



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 08:21 AM GMT

PDB ID : 3EAO
Title : Crystal structure of recombinant rat selenoprotein thioredoxin reductase 1 with oxidized C-terminal tail
Authors : Sandalova, T.; Cheng, Q.; Lindqvist, Y.; Arner, E.
Deposited on : 2008-08-26
Resolution : 3.10 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (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

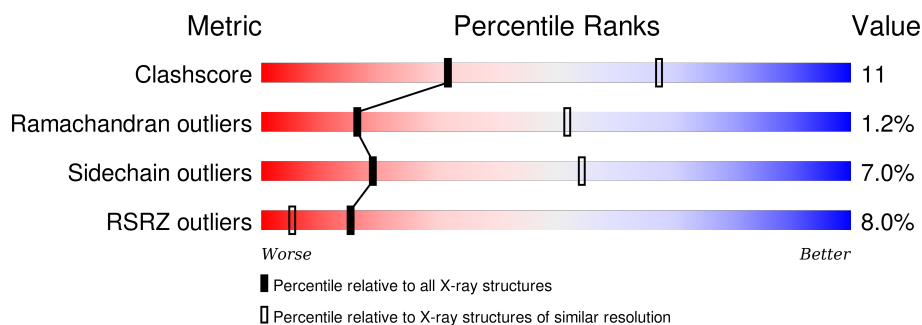
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1222 (3.14-3.06)
Ramachandran outliers	100387	1174 (3.14-3.06)
Sidechain outliers	100360	1174 (3.14-3.06)
RSRZ outliers	91569	1119 (3.14-3.06)

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	499	<div> <div>6%</div> <div>74%</div> <div>22%</div> <div>...</div> </div>
1	B	499	<div> <div>5%</div> <div>76%</div> <div>20%</div> <div>...</div> </div>
1	C	499	<div> <div>6%</div> <div>72%</div> <div>22%</div> <div>...</div> </div>
1	D	499	<div> <div>%</div> <div>73%</div> <div>22%</div> <div>...</div> </div>
1	E	499	<div> <div>4%</div> <div>75%</div> <div>21%</div> <div>...</div> </div>
1	F	499	<div> <div>24%</div> <div>68%</div> <div>27%</div> <div>..</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
3	NAP	F	601	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 23114 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.

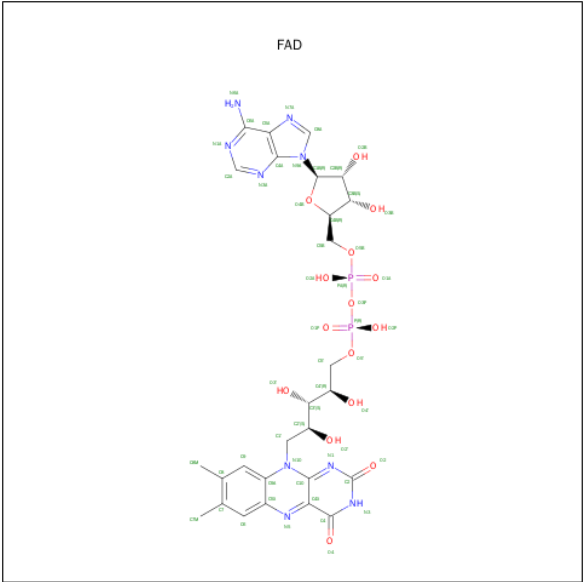
- Molecule 1 is a protein called Thioredoxin reductase 1, cytoplasmic.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	489	Total	C	N	O	S	Se	0	0	0
			3762	2390	636	714	21	1			
1	B	490	Total	C	N	O	S	Se	0	0	0
			3768	2393	637	716	21	1			
1	C	486	Total	C	N	O	S	Se	0	0	0
			3731	2368	633	708	21	1			
1	D	491	Total	C	N	O	S	Se	0	0	0
			3777	2399	639	717	21	1			
1	E	490	Total	C	N	O	S	Se	0	0	0
			3768	2393	637	716	21	1			
1	F	489	Total	C	N	O	S	Se	0	0	0
			3762	2390	636	714	21	1			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	52	ARG	ASN	CONFLICT	UNP O89049
A	53	TRP	GLY	CONFLICT	UNP O89049
B	52	ARG	ASN	CONFLICT	UNP O89049
B	53	TRP	GLY	CONFLICT	UNP O89049
C	52	ARG	ASN	CONFLICT	UNP O89049
C	53	TRP	GLY	CONFLICT	UNP O89049
D	52	ARG	ASN	CONFLICT	UNP O89049
D	53	TRP	GLY	CONFLICT	UNP O89049
E	52	ARG	ASN	CONFLICT	UNP O89049
E	53	TRP	GLY	CONFLICT	UNP O89049
F	52	ARG	ASN	CONFLICT	UNP O89049
F	53	TRP	GLY	CONFLICT	UNP O89049

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	E	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	F	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

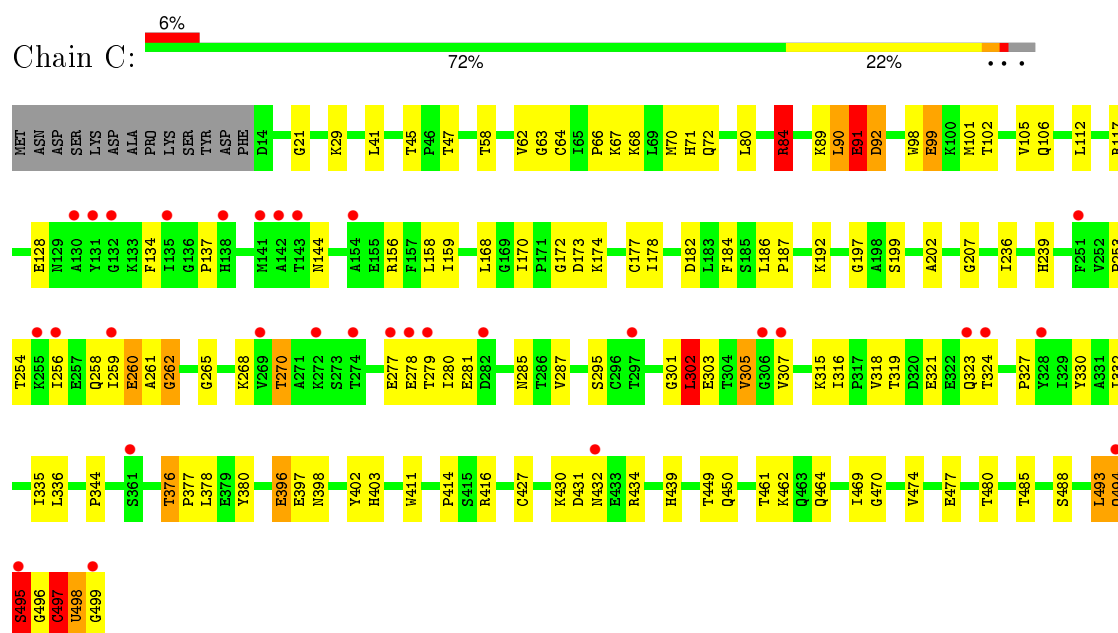
- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



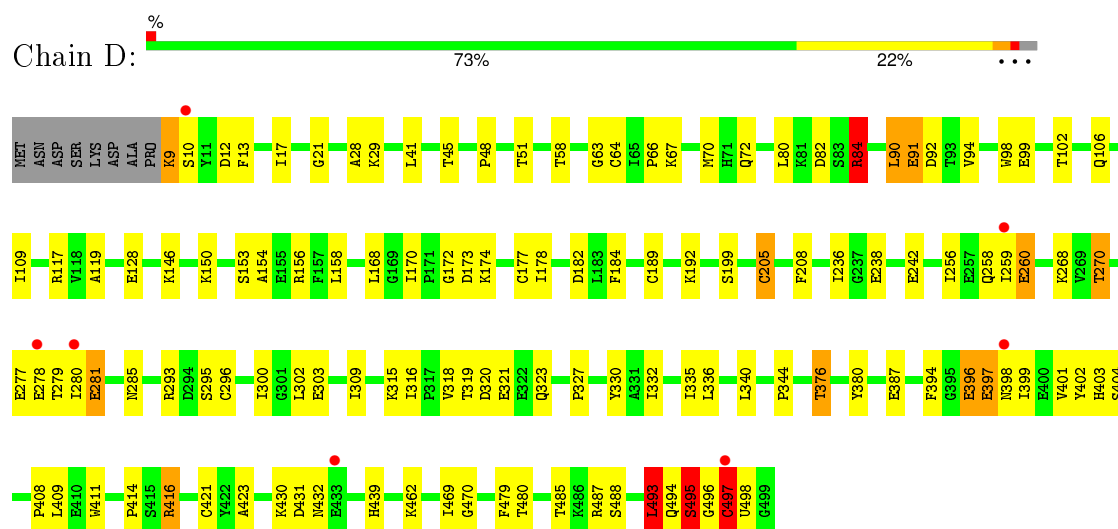
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			32	11	5	13	3		
3	B	1	Total	C	N	O	P	0	0
			32	11	5	13	3		
3	C	1	Total	C	N	O	P	0	0
			32	11	5	13	3		
3	D	1	Total	C	N	O	P	0	0
			32	11	5	13	3		
3	E	1	Total	C	N	O	P	0	0
			32	11	5	13	3		
3	F	1	Total	C	N	O	P	0	0
			32	11	5	13	3		

- Molecule 4 is water.

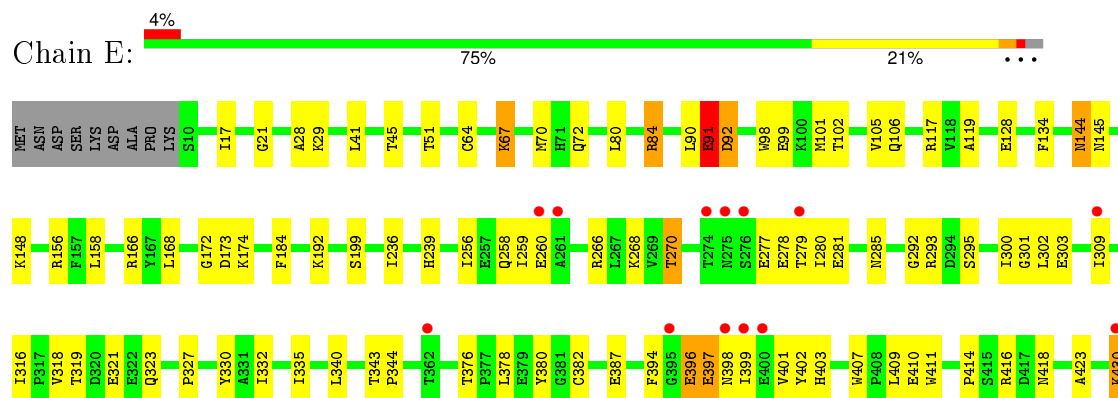
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	O	0	0
			3	3		
4	B	8	Total	O	0	0
			8	8		
4	C	5	Total	O	0	0
			5	5		
4	D	11	Total	O	0	0
			11	11		
4	E	4	Total	O	0	0
			4	4		
4	F	5	Total	O	0	0
			5	5		

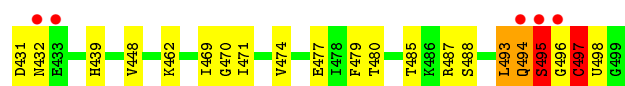


- Molecule 1: Thioredoxin reductase 1, cytoplasmic

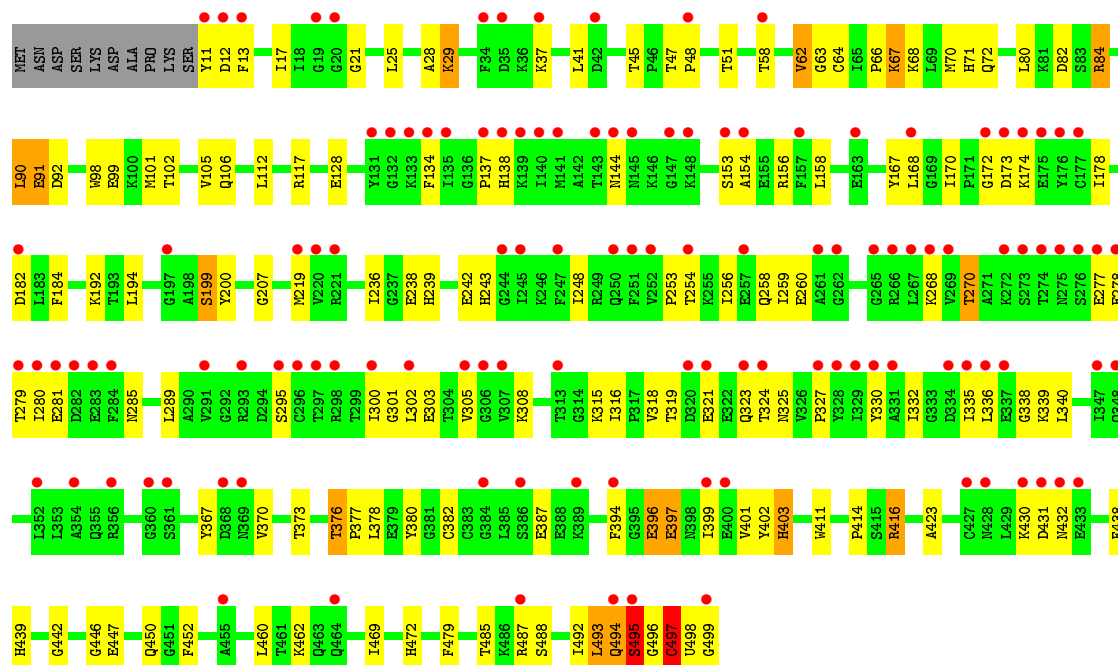


- Molecule 1: Thioredoxin reductase 1, cytoplasmic





- Molecule 1: Thioredoxin reductase 1, cytoplasmic



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	78.23Å 137.75Å 168.95Å 90.00° 94.00° 90.00°	Depositor
Resolution (Å)	30.00 – 3.10 64.35 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.3 (30.00-3.10) 99.3 (64.35-3.10)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.01 (at 3.13Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.256 , 0.289 0.260 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	63.3	Xtriage
Anisotropy	0.019	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 58.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 64479 reflections	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	23114	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.84% 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, 0.375 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: NAP, SEC, FAD

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	0.66	2/3829 (0.1%)	0.77	5/5177 (0.1%)
1	B	0.64	1/3835 (0.0%)	0.72	2/5185 (0.0%)
1	C	0.70	0/3796	0.80	6/5132 (0.1%)
1	D	0.69	5/3844 (0.1%)	0.76	4/5196 (0.1%)
1	E	0.67	0/3835	0.72	3/5185 (0.1%)
1	F	0.87	1/3829 (0.0%)	0.76	3/5177 (0.1%)
All	All	0.71	9/22968 (0.0%)	0.76	23/31052 (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	1
1	B	0	4
1	C	0	6
1	D	0	3
1	E	0	1
1	F	0	1
All	All	0	16

The worst 5 of 9 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	205	CYS	CB-SG	-6.33	1.71	1.82
1	D	189	CYS	CB-SG	-6.30	1.71	1.82
1	A	497	CYS	CB-SG	6.15	1.92	1.82
1	D	296	CYS	CB-SG	-5.33	1.73	1.81
1	F	497	CYS	CB-SG	5.33	1.91	1.82

The worst 5 of 23 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	84	ARG	NE-CZ-NH2	15.57	128.08	120.30
1	D	84	ARG	NE-CZ-NH1	-15.24	112.68	120.30
1	A	221	ARG	NE-CZ-NH2	14.32	127.46	120.30
1	A	221	ARG	NE-CZ-NH1	-14.06	113.27	120.30
1	C	84	ARG	NE-CZ-NH1	-13.93	113.34	120.30

There are no chirality outliers.

5 of 16 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	493	LEU	Peptide
1	B	10	SER	Peptide
1	B	197	GLY	Peptide
1	B	493	LEU	Peptide
1	B	91	GLU	Peptide

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	3762	0	3761	86	0
1	B	3768	0	3767	73	0
1	C	3731	0	3740	90	0
1	D	3777	0	3779	89	0
1	E	3768	0	3766	87	0
1	F	3762	0	3761	111	0
2	A	53	0	31	1	0
2	B	53	0	31	1	0
2	C	53	0	31	1	0
2	D	53	0	31	1	0
2	E	53	0	31	1	0
2	F	53	0	31	1	0
3	A	32	0	11	2	0
3	B	32	0	11	0	0
3	C	32	0	11	0	0
3	D	32	0	11	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	E	32	0	11	0	0
3	F	32	0	11	0	0
4	A	3	0	0	0	0
4	B	8	0	0	0	0
4	C	5	0	0	0	0
4	D	11	0	0	0	0
4	E	4	0	0	0	0
4	F	5	0	0	3	0
All	All	23114	0	22826	492	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 11.

The worst 5 of 492 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:497:CYS:SG	1:B:498:SEC:SE	2.55	1.15
1:A:493:LEU:HB3	1:A:494:GLN:HB2	1.23	1.11
1:B:197:GLY:O	1:B:202:ALA:HB1	1.55	1.06
1:F:248:ILE:HA	4:F:602:HOH:O	1.56	1.04
1:E:493:LEU:HB3	1:E:494:GLN:HB2	1.45	0.99

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	486/499 (97%)	436 (90%)	44 (9%)	6 (1%)	16	52
1	B	487/499 (98%)	440 (90%)	42 (9%)	5 (1%)	19	58
1	C	483/499 (97%)	437 (90%)	42 (9%)	4 (1%)	24	63

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	488/499 (98%)	450 (92%)	31 (6%)	7 (1%)	14	48
1	E	487/499 (98%)	441 (91%)	40 (8%)	6 (1%)	16	52
1	F	486/499 (97%)	437 (90%)	43 (9%)	6 (1%)	16	52
All	All	2917/2994 (97%)	2641 (90%)	242 (8%)	34 (1%)	16	52

5 of 34 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	495	SER
1	B	493	LEU
1	B	495	SER
1	C	495	SER
1	D	493	LEU

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	404/413 (98%)	377 (93%)	27 (7%)	20	56
1	B	405/413 (98%)	376 (93%)	29 (7%)	18	53
1	C	401/413 (97%)	373 (93%)	28 (7%)	19	54
1	D	406/413 (98%)	377 (93%)	29 (7%)	18	54
1	E	405/413 (98%)	377 (93%)	28 (7%)	19	55
1	F	404/413 (98%)	376 (93%)	28 (7%)	19	55
All	All	2425/2478 (98%)	2256 (93%)	169 (7%)	19	54

5 of 169 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	318	VAL
1	D	99	GLU
1	F	279	THR
1	C	396	GLU

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Mol	Chain	Res	Type
1	D	10	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 46 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	439	HIS
1	D	145	ASN
1	F	323	GLN
1	C	494	GLN
1	D	85	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](#)

12 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	FAD	A	600	-	48,58,58	1.48	7 (14%)	54,89,89	2.01	7 (12%)
3	NAP	A	601	-	27,34,52	1.57	4 (14%)	35,53,80	2.07	2 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	FAD	B	600	-	48,58,58	1.69	6 (12%)	54,89,89	2.19	7 (12%)
3	NAP	B	601	-	27,34,52	1.37	4 (14%)	35,53,80	2.11	6 (17%)
2	FAD	C	600	-	48,58,58	1.93	9 (18%)	54,89,89	1.95	9 (16%)
3	NAP	C	601	-	27,34,52	1.83	5 (18%)	35,53,80	2.30	3 (8%)
2	FAD	D	600	-	48,58,58	1.54	9 (18%)	54,89,89	1.90	8 (14%)
3	NAP	D	601	-	27,34,52	1.31	3 (11%)	35,53,80	2.06	5 (14%)
2	FAD	E	600	-	48,58,58	1.69	9 (18%)	54,89,89	1.84	6 (11%)
3	NAP	E	601	-	27,34,52	1.49	4 (14%)	35,53,80	2.01	4 (11%)
2	FAD	F	600	-	48,58,58	2.10	7 (14%)	54,89,89	2.37	12 (22%)
3	NAP	F	601	-	27,34,52	1.99	4 (14%)	35,53,80	2.30	5 (14%)

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	FAD	A	600	-	-	0/30/50/50	0/6/6/6
3	NAP	A	601	-	-	0/20/40/67	0/3/3/5
2	FAD	B	600	-	-	0/30/50/50	0/6/6/6
3	NAP	B	601	-	-	0/20/40/67	0/3/3/5
2	FAD	C	600	-	-	0/30/50/50	0/6/6/6
3	NAP	C	601	-	-	0/20/40/67	0/3/3/5
2	FAD	D	600	-	-	0/30/50/50	0/6/6/6
3	NAP	D	601	-	-	0/20/40/67	0/3/3/5
2	FAD	E	600	-	-	0/30/50/50	0/6/6/6
3	NAP	E	601	-	-	0/20/40/67	0/3/3/5
2	FAD	F	600	-	-	0/30/50/50	0/6/6/6
3	NAP	F	601	-	-	0/20/40/67	0/3/3/5

The worst 5 of 71 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	601	NAP	P2B-O3X	-2.51	1.45	1.54
2	D	600	FAD	PA-O2A	-2.42	1.44	1.54
2	E	600	FAD	PA-O2A	-2.37	1.44	1.54
2	D	600	FAD	C5A-N7A	-2.23	1.31	1.39
3	B	601	NAP	P2B-O3X	-2.22	1.46	1.54

The worst 5 of 74 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	C	601	NAP	N3A-C2A-N1A	-12.02	119.69	128.89
2	F	600	FAD	N3A-C2A-N1A	-11.78	119.87	128.89
3	F	601	NAP	N3A-C2A-N1A	-11.76	119.89	128.89
2	B	600	FAD	N3A-C2A-N1A	-11.60	120.01	128.89
3	A	601	NAP	N3A-C2A-N1A	-10.56	120.81	128.89

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	600	FAD	1	0
3	A	601	NAP	2	0
2	B	600	FAD	1	0
2	C	600	FAD	1	0
2	D	600	FAD	1	0
2	E	600	FAD	1	0
2	F	600	FAD	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	488/499 (97%)	0.44	30 (6%) 25 10	23, 55, 69, 96	0
1	B	489/499 (97%)	0.34	25 (5%) 32 13	23, 55, 69, 96	0
1	C	485/499 (97%)	0.55	31 (6%) 23 9	23, 55, 68, 96	0
1	D	490/499 (98%)	0.36	7 (1%) 78 60	23, 55, 68, 96	0
1	E	489/499 (97%)	0.42	18 (3%) 45 22	23, 55, 68, 96	0
1	F	488/499 (97%)	1.09	122 (25%) 1 0	23, 55, 69, 96	0
All	All	2929/2994 (97%)	0.53	233 (7%) 15 5	23, 55, 69, 96	0

The worst 5 of 233 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	143	THR	6.3
1	F	278	GLU	5.0
1	F	282	ASP	4.9
1	A	132	GLY	4.9
1	F	177	CYS	4.9

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	A	601	32/48	0.90	0.34	1.17	64,73,87,89	0
2	FAD	E	600	53/53	0.95	0.30	0.75	54,57,61,61	0
2	FAD	B	600	53/53	0.95	0.28	0.61	54,57,61,62	0
3	NAP	E	601	32/48	0.87	0.30	0.57	64,73,87,89	0
3	NAP	C	601	32/48	0.83	0.37	0.54	64,73,87,89	0
2	FAD	D	600	53/53	0.96	0.28	0.52	54,57,61,62	0
2	FAD	C	600	53/53	0.89	0.33	0.48	54,57,61,62	0
3	NAP	D	601	32/48	0.89	0.32	0.46	64,73,87,89	0
2	FAD	F	600	53/53	0.83	0.39	0.33	54,57,61,62	0
3	NAP	F	601	32/48	0.79	0.45	0.22	64,73,87,89	0
3	NAP	B	601	32/48	0.86	0.31	0.11	64,73,87,89	0
2	FAD	A	600	53/53	0.94	0.27	-0.21	54,57,61,62	0

6.5 Other polymers [i](#)

There are no such residues in this entry.