



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 8S5R
Title : Structure of the Chlamydia pneumoniae effector SemD
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Deposited on : 2024-02-25
Resolution : 2.10 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.37.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.37.1

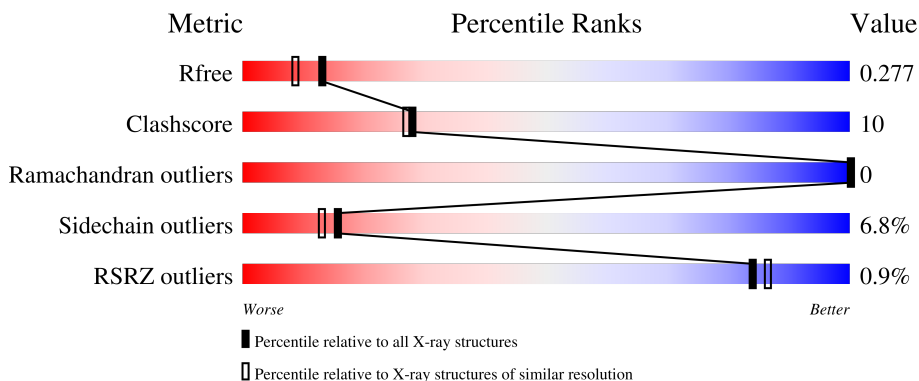
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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))
R_{free}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	A	322	
1	B	322	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3776 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 Effector SemD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	B	231	1782	1107	324	345	6	0	0	0
1	A	228	1778	1105	326	341	6	0	2	0

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-74	MET	-	initiating methionine	UNP Q9Z7M7
B	7	LEU	-	insertion	UNP Q9Z7M7
B	8	LYS	-	insertion	UNP Q9Z7M7
B	9	ASP	-	insertion	UNP Q9Z7M7
B	10	ALA	-	insertion	UNP Q9Z7M7
B	11	VAL	-	insertion	UNP Q9Z7M7
B	74	PRO	THR	conflict	UNP Q9Z7M7
B	75	VAL	SER	conflict	UNP Q9Z7M7
B	76	PRO	SER	conflict	UNP Q9Z7M7
B	77	GLN	THR	conflict	UNP Q9Z7M7
B	78	ALA	GLY	conflict	UNP Q9Z7M7
B	247	ALA	-	expression tag	UNP Q9Z7M7
A	-74	MET	-	initiating methionine	UNP Q9Z7M7
A	7	LEU	-	insertion	UNP Q9Z7M7
A	8	LYS	-	insertion	UNP Q9Z7M7
A	9	ASP	-	insertion	UNP Q9Z7M7
A	10	ALA	-	insertion	UNP Q9Z7M7
A	11	VAL	-	insertion	UNP Q9Z7M7
A	74	PRO	THR	conflict	UNP Q9Z7M7
A	75	VAL	SER	conflict	UNP Q9Z7M7
A	76	PRO	SER	conflict	UNP Q9Z7M7
A	77	GLN	THR	conflict	UNP Q9Z7M7
A	78	ALA	GLY	conflict	UNP Q9Z7M7
A	247	ALA	-	expression tag	UNP Q9Z7M7

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	110	Total 110	O 110	0	0
2	A	106	Total 106	O 106	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	47.00Å 51.96Å 58.63Å 88.33° 105.89° 117.43°	Depositor
Resolution (Å)	45.82 – 2.10 45.82 – 1.90	Depositor EDS
% Data completeness (in resolution range)	93.7 (45.82-2.10) 82.5 (45.82-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	-0.18 (at 1.89Å)	Xtrriage
Refinement program	PHENIX 1.19.2_4158, PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.218 , 0.278 0.218 , 0.277	Depositor DCC
R_{free} test set	1922 reflections (5.52%)	wwPDB-VP
Wilson B-factor (Å ²)	26.2	Xtrriage
Anisotropy	0.127	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 39.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3776	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.31% 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 [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	0.56	1/1807 (0.1%)	0.73	2/2449 (0.1%)
1	B	0.50	1/1812 (0.1%)	0.72	1/2459 (0.0%)
All	All	0.53	2/3619 (0.1%)	0.72	3/4908 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	24	LYS	CE-NZ	-7.33	1.30	1.49
1	B	218	GLU	CD-OE1	6.22	1.32	1.25

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	61	GLN	CA-CB-CG	8.35	131.77	113.40
1	B	14	ASP	CB-CG-OD2	-6.27	112.66	118.30
1	A	61	GLN	CB-CA-C	-6.20	98.01	110.40

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	1778	0	1774	45	0
1	B	1782	0	1769	28	0
2	A	106	0	0	12	3

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	110	0	0	7	5
All	All	3776	0	3543	70	5

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 10.

All (70) 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:247:ALA:O	2:A:301:HOH:O	1.83	0.95
1:B:11:VAL:HA	1:A:29:VAL:HG21	1.61	0.82
1:A:211:ASP:HB3	2:A:318:HOH:O	1.80	0.82
1:B:214:ARG:O	1:B:218:GLU:OE1	1.97	0.82
1:A:58:ARG:NH2	2:A:304:HOH:O	2.17	0.77
1:A:12:LEU:N	2:A:307:HOH:O	2.20	0.74
1:A:156:THR:O	2:A:302:HOH:O	2.05	0.73
1:A:205:GLY:O	2:A:303:HOH:O	2.08	0.70
1:B:68:SER:OG	2:B:301:HOH:O	2.11	0.69
1:A:93:GLN:HG2	1:A:161:ARG:CZ	2.25	0.67
1:A:58:ARG:NH2	2:A:308:HOH:O	2.30	0.65
1:B:47:ASP:OD2	2:B:302:HOH:O	2.14	0.65
1:B:114:VAL:O	1:B:118:VAL:HG13	1.98	0.63
1:B:142:ARG:HH11	1:B:142:ARG:HG3	1.64	0.63
1:B:12:LEU:HG	1:B:14:ASP:HB2	1.82	0.62
1:A:91:VAL:HG12	1:A:162:LEU:HD11	1.83	0.61
1:A:19:HIS:O	1:A:23:GLN:HG2	2.02	0.60
1:A:70:LEU:O	2:A:305:HOH:O	2.17	0.58
1:B:16:VAL:HG13	1:A:228:VAL:HG21	1.86	0.56
1:B:167:ASP:O	1:B:171:GLU:HG3	2.06	0.55
1:A:65:THR:O	1:A:69:MET:HG3	2.07	0.55
1:B:197:THR:O	1:B:201:ILE:HG12	2.08	0.54
1:B:139:LEU:HD11	1:B:162:LEU:HD12	1.90	0.53
1:B:142:ARG:HH11	1:B:142:ARG:CG	2.21	0.53
1:A:154:ASP:OD2	2:A:306:HOH:O	2.18	0.53
1:B:71:ILE:O	2:B:305:HOH:O	2.19	0.53
1:A:27:LYS:HG3	1:A:28:LYS:N	2.25	0.52
1:B:29:VAL:HG12	1:B:226:MET:HE2	1.92	0.52
1:B:36:LYS:NZ	2:B:303:HOH:O	2.15	0.51
1:B:61:GLN:O	1:B:65:THR:HG23	2.11	0.51
1:A:106:LYS:HG3	1:A:107:PRO:HD2	1.92	0.51
1:A:209:THR:O	1:A:213:VAL:HG23	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:13:LYS:H	1:A:13:LYS:HD2	1.76	0.50
1:A:13:LYS:HD2	2:A:322:HOH:O	2.12	0.49
1:A:161:ARG:HD3	2:A:381:HOH:O	2.12	0.49
1:A:56:ILE:HG23	1:A:124:ILE:HD11	1.94	0.49
1:A:62:GLN:HG3	1:A:66:GLU:OE2	2.12	0.49
1:A:140:GLU:O	1:A:143:VAL:HG12	2.13	0.49
1:A:18:SER:O	1:A:21:GLU:HG3	2.13	0.48
1:B:88:LYS:O	1:B:92:THR:HG23	2.14	0.48
1:B:194:SER:O	1:B:197:THR:HG23	2.13	0.48
1:A:203:ASN:HD22	1:A:207:ASN:HB3	1.80	0.46
1:A:214:ARG:NH2	2:A:318:HOH:O	2.49	0.45
1:A:71:ILE:HD11	1:A:182:ILE:CD1	2.47	0.45
1:B:155:TYR:O	1:B:159:ILE:HG23	2.17	0.45
1:A:91:VAL:O	1:A:95:VAL:HG12	2.16	0.45
1:B:247:ALA:N	2:B:311:HOH:O	2.45	0.45
1:A:87:ALA:HB1	1:A:166:LEU:HD12	1.98	0.45
1:A:13:LYS:H	1:A:13:LYS:CD	2.30	0.44
1:A:71:ILE:HD11	1:A:182:ILE:HD12	1.99	0.44
1:B:63:THR:HG21	1:B:123:GLY:O	2.19	0.43
1:B:68:SER:N	1:B:182:ILE:HD11	2.32	0.43
1:B:193:HIS:O	1:B:197:THR:HG22	2.18	0.43
1:A:131:PRO:HA	1:A:170:ARG:HH21	1.84	0.43
1:A:68:SER:HA	1:A:182:ILE:HD11	2.02	0.42
1:A:131:PRO:HA	1:A:170:ARG:NH2	2.35	0.42
1:B:151:SER:OG	1:B:152:ASP:N	2.52	0.42
1:A:156:THR:O	1:A:160:ALA:HB2	2.20	0.42
1:B:228:VAL:HG21	1:A:16:VAL:HG13	2.01	0.42
1:A:24:LYS:HD3	1:A:27:LYS:HE3	2.02	0.42
1:B:92:THR:HG22	2:B:327:HOH:O	2.19	0.41
1:A:226:MET:HE2	1:A:226:MET:HB2	1.71	0.41
1:A:120:ILE:HA	1:A:188:LEU:HD13	2.03	0.41
1:A:34:GLN:O	1:A:38:THR:HG23	2.21	0.41
1:A:13:LYS:CD	1:A:13:LYS:N	2.84	0.41
1:A:87:ALA:O	1:A:91:VAL:HG13	2.19	0.41
1:A:88:LYS:O	1:A:91:VAL:HG22	2.21	0.41
1:B:114:VAL:HG22	1:B:155:TYR:CZ	2.55	0.41
1:B:197:THR:HB	2:B:369:HOH:O	2.21	0.40
1:A:24:LYS:O	1:A:27:LYS:HG2	2.21	0.40

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:399:HOH:O	2:A:391:HOH:O[1_556]	1.99	0.21
2:B:386:HOH:O	2:A:367:HOH:O[1_545]	2.01	0.19
2:B:390:HOH:O	2:B:391:HOH:O[1_655]	2.02	0.18
2:B:394:HOH:O	2:A:396:HOH:O[1_556]	2.13	0.07
2:B:375:HOH:O	2:B:391:HOH:O[1_655]	2.19	0.01

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	224/322 (70%)	223 (100%)	1 (0%)	0	100	100
1	B	227/322 (70%)	226 (100%)	1 (0%)	0	100	100
All	All	451/644 (70%)	449 (100%)	2 (0%)	0	100	100

There are no Ramachandran outliers to report.

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	193/262 (74%)	186 (96%)	7 (4%)	35	36
1	B	193/262 (74%)	174 (90%)	19 (10%)	8	5
All	All	386/524 (74%)	360 (93%)	26 (7%)	16	13

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	12	LEU
1	B	13	LYS
1	B	14	ASP
1	B	18	SER
1	B	21	GLU
1	B	27	LYS
1	B	51	HIS
1	B	63	THR
1	B	80	VAL
1	B	92	THR
1	B	100	LYS
1	B	110	ASP
1	B	118	VAL
1	B	142	ARG
1	B	150	ASP
1	B	168	ARG
1	B	197	THR
1	B	221	ARG
1	B	228	VAL
1	A	13	LYS
1	A	51	HIS
1	A	54	ARG
1	A	63	THR
1	A	142	ARG
1	A	161	ARG
1	A	228	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	104	ASN
1	B	203	ASN
1	A	203	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 monosaccharides 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 [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	228/322 (70%)	-0.07	4 (1%) 68 72	25, 37, 61, 91	0
1	B	231/322 (71%)	-0.23	0 100 100	21, 35, 55, 72	0
All	All	459/644 (71%)	-0.15	4 (0%) 84 86	21, 36, 58, 91	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	108	GLY	5.5
1	A	107	PRO	3.3
1	A	148	GLY	3.2
1	A	110	ASP	2.6

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 monosaccharides 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.