



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:20 PM UTC

PDB ID : 2XEM / pdb_00002xem
Title : Induced-fit and allosteric effects upon polyene binding revealed by crystal structures of the Dynemicin thioesterase
Authors : Liew, C.W.; Sharff, A.; Kotaka, M.; Kong, R.; Bricogne, G.; Liang, Z.X.; Lescar, J.
Deposited on : 2010-05-17
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 : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
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

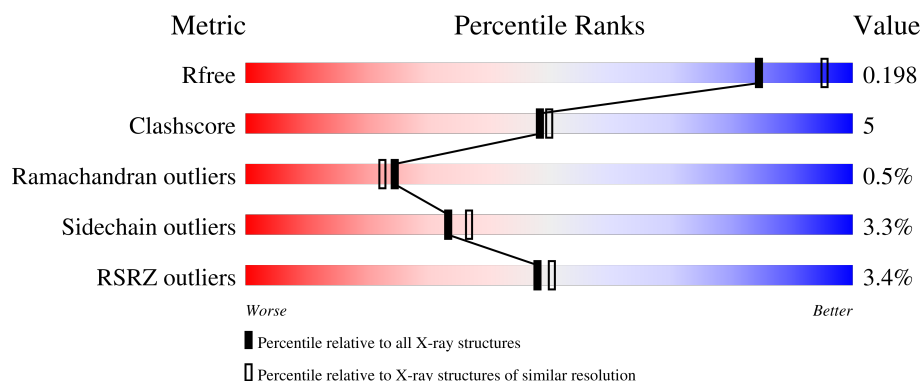
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (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	150	2% 83% 13% ..
1	B	150	3% 77% 11% 9%
1	C	150	3% 87% 6% 7%
1	D	150	5% 84% 12% ..

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
2	SSV	B	1145	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5033 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 DYNE7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	145	Total	C	N	O	S	0	2	0
			1155	729	213	203	10			
1	B	136	Total	C	N	O	S	0	0	0
			1062	668	195	191	8			
1	C	139	Total	C	N	O	S	0	1	0
			1099	693	203	195	8			
1	D	146	Total	C	N	O	S	0	1	0
			1155	728	214	204	9			

There are 24 discrepancies between the modelled and reference sequences:

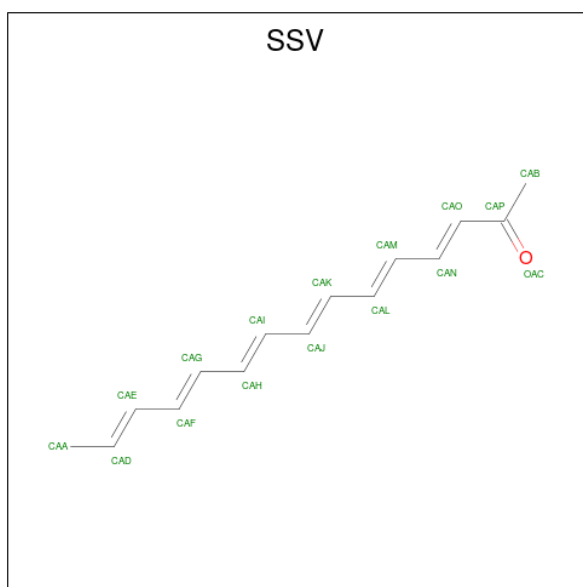
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	HIS	-	expression tag	UNP Q84HI7
A	2	HIS	-	expression tag	UNP Q84HI7
A	3	HIS	-	expression tag	UNP Q84HI7
A	4	HIS	-	expression tag	UNP Q84HI7
A	5	HIS	-	expression tag	UNP Q84HI7
A	6	HIS	-	expression tag	UNP Q84HI7
B	1	HIS	-	expression tag	UNP Q84HI7
B	2	HIS	-	expression tag	UNP Q84HI7
B	3	HIS	-	expression tag	UNP Q84HI7
B	4	HIS	-	expression tag	UNP Q84HI7
B	5	HIS	-	expression tag	UNP Q84HI7
B	6	HIS	-	expression tag	UNP Q84HI7
C	1	HIS	-	expression tag	UNP Q84HI7
C	2	HIS	-	expression tag	UNP Q84HI7
C	3	HIS	-	expression tag	UNP Q84HI7
C	4	HIS	-	expression tag	UNP Q84HI7
C	5	HIS	-	expression tag	UNP Q84HI7
C	6	HIS	-	expression tag	UNP Q84HI7
D	1	HIS	-	expression tag	UNP Q84HI7
D	2	HIS	-	expression tag	UNP Q84HI7
D	3	HIS	-	expression tag	UNP Q84HI7

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Chain	Residue	Modelled	Actual	Comment	Reference
D	4	HIS	-	expression tag	UNP Q84HI7
D	5	HIS	-	expression tag	UNP Q84HI7
D	6	HIS	-	expression tag	UNP Q84HI7

- Molecule 2 is (3E,5E,7E,9E,11E,13E)-pentadeca-3,5,7,9,11,13-hexaen-2-one (CCD ID: SSV) (formula: C₁₅H₁₈O).



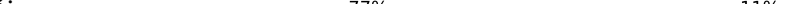
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			16	15	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	149	Total	O	0	0
			149	149		
3	B	110	Total	O	0	0
			110	110		
3	C	150	Total	O	0	0
			150	150		
3	D	137	Total	O	0	0
			137	137		

- Molecule 1: DYNE7



- Chain B:  3% 77% 11% 9%



- Chain C:  3% 87% 6% 7%



- Chain D: 5% 84% 12%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.47Å 65.39Å 94.41Å 90.00° 99.85° 90.00°	Depositor
Resolution (Å)	14.84 – 2.10 14.84 – 2.10	Depositor EDS
% Data completeness (in resolution range)	96.9 (14.84-2.10) 96.6 (14.84-2.10)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.26 (at 2.10Å)	Xtriage
Refinement program	BUSTER 2.11.0	Depositor
R, R_{free}	0.161 , 0.194 0.165 , 0.198	Depositor DCC
R_{free} test set	2008 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	26.0	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 63.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5033	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.59% 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: SSV

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	0.81	0/1194	1.17	4/1623 (0.2%)
1	B	0.85	0/1090	1.17	2/1482 (0.1%)
1	C	0.82	0/1132	1.07	1/1539 (0.1%)
1	D	0.83	0/1191	1.12	2/1620 (0.1%)
All	All	0.83	0/4607	1.13	9/6264 (0.1%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	78	PHE	CA-CB-CG	-8.88	104.92	113.80
1	A	9	ASP	CA-C-N	7.71	135.78	122.82
1	A	9	ASP	C-N-CA	7.71	135.78	122.82
1	A	80	GLU	CB-CG-CD	6.01	122.82	112.60
1	C	78	PHE	CA-CB-CG	5.12	118.92	113.80
1	B	37	GLN	CA-C-N	5.05	125.55	119.94
1	B	37	GLN	C-N-CA	5.05	125.55	119.94
1	D	139	ARG	CA-C-N	5.00	127.29	120.54
1	D	139	ARG	C-N-CA	5.00	127.29	120.54

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	1155	0	1103	16	0
1	B	1062	0	1008	20	0
1	C	1099	0	1049	10	0
1	D	1155	0	1099	12	0
2	B	16	0	18	8	0
3	A	149	0	0	2	0
3	B	110	0	0	2	0
3	C	150	0	0	1	0
3	D	137	0	0	0	0
All	All	5033	0	4277	48	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 5.

All (48) 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:8:PRO:HG3	1:A:12:VAL:HG21	1.41	1.01
1:B:142:LEU:HD23	2:B:1145:SSV:HAE	1.52	0.89
1:B:85:MET:HE3	2:B:1145:SSV:HAF	1.58	0.86
1:A:68:HIS:HD2	1:C:68:HIS:HD2	1.25	0.84
1:B:68:HIS:HD2	1:D:68:HIS:HD2	1.31	0.79
1:A:68:HIS:CD2	1:C:68:HIS:HD2	2.01	0.78
1:A:68:HIS:HD2	1:C:68:HIS:CD2	2.00	0.78
1:C:13:HIS:HD2	1:C:43:HIS:HD2	1.28	0.78
1:B:85:MET:CE	2:B:1145:SSV:HAF	2.17	0.75
1:B:68:HIS:HD2	1:D:68:HIS:CD2	2.07	0.73
1:B:68:HIS:CD2	1:D:68:HIS:HD2	2.07	0.73
1:B:124:ARG:HB3	1:B:129:LEU:HB3	1.71	0.73
1:C:13:HIS:HD2	1:C:43:HIS:CD2	2.07	0.72
1:B:41:ARG:HB2	2:B:1145:SSV:HAJ	1.76	0.68
1:B:142:LEU:HD21	2:B:1145:SSV:HAG	1.75	0.68
1:A:6:HIS:HA	3:A:2008:HOH:O	1.94	0.67
1:C:123:ARG:HG3	1:C:132:VAL:HG11	1.76	0.67
1:B:86:MET:HG3	3:B:2082:HOH:O	1.94	0.66
1:B:85:MET:HE3	2:B:1145:SSV:HAD	1.77	0.66
1:A:8:PRO:CG	1:A:12:VAL:HG21	2.21	0.65
1:B:87:LEU:HD13	1:B:142:LEU:HD22	1.78	0.65
1:B:90:LEU:HD12	1:B:142:LEU:HD13	1.79	0.64
1:A:13:HIS:HD2	1:A:43:HIS:ND1	1.97	0.62
1:D:8:PRO:HG2	1:D:86:MET:HE1	1.80	0.62
1:D:122:MET:HE2	1:D:131:PRO:HA	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:13:HIS:CD2	1:C:43:HIS:HD2	2.17	0.57
1:A:20:ASP:OD1	1:B:39:HIS:HE1	1.89	0.55
1:A:144:ARG:HD3	1:A:145:TYR:CZ	2.42	0.55
1:A:19:PHE:CD2	1:D:24:LEU:HD21	2.41	0.55
1:A:8:PRO:HG3	1:A:12:VAL:CG2	2.29	0.54
1:B:52:VAL:HG21	1:B:138:LEU:HB2	1.89	0.54
1:B:62:LEU:HB2	2:B:1145:SSV:HAO	1.88	0.54
1:A:19:PHE:CE2	1:D:24:LEU:HD21	2.44	0.51
1:D:13:HIS:ND1	1:D:43:HIS:HD2	2.08	0.51
1:C:13:HIS:HE1	3:C:2015:HOH:O	1.94	0.50
1:B:45:LEU:HB3	1:B:53:MET:HG2	1.96	0.48
1:A:144:ARG:HD3	1:A:145:TYR:CE2	2.49	0.47
1:B:13:HIS:HE1	3:B:2015:HOH:O	1.99	0.46
1:C:52:VAL:HG21	1:C:138:LEU:HB2	1.98	0.46
1:D:15:HIS:CE1	1:D:36:TRP:CZ2	3.04	0.45
1:D:123:ARG:HD2	1:D:132:VAL:CG2	2.45	0.45
1:A:52:VAL:HG21	1:A:138:LEU:HB2	2.00	0.44
1:A:63:VAL:HB	1:A:122:MET:HE3	2.01	0.43
1:C:20:ASP:OD1	1:D:39:HIS:HE1	2.01	0.43
1:A:104:VAL:HG12	3:A:2008:HOH:O	2.17	0.43
1:D:52:VAL:HG21	1:D:138:LEU:HB2	2.00	0.42
1:B:85:MET:HE3	2:B:1145:SSV:CAF	2.40	0.41
1:B:124:ARG:HB3	1:B:129:LEU:CB	2.45	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	145/150 (97%)	144 (99%)	0	1 (1%)	18	15
1	B	134/150 (89%)	127 (95%)	5 (4%)	2 (2%)	8	4

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	138/150 (92%)	135 (98%)	3 (2%)	0	100	100
1	D	145/150 (97%)	143 (99%)	2 (1%)	0	100	100
All	All	562/600 (94%)	549 (98%)	10 (2%)	3 (0%)	24	22

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	92	GLY
1	B	92	GLY
1	B	59	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	
1	A	119/121 (98%)	116 (98%)	3 (2%)	42	48
1	B	108/121 (89%)	102 (94%)	6 (6%)	19	18
1	C	112/121 (93%)	112 (100%)	0	100	100
1	D	118/121 (98%)	112 (95%)	6 (5%)	21	21
All	All	457/484 (94%)	442 (97%)	15 (3%)	33	37

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	HIS
1	A	90	LEU
1	A	127	HIS
1	B	45	LEU
1	B	53	MET
1	B	91	ASP
1	B	94	ARG
1	B	124	ARG
1	B	129	LEU

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Mol	Chain	Res	Type
1	D	27	ASN
1	D	89	ARG
1	D	90	LEU
1	D	122	MET
1	D	123	ARG
1	D	133	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	HIS
1	A	23	ASN
1	A	27	ASN
1	A	68	HIS
1	A	114	GLN
1	B	13	HIS
1	B	39	HIS
1	B	48	HIS
1	B	68	HIS
1	B	127	HIS
1	C	13	HIS
1	C	39	HIS
1	C	43	HIS
1	C	68	HIS
1	D	15	HIS
1	D	27	ASN
1	D	39	HIS
1	D	43	HIS
1	D	68	HIS
1	D	114	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

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	$\# Z > 2$
2	SSV	B	1145	-	15,15,15	1.61	2 (13%)	15,15,15	2.12	4 (26%)

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
2	SSV	B	1145	-	-	0/13/13/13	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1145	SSV	CAO-CAP	-3.42	1.39	1.47
2	B	1145	SSV	CAE-CAD	2.22	1.39	1.33

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1145	SSV	CAJ-CAI-CAH	-4.41	114.52	124.72
2	B	1145	SSV	CAL-CAK-CAJ	-3.77	116.01	124.72
2	B	1145	SSV	CAM-CAN-CAO	-3.29	116.87	124.65
2	B	1145	SSV	CAH-CAG-CAF	-3.22	117.27	124.72

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1145	SSV	8	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	145/150 (96%)	-0.41	3 (2%) 63 66	16, 28, 63, 87	2 (1%)
1	B	136/150 (90%)	-0.26	5 (3%) 45 47	17, 29, 83, 104	0
1	C	139/150 (92%)	-0.43	4 (2%) 53 56	18, 27, 58, 83	1 (0%)
1	D	146/150 (97%)	-0.33	7 (4%) 35 38	17, 28, 72, 91	1 (0%)
All	All	566/600 (94%)	-0.36	19 (3%) 48 50	16, 28, 67, 104	4 (0%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	125	ALA	4.7
1	A	148	VAL	3.5
1	C	147	VAL	3.3
1	C	127	HIS	2.9
1	B	127	HIS	2.9
1	D	125	ALA	2.8
1	B	129	LEU	2.7
1	D	4	HIS	2.6
1	B	124	ARG	2.6
1	C	126	GLY	2.6
1	A	127	HIS	2.5
1	C	9	ASP	2.4
1	B	142	LEU	2.4
1	D	90	LEU	2.3
1	A	57	ALA	2.2
1	D	7	MET	2.1
1	D	127	HIS	2.1
1	D	5	HIS	2.0
1	D	57	ALA	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($\text{\AA}^2$)	Q<0.9
2	SSV	B	1145	16/16	0.76	0.16	63,66,68,68	0

6.5 Other polymers [i](#)

There are no such residues in this entry.