Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 05:54 PM UTC

PDB ID	:	2OBK / pdb_00002obk
Title	:	X-Ray structure of the putative Se binding protein from Pseudomonas fluorescens. Northeast Structural Genomics Consortium target PlR6.
Authors	:	Kuzin, A.P.; Su, M.; Seetharaman, J.; Chen, C.X.; Fang, Y.; Cunningham, K.; Ma, L.C.; Xiao, R.; Liu, J.; Baran, M.C.; Acton, T.B.; Rost, B.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on	:	2006-12-19
Resolution	:	2.70 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

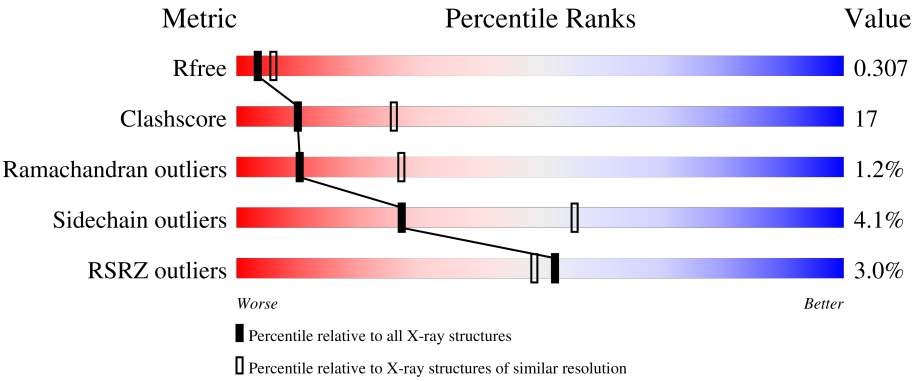
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

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 2.70 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}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	103	55% 22% 18%
1	B	103	52% 28% 17%
1	C	103	59% 17% 20%
1	D	103	40% 39% 18%

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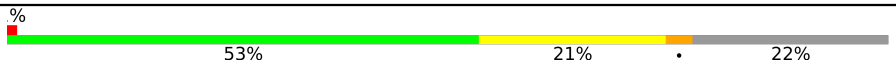
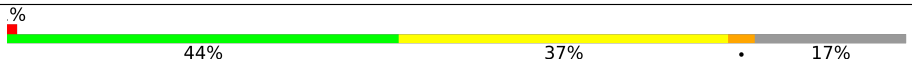
Mol	Chain	Length	Quality of chain
1	E	103	
1	F	103	
1	G	103	
1	H	103	

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 5443 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 SelT/selW/selH selenoprotein domain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	84	Total	C	N	O	S	0	0	0
			670	423	119	125	3			
1	B	85	Total	C	N	O	S	0	0	0
			678	429	120	126	3			
1	C	82	Total	C	N	O	S	0	0	0
			651	413	114	121	3			
1	D	84	Total	C	N	O	S	0	0	0
			670	423	119	125	3			
1	E	80	Total	C	N	O	S	0	0	0
			631	402	109	117	3			
1	F	86	Total	C	N	O	S	0	0	0
			682	431	121	127	3			
1	G	84	Total	C	N	O	S	0	0	0
			670	423	119	125	3			
1	H	83	Total	C	N	O	S	0	0	0
			659	417	115	124	3			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	modified residue	UNP Q4KGC5
A	96	LEU	-	cloning artifact	UNP Q4KGC5
A	97	GLU	-	cloning artifact	UNP Q4KGC5
A	98	HIS	-	cloning artifact	UNP Q4KGC5
A	99	HIS	-	cloning artifact	UNP Q4KGC5
A	100	HIS	-	cloning artifact	UNP Q4KGC5
A	101	HIS	-	cloning artifact	UNP Q4KGC5
A	102	HIS	-	cloning artifact	UNP Q4KGC5
A	103	HIS	-	cloning artifact	UNP Q4KGC5
B	1	MSE	MET	modified residue	UNP Q4KGC5
B	96	LEU	-	cloning artifact	UNP Q4KGC5
B	97	GLU	-	cloning artifact	UNP Q4KGC5
B	98	HIS	-	cloning artifact	UNP Q4KGC5

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Chain	Residue	Modelled	Actual	Comment	Reference
B	99	HIS	-	cloning artifact	UNP Q4KGC5
B	100	HIS	-	cloning artifact	UNP Q4KGC5
B	101	HIS	-	cloning artifact	UNP Q4KGC5
B	102	HIS	-	cloning artifact	UNP Q4KGC5
B	103	HIS	-	cloning artifact	UNP Q4KGC5
C	1	MSE	MET	modified residue	UNP Q4KGC5
C	96	LEU	-	cloning artifact	UNP Q4KGC5
C	97	GLU	-	cloning artifact	UNP Q4KGC5
C	98	HIS	-	cloning artifact	UNP Q4KGC5
C	99	HIS	-	cloning artifact	UNP Q4KGC5
C	100	HIS	-	cloning artifact	UNP Q4KGC5
C	101	HIS	-	cloning artifact	UNP Q4KGC5
C	102	HIS	-	cloning artifact	UNP Q4KGC5
C	103	HIS	-	cloning artifact	UNP Q4KGC5
D	1	MSE	MET	modified residue	UNP Q4KGC5
D	96	LEU	-	cloning artifact	UNP Q4KGC5
D	97	GLU	-	cloning artifact	UNP Q4KGC5
D	98	HIS	-	cloning artifact	UNP Q4KGC5
D	99	HIS	-	cloning artifact	UNP Q4KGC5
D	100	HIS	-	cloning artifact	UNP Q4KGC5
D	101	HIS	-	cloning artifact	UNP Q4KGC5
D	102	HIS	-	cloning artifact	UNP Q4KGC5
D	103	HIS	-	cloning artifact	UNP Q4KGC5
E	1	MSE	MET	modified residue	UNP Q4KGC5
E	96	LEU	-	cloning artifact	UNP Q4KGC5
E	97	GLU	-	cloning artifact	UNP Q4KGC5
E	98	HIS	-	cloning artifact	UNP Q4KGC5
E	99	HIS	-	cloning artifact	UNP Q4KGC5
E	100	HIS	-	cloning artifact	UNP Q4KGC5
E	101	HIS	-	cloning artifact	UNP Q4KGC5
E	102	HIS	-	cloning artifact	UNP Q4KGC5
E	103	HIS	-	cloning artifact	UNP Q4KGC5
F	1	MSE	MET	modified residue	UNP Q4KGC5
F	96	LEU	-	cloning artifact	UNP Q4KGC5
F	97	GLU	-	cloning artifact	UNP Q4KGC5
F	98	HIS	-	cloning artifact	UNP Q4KGC5
F	99	HIS	-	cloning artifact	UNP Q4KGC5
F	100	HIS	-	cloning artifact	UNP Q4KGC5
F	101	HIS	-	cloning artifact	UNP Q4KGC5
F	102	HIS	-	cloning artifact	UNP Q4KGC5
F	103	HIS	-	cloning artifact	UNP Q4KGC5
G	1	MSE	MET	modified residue	UNP Q4KGC5

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Chain	Residue	Modelled	Actual	Comment	Reference
G	96	LEU	-	cloning artifact	UNP Q4KGC5
G	97	GLU	-	cloning artifact	UNP Q4KGC5
G	98	HIS	-	cloning artifact	UNP Q4KGC5
G	99	HIS	-	cloning artifact	UNP Q4KGC5
G	100	HIS	-	cloning artifact	UNP Q4KGC5
G	101	HIS	-	cloning artifact	UNP Q4KGC5
G	102	HIS	-	cloning artifact	UNP Q4KGC5
G	103	HIS	-	cloning artifact	UNP Q4KGC5
H	1	MSE	MET	modified residue	UNP Q4KGC5
H	96	LEU	-	cloning artifact	UNP Q4KGC5
H	97	GLU	-	cloning artifact	UNP Q4KGC5
H	98	HIS	-	cloning artifact	UNP Q4KGC5
H	99	HIS	-	cloning artifact	UNP Q4KGC5
H	100	HIS	-	cloning artifact	UNP Q4KGC5
H	101	HIS	-	cloning artifact	UNP Q4KGC5
H	102	HIS	-	cloning artifact	UNP Q4KGC5
H	103	HIS	-	cloning artifact	UNP Q4KGC5

- Molecule 2 is water.

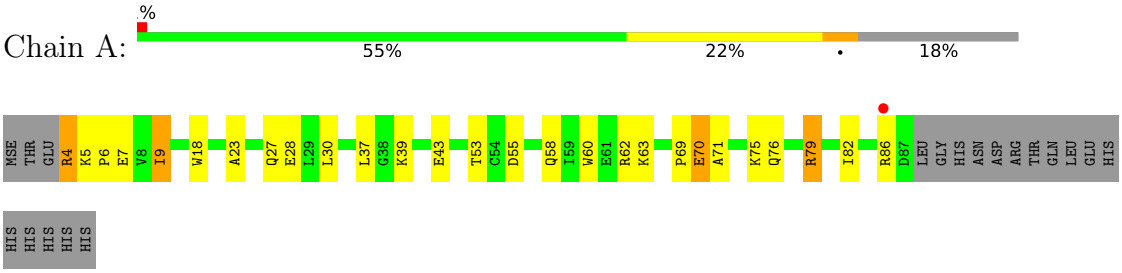
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	27	Total O 27 27	0	0
2	B	20	Total O 20 20	0	0
2	C	8	Total O 8 8	0	0
2	D	17	Total O 17 17	0	0
2	E	27	Total O 27 27	0	0
2	F	13	Total O 13 13	0	0
2	G	7	Total O 7 7	0	0
2	H	13	Total O 13 13	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry 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.

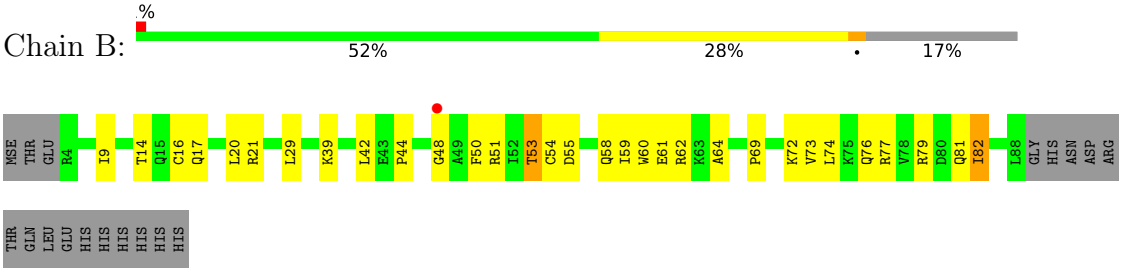
• Molecule 1: SelT/selW/selH selenoprotein domain

Chain A:



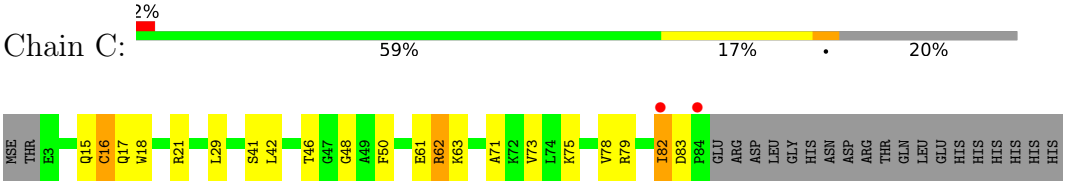
• Molecule 1: SelT/selW/selH selenoprotein domain

Chain B:



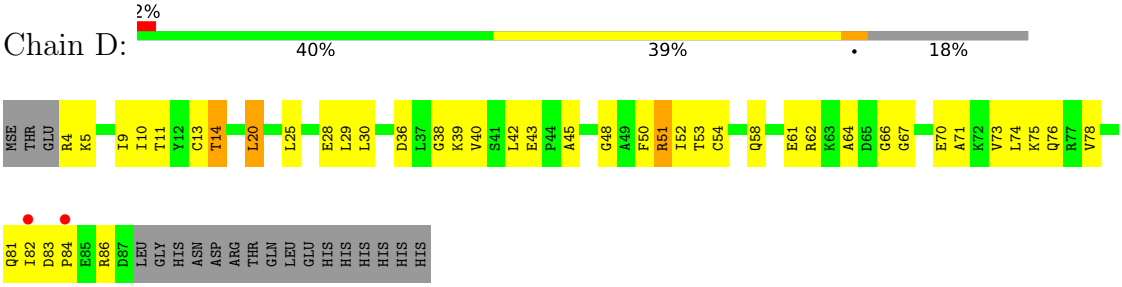
• Molecule 1: SelT/selW/selH selenoprotein domain

Chain C:



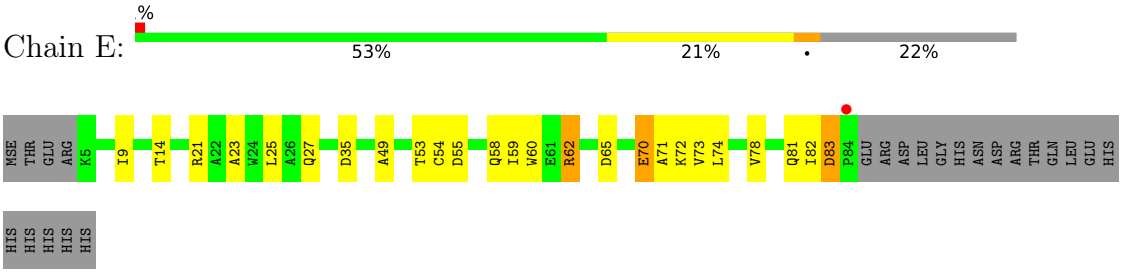
• Molecule 1: SelT/selW/selH selenoprotein domain

Chain D:



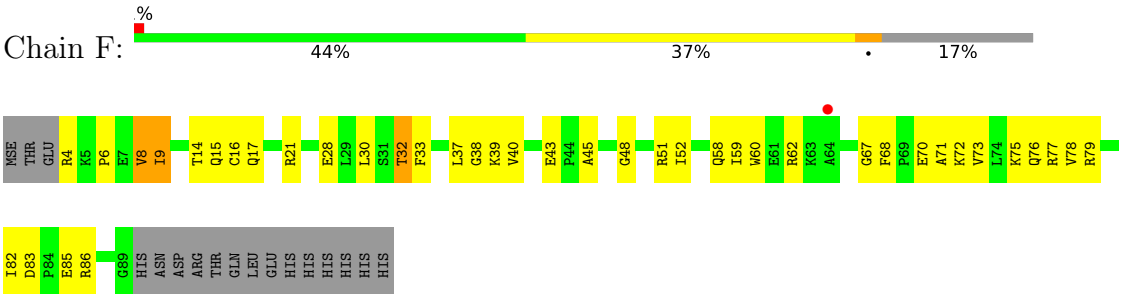
• Molecule 1: SelT/selW/selH selenoprotein domain

Chain E:



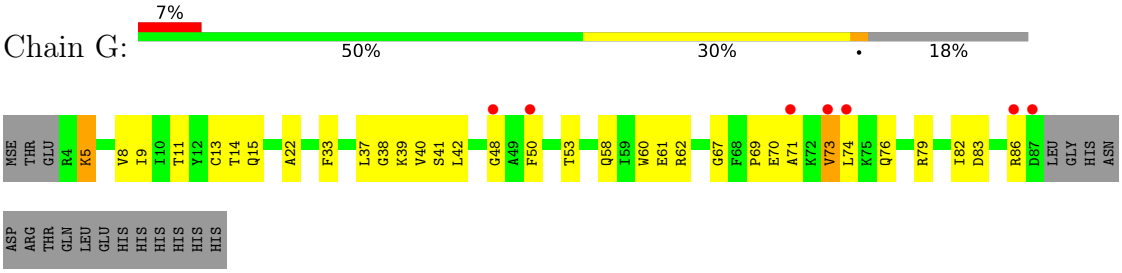
• Molecule 1: SelT/selW/selH selenoprotein domain

Chain F:



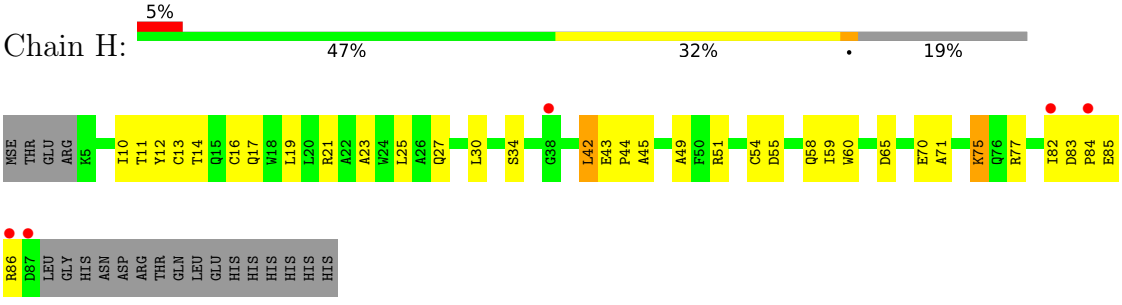
• Molecule 1: SelT/selW/selH selenoprotein domain

Chain G:



• Molecule 1: SelT/selW/selH selenoprotein domain

Chain H:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	54.31Å 112.25Å 148.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.98 – 2.70 19.98 – 2.70	Depositor EDS
% Data completeness (in resolution range)	78.8 (19.98-2.70) 87.7 (19.98-2.70)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.92 (at 2.69Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.215 , 0.292 0.232 , 0.307	Depositor DCC
R_{free} test set	2288 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å^2)	41.2	Xtriage
Anisotropy	0.794	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 51.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	5443	wwPDB-VP
Average B, all atoms (Å^2)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 20.52 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.5427e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

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	0.51	0/683	0.97	2/924 (0.2%)
1	B	0.50	0/691	0.98	3/935 (0.3%)
1	C	0.44	0/664	0.95	3/899 (0.3%)
1	D	0.47	0/683	0.97	5/924 (0.5%)
1	E	0.49	0/644	0.99	3/873 (0.3%)
1	F	0.46	0/695	0.96	2/940 (0.2%)
1	G	0.38	0/683	0.99	4/924 (0.4%)
1	H	0.43	0/672	0.87	2/910 (0.2%)
All	All	0.46	0/5415	0.96	24/7329 (0.3%)

There are no bond length outliers.

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	G	67	GLY	N-CA-C	7.06	118.69	111.56
1	B	82	ILE	N-CA-C	6.82	117.42	111.90
1	G	5	LYS	CA-C-N	6.59	126.61	119.89
1	G	5	LYS	C-N-CA	6.59	126.61	119.89
1	F	70	GLU	N-CA-C	-6.56	101.70	110.55
1	F	86	ARG	N-CA-C	6.14	118.73	108.96
1	A	70	GLU	N-CA-C	-6.12	100.59	110.32
1	G	73	VAL	N-CA-C	5.92	116.38	110.23
1	C	16	CYS	N-CA-C	-5.75	106.22	113.18
1	B	82	ILE	CB-CA-C	-5.74	105.79	111.30
1	A	28	GLU	N-CA-C	-5.70	104.97	111.07
1	D	70	GLU	N-CA-C	-5.70	101.26	110.32
1	E	70	GLU	N-CA-C	-5.51	101.56	110.32
1	C	82	ILE	N-CA-C	5.50	120.78	109.34
1	D	5	LYS	CA-C-N	5.47	125.48	119.90
1	D	5	LYS	C-N-CA	5.47	125.48	119.90
1	D	66	GLY	N-CA-C	5.28	122.64	115.30
1	D	14	THR	N-CA-C	5.25	116.69	111.07
1	C	73	VAL	N-CA-C	5.24	115.38	110.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	H	14	THR	N-CA-C	5.24	116.67	111.07
1	B	59	ILE	N-CA-C	-5.21	108.20	113.47
1	H	70	GLU	N-CA-C	-5.17	102.82	110.48
1	E	14	THR	N-CA-C	5.07	116.66	111.03
1	E	59	ILE	N-CA-C	-5.02	108.23	113.10

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	670	0	665	20	0
1	B	678	0	676	22	0
1	C	651	0	648	16	0
1	D	670	0	665	32	0
1	E	631	0	629	17	0
1	F	682	0	679	26	0
1	G	670	0	665	29	0
1	H	659	0	652	31	0
2	A	27	0	0	2	0
2	B	20	0	0	3	0
2	C	8	0	0	2	0
2	D	17	0	0	1	0
2	E	27	0	0	5	0
2	F	13	0	0	0	0
2	G	7	0	0	0	0
2	H	13	0	0	0	0
All	All	5443	0	5279	183	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 17.

All (183) 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:E:35:ASP:HB3	2:E:113:HOH:O	1.73	0.88
1:B:51:ARG:HH21	1:B:58:GLN:HE21	1.15	0.86
1:E:21:ARG:HD2	2:E:125:HOH:O	1.81	0.79
1:F:4:ARG:O	1:F:4:ARG:HD3	1.83	0.77
1:G:9:ILE:HG13	1:G:41:SER:HB2	1.68	0.76
1:B:51:ARG:NH2	1:B:58:GLN:HE21	1.84	0.75
1:A:23:ALA:O	1:A:27:GLN:HG3	1.88	0.74
1:F:51:ARG:HD2	1:F:58:GLN:OE1	1.88	0.73
1:F:28:GLU:O	1:F:32:THR:HG22	1.88	0.73
1:B:9:ILE:HB	1:B:53:THR:HG23	1.70	0.73
1:G:76:GLN:HG2	1:G:79:ARG:HH11	1.54	0.71
1:A:53:THR:HG22	1:A:58:GLN:HA	1.74	0.68
1:C:21:ARG:HD2	2:C:104:HOH:O	1.93	0.68
1:E:82:ILE:O	1:E:83:ASP:HB2	1.94	0.68
1:G:83:ASP:OD1	1:G:86:ARG:HB2	1.94	0.68
1:D:82:ILE:HG13	1:D:83:ASP:H	1.58	0.67
1:F:14:THR:HG23	1:F:15:GLN:H	1.60	0.67
1:B:60:TRP:CE3	1:B:69:PRO:HB3	2.31	0.65
1:G:73:VAL:HG13	1:G:74:LEU:H	1.61	0.65
1:C:15:GLN:H	1:C:15:GLN:CD	2.06	0.64
1:A:79:ARG:HH11	1:A:79:ARG:HB3	1.62	0.64
1:C:62:ARG:HH11	1:C:62:ARG:HB3	1.63	0.64
1:F:62:ARG:HG3	1:F:67:GLY:HA2	1.79	0.63
1:C:79:ARG:O	1:C:82:ILE:HG13	1.97	0.63
1:H:51:ARG:HE	1:H:58:GLN:HE21	1.46	0.62
1:A:4:ARG:HD3	1:A:5:LYS:H	1.63	0.62
1:H:82:ILE:O	1:H:84:PRO:HD3	2.00	0.62
1:E:9:ILE:HG12	1:E:53:THR:OG1	2.00	0.61
1:F:14:THR:HG22	1:F:45:ALA:O	2.00	0.61
1:H:11:THR:HG22	1:H:45:ALA:HB2	1.82	0.61
1:G:73:VAL:HG13	1:G:74:LEU:N	2.16	0.61
1:B:51:ARG:NH2	1:B:58:GLN:HB2	2.16	0.60
1:F:9:ILE:HD11	1:F:43:GLU:OE2	2.01	0.60
1:H:12:TYR:OH	1:H:19:LEU:HD13	2.01	0.60
1:A:76:GLN:HG2	1:A:79:ARG:HH12	1.67	0.60
1:F:73:VAL:O	1:F:77:ARG:HG3	2.03	0.59
1:C:18:TRP:HA	2:C:104:HOH:O	2.03	0.59
1:D:11:THR:HG22	1:D:45:ALA:HB2	1.84	0.59
1:G:39:LYS:HB3	1:H:43:GLU:HB3	1.85	0.59
1:D:84:PRO:HD3	2:D:113:HOH:O	2.03	0.59
1:H:71:ALA:O	1:H:75:LYS:HB2	2.03	0.58
1:D:61:GLU:HG3	1:D:64:ALA:H	1.68	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:58:GLN:NE2	1:G:61:GLU:HB2	2.19	0.58
1:F:16:CYS:SG	1:F:48:GLY:HA2	2.45	0.56
1:E:70:GLU:HB2	1:E:73:VAL:HG23	1.87	0.56
1:D:78:VAL:O	1:D:82:ILE:HG12	2.05	0.56
1:E:71:ALA:HB3	1:H:21:ARG:NH2	2.21	0.56
1:E:25:LEU:HD22	1:E:74:LEU:HD22	1.88	0.56
1:B:72:LYS:O	1:B:76:GLN:HG3	2.06	0.56
1:B:73:VAL:O	1:B:77:ARG:HG3	2.05	0.56
1:D:51:ARG:HH11	1:D:58:GLN:HB2	1.72	0.55
1:H:54:CYS:CB	1:H:59:ILE:HD11	2.36	0.55
1:A:9:ILE:HD11	1:A:43:GLU:HB3	1.89	0.54
1:C:16:CYS:O	1:C:17:GLN:HB2	2.08	0.54
1:A:7:GLU:HG2	1:A:39:LYS:HG2	1.89	0.54
1:A:18:TRP:HA	2:A:106:HOH:O	2.07	0.53
1:D:82:ILE:HG13	1:D:83:ASP:N	2.22	0.53
1:F:62:ARG:HA	1:F:67:GLY:O	2.09	0.53
1:F:14:THR:HG23	1:F:15:GLN:N	2.24	0.53
1:F:60:TRP:NE1	1:F:67:GLY:O	2.41	0.53
1:G:22:ALA:HB1	1:G:42:LEU:HD22	1.89	0.53
1:H:54:CYS:HB2	1:H:59:ILE:HD11	1.91	0.53
1:B:16:CYS:SG	1:B:48:GLY:HA2	2.50	0.51
1:H:21:ARG:HH11	1:H:21:ARG:HG3	1.76	0.51
1:D:25:LEU:HD22	1:D:74:LEU:HD23	1.93	0.51
1:H:59:ILE:HG23	1:H:77:ARG:HB3	1.91	0.51
1:A:86:ARG:HG2	1:A:86:ARG:HH11	1.76	0.50
1:E:53:THR:HG22	1:E:58:GLN:HA	1.92	0.50
1:E:54:CYS:O	1:E:55:ASP:HB2	2.09	0.50
1:B:51:ARG:HH21	1:B:58:GLN:NE2	1.95	0.50
1:E:9:ILE:O	1:E:9:ILE:HG13	2.11	0.50
1:F:8:VAL:HG22	1:F:40:VAL:HG13	1.94	0.50
1:B:21:ARG:HG3	2:B:104:HOH:O	2.12	0.50
1:H:43:GLU:O	1:H:43:GLU:HG3	2.10	0.50
1:B:76:GLN:HB3	2:B:106:HOH:O	2.12	0.49
1:E:78:VAL:O	1:E:82:ILE:HG13	2.11	0.49
1:A:4:ARG:CD	1:A:5:LYS:H	2.25	0.49
1:B:51:ARG:HE	1:B:58:GLN:NE2	2.10	0.49
1:D:71:ALA:O	1:D:75:LYS:HG3	2.13	0.49
1:G:8:VAL:C	1:G:9:ILE:HD12	2.38	0.49
1:D:11:THR:O	1:D:50:PHE:HA	2.11	0.49
1:A:9:ILE:HD11	1:A:43:GLU:CB	2.43	0.49
1:D:29:LEU:HD21	1:D:74:LEU:HG	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:11:THR:HG23	1:H:43:GLU:HG3	1.93	0.49
1:F:16:CYS:O	1:F:17:GLN:HB2	2.13	0.48
1:D:10:ILE:HG12	1:D:52:ILE:HG12	1.94	0.48
1:B:61:GLU:HB3	1:B:64:ALA:HB3	1.95	0.48
1:D:9:ILE:HD11	1:D:43:GLU:HB2	1.95	0.48
1:D:13:CYS:HB2	1:D:48:GLY:HA2	1.96	0.48
1:B:54:CYS:O	1:B:55:ASP:HB2	2.13	0.48
1:C:41:SER:HB3	1:D:39:LYS:HD2	1.95	0.47
1:F:79:ARG:HG2	1:F:83:ASP:O	2.13	0.47
1:G:11:THR:O	1:G:50:PHE:HA	2.14	0.47
1:C:71:ALA:O	1:C:75:LYS:HG3	2.15	0.47
1:D:51:ARG:NH1	1:D:58:GLN:HB2	2.28	0.47
1:B:54:CYS:SG	1:B:82:ILE:HD11	2.55	0.46
1:F:38:GLY:O	1:F:39:LYS:HB2	2.15	0.46
1:H:59:ILE:CG2	1:H:77:ARG:HB3	2.46	0.46
1:H:60:TRP:HA	1:H:65:ASP:OD2	2.15	0.46
1:B:29:LEU:HD21	1:B:74:LEU:HG	1.97	0.46
1:C:29:LEU:HD22	1:C:78:VAL:HG21	1.97	0.46
1:D:86:ARG:NH1	1:D:86:ARG:HB3	2.31	0.46
1:E:23:ALA:O	1:E:27:GLN:HG3	2.15	0.46
1:G:8:VAL:O	1:G:40:VAL:HA	2.16	0.46
1:A:86:ARG:HG2	1:A:86:ARG:NH1	2.31	0.46
1:B:39:LYS:HB3	1:B:39:LYS:HE2	1.77	0.46
1:F:33:PHE:HB3	1:F:37:LEU:HG	1.98	0.46
1:G:76:GLN:CG	1:G:79:ARG:HH11	2.26	0.45
1:D:9:ILE:HG22	1:D:53:THR:OG1	2.16	0.45
1:G:76:GLN:HA	1:G:79:ARG:HB3	1.98	0.45
1:D:53:THR:HG22	1:D:58:GLN:HA	1.98	0.45
1:C:61:GLU:OE1	1:C:63:LYS:HB2	2.16	0.45
2:E:124:HOH:O	1:F:14:THR:HG21	2.16	0.45
1:E:54:CYS:HB3	1:E:81:GLN:HG2	1.98	0.45
1:H:13:CYS:HB2	1:H:49:ALA:H	1.82	0.45
1:G:8:VAL:O	1:G:9:ILE:HD12	2.17	0.45
1:D:62:ARG:HA	1:D:67:GLY:O	2.17	0.45
1:F:8:VAL:HG13	1:F:37:LEU:HD22	1.98	0.45
1:G:82:ILE:HG22	1:G:82:ILE:O	2.16	0.45
1:A:63:LYS:HB2	1:A:63:LYS:NZ	2.32	0.44
1:A:70:GLU:HG3	2:A:120:HOH:O	2.17	0.44
1:D:25:LEU:HD22	1:D:74:LEU:CD2	2.48	0.44
1:F:52:ILE:O	1:F:59:ILE:HG12	2.18	0.44
1:B:50:PHE:HB3	1:B:62:ARG:HB2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:50:PHE:HB3	1:C:62:ARG:HB2	2.00	0.44
1:G:83:ASP:CG	1:G:86:ARG:HB2	2.43	0.44
1:D:28:GLU:HB3	1:D:75:LYS:HZ3	1.83	0.44
1:B:16:CYS:O	1:B:17:GLN:HB2	2.18	0.43
1:G:42:LEU:HB2	1:H:30:LEU:HD11	2.00	0.43
1:H:83:ASP:OD1	1:H:86:ARG:HB2	2.18	0.43
1:A:71:ALA:O	1:A:75:LYS:HG3	2.18	0.43
1:D:9:ILE:HD11	1:D:43:GLU:CB	2.49	0.43
1:G:60:TRP:NE1	1:G:69:PRO:HD3	2.32	0.43
1:F:72:LYS:O	1:F:76:GLN:HG3	2.19	0.43
1:G:61:GLU:O	1:G:62:ARG:C	2.61	0.43
1:D:4:ARG:HH11	1:D:4:ARG:HG2	1.84	0.43
1:A:9:ILE:CG2	1:A:53:THR:OG1	2.67	0.43
1:G:38:GLY:O	1:H:44:PRO:HD2	2.18	0.43
1:H:54:CYS:HB3	1:H:59:ILE:HD11	2.00	0.43
1:G:33:PHE:HD2	1:G:37:LEU:HD11	1.84	0.42
1:D:54:CYS:SG	1:D:81:GLN:HG2	2.59	0.42
1:F:30:LEU:HD21	1:F:40:VAL:HG23	2.01	0.42
1:B:81:GLN:HG2	2:B:118:HOH:O	2.19	0.42
1:C:29:LEU:CD2	1:C:78:VAL:HG21	2.50	0.42
1:F:21:ARG:NH1	1:F:68:PHE:CZ	2.88	0.42
1:F:6:PRO:HG3	1:F:82:ILE:HG21	2.02	0.42
1:H:16:CYS:O	1:H:17:GLN:HB2	2.20	0.42
1:E:49:ALA:HA	2:E:120:HOH:O	2.20	0.42
1:D:73:VAL:O	1:D:76:GLN:HB2	2.19	0.42
1:H:54:CYS:O	1:H:55:ASP:HB2	2.20	0.42
1:D:86:ARG:HB3	1:D:86:ARG:CZ	2.50	0.41
1:G:70:GLU:H	1:G:73:VAL:HG11	1.85	0.41
1:H:23:ALA:O	1:H:27:GLN:HG3	2.20	0.41
1:D:38:GLY:O	1:D:39:LYS:HB3	2.20	0.41
1:F:71:ALA:O	1:F:75:LYS:HG3	2.20	0.41
1:H:83:ASP:CG	1:H:86:ARG:HB2	2.45	0.41
1:E:70:GLU:HG3	2:E:119:HOH:O	2.19	0.41
1:G:39:LYS:HB3	1:H:43:GLU:CB	2.48	0.41
1:B:14:THR:HG21	1:B:44:PRO:HB3	2.03	0.41
1:C:15:GLN:HE22	1:C:46:THR:HG22	1.86	0.41
1:D:30:LEU:HD21	1:D:40:VAL:HG23	2.02	0.41
1:G:14:THR:HG21	1:H:34:SER:HB2	2.03	0.41
1:G:53:THR:HG22	1:G:58:GLN:HA	2.03	0.41
1:C:82:ILE:HD12	1:C:83:ASP:N	2.36	0.41
1:A:6:PRO:HG2	1:A:37:LEU:HA	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:70:GLU:OE1	1:E:72:LYS:HE2	2.20	0.41
1:H:43:GLU:HA	1:H:44:PRO:HD3	1.99	0.41
1:A:60:TRP:CE3	1:A:69:PRO:HB3	2.55	0.41
1:D:20:LEU:HD12	1:D:20:LEU:HA	1.92	0.41
1:F:78:VAL:O	1:F:82:ILE:HG12	2.20	0.41
1:G:14:THR:CG2	1:H:34:SER:HB2	2.51	0.41
1:G:70:GLU:HG3	1:G:73:VAL:HG12	2.02	0.41
1:A:6:PRO:HA	1:A:55:ASP:OD1	2.21	0.41
1:A:30:LEU:HD22	1:B:44:PRO:HG3	2.02	0.40
1:H:21:ARG:HG3	1:H:21:ARG:NH1	2.36	0.40
1:H:25:LEU:HD23	1:H:25:LEU:HA	1.92	0.40
1:C:42:LEU:HB2	1:D:30:LEU:HD11	2.04	0.40
1:G:33:PHE:HB2	1:G:37:LEU:HD12	2.03	0.40
1:C:15:GLN:NE2	1:C:46:THR:HG22	2.37	0.40
1:D:11:THR:HA	1:D:43:GLU:O	2.22	0.40
1:H:10:ILE:HB	1:H:42:LEU:HD12	2.03	0.40
1:E:60:TRP:HA	1:E:65:ASP:OD2	2.21	0.40
1:G:50:PHE:HB3	1:G:62:ARG:HB2	2.03	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	82/103 (80%)	71 (87%)	9 (11%)	2 (2%)	4 12
1	B	83/103 (81%)	78 (94%)	5 (6%)	0	100 100
1	C	80/103 (78%)	63 (79%)	16 (20%)	1 (1%)	9 25
1	D	82/103 (80%)	76 (93%)	6 (7%)	0	100 100
1	E	78/103 (76%)	71 (91%)	5 (6%)	2 (3%)	4 11
1	F	84/103 (82%)	73 (87%)	11 (13%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	G	82/103 (80%)	62 (76%)	17 (21%)	3 (4%)	2 6
1	H	81/103 (79%)	76 (94%)	5 (6%)	0	100 100
All	All	652/824 (79%)	570 (87%)	74 (11%)	8 (1%)	10 27

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	83	ASP
1	G	71	ALA
1	A	62	ARG
1	E	62	ARG
1	G	5	LYS
1	C	48	GLY
1	G	48	GLY
1	A	82	ILE

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	71/88 (81%)	68 (96%)	3 (4%)	26 55
1	B	72/88 (82%)	68 (94%)	4 (6%)	19 44
1	C	69/88 (78%)	68 (99%)	1 (1%)	59 82
1	D	71/88 (81%)	66 (93%)	5 (7%)	14 34
1	E	67/88 (76%)	66 (98%)	1 (2%)	57 81
1	F	72/88 (82%)	68 (94%)	4 (6%)	19 44
1	G	71/88 (81%)	69 (97%)	2 (3%)	38 68
1	H	70/88 (80%)	67 (96%)	3 (4%)	26 54
All	All	563/704 (80%)	540 (96%)	23 (4%)	27 56

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ARG
1	A	9	ILE
1	A	79	ARG
1	B	20	LEU
1	B	42	LEU
1	B	53	THR
1	B	79	ARG
1	C	62	ARG
1	D	14	THR
1	D	20	LEU
1	D	36	ASP
1	D	42	LEU
1	D	51	ARG
1	E	62	ARG
1	F	8	VAL
1	F	9	ILE
1	F	32	THR
1	F	85	GLU
1	G	13	CYS
1	G	15	GLN
1	H	42	LEU
1	H	75	LYS
1	H	85	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	76	GLN
1	A	81	GLN
1	B	15	GLN
1	B	58	GLN
1	B	81	GLN
1	C	15	GLN
1	C	81	GLN
1	D	76	GLN
1	D	81	GLN
1	E	15	GLN
1	E	27	GLN
1	E	81	GLN
1	F	17	GLN
1	F	76	GLN
1	F	81	GLN

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Mol	Chain	Res	Type
1	G	17	GLN
1	G	81	GLN
1	H	58	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 oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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^{th} 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	84/103 (81%)	-0.12	1 (1%)	76	75	12, 33, 60, 78	0
1	B	85/103 (82%)	-0.16	1 (1%)	76	75	20, 34, 47, 70	0
1	C	82/103 (79%)	0.20	2 (2%)	59	56	25, 45, 67, 83	0
1	D	84/103 (81%)	0.22	2 (2%)	59	56	15, 41, 78, 111	0
1	E	80/103 (77%)	-0.16	1 (1%)	75	73	19, 35, 50, 72	0
1	F	86/103 (83%)	0.13	1 (1%)	76	75	26, 42, 59, 72	0
1	G	84/103 (81%)	0.99	7 (8%)	17	14	38, 69, 106, 122	0
1	H	83/103 (80%)	0.58	5 (6%)	27	24	17, 57, 105, 122	0
All	All	668/824 (81%)	0.21	20 (2%)	52	49	12, 42, 79, 122	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	87	ASP	3.3
1	H	84	PRO	3.0
1	F	64	ALA	3.0
1	C	82	ILE	2.9
1	D	82	ILE	2.7
1	G	87	ASP	2.7
1	E	84	PRO	2.5
1	H	82	ILE	2.5
1	G	71	ALA	2.4
1	D	84	PRO	2.4
1	C	84	PRO	2.4
1	G	48	GLY	2.3
1	H	38	GLY	2.3
1	G	74	LEU	2.2
1	A	86	ARG	2.2
1	B	48	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	H	86	ARG	2.0
1	G	50	PHE	2.0
1	G	86	ARG	2.0
1	G	73	VAL	2.0

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 oligosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.