UNP Q31QF3

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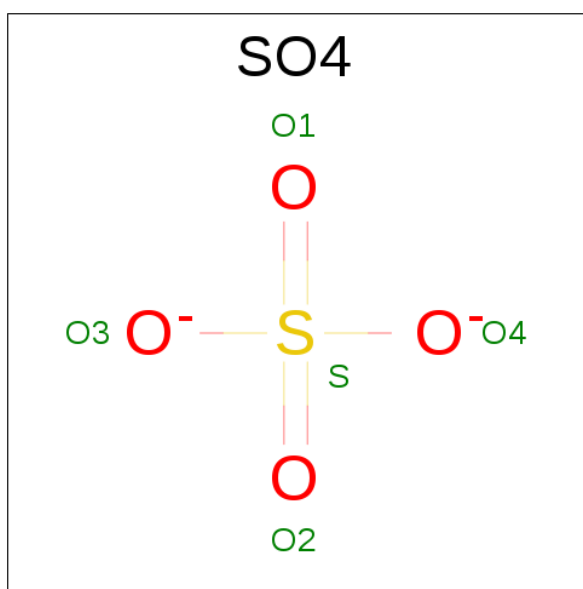
Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	GLY	-	EXPRESSION TAG	UNP Q31QF3
B	-17	SER	-	EXPRESSION TAG	UNP Q31QF3
B	-16	SER	-	EXPRESSION TAG	UNP Q31QF3
B	-15	HIS	-	EXPRESSION TAG	UNP Q31QF3
B	-14	HIS	-	EXPRESSION TAG	UNP Q31QF3
B	-13	HIS	-	EXPRESSION TAG	UNP Q31QF3
B	-12	HIS	-	EXPRESSION TAG	UNP Q31QF3
B	-11	HIS	-	EXPRESSION TAG	UNP Q31QF3
B	-10	HIS	-	EXPRESSION TAG	UNP Q31QF3
B	-9	SER	-	EXPRESSION TAG	UNP Q31QF3
B	-8	SER	-	EXPRESSION TAG	UNP Q31QF3
B	-7	GLY	-	EXPRESSION TAG	UNP Q31QF3
B	-6	LEU	-	EXPRESSION TAG	UNP Q31QF3
B	-5	VAL	-	EXPRESSION TAG	UNP Q31QF3
B	-4	PRO	-	EXPRESSION TAG	UNP Q31QF3
B	-3	ARG	-	EXPRESSION TAG	UNP Q31QF3
B	-2	GLY	-	EXPRESSION TAG	UNP Q31QF3
B	-1	SER	-	EXPRESSION TAG	UNP Q31QF3
B	0	HIS	-	EXPRESSION TAG	UNP Q31QF3
C	-19	MET	-	EXPRESSION TAG	UNP Q31QF3
C	-18	GLY	-	EXPRESSION TAG	UNP Q31QF3
C	-17	SER	-	EXPRESSION TAG	UNP Q31QF3
C	-16	SER	-	EXPRESSION TAG	UNP Q31QF3
C	-15	HIS	-	EXPRESSION TAG	UNP Q31QF3
C	-14	HIS	-	EXPRESSION TAG	UNP Q31QF3
C	-13	HIS	-	EXPRESSION TAG	UNP Q31QF3
C	-12	HIS	-	EXPRESSION TAG	UNP Q31QF3
C	-11	HIS	-	EXPRESSION TAG	UNP Q31QF3
C	-10	HIS	-	EXPRESSION TAG	UNP Q31QF3
C	-9	SER	-	EXPRESSION TAG	UNP Q31QF3
C	-8	SER	-	EXPRESSION TAG	UNP Q31QF3
C	-7	GLY	-	EXPRESSION TAG	UNP Q31QF3
C	-6	LEU	-	EXPRESSION TAG	UNP Q31QF3
C	-5	VAL	-	EXPRESSION TAG	UNP Q31QF3
C	-4	PRO	-	EXPRESSION TAG	UNP Q31QF3
C	-3	ARG	-	EXPRESSION TAG	UNP Q31QF3
C	-2	GLY	-	EXPRESSION TAG	UNP Q31QF3
C	-1	SER	-	EXPRESSION TAG	UNP Q31QF3
C	0	HIS	-	EXPRESSION TAG	UNP Q31QF3
D	-19	MET	-	EXPRESSION TAG	UNP Q31QF3
D	-18	GLY	-	EXPRESSION TAG	UNP Q31QF3
D	-17	SER	-	EXPRESSION TAG	UNP Q31QF3

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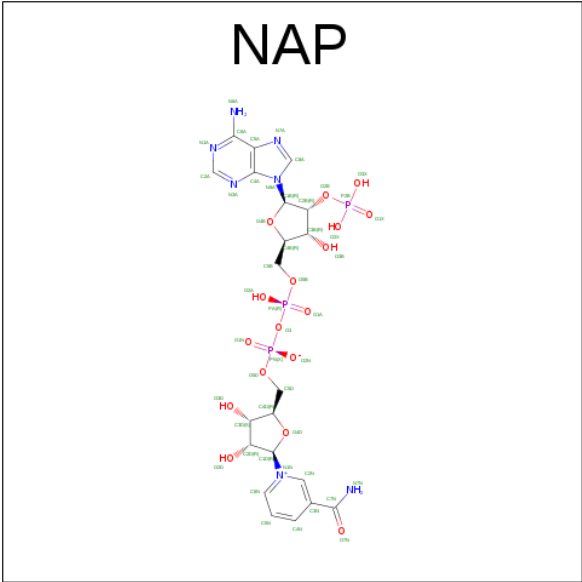
Chain	Residue	Modelled	Actual	Comment	Reference
D	-16	SER	-	EXPRESSION TAG	UNP Q31QF3
D	-15	HIS	-	EXPRESSION TAG	UNP Q31QF3
D	-14	HIS	-	EXPRESSION TAG	UNP Q31QF3
D	-13	HIS	-	EXPRESSION TAG	UNP Q31QF3
D	-12	HIS	-	EXPRESSION TAG	UNP Q31QF3
D	-11	HIS	-	EXPRESSION TAG	UNP Q31QF3
D	-10	HIS	-	EXPRESSION TAG	UNP Q31QF3
D	-9	SER	-	EXPRESSION TAG	UNP Q31QF3
D	-8	SER	-	EXPRESSION TAG	UNP Q31QF3
D	-7	GLY	-	EXPRESSION TAG	UNP Q31QF3
D	-6	LEU	-	EXPRESSION TAG	UNP Q31QF3
D	-5	VAL	-	EXPRESSION TAG	UNP Q31QF3
D	-4	PRO	-	EXPRESSION TAG	UNP Q31QF3
D	-3	ARG	-	EXPRESSION TAG	UNP Q31QF3
D	-2	GLY	-	EXPRESSION TAG	UNP Q31QF3
D	-1	SER	-	EXPRESSION TAG	UNP Q31QF3
D	0	HIS	-	EXPRESSION TAG	UNP Q31QF3

- Molecule 2 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



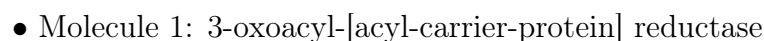
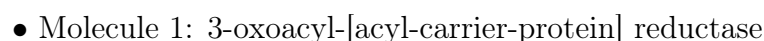
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	C	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	D	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

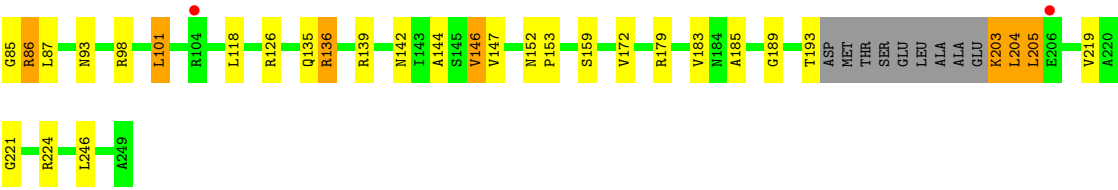
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	65	Total	O	0	0
			65	65		
4	B	58	Total	O	0	0
			58	58		
4	C	38	Total	O	0	0
			38	38		
4	D	27	Total	O	0	0
			27	27		

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- Molecule 1: 3-oxoacyl-[acyl-carrier-protein] reductase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	63.98Å 116.92Å 125.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.78 – 2.38 42.78 – 2.38	Depositor EDS
% Data completeness (in resolution range)	99.8 (42.78-2.38) 99.8 (42.78-2.38)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.88 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.164 , 0.216 0.161 , 0.212	Depositor DCC
R_{free} test set	1914 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	30.1	Xtriage
Anisotropy	0.108	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 35.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7259	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.46% 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: NAP, SO4

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	1.06	2/1785 (0.1%)	1.00	4/2421 (0.2%)
1	B	1.07	1/1739 (0.1%)	0.95	0/2358
1	C	1.02	1/1744 (0.1%)	0.87	0/2363
1	D	1.01	1/1719 (0.1%)	0.92	4/2330 (0.2%)
All	All	1.04	5/6987 (0.1%)	0.94	8/9472 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	74	GLU	CG-CD	6.39	1.61	1.51
1	A	78	ALA	CA-CB	6.24	1.65	1.52
1	C	51	ALA	CA-CB	6.15	1.65	1.52
1	B	94	ALA	CA-CB	5.48	1.64	1.52
1	D	48	GLU	CG-CD	5.20	1.59	1.51

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	126	ARG	NE-CZ-NH2	-6.85	116.88	120.30
1	A	126	ARG	NE-CZ-NH1	6.60	123.60	120.30
1	A	67	VAL	CB-CA-C	5.52	121.89	111.40
1	D	139	ARG	NE-CZ-NH1	5.30	122.95	120.30
1	D	66	ASP	CB-CG-OD1	5.18	122.97	118.30
1	D	126	ARG	NE-CZ-NH2	-5.18	117.71	120.30
1	A	224	ARG	NE-CZ-NH2	-5.09	117.75	120.30
1	D	126	ARG	NE-CZ-NH1	5.06	122.83	120.30

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	1768	0	1832	23	0
1	B	1723	0	1785	19	0
1	C	1728	0	1792	24	0
1	D	1703	0	1773	23	0
2	A	5	0	0	0	0
3	B	48	0	25	4	0
3	C	48	0	25	7	0
3	D	48	0	25	4	0
4	A	65	0	0	1	0
4	B	58	0	0	0	0
4	C	38	0	0	0	0
4	D	27	0	0	0	0
All	All	7259	0	7257	84	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 6.

All (84) 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:B:87:LEU:H	1:B:135:GLN:HE22	1.19	0.87
1:C:87:LEU:H	1:C:135:GLN:HE22	1.20	0.85
1:C:210:LEU:HD12	1:C:244:GLY:HA2	1.59	0.84
1:A:87:LEU:H	1:A:135:GLN:HE22	1.27	0.80
1:C:107:ARG:HH22	1:C:111:GLN:HG2	1.54	0.72
1:B:15:GLY:H	1:B:38:ASN:HD22	1.40	0.69
1:A:199:LEU:H	1:A:199:LEU:HD22	1.59	0.68
1:C:20:ILE:HB	3:C:301:NAP:H51N	1.74	0.67
1:D:15:GLY:H	1:D:38:ASN:HD22	1.43	0.67
1:A:15:GLY:H	1:A:38:ASN:HD22	1.46	0.63
1:C:15:GLY:H	1:C:38:ASN:HD22	1.46	0.62
1:A:92:ASN:HD21	1:A:124:CYS:HB2	1.65	0.62
1:B:189:GLY:O	3:B:301:NAP:H4N	2.00	0.62
1:C:193:THR:HG23	1:C:194:ASP:N	2.18	0.58
1:D:203:LYS:O	1:D:205:LEU:N	2.35	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:87:LEU:H	1:D:135:GLN:HE22	1.52	0.57
1:C:7:THR:O	1:C:8:ASP:HB2	2.04	0.56
1:B:204:LEU:N	1:B:205:LEU:HB3	2.20	0.56
1:C:15:GLY:HA2	3:C:301:NAP:H1B	1.88	0.53
1:D:69:GLN:HB2	1:D:72:GLU:HB2	1.91	0.53
1:C:146:VAL:HG21	1:C:246:LEU:HD23	1.91	0.53
1:D:189:GLY:O	3:D:301:NAP:H4N	2.08	0.53
1:A:130:LYS:HG3	1:B:103:LEU:HD23	1.92	0.52
1:D:142:ASN:O	1:D:185:ALA:HA	2.09	0.52
1:D:7:THR:O	1:D:8:ASP:HB2	2.09	0.52
1:A:146:VAL:HG21	1:A:246:LEU:HD23	1.92	0.51
1:B:14:THR:O	1:B:93:ASN:HB3	2.11	0.51
1:C:189:GLY:O	3:C:301:NAP:H4N	2.11	0.50
1:D:147:VAL:HG12	1:D:159:SER:HB3	1.93	0.50
1:D:172:VAL:HG12	1:D:183:VAL:HG21	1.94	0.50
1:D:85:GLY:O	1:D:86:ARG:HB3	2.13	0.49
1:A:199:LEU:CD2	1:A:199:LEU:H	2.25	0.49
1:A:201:ALA:HB1	1:A:213:TYR:CE1	2.49	0.48
1:B:210:LEU:HD12	1:B:244:GLY:HA2	1.96	0.48
1:D:14:THR:HA	1:D:38:ASN:HB3	1.95	0.48
1:C:155:GLN:HB3	1:C:158:TYR:HB3	1.96	0.48
1:A:147:VAL:HG12	1:A:159:SER:HB3	1.96	0.47
1:C:14:THR:O	1:C:93:ASN:HB3	2.15	0.47
1:D:146:VAL:HG11	1:D:246:LEU:HD23	1.96	0.47
1:B:38:ASN:HD21	3:B:301:NAP:H2A	1.79	0.47
1:D:42:SER:OG	1:D:45:ALA:HB3	2.14	0.47
1:A:15:GLY:H	1:A:38:ASN:ND2	2.10	0.47
1:A:190:PHE:HB3	1:A:195:MET:SD	2.55	0.46
1:A:195:MET:HG2	1:A:198:GLU:HG3	1.97	0.46
1:A:130:LYS:HG2	1:B:103:LEU:CD2	2.46	0.45
1:A:7:THR:O	1:A:8:ASP:HB2	2.16	0.45
1:C:144:ALA:O	3:C:301:NAP:H6N	2.16	0.45
1:B:144:ALA:O	3:B:301:NAP:H6N	2.17	0.45
1:C:15:GLY:CA	3:C:301:NAP:H1B	2.47	0.45
1:C:193:THR:HG21	3:C:301:NAP:O1N	2.17	0.45
1:D:66:ASP:OD2	1:D:69:GLN:HG2	2.17	0.45
1:A:179:ARG:NE	4:A:461:HOH:O	2.50	0.45
1:B:90:LEU:HD22	1:B:128:ALA:HB2	1.99	0.45
1:A:207:VAL:O	1:A:209:PRO:HD3	2.17	0.45
1:C:38:ASN:HD21	3:C:301:NAP:H2A	1.82	0.44
1:C:107:ARG:NH2	1:C:111:GLN:HG2	2.27	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:ALA:O	1:A:64:LYS:HE3	2.18	0.44
1:A:92:ASN:HD21	1:A:124:CYS:CB	2.31	0.44
1:C:136:ARG:HG2	1:C:136:ARG:HH11	1.83	0.44
1:A:91:VAL:HA	1:A:141:ILE:O	2.18	0.44
1:D:93:ASN:OD1	3:D:301:NAP:H4D	2.17	0.44
1:B:14:THR:HA	1:B:38:ASN:HB3	2.00	0.43
1:C:92:ASN:HD21	1:C:124:CYS:HB2	1.83	0.43
1:C:130:LYS:HE2	1:C:130:LYS:HB2	1.68	0.43
1:B:231:ALA:HB2	1:C:221:GLY:HA3	2.01	0.43
1:C:136:ARG:HH11	1:C:136:ARG:CG	2.32	0.43
1:D:101:LEU:HD12	1:D:101:LEU:HA	1.88	0.42
1:B:146:VAL:HG21	1:B:246:LEU:HD23	2.01	0.42
1:D:15:GLY:HA2	3:D:301:NAP:H1B	2.01	0.42
1:A:231:ALA:HB2	1:D:221:GLY:HA3	2.01	0.42
1:A:130:LYS:CG	1:B:103:LEU:CD2	2.98	0.41
1:B:15:GLY:HA2	3:B:301:NAP:H1B	2.02	0.41
1:A:4:LEU:HD23	1:A:4:LEU:HA	1.93	0.41
1:B:20:ILE:HG13	1:B:193:THR:HG21	2.03	0.41
1:D:203:LYS:HG2	1:D:204:LEU:HB2	2.02	0.41
1:B:204:LEU:N	1:B:204:LEU:HD13	2.35	0.41
1:D:136:ARG:HD3	1:D:179:ARG:HD2	2.02	0.41
1:C:142:ASN:O	1:C:185:ALA:HA	2.21	0.41
1:C:106:LYS:HA	1:C:106:LYS:HD3	1.90	0.41
1:D:86:ARG:HB2	1:D:135:GLN:NE2	2.36	0.41
1:A:146:VAL:HB	1:A:247:VAL:HG21	2.02	0.40
1:D:144:ALA:O	3:D:301:NAP:H6N	2.21	0.40
1:D:152:ASN:HA	1:D:153:PRO:HD3	1.98	0.40
1:B:99:ASP:OD1	1:B:155:GLN:HG2	2.21	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	244/269 (91%)	233 (96%)	11 (4%)	0	100	100
1	B	236/269 (88%)	224 (95%)	11 (5%)	1 (0%)	38	51
1	C	236/269 (88%)	222 (94%)	13 (6%)	1 (0%)	38	51
1	D	233/269 (87%)	218 (94%)	14 (6%)	1 (0%)	38	51
All	All	949/1076 (88%)	897 (94%)	49 (5%)	3 (0%)	44	59

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	204	LEU
1	B	205	LEU
1	C	194	ASP

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	172/191 (90%)	159 (92%)	13 (8%)	15	22
1	B	168/191 (88%)	160 (95%)	8 (5%)	30	44
1	C	168/191 (88%)	158 (94%)	10 (6%)	22	33
1	D	165/191 (86%)	149 (90%)	16 (10%)	9	12
All	All	673/764 (88%)	626 (93%)	47 (7%)	18	25

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	ARG
1	A	26	LEU
1	A	28	LEU
1	A	47	ASP
1	A	67	VAL
1	A	83	ARG
1	A	101	LEU
1	A	118	LEU

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Mol	Chain	Res	Type
1	A	136	ARG
1	A	198	GLU
1	A	199	LEU
1	A	202	GLU
1	A	204	LEU
1	B	61	PHE
1	B	73	VAL
1	B	82	GLU
1	B	103	LEU
1	B	104	ARG
1	B	107	ARG
1	B	196	THR
1	B	204	LEU
1	C	26	LEU
1	C	47	ASP
1	C	107	ARG
1	C	136	ARG
1	C	193	THR
1	C	194	ASP
1	C	195	MET
1	C	203	LYS
1	C	204	LEU
1	C	205	LEU
1	D	7	THR
1	D	17	SER
1	D	42	SER
1	D	64	LYS
1	D	74	GLU
1	D	86	ARG
1	D	98	ARG
1	D	101	LEU
1	D	118	LEU
1	D	136	ARG
1	D	146	VAL
1	D	193	THR
1	D	203	LYS
1	D	205	LEU
1	D	219	VAL
1	D	224	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	ASN
1	A	92	ASN
1	A	111	GLN
1	A	135	GLN
1	A	241	ASN
1	B	38	ASN
1	B	111	GLN
1	B	135	GLN
1	C	38	ASN
1	C	92	ASN
1	C	135	GLN
1	D	38	ASN
1	D	135	GLN
1	D	241	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

4 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	SO4	A	301	-	4,4,4	0.18	0	6,6,6	0.68	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAP	B	301	-	44,52,52	1.64	9 (20%)	51,80,80	2.14	12 (23%)
3	NAP	C	301	-	44,52,52	1.65	8 (18%)	51,80,80	2.03	9 (17%)
3	NAP	D	301	-	44,52,52	1.66	12 (27%)	51,80,80	2.10	9 (17%)

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	SO4	A	301	-	-	0/0/0/0	0/0/0/0
3	NAP	B	301	-	-	0/27/67/67	0/5/5/5
3	NAP	C	301	-	-	0/27/67/67	0/5/5/5
3	NAP	D	301	-	-	0/27/67/67	0/5/5/5

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	301	NAP	P2B-O3X	2.04	1.63	1.54
3	B	301	NAP	PA-O2A	2.05	1.65	1.55
3	D	301	NAP	P2B-O3X	2.11	1.63	1.54
3	B	301	NAP	O3B-C3B	2.14	1.47	1.43
3	D	301	NAP	C5B-C4B	2.17	1.58	1.51
3	C	301	NAP	C6N-N1N	2.19	1.41	1.35
3	D	301	NAP	C7N-N7N	2.22	1.37	1.33
3	D	301	NAP	C3N-C7N	2.24	1.54	1.50
3	B	301	NAP	C5B-C4B	2.34	1.59	1.51
3	C	301	NAP	PA-O5B	2.41	1.69	1.59
3	D	301	NAP	C2D-C1D	2.42	1.57	1.53
3	D	301	NAP	PA-O2A	2.46	1.67	1.55
3	C	301	NAP	C2D-C1D	2.47	1.57	1.53
3	B	301	NAP	O4D-C1D	2.52	1.44	1.41
3	B	301	NAP	C3N-C7N	2.76	1.54	1.50
3	D	301	NAP	O4B-C1B	2.79	1.45	1.41
3	D	301	NAP	P2B-O2X	2.80	1.66	1.54
3	D	301	NAP	C6N-N1N	2.88	1.42	1.35
3	D	301	NAP	PA-O5B	3.12	1.72	1.59
3	C	301	NAP	PA-O2A	3.29	1.71	1.55
3	D	301	NAP	P2B-O2B	3.34	1.65	1.59
3	B	301	NAP	P2B-O2B	3.39	1.65	1.59
3	B	301	NAP	O4B-C4B	3.40	1.52	1.45
3	C	301	NAP	C3N-C7N	3.52	1.56	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	301	NAP	C6N-N1N	3.74	1.45	1.35
3	B	301	NAP	O4B-C1B	4.06	1.46	1.41
3	D	301	NAP	O4D-C1D	4.19	1.47	1.41
3	C	301	NAP	O4B-C1B	4.68	1.47	1.41
3	C	301	NAP	O4D-C1D	5.40	1.48	1.41

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	301	NAP	N3A-C2A-N1A	-9.39	120.68	128.86
3	B	301	NAP	N3A-C2A-N1A	-9.00	121.02	128.86
3	C	301	NAP	N3A-C2A-N1A	-8.37	121.57	128.86
3	C	301	NAP	O2A-PA-O5B	-5.91	80.21	108.14
3	D	301	NAP	O2A-PA-O5B	-5.73	81.09	108.14
3	C	301	NAP	O5B-PA-O1A	-5.71	86.20	109.25
3	D	301	NAP	O5B-PA-O1A	-5.63	86.54	109.25
3	B	301	NAP	O2A-PA-O5B	-5.24	83.39	108.14
3	B	301	NAP	O5B-PA-O1A	-5.14	88.51	109.25
3	B	301	NAP	C4A-C5A-N7A	-3.10	106.42	109.41
3	C	301	NAP	C3N-C7N-N7N	-3.03	114.31	117.77
3	B	301	NAP	C3N-C7N-N7N	-2.97	114.39	117.77
3	B	301	NAP	O4D-C4D-C5D	-2.73	100.18	109.40
3	D	301	NAP	C4D-O4D-C1D	-2.61	106.99	109.77
3	B	301	NAP	C4D-O4D-C1D	-2.54	107.06	109.77
3	D	301	NAP	O4B-C1B-C2B	-2.47	102.27	106.59
3	C	301	NAP	C4A-C5A-N7A	-2.36	107.13	109.41
3	D	301	NAP	C4A-C5A-N7A	-2.34	107.15	109.41
3	C	301	NAP	O4B-C1B-C2B	-2.27	102.62	106.59
3	C	301	NAP	O3X-P2B-O2X	-2.23	98.62	107.61
3	D	301	NAP	O3X-P2B-O2X	-2.21	98.69	107.61
3	D	301	NAP	O7N-C7N-C3N	-2.11	117.15	119.62
3	B	301	NAP	C3N-C2N-N1N	-2.00	118.41	120.43
3	C	301	NAP	O7N-C7N-N7N	2.01	125.44	122.58
3	B	301	NAP	O4B-C4B-C5B	2.20	116.82	109.40
3	B	301	NAP	O7N-C7N-N7N	2.31	125.87	122.58
3	B	301	NAP	C2N-C3N-C4N	2.46	121.07	118.26
3	C	301	NAP	O2A-PA-O1A	2.59	125.66	112.28
3	D	301	NAP	O7N-C7N-N7N	2.69	126.40	122.58
3	B	301	NAP	O3X-P2B-O1X	2.82	121.52	110.50

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	301	NAP	4	0
3	C	301	NAP	7	0
3	D	301	NAP	4	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	246/269 (91%)	-0.26	8 (3%) 47 49	10, 22, 43, 71	0
1	B	240/269 (89%)	-0.32	5 (2%) 64 65	10, 24, 49, 61	0
1	C	240/269 (89%)	-0.23	6 (2%) 58 59	14, 26, 46, 80	0
1	D	237/269 (88%)	0.04	12 (5%) 29 31	15, 32, 57, 68	0
All	All	963/1076 (89%)	-0.19	31 (3%) 48 50	10, 25, 53, 80	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	249	ALA	5.0
1	D	55	ALA	4.5
1	C	203	LYS	3.8
1	A	200	ALA	3.6
1	D	54	ALA	3.4
1	D	4	LEU	3.3
1	D	42	SER	3.2
1	C	195	MET	3.1
1	A	204	LEU	3.1
1	B	197	SER	3.0
1	D	57	GLY	2.8
1	D	53	ILE	2.8
1	C	55	ALA	2.8
1	D	52	ALA	2.8
1	C	194	ASP	2.6
1	A	56	ALA	2.6
1	B	193	THR	2.5
1	A	206	GLU	2.5
1	A	202	GLU	2.5
1	D	206	GLU	2.5
1	A	199	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	206	GLU	2.4
1	D	104	ARG	2.3
1	B	213	TYR	2.3
1	A	203	LYS	2.3
1	B	206	GLU	2.3
1	D	56	ALA	2.2
1	A	247	VAL	2.2
1	D	58	GLY	2.2
1	B	194	ASP	2.1
1	D	39	TYR	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 carbohydrates 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. 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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	NAP	D	301	48/48	0.91	0.20	1.24	41,53,64,66	0
2	SO4	A	301	5/5	0.94	0.13	0.41	55,56,60,61	0
3	NAP	C	301	48/48	0.93	0.15	0.35	27,51,58,59	0
3	NAP	B	301	48/48	0.97	0.14	0.34	17,31,36,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.