



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 03:45 pm GMT

PDB ID : 4OGE
Title : Crystal structure of the Type II-C Cas9 enzyme from *Actinomyces naeslundii*
Authors : Jiang, F.; Ma, E.; Lin, S.; Doudna, J.A.
Deposited on : 2014-01-15
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

<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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
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) : recalc28949

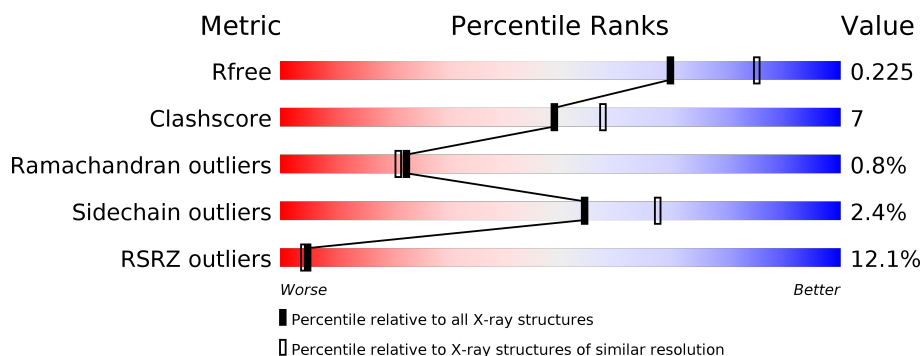
1 Overall quality at a glance

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}	100719	4002 (2.20-2.20)
Clashscore	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)
RSRZ outliers	101464	4033 (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 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	1101	

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	MG	A	1203	-	-	-	X
4	SPD	A	1205	-	-	X	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	SPD	A	1206	-	-	-	X

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7799 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 HNH endonuclease domain protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	977	Total	C	N	O	S	0	0	0
			7432	4620	1392	1391	29			

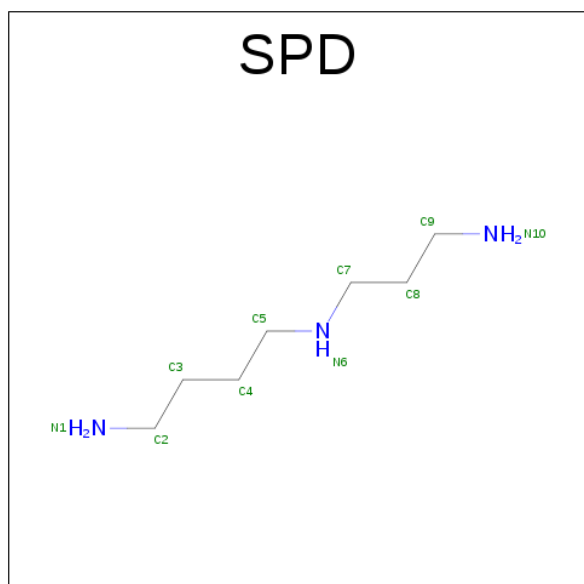
- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total	Mg	0	0
			3	3		

- Molecule 4 is SPERMIDINE (three-letter code: SPD) (formula: C₇H₁₉N₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	N	0	0
			10	7	3		
4	A	1	Total	C	N	0	0
			10	7	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	343	Total	O	0	0
			343	343		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	75.42Å 133.03Å 80.69Å 90.00° 96.22° 90.00°	Depositor
Resolution (Å)	68.69 – 2.20 68.69 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.9 (68.69-2.20) 97.9 (68.69-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.20Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.186 , 0.226 0.185 , 0.225	Depositor DCC
R_{free} test set	3926 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	39.4	Xtriage
Anisotropy	0.245	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 50.1	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	0.95	EDS
Total number of atoms	7799	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.37% 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: ZN, SPD, MG

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	0.41	0/7575	0.56	0/10287

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

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	7432	0	7179	99	0
2	A	1	0	0	0	0
3	A	3	0	0	0	0
4	A	20	0	38	12	0
5	A	343	0	0	6	0
All	All	7799	0	7217	99	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 7.

All (99) 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:A:1031:THR:HG23	1:A:1033:SER:H	1.47	0.80
1:A:757:ARG:HB2	4:A:1205:SPD:H72	1.63	0.80
1:A:632:ILE:HD12	1:A:658:ILE:HG23	1.70	0.72
1:A:757:ARG:HH21	4:A:1205:SPD:C4	2.04	0.70
1:A:1055:ASN:ND2	1:A:1059:GLU:OE2	2.24	0.70
1:A:710:ARG:NH1	5:A:1584:HOH:O	2.23	0.70
1:A:857:CYS:O	1:A:861:ARG:HG2	1.92	0.70
1:A:539:GLN:HG3	1:A:548:ILE:HD11	1.72	0.70
1:A:757:ARG:HH21	4:A:1205:SPD:H41	1.57	0.70
1:A:319:ASP:OD2	1:A:652:ARG:NH2	2.26	0.68
1:A:1066:ALA:HB3	1:A:1069:VAL:HG13	1.76	0.67
1:A:378:TRP:O	1:A:386:ARG:NH1	2.29	0.66
1:A:71:ARG:HB3	1:A:397:PRO:HG3	1.76	0.66
1:A:1025:ARG:NH2	5:A:1357:HOH:O	2.29	0.64
1:A:660:ARG:NH2	1:A:667:ASP:OD2	2.29	0.64
1:A:382:ASP:OD1	1:A:382:ASP:N	2.31	0.64
1:A:640:LYS:HE3	1:A:647:SER:HB3	1.83	0.61
1:A:761:ARG:HD3	1:A:773:TRP:CZ2	2.36	0.61
1:A:500:GLU:HG3	1:A:501:VAL:HG13	1.83	0.60
1:A:925:SER:HA	4:A:1206:SPD:H72	1.83	0.60
1:A:952:LEU:O	1:A:955:ARG:HG3	2.01	0.60
1:A:95:PHE:HB3	1:A:143:ILE:HD11	1.84	0.59
1:A:757:ARG:CB	4:A:1205:SPD:H72	2.32	0.59
1:A:477:SER:O	1:A:480:ARG:HG2	2.01	0.59
1:A:301:ASN:ND2	5:A:1417:HOH:O	2.35	0.59
1:A:628:VAL:O	1:A:632:ILE:HG12	2.02	0.58
1:A:251:PRO:HG2	1:A:256:VAL:HG11	1.84	0.58
1:A:1025:ARG:NH1	5:A:1534:HOH:O	2.36	0.58
1:A:410:PRO:HA	1:A:413:ASP:HB2	1.85	0.58
1:A:754:LEU:O	4:A:1205:SPD:H81	2.04	0.57
1:A:242:LEU:O	1:A:244:ALA:N	2.38	0.56
1:A:352:ARG:NH2	1:A:683:ASN:OD1	2.36	0.56
1:A:507:VAL:CG1	1:A:511:PHE:HB2	2.36	0.56
1:A:97:ILE:HG13	1:A:98:PRO:HD3	1.88	0.55
1:A:708:ALA:HB1	4:A:1205:SPD:H82	1.87	0.55
1:A:242:LEU:O	1:A:245:VAL:N	2.36	0.54
1:A:638:TRP:HB2	1:A:654:LYS:HD3	1.89	0.54
1:A:421:HIS:NE2	1:A:512:THR:O	2.41	0.54
1:A:925:SER:HA	4:A:1206:SPD:C7	2.37	0.54
1:A:258:ARG:HE	1:A:260:ALA:HA	1.74	0.53
1:A:446:ASP:OD2	1:A:454:ARG:HD2	2.08	0.53
1:A:279:GLU:OE1	1:A:435:THR:HG22	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:381:ALA:HB3	1:A:386:ARG:HD2	1.89	0.52
1:A:699:MET:HG2	1:A:701:TYR:CZ	2.44	0.52
1:A:383:SER:HA	1:A:386:ARG:HD3	1.92	0.52
1:A:725:LYS:N	1:A:726:GLY:HA3	2.25	0.51
1:A:401:GLU:O	1:A:405:ILE:HG12	2.10	0.51
1:A:757:ARG:HH21	4:A:1205:SPD:H42	1.75	0.50
1:A:849:ARG:NH2	1:A:904:ASP:OD2	2.40	0.50
1:A:674:SER:O	1:A:702:ARG:NH1	2.45	0.50
1:A:91:ARG:HA	1:A:95:PHE:O	2.13	0.49
1:A:666:GLU:OE2	5:A:1387:HOH:O	2.20	0.48
1:A:845:GLN:HB3	1:A:907:GLU:O	2.14	0.47
1:A:403:ALA:HB1	1:A:406:ILE:HB	1.95	0.47
1:A:257:SER:HA	1:A:258:ARG:HA	1.45	0.47
1:A:509:ASP:N	1:A:510:GLY:HA3	2.30	0.47
1:A:507:VAL:HG12	1:A:511:PHE:HB2	1.95	0.47
1:A:65:LEU:O	1:A:69:ALA:HB3	2.14	0.47
1:A:237:VAL:O	1:A:241:LEU:HG	2.15	0.47
1:A:672:GLU:H	1:A:672:GLU:CD	2.18	0.47
1:A:258:ARG:HB2	1:A:258:ARG:HH11	1.81	0.46
1:A:12:LEU:HB2	1:A:499:PRO:HA	1.97	0.46
1:A:794:MET:O	1:A:798:THR:HG23	2.16	0.46
1:A:1031:THR:HG22	1:A:1034:LYS:HB2	1.98	0.46
1:A:1013:ARG:O	1:A:1016:GLU:HB2	2.16	0.45
1:A:378:TRP:CZ2	1:A:386:ARG:HG2	2.51	0.45
1:A:371:ILE:HD11	1:A:416:LYS:O	2.16	0.45
1:A:293:ARG:HB2	1:A:341:ARG:HB2	1.98	0.45
1:A:715:ILE:HB	1:A:798:THR:HG21	1.98	0.45
1:A:258:ARG:HH21	1:A:261:PRO:HD3	1.81	0.45
1:A:880:ARG:NH1	1:A:885:GLU:HB2	2.31	0.45
1:A:89:VAL:O	1:A:93:LEU:HB2	2.16	0.45
1:A:926:ILE:O	4:A:1206:SPD:H52	2.17	0.45
1:A:229:ALA:HA	1:A:230:ARG:HA	1.54	0.44
1:A:431:ARG:O	1:A:435:THR:HG23	2.18	0.43
1:A:758:SER:HB2	4:A:1205:SPD:H92	1.99	0.43
1:A:1057:ILE:O	1:A:1062:GLY:HA2	2.19	0.43
1:A:315:PHE:CD1	1:A:332:LYS:HE3	2.54	0.43
1:A:712:ALA:HB2	4:A:1205:SPD:H42	2.01	0.43
1:A:656:GLU:O	1:A:660:ARG:HG3	2.19	0.43
1:A:232:GLY:HA3	1:A:233:VAL:HA	1.59	0.43
1:A:416:LYS:O	1:A:419:SER:HB3	2.19	0.42
1:A:604:ARG:C	1:A:604:ARG:HD2	2.39	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:VAL:HG13	1:A:158:SER:OG	2.20	0.42
1:A:298:LYS:HA	1:A:299:GLY:HA2	1.64	0.42
1:A:392:TYR:OH	1:A:403:ALA:HB2	2.19	0.42
1:A:1079:ARG:HG3	1:A:1096:THR:O	2.19	0.42
1:A:401:GLU:O	1:A:405:ILE:N	2.52	0.42
1:A:483:LYS:O	1:A:487:ARG:HG2	2.19	0.42
1:A:419:SER:O	1:A:421:HIS:ND1	2.53	0.41
1:A:596:GLY:HA3	1:A:656:GLU:HG2	2.03	0.41
1:A:166:SER:HA	1:A:167:PRO:HD3	1.87	0.41
1:A:820:LEU:HD22	1:A:948:THR:HG23	2.03	0.41
1:A:848:ASP:HB3	1:A:909:PRO:HG3	2.02	0.41
1:A:471:ALA:HA	1:A:472:PRO:HD3	1.87	0.41
1:A:271:ARG:HG2	5:A:1552:HOH:O	2.21	0.40
1:A:417:LEU:HA	1:A:417:LEU:HD23	1.74	0.40
1:A:389:MET:O	1:A:393:LEU:HG	2.21	0.40
1:A:994:VAL:O	1:A:997:ASP:HB2	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	965/1101 (88%)	914 (95%)	43 (4%)	8 (1%)	22	21

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	243	ARG
1	A	508	ARG
1	A	407	ALA
1	A	414	GLN
1	A	923	GLY

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Mol	Chain	Res	Type
1	A	162	PRO
1	A	231	GLN
1	A	245	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	A	759/931 (82%)	741 (98%)	18 (2%)	54	67

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	79	ARG
1	A	86	LEU
1	A	90	LEU
1	A	245	VAL
1	A	246	PHE
1	A	258	ARG
1	A	382	ASP
1	A	401	GLU
1	A	402	CYS
1	A	408	GLU
1	A	475	ASN
1	A	512	THR
1	A	561	LEU
1	A	601	VAL
1	A	640	LYS
1	A	867	GLU
1	A	997	ASP
1	A	1042	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 ⓘ

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

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	$\# Z > 2$
4	SPD	A	1205	-	9,9,9	0.47	0	8,8,8	1.41	2 (25%)
4	SPD	A	1206	-	9,9,9	0.45	0	8,8,8	0.76	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
4	SPD	A	1205	-	-	0/7/7/7	0/0/0/0
4	SPD	A	1206	-	-	0/7/7/7	0/0/0/0

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1205	SPD	C4-C5-N6	-2.52	105.62	112.06
4	A	1205	SPD	C7-C8-C9	-2.16	105.64	114.19

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1205	SPD	9	0
4	A	1206	SPD	3	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, 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	977/1101 (88%)	0.73	118 (12%) 5 4	20, 42, 147, 199	0

All (118) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	163	ALA	12.7
1	A	227	ILE	11.0
1	A	233	VAL	10.9
1	A	231	GLN	10.4
1	A	161	SER	9.7
1	A	143	ILE	9.5
1	A	169	MET	9.3
1	A	160	LEU	9.2
1	A	149	TRP	8.9
1	A	226	LYS	8.8
1	A	98	PRO	8.8
1	A	97	ILE	7.8
1	A	162	PRO	7.8
1	A	65	LEU	7.7
1	A	257	SER	7.5
1	A	138	MET	7.4
1	A	245	VAL	7.4
1	A	244	ALA	7.3
1	A	168	PHE	7.2
1	A	139	ALA	6.8
1	A	151	ASN	6.8
1	A	152	PRO	6.8
1	A	144	ALA	6.8
1	A	230	ARG	6.6
1	A	228	CYS	6.2
1	A	823	GLY	6.2
1	A	234	SER	6.2

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Mol	Chain	Res	Type	RSRZ
1	A	237	VAL	6.1
1	A	511	PHE	5.9
1	A	89	VAL	5.8
1	A	249	ASP	5.8
1	A	822	ASP	5.7
1	A	239	LYS	5.7
1	A	248	ALA	5.5
1	A	826	HIS	5.5
1	A	235	PRO	5.5
1	A	415	ALA	5.5
1	A	155	LYS	5.4
1	A	73	ARG	5.3
1	A	140	VAL	5.3
1	A	827	THR	5.3
1	A	76	LEU	5.2
1	A	238	CYS	5.1
1	A	69	ALA	5.0
1	A	153	TYR	4.9
1	A	251	PRO	4.8
1	A	824	ASN	4.8
1	A	66	SER	4.7
1	A	167	PRO	4.7
1	A	241	LEU	4.5
1	A	164	GLU	4.4
1	A	236	ASP	4.2
1	A	509	ASP	4.2
1	A	86	LEU	4.2
1	A	170	LYS	4.2
1	A	80	ARG	4.2
1	A	142	HIS	4.1
1	A	77	HIS	4.1
1	A	254	SER	4.1
1	A	141	ARG	4.0
1	A	78	HIS	3.9
1	A	146	HIS	3.9
1	A	521	ASP	3.8
1	A	242	LEU	3.7
1	A	87	ASP	3.7
1	A	419	SER	3.6
1	A	518	ASP	3.6
1	A	829	ASN	3.6
1	A	240	GLN	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	253	GLY	3.5
1	A	510	GLY	3.5
1	A	229	ALA	3.5
1	A	137	SER	3.5
1	A	825	ALA	3.4
1	A	81	THR	3.4
1	A	68	ILE	3.4
1	A	75	LEU	3.4
1	A	250	SER	3.3
1	A	96	PRO	3.2
1	A	165	GLU	3.2
1	A	93	LEU	3.1
1	A	232	GLY	3.1
1	A	507	VAL	3.1
1	A	95	PHE	3.0
1	A	258	ARG	3.0
1	A	148	GLY	3.0
1	A	88	GLU	3.0
1	A	159	LEU	2.9
1	A	85	GLN	2.9
1	A	298	LYS	2.9
1	A	299	GLY	2.9
1	A	512	THR	2.8
1	A	156	VAL	2.8
1	A	517	ALA	2.8
1	A	83	LEU	2.8
1	A	90	LEU	2.6
1	A	405	ILE	2.6
1	A	256	VAL	2.5
1	A	828	VAL	2.5
1	A	404	GLU	2.5
1	A	154	SER	2.4
1	A	71	ARG	2.4
1	A	94	GLY	2.4
1	A	402	CYS	2.4
1	A	300	GLU	2.4
1	A	158	SER	2.4
1	A	516	MET	2.4
1	A	92	ASP	2.3
1	A	416	LYS	2.3
1	A	411	GLU	2.2
1	A	74	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	67	GLY	2.1
1	A	392	TYR	2.1
1	A	410	PRO	2.1
1	A	963	VAL	2.1
1	A	70	ARG	2.0
1	A	84	GLN	2.0
1	A	147	ARG	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 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
4	SPD	A	1206	10/10	0.72	0.35	6.98	41,70,89,90	0
4	SPD	A	1205	10/10	0.85	0.22	3.55	26,45,49,53	0
3	MG	A	1203	1/1	0.90	0.19	2.68	53,53,53,53	0
3	MG	A	1204	1/1	0.95	0.19	1.90	14,14,14,14	0
2	ZN	A	1201	1/1	0.99	0.19	1.79	28,28,28,28	0
3	MG	A	1202	1/1	0.95	0.17	-	19,19,19,19	0

6.5 Other polymers [i](#)

There are no such residues in this entry.