



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 23, 2017 – 07:59 PM EDT

PDB ID : 3GTY
Title : Promiscuous Substrate Recognition in Folding and Assembly Activities of the Trigger Factor Chaperone
Authors : Martinez-Hackert, E.; Hendrickson, W.A.
Deposited on : unknown
Resolution : 3.40 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
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030345
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)	:	rb-20030345

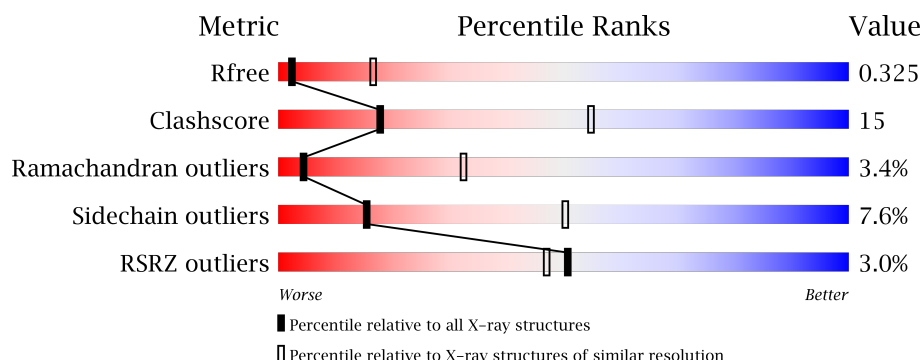
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 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	1679 (3.50-3.30)
Clashscore	112137	1832 (3.50-3.30)
Ramachandran outliers	110173	1789 (3.50-3.30)
Sidechain outliers	110143	1789 (3.50-3.30)
RSRZ outliers	101464	1709 (3.50-3.30)

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	X	433	
2	S	149	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4610 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 Trigger factor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	X	412	Total	C	N	O	S	0	0	0
			3405	2147	575	677	6			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
X	426	LEU	-	EXPRESSION TAG	UNP Q9WZF8
X	427	GLU	-	EXPRESSION TAG	UNP Q9WZF8
X	428	HIS	-	EXPRESSION TAG	UNP Q9WZF8
X	429	HIS	-	EXPRESSION TAG	UNP Q9WZF8
X	430	HIS	-	EXPRESSION TAG	UNP Q9WZF8
X	431	HIS	-	EXPRESSION TAG	UNP Q9WZF8
X	432	HIS	-	EXPRESSION TAG	UNP Q9WZF8
X	433	HIS	-	EXPRESSION TAG	UNP Q9WZF8

- Molecule 2 is a protein called 30S ribosomal protein S7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	S	149	Total	C	N	O	S	0	0	0
			1205	767	226	208	4			

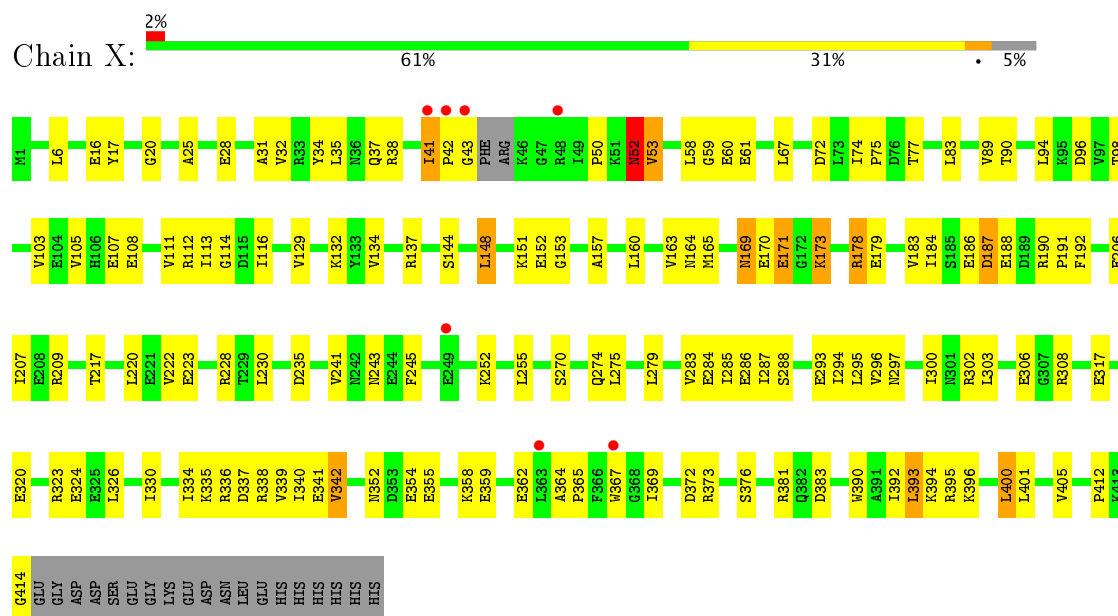
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
S	7	MET	-	EXPRESSION TAG	UNP P38526
S	8	GLY	-	EXPRESSION TAG	UNP P38526

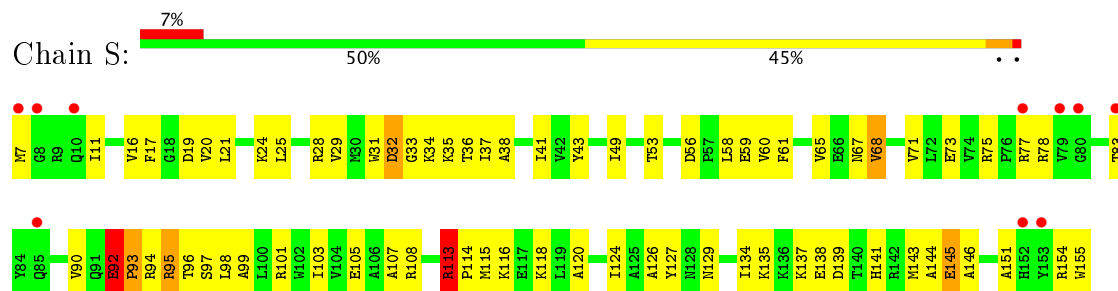
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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 ($\text{RSRZ} > 2$). 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: Trigger factor



• Molecule 2: 30S ribosomal protein S7



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	94.16 Å 94.16 Å 193.07 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	12.00 – 3.40 39.08 – 3.31	Depositor EDS
% Data completeness (in resolution range)	100.0 (12.00-3.40) 97.0 (39.08-3.31)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	63.10 (at 3.32 Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.257 , 0.329 0.250 , 0.325	Depositor DCC
R_{free} test set	620 reflections (5.46%)	DCC
Wilson B-factor (Å ²)	83.5	Xtriage
Anisotropy	0.240	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 1.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	4610	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.15% 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 [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	X	0.77	4/3443 (0.1%)	0.72	0/4617
2	S	0.59	1/1229 (0.1%)	0.61	0/1655
All	All	0.73	5/4672 (0.1%)	0.69	0/6272

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	X	43	GLY	C-O	7.46	1.35	1.23
2	S	145	GLU	CD-OE1	6.43	1.32	1.25
1	X	317	GLU	CD-OE1	6.23	1.32	1.25
1	X	107	GLU	CG-CD	5.35	1.59	1.51
1	X	37	GLN	CD-OE1	5.10	1.35	1.24

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	X	3405	0	3480	85	0
2	S	1205	0	1257	62	0
All	All	4610	0	4737	142	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 15.

All (142) 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:X:169:ASN:HD22	1:X:170:GLU:H	0.97	0.94
1:X:169:ASN:HD22	1:X:170:GLU:N	1.73	0.85
2:S:92:GLU:HB3	2:S:93:PRO:HD2	1.57	0.85
2:S:95:ARG:HH12	2:S:96:THR:HB	1.42	0.84
2:S:120:ALA:O	2:S:124:ILE:HG22	1.79	0.82
1:X:111:VAL:HA	1:X:285:ILE:HG22	1.60	0.81
2:S:135:LYS:O	2:S:139:ASP:HB2	1.85	0.76
2:S:107:ALA:O	2:S:118:LYS:HD3	1.87	0.74
2:S:16:VAL:HB	2:S:43:TYR:HE1	1.51	0.74
2:S:95:ARG:NH1	2:S:96:THR:HB	2.02	0.74
2:S:92:GLU:HB3	2:S:93:PRO:CD	2.17	0.74
1:X:381:ARG:HH12	2:S:145:GLU:HB3	1.54	0.73
1:X:178:ARG:HH11	1:X:178:ARG:HG3	1.54	0.72
2:S:21:LEU:HD21	2:S:65:VAL:HG11	1.70	0.72
2:S:16:VAL:HB	2:S:43:TYR:CE1	2.25	0.72
1:X:306:GLU:HB2	1:X:308:ARG:HG3	1.72	0.70
1:X:338:ARG:O	1:X:341:GLU:HG2	1.92	0.69
1:X:52:ASN:HD22	1:X:52:ASN:C	1.97	0.68
1:X:183:VAL:HG12	1:X:228:ARG:HD2	1.76	0.68
1:X:17:TYR:CZ	1:X:74:ILE:HG23	2.30	0.66
1:X:338:ARG:HD2	1:X:341:GLU:OE2	1.95	0.66
2:S:29:VAL:HG13	2:S:38:ALA:HB1	1.76	0.66
1:X:151:LYS:HG2	1:X:153:GLY:H	1.61	0.65
2:S:113:ARG:HB2	2:S:114:PRO:HD3	1.79	0.65
1:X:230:LEU:HD12	1:X:230:LEU:H	1.61	0.64
2:S:143:MET:HA	2:S:146:ALA:HB3	1.78	0.64
1:X:72:ASP:O	1:X:75:PRO:HD2	1.97	0.64
1:X:326:LEU:O	1:X:330:ILE:HG13	1.99	0.62
1:X:393:LEU:O	1:X:396:LYS:HB2	1.99	0.62
2:S:126:ALA:HA	2:S:134:ILE:HD11	1.83	0.61
1:X:283:VAL:HG21	1:X:339:VAL:HG21	1.82	0.60
1:X:340:ILE:HD13	1:X:394:LYS:HG3	1.84	0.60
2:S:94:ARG:HG3	2:S:94:ARG:HH11	1.67	0.59
2:S:19:ASP:OD2	2:S:21:LEU:HB3	2.03	0.58
1:X:381:ARG:HE	1:X:383:ASP:HB3	1.68	0.58
1:X:293:GLU:HG2	1:X:323:ARG:NH2	2.18	0.57
1:X:169:ASN:HB2	1:X:173:LYS:O	2.05	0.57
1:X:165:MET:HG3	1:X:220:LEU:HD12	1.87	0.56
2:S:113:ARG:HE	2:S:113:ARG:N	2.03	0.56
1:X:112:ARG:HH12	1:X:286:GLU:HG2	1.71	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:191:PRO:O	1:X:192:PHE:HB2	2.06	0.56
1:X:354:GLU:O	1:X:358:LYS:HG3	2.05	0.56
2:S:143:MET:HA	2:S:146:ALA:CB	2.35	0.55
1:X:20:GLY:HA2	1:X:98:THR:HA	1.88	0.55
1:X:137:ARG:HG2	1:X:255:LEU:HD22	1.89	0.54
1:X:20:GLY:CA	1:X:98:THR:HA	2.38	0.54
2:S:68:VAL:HG21	2:S:103:ILE:HD11	1.91	0.53
1:X:165:MET:HG3	1:X:220:LEU:CD1	2.39	0.52
1:X:414:GLY:C	2:S:113:ARG:HB3	2.29	0.52
1:X:392:ILE:O	1:X:395:ARG:HG2	2.08	0.52
1:X:163:VAL:HG12	1:X:222:VAL:HA	1.91	0.52
2:S:155:TRP:CE3	2:S:155:TRP:HA	2.45	0.52
1:X:83:LEU:HD21	1:X:105:VAL:HG13	1.92	0.51
2:S:35:LYS:HG2	2:S:36:THR:H	1.74	0.51
1:X:338:ARG:HA	1:X:341:GLU:HG2	1.93	0.51
2:S:113:ARG:HB2	2:S:114:PRO:CD	2.41	0.50
1:X:89:VAL:HG22	1:X:103:VAL:HG12	1.93	0.50
1:X:270:SER:O	1:X:274:GLN:HG2	2.12	0.50
2:S:93:PRO:O	2:S:96:THR:HG22	2.12	0.50
1:X:157:ALA:HA	1:X:184:ILE:HB	1.92	0.50
1:X:165:MET:SD	1:X:220:LEU:HD11	2.52	0.49
2:S:61:PHE:O	2:S:65:VAL:HG12	2.12	0.49
1:X:300:ILE:O	1:X:303:LEU:HB2	2.12	0.49
1:X:28:GLU:O	1:X:31:ALA:HB3	2.13	0.49
1:X:337:ASP:HA	1:X:340:ILE:HD12	1.94	0.49
1:X:364:ALA:N	1:X:365:PRO:HD2	2.27	0.49
2:S:77:ARG:HG2	2:S:78:ARG:H	1.78	0.49
1:X:164:ASN:HB2	1:X:179:GLU:HG2	1.95	0.49
1:X:17:TYR:CE1	1:X:74:ILE:HG23	2.48	0.48
1:X:59:GLY:O	1:X:60:GLU:HB3	2.14	0.48
2:S:28:ARG:HH11	2:S:28:ARG:HG3	1.78	0.48
2:S:49:ILE:HD11	2:S:120:ALA:HA	1.96	0.48
2:S:124:ILE:O	2:S:127:TYR:HB3	2.13	0.48
2:S:11:ILE:HD11	2:S:24:LYS:HG2	1.95	0.47
2:S:56:ASP:OD2	2:S:59:GLU:HG2	2.14	0.47
2:S:73:GLU:HG2	2:S:90:VAL:CG2	2.44	0.47
2:S:75:ARG:NH2	2:S:94:ARG:HH22	2.13	0.47
1:X:352:ASN:OD1	1:X:355:GLU:HB2	2.15	0.46
1:X:367:TRP:CD1	2:S:143:MET:HB3	2.50	0.46
1:X:134:VAL:HG11	1:X:252:LYS:HG2	1.97	0.46
2:S:41:ILE:HG23	2:S:116:LYS:HA	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:369:ILE:HD12	1:X:373:ARG:HB3	1.99	0.45
1:X:320:GLU:O	1:X:324:GLU:HB2	2.17	0.45
1:X:294:ILE:O	1:X:297:ASN:N	2.49	0.45
2:S:108:ARG:HG2	2:S:115:MET:HE3	1.99	0.45
2:S:65:VAL:HA	2:S:68:VAL:CG2	2.47	0.45
2:S:92:GLU:CB	2:S:93:PRO:CD	2.90	0.45
2:S:101:ARG:O	2:S:105:GLU:HB2	2.16	0.45
1:X:52:ASN:ND2	1:X:52:ASN:C	2.69	0.45
2:S:141:HIS:O	2:S:144:ALA:HB3	2.17	0.45
1:X:28:GLU:O	1:X:32:VAL:HG23	2.17	0.44
1:X:293:GLU:O	1:X:296:VAL:HB	2.16	0.44
2:S:113:ARG:NE	2:S:114:PRO:HD3	2.32	0.44
1:X:60:GLU:O	1:X:61:GLU:C	2.56	0.44
2:S:60:VAL:HG13	2:S:127:TYR:CD1	2.53	0.44
1:X:169:ASN:HB3	1:X:171:GLU:OE1	2.18	0.44
1:X:275:LEU:HD13	1:X:405:VAL:HG11	2.00	0.44
1:X:283:VAL:HG12	1:X:284:GLU:N	2.32	0.44
1:X:394:LYS:C	1:X:396:LYS:H	2.20	0.44
1:X:52:ASN:O	1:X:53:VAL:C	2.56	0.43
1:X:412:PRO:HA	2:S:113:ARG:HD3	2.00	0.43
2:S:71:VAL:O	2:S:90:VAL:N	2.47	0.43
1:X:334:ILE:O	1:X:335:LYS:C	2.57	0.43
2:S:134:ILE:O	2:S:137:LYS:HB3	2.19	0.43
1:X:279:LEU:O	1:X:283:VAL:HG23	2.19	0.43
2:S:92:GLU:O	2:S:94:ARG:N	2.51	0.43
1:X:206:GLU:C	1:X:207:ILE:HG13	2.38	0.43
2:S:113:ARG:CB	2:S:114:PRO:HD3	2.46	0.42
2:S:68:VAL:HG23	2:S:99:ALA:HB1	1.99	0.42
2:S:77:ARG:HG2	2:S:78:ARG:N	2.33	0.42
2:S:17:PHE:HB3	2:S:58:LEU:HD22	2.00	0.42
2:S:25:LEU:HD22	2:S:61:PHE:CZ	2.54	0.42
1:X:338:ARG:O	1:X:342:VAL:HG23	2.18	0.42
2:S:151:ALA:O	2:S:154:ARG:HB2	2.20	0.42
1:X:132:LYS:HA	1:X:132:LYS:HD3	1.71	0.42
1:X:178:ARG:HH11	1:X:178:ARG:CG	2.29	0.42
1:X:359:GLU:O	1:X:362:GLU:HB2	2.20	0.42
1:X:67:LEU:HD21	1:X:94:LEU:HB2	2.02	0.42
1:X:148:LEU:HD23	1:X:160:LEU:HD13	2.01	0.41
2:S:94:ARG:NH1	2:S:98:LEU:HD13	2.35	0.41
1:X:74:ILE:H	1:X:75:PRO:CD	2.34	0.41
2:S:29:VAL:O	2:S:108:ARG:NH2	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:116:ILE:HG12	1:X:401:LEU:HD22	2.01	0.41
1:X:112:ARG:NH1	1:X:286:GLU:HG2	2.35	0.41
1:X:74:ILE:N	1:X:75:PRO:CD	2.84	0.41
1:X:41:ILE:HB	1:X:42:PRO:CD	2.51	0.41
2:S:20:VAL:O	2:S:24:LYS:HG3	2.20	0.41
1:X:6:LEU:CD1	1:X:16:GLU:HB2	2.51	0.41
2:S:31:TRP:O	2:S:32:ASP:HB3	2.20	0.41
2:S:75:ARG:HA	2:S:75:ARG:HD3	1.92	0.41
1:X:412:PRO:HB3	2:S:113:ARG:HG2	2.02	0.41
1:X:25:ALA:HA	1:X:28:GLU:HB2	2.02	0.41
1:X:6:LEU:HD11	1:X:16:GLU:HB2	2.02	0.41
1:X:191:PRO:O	1:X:192:PHE:CB	2.68	0.40
1:X:111:VAL:HG21	1:X:342:VAL:HG21	2.03	0.40
1:X:34:TYR:O	1:X:38:ARG:HG2	2.22	0.40
1:X:400:LEU:HD22	1:X:400:LEU:HA	1.91	0.40
2:S:138:GLU:HA	2:S:141:HIS:HB3	2.04	0.40
2:S:34:LYS:HB3	2:S:37:ILE:HD12	2.03	0.40
1:X:58:LEU:HA	1:X:58:LEU:HD23	1.94	0.40
2:S:67:ASN:O	2:S:137:LYS:HE3	2.21	0.40
1:X:178:ARG:NH1	1:X:178:ARG:HG3	2.28	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	X	408/433 (94%)	331 (81%)	63 (15%)	14 (3%)	4	32
2	S	147/149 (99%)	117 (80%)	25 (17%)	5 (3%)	4	32
All	All	555/582 (95%)	448 (81%)	88 (16%)	19 (3%)	4	32

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	X	53	VAL
1	X	113	ILE
1	X	186	GLU
1	X	173	LYS
1	X	243	ASN
2	S	92	GLU
2	S	113	ARG
1	X	50	PRO
1	X	52	ASN
1	X	187	ASP
1	X	295	LEU
1	X	90	THR
1	X	114	GLY
1	X	302	ARG
1	X	336	ARG
2	S	33	GLY
2	S	93	PRO
2	S	83	THR
1	X	342	VAL

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	X	379/398 (95%)	350 (92%)	29 (8%)	15	49
2	S	124/124 (100%)	115 (93%)	9 (7%)	16	52
All	All	503/522 (96%)	465 (92%)	38 (8%)	15	50

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	X	35	LEU
1	X	41	ILE
1	X	52	ASN
1	X	77	THR
1	X	96	ASP

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Mol	Chain	Res	Type
1	X	108	GLU
1	X	129	VAL
1	X	144	SER
1	X	148	LEU
1	X	152	GLU
1	X	169	ASN
1	X	171	GLU
1	X	178	ARG
1	X	187	ASP
1	X	188	GLU
1	X	190	ARG
1	X	209	ARG
1	X	217	THR
1	X	223	GLU
1	X	235	ASP
1	X	241	VAL
1	X	245	PHE
1	X	287	ILE
1	X	288	SER
1	X	372	ASP
1	X	376	SER
1	X	390	TRP
1	X	393	LEU
1	X	400	LEU
2	S	7	MET
2	S	32	ASP
2	S	53	THR
2	S	68	VAL
2	S	92	GLU
2	S	95	ARG
2	S	97	SER
2	S	113	ARG
2	S	129	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	X	52	ASN
1	X	169	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](#)

There are no ligands in this entry.

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	X	412/433 (95%)	-0.09	7 (1%) 70 66	2, 35, 87, 91	0
2	S	149/149 (100%)	0.44	10 (6%) 19 19	42, 71, 122, 139	0
All	All	561/582 (96%)	0.05	17 (3%) 51 47	2, 49, 91, 139	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	S	8	GLY	9.7
2	S	7	MET	8.4
1	X	41	ILE	5.5
2	S	80	GLY	4.7
1	X	42	PRO	3.4
2	S	10	GLN	3.2
2	S	79	VAL	3.2
2	S	153	TYR	3.1
2	S	83	THR	3.1
1	X	48	ARG	3.0
2	S	152	HIS	2.9
1	X	43	GLY	2.8
1	X	363	LEU	2.3
2	S	85	GLN	2.3
2	S	77	ARG	2.3
1	X	249	GLU	2.1
1	X	367	TRP	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 [i](#)

There are no carbohydrates 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.