



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:52 PM GMT

PDB ID : 1YW4
Title : Crystal Structure of the Succinylglutamate Desuccinylase from *Chromobacterium violaceum*, Northeast Structural Genomics Target CvR22.
Authors : Forouhar, F.; Abashidze, M.; Conover, K.; Acton, T.B.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2005-02-17
Resolution : 2.00 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

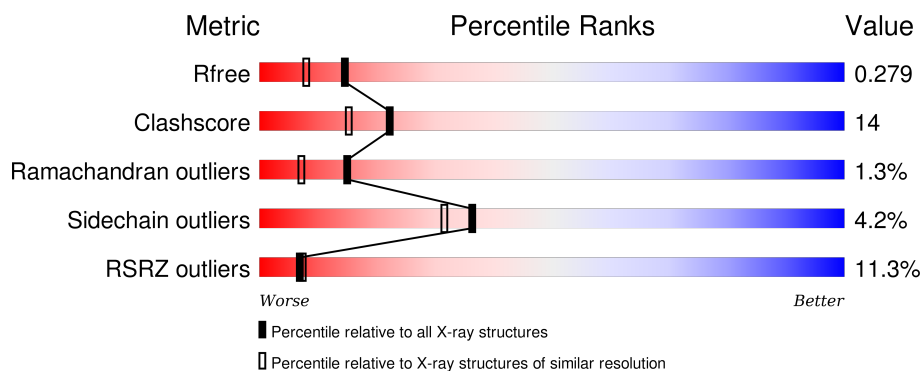
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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	341	<div> <div>10%</div> <div>67%</div> <div>24%</div> <div>• 6%</div> </div>
1	B	341	<div> <div>10%</div> <div>65%</div> <div>23%</div> <div>• 9%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5277 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 Succinylglutamate desuccinylase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	319	Total	C	N	O	S	Se	0	0	0
			2465	1563	442	451	5	4			
1	B	309	Total	C	N	O	S	Se	0	0	0
			2388	1511	426	442	5	4			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	MODIFIED RESIDUE	UNP Q7NU26
A	69	MSE	MET	MODIFIED RESIDUE	UNP Q7NU26
A	87	MSE	MET	MODIFIED RESIDUE	UNP Q7NU26
A	105	MSE	MET	MODIFIED RESIDUE	UNP Q7NU26
A	289	MSE	MET	MODIFIED RESIDUE	UNP Q7NU26
A	334	LEU	-	EXPRESSION TAG	UNP Q7NU26
A	335	GLU	-	EXPRESSION TAG	UNP Q7NU26
A	336	HIS	-	EXPRESSION TAG	UNP Q7NU26
A	337	HIS	-	EXPRESSION TAG	UNP Q7NU26
A	338	HIS	-	EXPRESSION TAG	UNP Q7NU26
A	339	HIS	-	EXPRESSION TAG	UNP Q7NU26
A	340	HIS	-	EXPRESSION TAG	UNP Q7NU26
A	341	HIS	-	EXPRESSION TAG	UNP Q7NU26
B	1	MSE	MET	MODIFIED RESIDUE	UNP Q7NU26
B	69	MSE	MET	MODIFIED RESIDUE	UNP Q7NU26
B	87	MSE	MET	MODIFIED RESIDUE	UNP Q7NU26
B	105	MSE	MET	MODIFIED RESIDUE	UNP Q7NU26
B	289	MSE	MET	MODIFIED RESIDUE	UNP Q7NU26
B	334	LEU	-	EXPRESSION TAG	UNP Q7NU26
B	335	GLU	-	EXPRESSION TAG	UNP Q7NU26
B	336	HIS	-	EXPRESSION TAG	UNP Q7NU26
B	337	HIS	-	EXPRESSION TAG	UNP Q7NU26
B	338	HIS	-	EXPRESSION TAG	UNP Q7NU26
B	339	HIS	-	EXPRESSION TAG	UNP Q7NU26
B	340	HIS	-	EXPRESSION TAG	UNP Q7NU26

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Chain	Residue	Modelled	Actual	Comment	Reference
B	341	HIS	-	EXPRESSION TAG	UNP Q7NU26

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

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

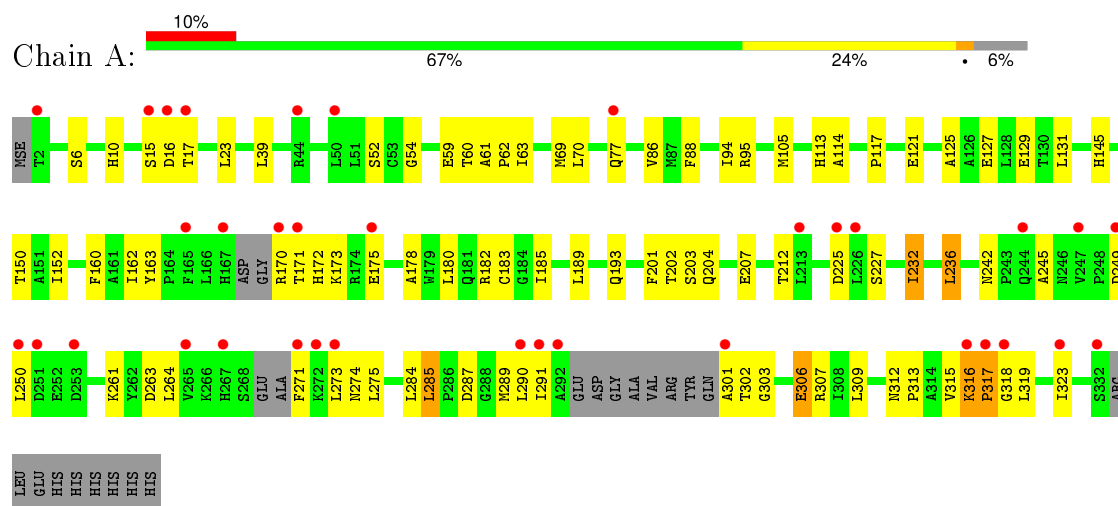
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	241	Total	O	0	0
			241	241		
3	B	181	Total	O	0	0
			181	181		

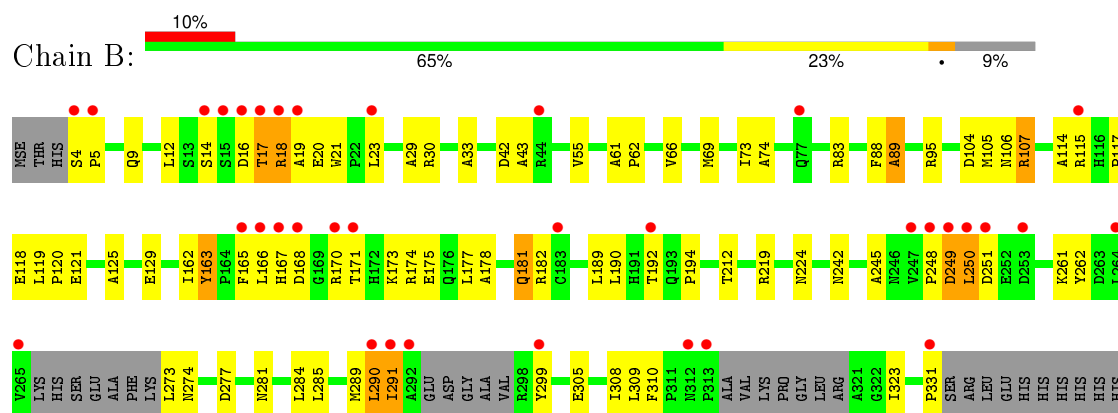
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Succinylglutamate desuccinylase



• Molecule 1: Succinylglutamate desuccinylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	62.95Å 39.44Å 131.59Å 90.00° 94.74° 90.00°	Depositor
Resolution (Å)	29.95 – 2.00 29.95 – 2.00	Depositor EDS
% Data completeness (in resolution range)	90.9 (29.95-2.00) 98.1 (29.95-2.00)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.79 (at 2.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.221 , 0.262 0.236 , 0.279	Depositor DCC
R_{free} test set	4230 reflections (9.74%)	DCC
Wilson B-factor (Å ²)	20.1	Xtriage
Anisotropy	0.346	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 53.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 83778 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5277	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.18% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.34	0/2517	0.57	1/3407 (0.0%)
1	B	0.34	0/2436	0.59	3/3298 (0.1%)
All	All	0.34	0/4953	0.58	4/6705 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	83	ARG	N-CA-C	-6.02	94.74	111.00
1	A	88	PHE	N-CA-C	-5.64	95.76	111.00
1	B	331	PRO	N-CA-CB	5.61	110.03	103.30
1	B	88	PHE	N-CA-C	-5.47	96.23	111.00

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	2465	0	2427	66	0
1	B	2388	0	2331	67	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	241	0	0	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	181	0	0	7	0
All	All	5277	0	4758	133	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 14.

All (133) 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:B:309:LEU:HD23	1:B:323:ILE:HD11	1.60	0.84
1:A:178:ALA:HB2	1:A:250:LEU:HB2	1.58	0.83
1:B:18:ARG:HH12	1:B:33:ALA:HA	1.45	0.82
1:A:317:PRO:HG2	1:A:319:LEU:H	1.45	0.81
1:B:181:GLN:HG2	1:B:250:LEU:HD23	1.62	0.79
1:A:203:SER:O	1:A:207:GLU:HA	1.86	0.76
1:A:315:VAL:C	1:A:317:PRO:HD3	2.11	0.70
1:A:261:LYS:HZ1	1:A:301:ALA:N	1.90	0.69
1:B:106:ASN:C	1:B:107:ARG:HD2	2.13	0.69
1:A:171:THR:HG22	1:A:172:HIS:H	1.57	0.68
1:A:61:ALA:HB3	1:A:62:PRO:HD3	1.76	0.68
1:A:317:PRO:HG2	1:A:318:GLY:H	1.59	0.67
1:B:242:ASN:ND2	1:B:245:ALA:HB2	2.10	0.67
1:A:127:GLU:O	1:A:131:LEU:HD13	1.96	0.66
1:B:273:LEU:HB3	1:B:291:ILE:HG21	1.79	0.65
1:A:16:ASP:HB2	3:A:619:HOH:O	1.97	0.65
1:A:290:LEU:HD13	1:A:302:THR:HB	1.80	0.64
1:A:23:LEU:HD21	1:A:39:LEU:HD23	1.82	0.61
1:B:107:ARG:HD2	1:B:107:ARG:N	2.16	0.61
1:B:182:ARG:HB3	3:B:597:HOH:O	1.99	0.61
1:B:281:ASN:ND2	1:B:310:PHE:H	1.98	0.60
1:B:14:SER:HB3	3:B:673:HOH:O	2.02	0.60
1:A:232:ILE:HD12	1:A:236:LEU:HD22	1.84	0.60
1:A:160:PHE:HE1	1:A:162:ILE:HD11	1.66	0.59
1:A:6:SER:HA	3:A:633:HOH:O	2.03	0.58
1:A:309:LEU:HB3	1:A:323:ILE:HG13	1.86	0.58
1:A:274:ASN:C	1:A:275:LEU:HD22	2.24	0.58
1:A:145:HIS:HD2	1:A:202:THR:OG1	1.86	0.57
1:A:114:ALA:O	1:A:117:PRO:HD3	2.05	0.56
1:B:194:PRO:HG2	3:B:621:HOH:O	2.04	0.56
1:B:12:LEU:O	1:B:95:ARG:HD3	2.05	0.56
1:A:10:HIS:CE1	1:A:15:SER:HB2	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:LEU:O	1:A:183:CYS:HB3	2.06	0.56
1:B:290:LEU:HD12	1:B:291:ILE:HG12	1.87	0.56
1:B:17:THR:HG23	1:B:18:ARG:N	2.21	0.56
1:A:16:ASP:O	1:A:17:THR:HG23	2.06	0.55
1:B:274:ASN:HD22	1:B:291:ILE:HB	1.71	0.55
1:B:178:ALA:HB2	1:B:250:LEU:HB2	1.89	0.54
1:B:106:ASN:HB3	1:B:107:ARG:NH1	2.23	0.54
1:B:167:HIS:O	1:B:170:ARG:HG3	2.08	0.54
1:B:4:SER:N	1:B:5:PRO:HD2	2.23	0.54
1:A:59:GLU:HB3	1:A:150:THR:OG1	2.08	0.53
1:A:175:GLU:HB2	3:A:689:HOH:O	2.07	0.53
1:B:61:ALA:HB3	1:B:62:PRO:HD3	1.89	0.53
1:B:23:LEU:HD13	1:B:74:ALA:HA	1.91	0.53
1:B:261:LYS:HG3	1:B:262:TYR:CD1	2.43	0.52
1:A:225:ASP:OD1	1:A:227:SER:HB3	2.09	0.52
1:A:242:ASN:ND2	1:A:245:ALA:HB2	2.25	0.51
1:A:274:ASN:O	1:A:275:LEU:HD13	2.11	0.51
1:B:20:GLU:HG3	1:B:30:ARG:HG2	1.93	0.51
1:A:70:LEU:HD21	1:A:86:VAL:HG21	1.92	0.51
1:B:115:ARG:HG3	1:B:115:ARG:HH11	1.74	0.51
1:A:95:ARG:HD3	3:A:597:HOH:O	2.10	0.51
1:B:18:ARG:HH12	1:B:33:ALA:CA	2.20	0.50
1:B:42:ASP:HB2	3:B:554:HOH:O	2.11	0.50
1:B:178:ALA:HB1	1:B:249:ASP:O	2.11	0.50
1:B:285:LEU:HD12	1:B:289:MSE:SE	2.62	0.50
1:B:290:LEU:CD1	1:B:291:ILE:HG12	2.42	0.50
1:A:285:LEU:HG	1:A:289:MSE:HE2	1.94	0.50
1:B:16:ASP:OD2	1:B:19:ALA:N	2.44	0.49
1:B:175:GLU:HG3	3:B:636:HOH:O	2.11	0.49
1:A:204:GLN:HG3	3:A:590:HOH:O	2.13	0.49
1:B:284:LEU:HD11	1:B:305:GLU:HA	1.95	0.49
1:A:54:GLY:HA3	1:A:94:ILE:HD11	1.95	0.49
1:A:113:HIS:HD2	3:A:653:HOH:O	1.96	0.49
1:A:315:VAL:HG23	1:A:316:LYS:N	2.28	0.48
1:B:21:TRP:HB2	1:B:29:ALA:HB3	1.94	0.48
1:A:317:PRO:HG2	1:A:318:GLY:N	2.28	0.48
1:B:171:THR:O	1:B:173:LYS:HD2	2.13	0.48
1:B:104:ASP:OD1	1:B:107:ARG:HD3	2.13	0.48
1:B:219:ARG:HB2	1:B:224:ASN:ND2	2.28	0.48
1:B:178:ALA:CB	1:B:250:LEU:HB2	2.44	0.48
1:A:113:HIS:HE1	1:A:129:GLU:OE1	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:VAL:HA	1:B:69:MSE:HE3	1.96	0.48
1:B:125:ALA:O	1:B:129:GLU:HG3	2.14	0.47
1:B:273:LEU:HD23	1:B:291:ILE:HG21	1.97	0.47
1:B:23:LEU:CD1	1:B:74:ALA:HA	2.45	0.47
1:B:17:THR:CG2	1:B:18:ARG:N	2.77	0.47
1:B:18:ARG:NH1	1:B:33:ALA:HA	2.23	0.47
1:B:9:GLN:NE2	3:B:579:HOH:O	2.48	0.46
1:A:273:LEU:CD1	1:A:275:LEU:HB2	2.46	0.46
1:A:287:ASP:HA	1:A:302:THR:OG1	2.16	0.46
1:A:170:ARG:HH12	1:A:173:LYS:HD2	1.78	0.46
1:B:178:ALA:O	1:B:181:GLN:HG3	2.16	0.45
1:A:316:LYS:HB2	1:A:316:LYS:NZ	2.31	0.45
1:A:316:LYS:N	1:A:317:PRO:HD3	2.31	0.45
1:B:181:GLN:NE2	1:B:251:ASP:OD2	2.49	0.45
1:B:250:LEU:O	1:B:250:LEU:HD13	2.16	0.45
1:A:309:LEU:O	1:A:323:ILE:HG12	2.17	0.44
1:B:114:ALA:O	1:B:117:PRO:HD3	2.17	0.44
1:B:189:LEU:HD22	1:B:309:LEU:CD2	2.47	0.44
1:A:152:ILE:N	1:A:152:ILE:HD12	2.33	0.44
1:A:60:THR:HA	1:A:63:ILE:HD12	2.00	0.44
1:A:317:PRO:CG	1:A:318:GLY:N	2.81	0.44
1:B:181:GLN:HG3	1:B:182:ARG:N	2.33	0.43
1:A:271:PHE:HZ	1:A:313:PRO:HB2	1.82	0.43
1:B:309:LEU:HB3	1:B:323:ILE:HG13	2.00	0.43
1:A:264:LEU:HD22	1:A:291:ILE:HD12	2.00	0.43
1:A:69:MSE:SE	1:A:236:LEU:HB3	2.68	0.43
1:A:162:ILE:HD12	1:A:189:LEU:HD23	2.01	0.43
1:B:181:GLN:CG	1:B:250:LEU:HD23	2.41	0.43
1:A:312:ASN:O	1:A:315:VAL:HG22	2.19	0.43
1:B:248:PRO:O	1:B:249:ASP:O	2.37	0.43
1:B:163:TYR:HB3	1:B:190:LEU:HD23	2.00	0.43
1:B:289:MSE:HE2	1:B:291:ILE:O	2.19	0.42
1:A:289:MSE:HG3	1:A:289:MSE:O	2.18	0.42
1:A:162:ILE:HB	1:A:212:THR:HB	2.00	0.42
1:B:162:ILE:HB	1:B:212:THR:HB	2.01	0.42
1:B:55:VAL:HA	1:B:89:ALA:HB3	2.01	0.42
1:A:125:ALA:O	1:A:129:GLU:HG3	2.20	0.42
1:B:105:MSE:HE2	1:B:121:GLU:O	2.20	0.42
1:B:165:PHE:O	1:B:192:THR:HA	2.20	0.42
1:B:224:ASN:ND2	3:B:504:HOH:O	2.50	0.41
1:B:285:LEU:HD21	1:B:308:ILE:N	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:160:PHE:HA	1:A:185:ILE:HG23	2.02	0.41
1:B:66:VAL:HA	1:B:69:MSE:CE	2.50	0.41
1:A:290:LEU:HD21	1:A:301:ALA:O	2.20	0.41
1:A:315:VAL:HB	1:A:317:PRO:HD3	2.03	0.41
1:A:302:THR:HA	1:A:306:GLU:OE2	2.20	0.41
1:A:263:ASP:OD1	1:A:323:ILE:HG22	2.20	0.41
1:B:23:LEU:HD12	1:B:73:ILE:HG22	2.02	0.41
1:A:105:MSE:HE3	1:A:121:GLU:HG2	2.02	0.41
1:A:178:ALA:CB	1:A:250:LEU:HB2	2.40	0.41
1:A:250:LEU:O	1:A:250:LEU:HD13	2.21	0.41
1:B:16:ASP:CG	1:B:19:ALA:HB2	2.41	0.41
1:A:129:GLU:HB3	1:A:201:PHE:CZ	2.56	0.41
1:B:119:LEU:HA	1:B:120:PRO:HD3	1.96	0.41
1:B:18:ARG:HG3	1:B:18:ARG:NH1	2.36	0.40
1:A:317:PRO:CG	1:A:318:GLY:H	2.24	0.40
1:A:284:LEU:HA	1:A:307:ARG:HG2	2.03	0.40
1:A:232:ILE:CD1	1:A:236:LEU:HD22	2.49	0.40
1:A:52:SER:OG	1:A:145:HIS:HE1	2.04	0.40
1:B:165:PHE:O	1:B:166:LEU:HD23	2.21	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	
1	A	311/341 (91%)	286 (92%)	22 (7%)	3 (1%)	19	11
1	B	301/341 (88%)	279 (93%)	17 (6%)	5 (2%)	11	4
All	All	612/682 (90%)	565 (92%)	39 (6%)	8 (1%)	15	7

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	317	PRO
1	B	249	ASP
1	B	17	THR
1	B	89	ALA
1	A	249	ASP
1	B	43	ALA
1	A	303	GLY
1	B	291	ILE

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	255/268 (95%)	246 (96%)	9 (4%)	43	40
1	B	245/268 (91%)	233 (95%)	12 (5%)	31	25
All	All	500/536 (93%)	479 (96%)	21 (4%)	36	31

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	77	GLN
1	A	163	TYR
1	A	182	ARG
1	A	193	GLN
1	A	232	ILE
1	A	236	LEU
1	A	285	LEU
1	A	306	GLU
1	A	316	LYS
1	B	18	ARG
1	B	107	ARG
1	B	118	GLU
1	B	163	TYR
1	B	168	ASP
1	B	174	ARG
1	B	177	LEU

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Mol	Chain	Res	Type
1	B	181	GLN
1	B	250	LEU
1	B	277	ASP
1	B	290	LEU
1	B	299	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	ASN
1	A	102	ASN
1	A	113	HIS
1	A	145	HIS
1	A	181	GLN
1	A	193	GLN
1	A	242	ASN
1	A	274	ASN
1	B	110	ASN
1	B	176	GLN
1	B	191	HIS
1	B	196	ASN
1	B	224	ASN
1	B	274	ASN
1	B	281	ASN

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 2 ligands modelled in this entry, 2 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

There are no such residues in this entry.

5.8 Polymer linkage issues

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	315/341 (92%)	0.69	35 (11%) 7 8	10, 24, 51, 58	0
1	B	305/341 (89%)	0.62	35 (11%) 6 7	9, 23, 47, 56	0
All	All	620/682 (90%)	0.66	70 (11%) 7 7	9, 24, 50, 58	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	15	SER	10.4
1	A	332	SER	9.6
1	B	17	THR	9.6
1	A	301	ALA	8.0
1	B	292	ALA	7.7
1	B	14	SER	6.7
1	A	316	LYS	6.4
1	B	18	ARG	6.1
1	B	253	ASP	5.7
1	A	170	ARG	5.4
1	B	249	ASP	5.3
1	A	271	PHE	5.3
1	A	247	VAL	5.1
1	B	265	VAL	4.9
1	B	44	ARG	4.8
1	B	290	LEU	4.6
1	A	253	ASP	4.5
1	B	291	ILE	4.5
1	B	5	PRO	4.4
1	A	44	ARG	4.3
1	B	19	ALA	4.3
1	B	250	LEU	4.0
1	A	290	LEU	3.9
1	A	317	PRO	3.9

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Mol	Chain	Res	Type	RSRZ
1	B	264	LEU	3.8
1	B	248	PRO	3.8
1	A	291	ILE	3.7
1	B	251	ASP	3.6
1	A	167	HIS	3.6
1	B	168	ASP	3.6
1	A	292	ALA	3.5
1	B	165	PHE	3.5
1	B	171	THR	3.4
1	A	265	VAL	3.4
1	A	267	HIS	3.2
1	A	272	LYS	3.1
1	B	192	THR	3.1
1	A	16	ASP	2.9
1	A	171	THR	2.9
1	B	16	ASP	2.9
1	A	249	ASP	2.8
1	A	250	LEU	2.7
1	A	244	GLN	2.7
1	B	23	LEU	2.6
1	B	77	GLN	2.6
1	B	4	SER	2.6
1	A	2	THR	2.5
1	A	273	LEU	2.5
1	B	313	PRO	2.5
1	B	247	VAL	2.5
1	A	225	ASP	2.4
1	B	15	SER	2.4
1	A	318	GLY	2.4
1	A	17	THR	2.4
1	A	77	GLN	2.4
1	B	166	LEU	2.4
1	B	183	CYS	2.4
1	B	167	HIS	2.4
1	B	115	ARG	2.3
1	B	170	ARG	2.3
1	A	226	LEU	2.3
1	B	331	PRO	2.2
1	A	165	PHE	2.2
1	A	251	ASP	2.2
1	A	213	LEU	2.2
1	B	299	TYR	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	312	ASN	2.1
1	A	50	LEU	2.1
1	A	175	GLU	2.0
1	A	323	ILE	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(\AA^2)	Q<0.9
2	ZN	A	501	1/1	0.99	0.04	-2.56	28,28,28,28	0
2	ZN	B	502	1/1	0.98	0.08	-4.84	40,40,40,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.