



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:35 AM GMT

PDB ID : 3F8P  
Title : Structure of Sulfolobus solfataricus TrxR-B3  
Authors : Ruggiero, A.; Masullo, M.; Ruocco, M.R.; Arcari, P.; Zagari, A.; Vitagliano, L.  
Deposited on : 2008-11-13  
Resolution : 1.80 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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) : trunk26865

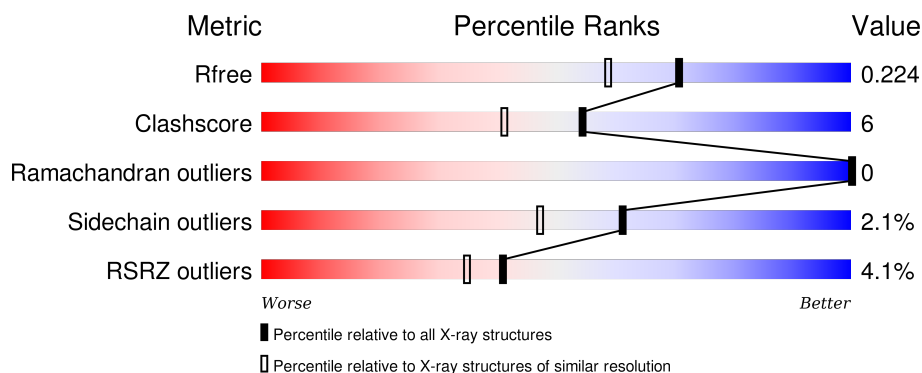
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 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))
$R_{free}$	91344	4533 (1.80-1.80)
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (1.80-1.80)

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  $\geq 3$ , 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	323	<div> <div>2%</div> <div>85% 10% 5%</div> </div>
1	B	323	<div> <div>4%</div> <div>85% 9% 5%</div> </div>
1	C	323	<div> <div>4%</div> <div>81% 13% 5%</div> </div>
1	D	323	<div> <div>6%</div> <div>83% 12% 5%</div> </div>

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	B	1501	-	-	-	X

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 9901 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 Thioredoxin reductase (TrxB-3).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	307	Total	C	N	O	S	0	0	0
			2359	1518	391	444	6			
1	B	306	Total	C	N	O	S	0	0	0
			2351	1514	389	442	6			
1	C	308	Total	C	N	O	S	0	0	0
			2368	1524	393	445	6			
1	D	310	Total	C	N	O	S	0	0	0
			2382	1530	397	449	6			

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula:  $C_{21}H_{27}N_7O_{14}P_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	D	1	Total	C	N	O	P	0	0
			26	10	5	9	2		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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	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	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is ACETIC ACID (three-letter code: ACY) (formula: C<sub>2</sub>H<sub>4</sub>O<sub>2</sub>).



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

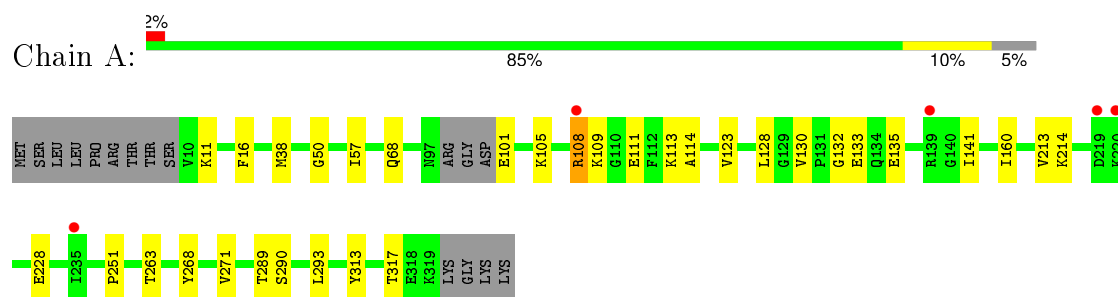
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	76	Total	O	0	0
			76	76		
5	B	73	Total	O	0	0
			73	73		
5	C	81	Total	O	0	0
			81	81		
5	D	70	Total	O	0	0
			70	70		

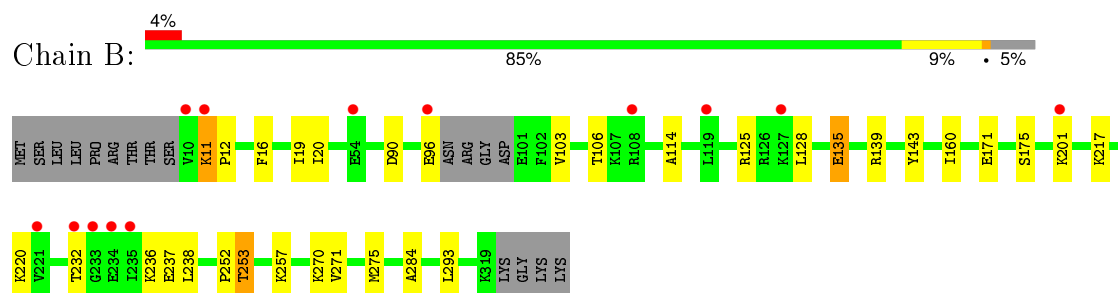
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

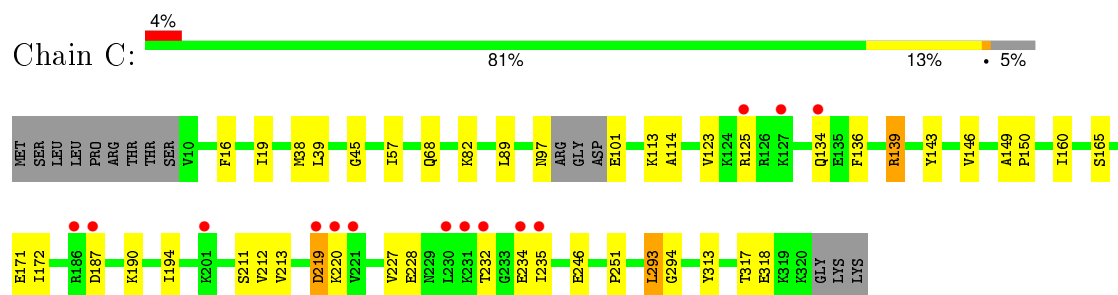
#### • Molecule 1: Thioredoxin reductase (TrxB-3)



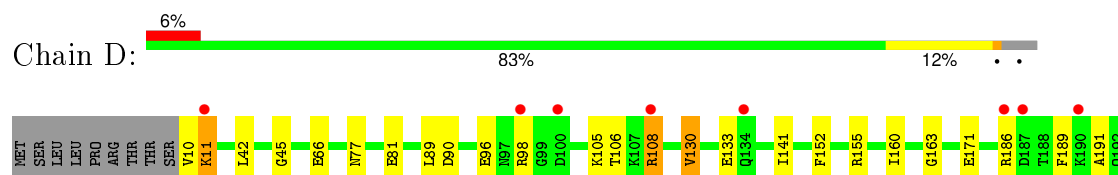
#### • Molecule 1: Thioredoxin reductase (TrxB-3)

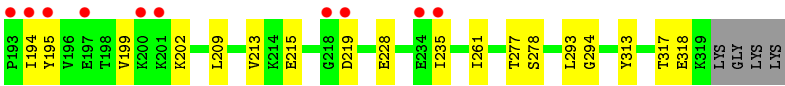


#### • Molecule 1: Thioredoxin reductase (TrxB-3)



#### • Molecule 1: Thioredoxin reductase (TrxB-3)







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	76.77Å 120.68Å 126.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 1.80 19.97 – 1.80	Depositor EDS
% Data completeness (in resolution range)	(Not available) (15.00-1.80) 98.1 (19.97-1.80)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.92 (at 1.80Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.200 , 0.237 0.199 , 0.224	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	17.2	Xtriage
Anisotropy	0.160	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.43 , 52.4	EDS
Estimated twinning fraction	0.013 for -h,l,k	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 108720 reflections (0.001%)	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	9901	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.02% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

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

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.69	0/2399	0.78	0/3242
1	B	0.64	0/2391	0.76	0/3231
1	C	0.69	0/2408	0.79	0/3253
1	D	0.63	0/2423	0.77	0/3275
All	All	0.66	0/9621	0.77	0/13001

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 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	2359	0	2417	33	0
1	B	2351	0	2411	23	0
1	C	2368	0	2430	33	0
1	D	2382	0	2438	44	0
2	A	27	0	12	1	0
2	B	27	0	12	0	0
2	C	27	0	12	0	0
2	D	26	0	12	0	0
3	A	15	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	10	0	0	0	0
3	D	5	0	0	0	0
4	B	4	0	3	0	0
5	A	76	0	0	0	0
5	B	73	0	0	0	0
5	C	81	0	0	0	0
5	D	70	0	0	2	0
All	All	9901	0	9747	124	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 (124) 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:D:261:ILE:HG23	1:D:277:THR:HG21	1.43	0.96
1:D:228:GLU:HB2	1:D:235:ILE:HD13	1.48	0.92
1:D:202:LYS:HE2	1:D:202:LYS:HA	1.52	0.92
1:D:261:ILE:HG23	1:D:277:THR:CG2	2.03	0.88
1:D:277:THR:HG22	1:D:278:SER:H	1.43	0.82
1:A:289:THR:HG22	1:A:290:SER:N	1.96	0.80
1:C:101:GLU:OE2	1:C:113:LYS:HD3	1.82	0.80
1:D:228:GLU:HB2	1:D:235:ILE:CD1	2.11	0.80
1:A:57:ILE:HD11	1:A:68:GLN:NE2	1.98	0.79
1:A:313:TYR:OH	1:B:171:GLU:HG2	1.83	0.79
1:D:98:ARG:HG2	1:D:98:ARG:O	1.84	0.78
1:A:271:VAL:HG21	1:A:289:THR:HG21	1.65	0.77
1:A:50:GLY:HA2	2:A:4001:NAD:O3B	1.85	0.77
1:B:217:LYS:HE2	1:B:237:GLU:OE1	1.85	0.77
1:A:271:VAL:HG22	1:A:289:THR:CG2	2.14	0.76
1:A:271:VAL:CG2	1:A:289:THR:HG21	2.16	0.76
1:A:132:GLY:HA2	1:A:135:GLU:HG2	1.68	0.75
1:D:277:THR:HG22	1:D:278:SER:N	2.03	0.73
1:D:90:ASP:OD1	1:D:108:ARG:HD2	1.88	0.73
1:D:90:ASP:OD2	1:D:106:THR:HG22	1.88	0.73
1:C:136:PHE:HA	1:C:139:ARG:HG3	1.71	0.72
1:D:277:THR:HG23	5:D:5021:HOH:O	1.89	0.72
1:D:261:ILE:CG2	1:D:277:THR:HG21	2.19	0.72
1:D:202:LYS:CA	1:D:202:LYS:HE2	2.19	0.72
1:C:293:LEU:HD13	1:C:294:GLY:N	2.06	0.71
1:C:160:ILE:CD1	1:C:213:VAL:HG21	2.20	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:271:VAL:CG2	1:A:289:THR:CG2	2.70	0.70
1:C:160:ILE:HD13	1:C:213:VAL:HG21	1.73	0.69
1:B:125:ARG:HD2	1:B:143:TYR:CD2	2.28	0.68
1:A:289:THR:HG22	1:A:290:SER:H	1.58	0.67
1:D:160:ILE:HD11	1:D:213:VAL:HG21	1.77	0.65
1:D:293:LEU:HD13	1:D:294:GLY:N	2.11	0.65
1:B:232:THR:O	1:B:232:THR:HG22	1.96	0.65
1:D:10:VAL:C	1:D:11:LYS:HD2	2.18	0.63
1:B:139:ARG:HD2	1:B:220:LYS:O	2.00	0.61
1:D:277:THR:CG2	5:D:5021:HOH:O	2.48	0.61
1:A:132:GLY:HA2	1:A:135:GLU:CG	2.31	0.59
1:D:90:ASP:OD2	1:D:106:THR:CG2	2.50	0.59
1:A:160:ILE:CD1	1:A:213:VAL:HG21	2.33	0.59
1:C:212:VAL:O	1:C:227:VAL:HG23	2.03	0.58
1:A:108:ARG:NE	1:A:109:LYS:HE3	2.17	0.58
1:C:194:ILE:HG12	1:D:318:GLU:HG2	1.85	0.58
1:D:10:VAL:HG11	1:D:42:LEU:HD13	1.87	0.57
1:D:77:ASN:O	1:D:81:GLU:HG2	2.04	0.57
1:D:228:GLU:CB	1:D:235:ILE:HD13	2.29	0.56
1:C:313:TYR:OH	1:D:171:GLU:HG3	2.05	0.56
1:A:101:GLU:OE2	1:A:113:LYS:HD3	2.07	0.55
1:A:289:THR:CG2	1:A:290:SER:N	2.67	0.55
1:D:163:GLY:C	1:D:191:ALA:HB2	2.26	0.55
1:C:57:ILE:HD11	1:C:68:GLN:HE21	1.72	0.55
1:C:136:PHE:CA	1:C:139:ARG:HG3	2.36	0.54
1:A:108:ARG:CD	1:A:109:LYS:HE3	2.37	0.54
1:C:82:LYS:NZ	1:D:66:GLU:OE2	2.41	0.54
1:B:11:LYS:NZ	1:B:11:LYS:HB2	2.22	0.54
1:B:135:GLU:N	1:B:135:GLU:OE1	2.41	0.53
1:A:105:LYS:HE2	1:A:111:GLU:OE2	2.08	0.53
1:D:98:ARG:O	1:D:98:ARG:CG	2.56	0.53
1:C:125:ARG:HD2	1:C:143:TYR:CD2	2.44	0.53
1:A:317:THR:HB	1:B:171:GLU:OE2	2.09	0.52
1:A:57:ILE:CD1	1:A:68:GLN:NE2	2.70	0.52
1:A:214:LYS:HE3	1:A:228:GLU:OE1	2.10	0.51
1:C:228:GLU:HB2	1:C:235:ILE:HD12	1.92	0.51
1:C:228:GLU:HB2	1:C:235:ILE:CD1	2.41	0.51
1:D:277:THR:CG2	1:D:278:SER:H	2.20	0.50
1:D:293:LEU:C	1:D:293:LEU:HD13	2.32	0.50
1:A:128:LEU:HG	1:A:130:VAL:HG13	1.92	0.50
1:D:96:GLU:OE2	1:D:105:LYS:HE3	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:160:ILE:HD11	1:A:213:VAL:HG21	1.94	0.50
1:A:108:ARG:HE	1:A:109:LYS:HE3	1.76	0.50
1:D:189:PHE:CZ	1:D:209:LEU:HD21	2.47	0.50
1:C:194:ILE:CG1	1:D:318:GLU:HG2	2.41	0.49
1:B:270:LYS:HD2	1:B:271:VAL:N	2.27	0.49
1:D:277:THR:CG2	1:D:278:SER:N	2.75	0.49
1:D:133:GLU:HA	1:D:141:ILE:CD1	2.42	0.49
1:B:253:THR:HG22	1:B:257:LYS:HE3	1.94	0.49
1:A:133:GLU:HA	1:A:141:ILE:CD1	2.43	0.49
1:C:125:ARG:HH11	1:C:125:ARG:HG3	1.78	0.49
1:D:195:TYR:O	1:D:199:VAL:HG23	2.12	0.48
1:C:136:PHE:C	1:C:139:ARG:HG3	2.33	0.48
1:C:146:VAL:HA	1:C:172:ILE:HD13	1.95	0.48
1:A:289:THR:CG2	1:A:290:SER:H	2.24	0.47
1:B:232:THR:CG2	1:B:232:THR:O	2.61	0.47
1:B:96:GLU:HB3	1:B:103:VAL:HB	1.96	0.47
1:D:90:ASP:CG	1:D:106:THR:HG22	2.34	0.47
1:C:171:GLU:HG3	1:D:313:TYR:OH	2.15	0.47
1:B:236:LYS:HE2	1:B:238:LEU:HD23	1.96	0.47
1:B:16:PHE:O	1:B:114:ALA:HA	2.14	0.47
1:C:318:GLU:HG2	1:D:194:ILE:HD11	1.95	0.47
1:A:123:VAL:HG12	1:A:251:PRO:HA	1.97	0.47
1:C:165:SER:OG	1:C:246:GLU:OE1	2.33	0.46
1:B:90:ASP:OD2	1:B:106:THR:HB	2.15	0.46
1:A:38:MET:HG2	1:B:175:SER:HB3	1.97	0.46
1:C:19:ILE:HD12	1:C:114:ALA:HB2	1.98	0.46
1:C:57:ILE:CD1	1:C:68:GLN:NE2	2.79	0.46
1:C:57:ILE:HD11	1:C:68:GLN:NE2	2.31	0.46
1:A:271:VAL:HG13	1:A:289:THR:HG23	1.99	0.45
1:A:293:LEU:HD23	1:A:293:LEU:C	2.37	0.45
1:D:160:ILE:CD1	1:D:213:VAL:HG21	2.44	0.45
1:C:219:ASP:HB3	1:C:220:LYS:H	1.50	0.45
1:B:19:ILE:HG13	1:B:114:ALA:HB2	1.99	0.45
1:C:232:THR:CG2	1:C:234:GLU:HB2	2.47	0.45
1:A:160:ILE:HD13	1:A:213:VAL:HG21	1.98	0.44
1:C:313:TYR:O	1:C:317:THR:HG23	2.18	0.44
1:A:16:PHE:O	1:A:114:ALA:HA	2.18	0.43
1:D:90:ASP:HB3	1:D:106:THR:HG21	2.00	0.43
1:B:217:LYS:CE	1:B:237:GLU:OE1	2.62	0.43
1:C:123:VAL:HG12	1:C:251:PRO:HA	2.00	0.43
1:C:211:SER:HB3	1:C:227:VAL:CG2	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:11:LYS:HG3	1:B:12:PRO:HD2	2.01	0.42
1:C:149:ALA:HB3	1:C:150:PRO:HD3	2.02	0.42
1:C:16:PHE:O	1:C:114:ALA:HA	2.21	0.41
1:D:313:TYR:O	1:D:317:THR:HG23	2.20	0.41
1:B:201:LYS:HE3	1:B:201:LYS:HB2	1.88	0.41
1:A:108:ARG:HG2	1:A:109:LYS:HG2	2.01	0.41
1:C:45:GLY:O	1:C:89:LEU:HA	2.21	0.41
1:D:130:VAL:CG2	1:D:215:GLU:HA	2.51	0.41
1:D:152:PHE:HA	1:D:155:ARG:HD3	2.02	0.41
1:D:81:GLU:HA	1:D:81:GLU:OE1	2.19	0.41
1:A:263:THR:HA	1:A:268:TYR:O	2.20	0.40
1:D:45:GLY:O	1:D:89:LEU:HA	2.21	0.40
1:B:128:LEU:HD21	1:B:160:ILE:HD13	2.02	0.40
1:C:39:LEU:HD11	1:C:313:TYR:HA	2.03	0.40
1:B:253:THR:CG2	1:B:257:LYS:HE3	2.51	0.40
1:B:275:MET:HB3	1:B:284:ALA:O	2.21	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	303/323 (94%)	296 (98%)	7 (2%)	0	100	100
1	B	302/323 (94%)	293 (97%)	9 (3%)	0	100	100
1	C	304/323 (94%)	299 (98%)	5 (2%)	0	100	100
1	D	308/323 (95%)	302 (98%)	6 (2%)	0	100	100
All	All	1217/1292 (94%)	1190 (98%)	27 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	253/267 (95%)	251 (99%)	2 (1%)	86	83
1	B	252/267 (94%)	246 (98%)	6 (2%)	57	41
1	C	254/267 (95%)	246 (97%)	8 (3%)	47	30
1	D	255/267 (96%)	250 (98%)	5 (2%)	63	49
All	All	1014/1068 (95%)	993 (98%)	21 (2%)	61	47

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	LYS
1	A	108	ARG
1	B	11	LYS
1	B	20	ILE
1	B	135	GLU
1	B	252	PRO
1	B	253	THR
1	B	293	LEU
1	C	38	MET
1	C	97	ASN
1	C	134	GLN
1	C	139	ARG
1	C	187	ASP
1	C	190	LYS
1	C	219	ASP
1	C	293	LEU
1	D	11	LYS
1	D	108	ARG
1	D	130	VAL
1	D	186	ARG
1	D	219	ASP

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	68	GLN
1	A	134	GLN
1	A	184	HIS
1	B	68	GLN
1	B	134	GLN
1	C	68	GLN
1	C	77	ASN
1	C	210	ASN
1	D	68	GLN
1	D	77	ASN
1	D	134	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

11 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	NAD	A	4001	-	23,29,48	1.42	5 (21%)	27,45,73	1.44	3 (11%)
3	SO4	A	4157	-	4,4,4	3.33	2 (50%)	6,6,6	1.07	0
3	SO4	A	4159	-	4,4,4	3.04	2 (50%)	6,6,6	0.89	0



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	SO4	A	4160	-	4,4,4	3.34	2 (50%)	6,6,6	0.89	0
4	ACY	B	1501	-	1,3,3	3.21	1 (100%)	0,3,3	0.00	-
2	NAD	B	4002	-	23,29,48	1.20	3 (13%)	27,45,73	0.85	1 (3%)
3	SO4	B	4158	-	4,4,4	3.18	2 (50%)	6,6,6	0.99	0
3	SO4	B	4161	-	4,4,4	3.30	2 (50%)	6,6,6	0.86	0
2	NAD	C	4003	-	23,29,48	1.04	3 (13%)	27,45,73	1.06	2 (7%)
2	NAD	D	4004	-	19,28,48	1.37	3 (15%)	21,42,73	1.23	2 (9%)
3	SO4	D	4162	-	4,4,4	3.23	2 (50%)	6,6,6	0.83	0

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	NAD	A	4001	-	-	0/12/32/62	0/3/3/5
3	SO4	A	4157	-	-	0/0/0/0	0/0/0/0
3	SO4	A	4159	-	-	0/0/0/0	0/0/0/0
3	SO4	A	4160	-	-	0/0/0/0	0/0/0/0
4	ACY	B	1501	-	-	0/0/0/0	0/0/0/0
2	NAD	B	4002	-	-	0/12/32/62	0/3/3/5
3	SO4	B	4158	-	-	0/0/0/0	0/0/0/0
3	SO4	B	4161	-	-	0/0/0/0	0/0/0/0
2	NAD	C	4003	-	-	0/12/32/62	0/3/3/5
2	NAD	D	4004	-	-	0/9/31/62	0/3/3/5
3	SO4	D	4162	-	-	0/0/0/0	0/0/0/0

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	4161	SO4	O3-S	-4.88	1.29	1.47
3	A	4157	SO4	O3-S	-4.79	1.30	1.47
3	A	4160	SO4	O3-S	-4.68	1.30	1.47
3	B	4158	SO4	O3-S	-4.44	1.31	1.47
3	A	4159	SO4	O3-S	-4.41	1.31	1.47
3	D	4162	SO4	O3-S	-4.35	1.31	1.47
2	A	4001	NAD	PN-O1N	-3.09	1.40	1.50
2	A	4001	NAD	C2B-C3B	-2.50	1.46	1.53
2	A	4001	NAD	O4B-C4B	-2.22	1.39	1.45
2	D	4004	NAD	O4B-C4B	-2.05	1.40	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	4002	NAD	O4B-C4B	-2.02	1.40	1.45
2	C	4003	NAD	PN-O5D	2.12	1.63	1.55
2	A	4001	NAD	C2A-N3A	2.16	1.36	1.32
2	C	4003	NAD	O3B-C3B	2.19	1.48	1.43
2	A	4001	NAD	PN-O5D	2.23	1.63	1.55
2	C	4003	NAD	C2A-N3A	2.33	1.36	1.32
2	B	4002	NAD	C4A-N3A	2.45	1.39	1.35
2	B	4002	NAD	PN-O5D	2.59	1.65	1.55
2	D	4004	NAD	O3B-C3B	2.87	1.49	1.43
4	B	1501	ACY	CH3-C	3.21	1.53	1.48
2	D	4004	NAD	C4A-N3A	3.70	1.41	1.35
3	A	4159	SO4	O1-S	4.06	1.61	1.47
3	B	4161	SO4	O1-S	4.30	1.61	1.47
3	B	4158	SO4	O1-S	4.43	1.62	1.47
3	D	4162	SO4	O1-S	4.53	1.62	1.47
3	A	4157	SO4	O1-S	4.56	1.62	1.47
3	A	4160	SO4	O1-S	4.68	1.63	1.47

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	4001	NAD	C4B-O4B-C1B	-4.33	104.96	109.72
2	A	4001	NAD	N3A-C2A-N1A	-2.91	126.67	128.89
2	C	4003	NAD	N3A-C2A-N1A	-2.71	126.82	128.89
2	B	4002	NAD	C4B-O4B-C1B	-2.29	107.21	109.72
2	C	4003	NAD	C2B-C1B-N9A	2.36	117.90	114.29
2	A	4001	NAD	C4A-C5A-N7A	2.43	111.71	109.48
2	D	4004	NAD	C4A-C5A-N7A	2.61	111.88	109.48
2	D	4004	NAD	C2B-C1B-N9A	2.90	118.72	114.29

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	4001	NAD	1	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	307/323 (95%)	0.01	5 (1%) 74 71	6, 18, 36, 45	0
1	B	306/323 (94%)	0.34	13 (4%) 40 34	8, 23, 42, 54	0
1	C	308/323 (95%)	0.11	14 (4%) 37 31	7, 18, 41, 57	0
1	D	310/323 (95%)	0.21	18 (5%) 26 21	7, 20, 46, 56	0
All	All	1231/1292 (95%)	0.16	50 (4%) 41 35	6, 20, 42, 57	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	194	ILE	5.9
1	B	11	LYS	4.5
1	C	127	LYS	4.1
1	B	127	LYS	4.0
1	D	200	LYS	4.0
1	D	201	LYS	3.8
1	C	235	ILE	3.7
1	B	235	ILE	3.6
1	C	187	ASP	3.6
1	D	98	ARG	3.5
1	C	219	ASP	3.5
1	A	220	LYS	3.2
1	D	11	LYS	3.1
1	C	134	GLN	3.1
1	B	108	ARG	3.1
1	D	186	ARG	3.1
1	B	10	VAL	3.0
1	A	235	ILE	3.0
1	C	220	LYS	3.0
1	C	234	GLU	2.9
1	D	219	ASP	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	119	LEU	2.9
1	B	201	LYS	2.9
1	A	219	ASP	2.8
1	B	96	GLU	2.8
1	B	234	GLU	2.7
1	C	186	ARG	2.7
1	D	190	LYS	2.6
1	C	232	THR	2.6
1	D	197	GLU	2.5
1	A	108	ARG	2.5
1	B	233	GLY	2.4
1	D	193	PRO	2.4
1	D	108	ARG	2.4
1	B	221	VAL	2.4
1	D	218	GLY	2.4
1	C	125	ARG	2.3
1	C	221	VAL	2.3
1	A	139	ARG	2.3
1	D	134	GLN	2.3
1	D	187	ASP	2.3
1	B	54	GLU	2.3
1	B	232	THR	2.3
1	D	195	TYR	2.2
1	C	230	LEU	2.2
1	C	231	LYS	2.2
1	D	235	ILE	2.2
1	C	201	LYS	2.2
1	D	100	ASP	2.1
1	D	234	GLU	2.1

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

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
4	ACY	B	1501	4/4	0.92	0.17	2.30	25,29,29,30	0
3	SO4	A	4160	5/5	0.91	0.21	1.01	56,56,57,57	0
3	SO4	A	4159	5/5	0.94	0.12	-0.05	47,48,49,50	0
3	SO4	D	4162	5/5	0.94	0.11	-0.38	41,42,43,43	0
2	NAD	D	4004	26/44	0.97	0.08	-0.49	10,14,18,18	0
3	SO4	B	4161	5/5	0.98	0.09	-0.49	46,47,48,49	0
2	NAD	B	4002	27/44	0.96	0.10	-0.61	14,23,26,32	0
2	NAD	A	4001	27/44	0.95	0.09	-0.64	10,14,18,35	0
2	NAD	C	4003	27/44	0.97	0.08	-0.83	10,12,16,21	0
3	SO4	A	4157	5/5	0.94	0.16	-	55,55,56,57	0
3	SO4	B	4158	5/5	0.97	0.16	-	44,44,45,48	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.