



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 07:04 PM UTC

PDB ID : 3DW9 / pdb_00003dw9
Title : SgrAI with cognate DNA and manganese bound
Authors : Dunten, P.W.; Horton, N.C.; Little, E.J.
Deposited on : 2008-07-21
Resolution : 2.20 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

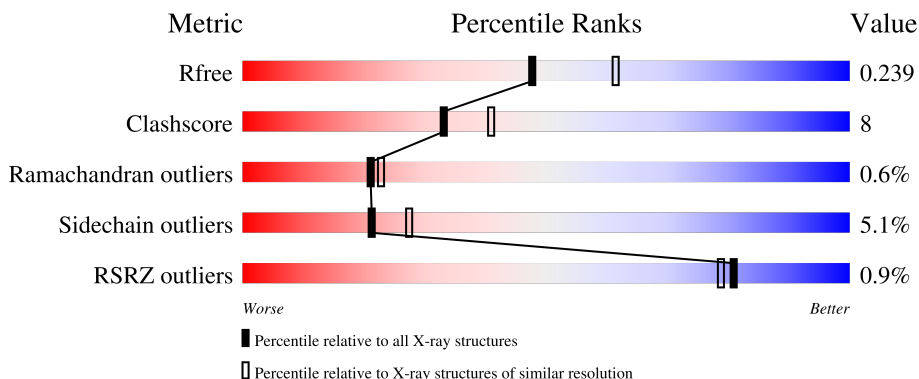
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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 $\leq 5\%$. 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	E	18	
1	F	18	
2	A	338	
2	B	338	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6335 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 DNA chain called DNA (5'-D(*DGP*DAP*DGP*DTP*DCP*DCP*DAP*DCP*DCP*DGP*DGP*DTP*DGP*DGP*DAP*DCP*DTP*DC)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	E	18	Total	C	N	O	P	0	0	0
			366	174	69	106	17			
1	F	18	Total	C	N	O	P	0	0	0
			366	174	69	106	17			

- Molecule 2 is a protein called SgrIR restriction enzyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	335	Total	C	N	O	S	0	0	0
			2639	1672	467	491	9			
2	B	334	Total	C	N	O	S	0	1	0
			2641	1672	468	492	9			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	63	ASP	ASN	cloning artifact	UNP Q9F6L0
B	63	ASP	ASN	cloning artifact	UNP Q9F6L0

- Molecule 3 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Mn	0	0
			2	2		
3	B	2	Total	Mn	0	0
			2	2		

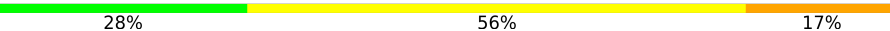
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	37	Total 37	O 37	0	0
4	F	30	Total 30	O 30	0	0
4	A	135	Total 135	O 135	0	0
4	B	117	Total 117	O 117	0	0

3 Residue-property plots

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: DNA (5'-D(*DGP*DAP*DGP*DTP*DCP*DCP*DAP*DCP*DCP*DGP*DGP*DT*P*DGP*DGP*DAP*DCP*DTP*DC)-3')

Chain E: 



- Molecule 1: DNA (5'-D(*DGP*DAP*DGP*DTP*DCP*DCP*DAP*DCP*DCP*DGP*DGP*DT*P*DGP*DGP*DAP*DCP*DTP*DC)-3')

Chain F: 



- Molecule 2: SgraIR restriction enzyme

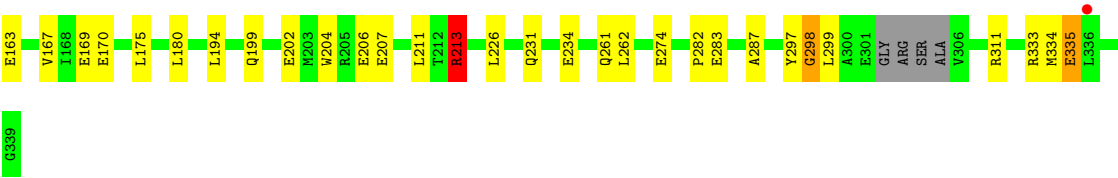
Chain A: 



- Molecule 2: SgraIR restriction enzyme

Chain B: 





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.26Å 117.10Å 63.70Å 90.00° 112.96° 90.00°	Depositor
Resolution (Å)	58.62 – 2.20 58.62 – 2.20	Depositor EDS
% Data completeness (in resolution range)	88.6 (58.62-2.20) 88.8 (58.62-2.20)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.5.0031	Depositor
R, R_{free}	0.174 , 0.240 0.182 , 0.239	Depositor DCC
R_{free} test set	1756 reflections (4.38%)	wwPDB-VP
Wilson B-factor (Å ²)	30.4	Xtriage
Anisotropy	0.709	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 36.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6335	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.31% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MN

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	E	0.91	0/410	1.79	13/631 (2.1%)
1	F	0.91	0/410	1.60	7/631 (1.1%)
2	A	1.17	4/2697 (0.1%)	1.13	5/3661 (0.1%)
2	B	1.16	2/2699 (0.1%)	1.08	3/3665 (0.1%)
All	All	1.14	6/6216 (0.1%)	1.21	28/8588 (0.3%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	51	ILE	N-CA	-6.39	1.41	1.46
2	A	54	PHE	CA-C	-6.18	1.45	1.52
2	A	70	ARG	C-O	5.96	1.30	1.24
2	B	108	ALA	CA-CB	5.80	1.62	1.53
2	B	287	ALA	N-CA	-5.63	1.38	1.46
2	A	92	ASN	N-CA	5.31	1.53	1.46

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	213	ARG	CA-C-N	-8.62	109.80	119.28
2	B	213	ARG	C-N-CA	-8.62	109.80	119.28
2	A	122	VAL	N-CA-C	-7.67	103.79	111.77
1	E	4	DT	P-O5'-C5'	-7.14	109.30	120.00
2	A	26	ALA	CA-C-N	6.99	127.54	120.14
2	A	26	ALA	C-N-CA	6.99	127.54	120.14
1	E	17	DT	P-O5'-C5'	6.97	130.46	120.00
1	F	10	DG	O3'-P-O5'	-6.78	93.84	104.00
1	E	3	DG	O4'-C1'-N9	6.43	118.04	108.40
1	E	3	DG	O5'-C5'-C4'	-6.27	101.39	110.80
1	F	15	DA	P-O3'-C3'	-6.25	110.83	120.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	12	DT	N1-C1'-C2'	6.11	122.66	113.50
2	A	231	GLN	CA-C-N	5.91	125.38	119.24
2	A	231	GLN	C-N-CA	5.91	125.38	119.24
1	E	13	DG	P-O3'-C3'	5.85	128.98	120.20
1	F	4	DT	O5'-C5'-C4'	-5.71	102.23	110.80
1	E	17	DT	O3'-P-O5'	-5.66	95.51	104.00
1	E	4	DT	C5'-C4'-O4'	-5.54	101.09	109.40
1	E	18	DC	C5'-C4'-C3'	-5.54	106.60	114.90
1	F	5	DC	C4'-C3'-O3'	-5.51	101.74	110.00
1	E	6	DC	C2'-C3'-O3'	5.49	119.73	111.50
1	E	7	DA	O4'-C1'-N9	-5.25	100.53	108.40
1	F	7	DA	C4'-C3'-O3'	-5.20	102.20	110.00
1	E	11	DG	O3'-P-O5'	-5.15	96.28	104.00
1	E	3	DG	O3'-P-O5'	-5.08	96.38	104.00
1	F	7	DA	P-O3'-C3'	5.05	127.77	120.20
1	F	4	DT	C1'-O4'-C4'	-5.05	102.13	109.70
2	B	261	GLN	N-CA-C	-5.01	105.53	111.69

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	366	0	203	5	0
1	F	366	0	203	1	0
2	A	2639	0	2596	51	0
2	B	2641	0	2594	40	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	135	0	0	1	0
4	B	117	0	0	9	0
4	E	37	0	0	0	0
4	F	30	0	0	0	0
All	All	6335	0	5596	90	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 8.

All (90) 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)
2:A:145:VAL:HG13	2:A:334:MET:HE2	1.52	0.90
2:A:145:VAL:CG1	2:A:334:MET:HE2	2.13	0.77
2:A:148:LEU:HD21	2:A:226:LEU:HD13	1.68	0.76
2:A:147:VAL:HG22	2:A:191:VAL:HG22	1.71	0.72
2:A:192:VAL:HG21	2:A:226:LEU:HD22	1.76	0.68
2:B:131:ARG:HD3	4:B:409:HOH:O	1.94	0.67
2:B:20:ILE:HD12	2:B:87:PHE:CE1	2.30	0.66
2:A:72:VAL:O	2:A:76:VAL:HG23	1.95	0.65
2:A:296:LEU:HD22	2:B:262:LEU:HD13	1.78	0.65
2:B:2:PRO:N	4:B:451:HOH:O	2.31	0.63
2:B:23:ILE:HD13	2:B:28:VAL:HG23	1.80	0.62
2:A:120:LEU:HD22	2:A:126:TRP:CE3	2.35	0.62
2:B:226:LEU:HD12	2:B:226:LEU:N	2.16	0.60
2:B:137:LEU:HD12	2:B:335:GLU:OE1	2.05	0.57
2:A:143:ARG:O	2:A:334:MET:HE3	2.05	0.56
2:A:312:GLU:OE2	2:A:325:ARG:NH1	2.30	0.56
2:B:143:ARG:O	2:B:334:MET:HE2	2.08	0.54
2:A:315:VAL:O	2:A:317:PRO:HD3	2.09	0.53
2:A:252:GLN:HB3	2:A:253:PRO:HD3	1.91	0.53
2:A:270:ALA:HB1	2:A:271:PRO:HD2	1.90	0.53
2:A:189:LEU:C	2:A:189:LEU:HD23	2.35	0.52
2:B:29:ARG:O	2:B:30:ASN:HB2	2.10	0.52
2:A:48:GLY:O	2:A:210:GLY:HA3	2.10	0.52
1:E:8:DC:H5''	4:B:347:HOH:O	2.11	0.51
2:A:52:PRO:O	2:A:109:VAL:HG22	2.10	0.51
1:F:11:DG:N3	2:B:92:ASN:HB3	2.25	0.51
2:A:126:TRP:CZ2	2:A:135:PRO:HD3	2.46	0.51
2:B:204:TRP:CH2	2:B:226:LEU:HD11	2.47	0.50
2:A:43:LEU:HD13	2:A:104:ILE:HG21	1.92	0.50
2:B:67:VAL:HG21	4:B:350:HOH:O	2.12	0.50
2:A:138:SER:O	2:A:143:ARG:NH1	2.45	0.50
2:B:213:ARG:HD2	4:B:404:HOH:O	2.12	0.49
1:E:11:DG:N3	2:A:92:ASN:HB3	2.26	0.49
2:A:20:ILE:HD12	2:A:87:PHE:HE1	1.78	0.49
2:B:72:VAL:HG23	4:B:361:HOH:O	2.11	0.49
2:B:120:LEU:CD2	2:B:135:PRO:HG2	2.43	0.49
2:A:254:LEU:O	2:A:258:ASN:ND2	2.41	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:115:ALA:HB3	2:B:207:GLU:HG2	1.95	0.48
2:B:24:ARG:HG2	2:B:34:GLN:HG2	1.94	0.48
2:A:232:PRO:HG3	2:A:268:LEU:HD13	1.96	0.48
2:A:163:GLU:O	2:A:167:VAL:HG23	2.14	0.47
2:A:272:LYS:HG3	2:A:307:HIS:CD2	2.48	0.47
2:A:132:TYR:O	2:A:133:SER:C	2.57	0.47
2:A:145:VAL:HG11	2:A:330:LEU:HD22	1.97	0.47
2:B:28:VAL:O	2:B:29:ARG:C	2.58	0.47
2:A:43:LEU:HD21	2:A:101:ILE:HG23	1.96	0.47
2:A:262:LEU:C	2:A:262:LEU:HD23	2.40	0.47
2:A:43:LEU:CD1	2:A:104:ILE:HG21	2.45	0.47
2:B:114:ALA:HB3	4:B:349:HOH:O	2.15	0.46
2:A:316:PRO:HA	2:A:317:PRO:HD2	1.79	0.46
2:A:185:SER:OG	2:A:186:THR:N	2.48	0.46
2:A:167:VAL:HB	2:A:268:LEU:HD21	1.97	0.46
2:A:37:LEU:HB2	2:A:97:VAL:HG21	1.98	0.45
2:B:20:ILE:HD12	2:B:87:PHE:HE1	1.78	0.45
2:B:274:GLU:HB3	2:B:311:ARG:HD2	1.98	0.45
2:B:121:MET:HA	2:B:121:MET:HE2	1.99	0.45
2:A:314:TYR:CE2	2:A:316:PRO:HB3	2.52	0.44
2:A:296:LEU:CD2	2:B:262:LEU:HD13	2.43	0.44
2:A:299:LEU:HD21	2:B:175:LEU:HD21	1.98	0.44
2:B:55:ALA:HB1	4:B:419:HOH:O	2.17	0.44
2:B:150:LEU:HD22	2:B:154:PHE:CE2	2.52	0.44
2:A:5:TYR:HB2	2:A:74:ASN:ND2	2.33	0.43
2:A:7:ILE:HD11	2:A:81:HIS:NE2	2.34	0.43
2:B:231:GLN:O	2:B:234:GLU:HB2	2.19	0.43
2:A:81:HIS:HE1	4:A:369:HOH:O	2.01	0.43
2:B:117:TRP:CH2	2:B:137:LEU:HD22	2.54	0.43
2:B:152:ARG:O	2:B:153:SER:HB2	2.18	0.43
2:A:126:TRP:CE2	2:A:135:PRO:HD3	2.54	0.42
2:A:62:MET:HE1	2:A:70:ARG:NH2	2.34	0.42
2:A:272:LYS:CG	2:A:307:HIS:CD2	3.02	0.42
2:B:297:TYR:O	2:B:298:GLY:C	2.62	0.42
2:A:306:VAL:HG12	2:B:180:LEU:HD21	2.01	0.42
2:B:163:GLU:O	2:B:167:VAL:HG23	2.20	0.42
2:A:65:ARG:HA	2:A:280:LEU:HD13	2.01	0.42
1:E:7:DA:H1'	1:E:8:DC:C6	2.54	0.42
2:A:20:ILE:HD12	2:A:87:PHE:CE1	2.55	0.42
2:B:23:ILE:CD1	2:B:28:VAL:HG23	2.48	0.42
2:B:97:VAL:HG12	2:B:101:ILE:HD12	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:70:ARG:HB3	2:B:71:PRO:HD3	2.00	0.41
2:B:108:ALA:HB2	2:B:211:LEU:HD21	2.03	0.41
2:A:7:ILE:HD12	2:A:7:ILE:HG21	1.90	0.41
2:A:44:ALA:O	2:A:48:GLY:CA	2.67	0.41
2:B:333:ARG:NH1	4:B:358:HOH:O	2.53	0.41
2:A:50:VAL:HG12	2:A:52:PRO:HD3	2.01	0.41
2:A:116:ARG:HB2	2:A:207:GLU:OE2	2.21	0.41
2:A:171:PHE:HZ	2:B:299:LEU:HD13	1.86	0.41
2:B:194:LEU:HD23	2:B:199:GLN:HA	2.03	0.41
2:B:282:PRO:HD2	2:B:283:GLU:OE1	2.21	0.41
1:E:9:DC:H2''	1:E:10:DG:N7	2.36	0.40
1:E:16:DC:H2'	1:E:17:DT:H71	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	A	331/338 (98%)	314 (95%)	15 (4%)	2 (1%)	21	23
2	B	331/338 (98%)	303 (92%)	26 (8%)	2 (1%)	21	23
All	All	662/676 (98%)	617 (93%)	41 (6%)	4 (1%)	21	23

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	29	ARG
2	A	197	GLU
2	A	199	GLN
2	B	298	GLY

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	
2	A	277/282 (98%)	262 (95%)	15 (5%)	20	25
2	B	278/282 (99%)	265 (95%)	13 (5%)	23	31
All	All	555/564 (98%)	527 (95%)	28 (5%)	21	28

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	6	SER
2	A	7	ILE
2	A	15	THR
2	A	122	VAL
2	A	129	GLN
2	A	133	SER
2	A	153	SER
2	A	202	GLU
2	A	213	ARG
2	A	226	LEU
2	A	285	THR
2	A	297	TYR
2	A	306	VAL
2	A	335	GLU
2	A	336	LEU
2	B	6	SER
2	B	8	GLU
2	B	30	ASN
2	B	49	GLU
2	B	122	VAL
2	B	137	LEU
2	B	153	SER
2	B	169	GLU
2	B	170	GLU
2	B	202	GLU
2	B	206	GLU
2	B	213	ARG

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Mol	Chain	Res	Type
2	B	335	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	30	ASN
2	A	74	ASN
2	A	199	GLN
2	A	231	GLN
2	A	307	HIS
2	B	36	GLN
2	B	74	ASN
2	B	231	GLN
2	B	338	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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(Å²)	Q<0.9	
1	E	18/18 (100%)	-0.84	0	100	100	30, 36, 58, 59	0
1	F	18/18 (100%)	-0.80	0	100	100	32, 37, 56, 58	0
2	A	335/338 (99%)	-0.02	4 (1%)	76	74	28, 47, 65, 73	0
2	B	334/338 (98%)	-0.08	2 (0%)	85	83	21, 46, 70, 77	1 (0%)
All	All	705/712 (99%)	-0.09	6 (0%)	81	79	21, 46, 67, 77	1 (0%)

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	302	GLY	3.7
2	B	29	ARG	2.5
2	B	336	LEU	2.4
2	A	123	GLY	2.3
2	A	285	THR	2.2
2	A	125	GLY	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](#)

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. 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	B-factors(Å ²)	Q<0.9
3	MN	A	341	1/1	0.92	0.07	34,34,34,34	1
3	MN	B	341	1/1	0.96	0.05	36,36,36,36	1
3	MN	A	340	1/1	0.99	0.02	34,34,34,34	0
3	MN	B	340	1/1	1.00	0.02	32,32,32,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.