



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:17 PM UTC

PDB ID : 1NZR / pdb_00001nzh
Title : CRYSTAL STRUCTURE OF THE AZURIN MUTANT NICKEL-TRP48MET FROM PSEUDOMONAS AERUGINOSA AT 2.2 ANGSTROMS RESOLUTION
Authors : Tsai, L.-C.; Sjolín, L.; Langer, V.; Bonander, N.; Karlsson, B.G.; Vanngard, T.; Hammann, C.; Nar, H.
Deposited on : 1994-12-09
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
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

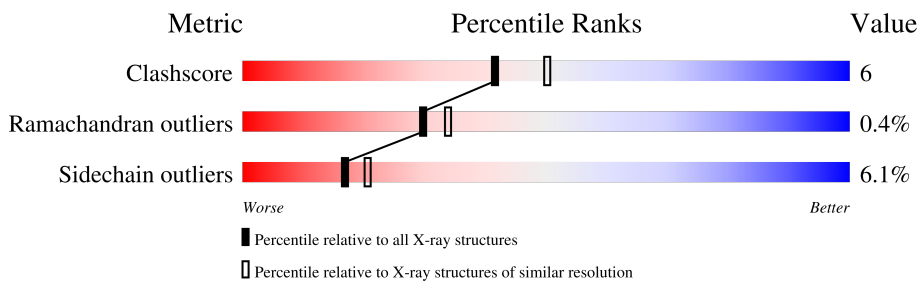
1 Overall quality at a glance

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))
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	128	
1	B	128	
1	C	128	
1	D	128	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5751 atoms, of which 1538 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 AZURIN.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	128	Total	C	H	N	O	S	0	0	0
			1183	601	216	163	193	10			
1	B	128	Total	C	H	N	O	S	0	0	0
			1183	601	216	163	193	10			
1	C	128	Total	C	H	N	O	S	0	0	0
			1183	601	216	163	193	10			
1	D	128	Total	C	H	N	O	S	0	0	0
			1183	601	216	163	193	10			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	48	MET	TRP	engineered mutation	UNP P00282
B	48	MET	TRP	engineered mutation	UNP P00282
C	48	MET	TRP	engineered mutation	UNP P00282
D	48	MET	TRP	engineered mutation	UNP P00282

- Molecule 2 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ni	0	0
			1	1		
2	B	1	Total	Ni	0	0
			1	1		
2	C	1	Total	Ni	0	0
			1	1		
2	D	1	Total	Ni	0	0
			1	1		

- Molecule 3 is NITRATE ION (CCD ID: NO3) (formula: NO₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	N	O	0	0
			4	1	3		

- Molecule 4 is water.

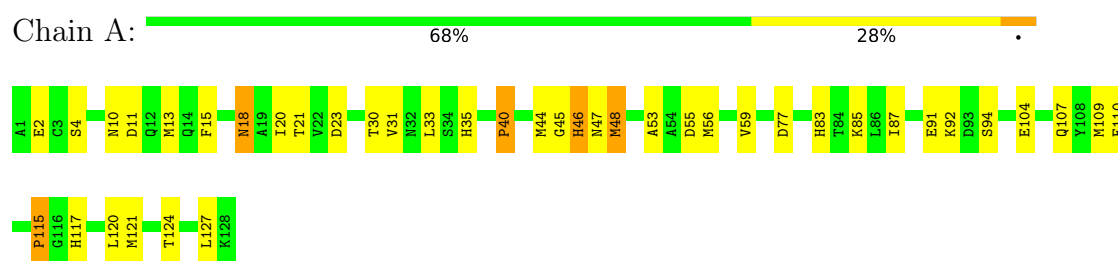
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	89	Total	H	O	0	0
			267	178	89		
4	B	92	Total	H	O	0	0
			276	184	92		
4	C	78	Total	H	O	0	0
			234	156	78		
4	D	78	Total	H	O	0	0
			234	156	78		

3 Residue-property plots

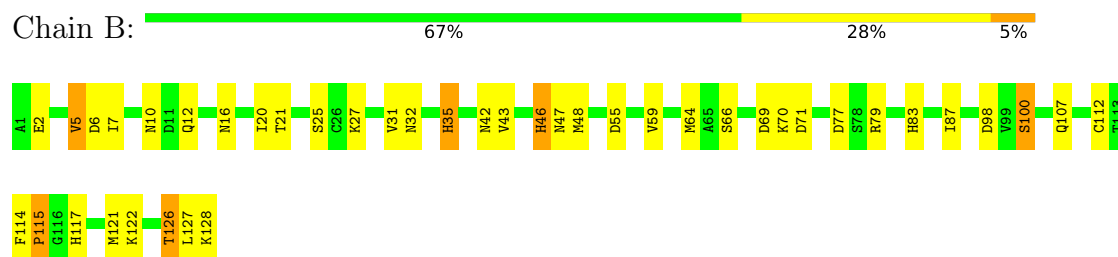
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

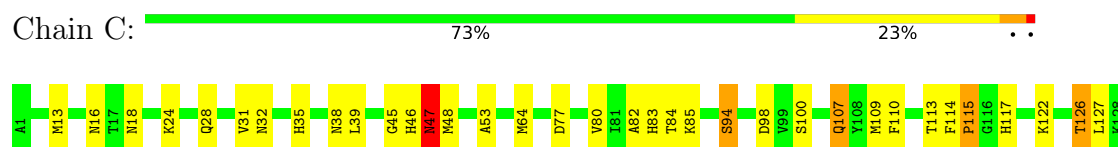
• Molecule 1: AZURIN



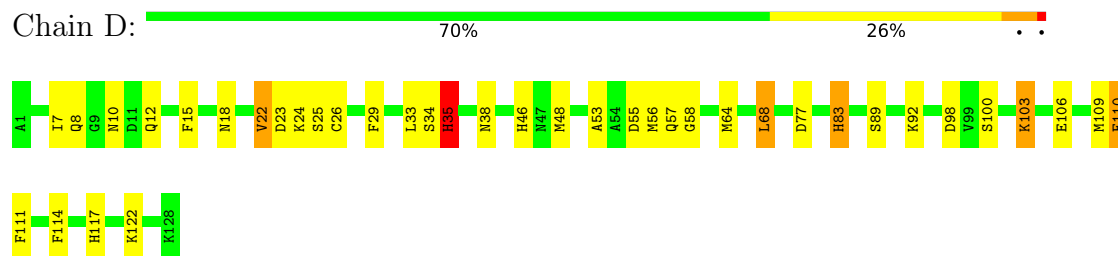
• Molecule 1: AZURIN



• Molecule 1: AZURIN



• Molecule 1: AZURIN



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	57.40Å 80.40Å 110.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.20	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.20)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.170 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5751	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

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: NO3, NI

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	1.05	6/982 (0.6%)	1.77	20/1322 (1.5%)
1	B	1.06	6/982 (0.6%)	1.82	22/1322 (1.7%)
1	C	1.12	7/982 (0.7%)	1.89	28/1322 (2.1%)
1	D	1.10	6/982 (0.6%)	1.84	26/1322 (2.0%)
All	All	1.08	25/3928 (0.6%)	1.83	96/5288 (1.8%)

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	35	HIS	CD2-NE2	-8.01	1.29	1.37
1	A	35	HIS	CD2-NE2	-7.27	1.29	1.37
1	D	117	HIS	CG-ND1	-6.80	1.30	1.38
1	D	46	HIS	CD2-NE2	-6.68	1.30	1.37
1	D	117	HIS	CD2-NE2	-6.66	1.30	1.37
1	A	83	HIS	CD2-NE2	-6.66	1.30	1.37
1	B	83	HIS	CD2-NE2	-6.61	1.30	1.37
1	C	117	HIS	CD2-NE2	-6.53	1.30	1.37
1	D	83	HIS	CD2-NE2	-6.51	1.30	1.37
1	C	83	HIS	CD2-NE2	-6.32	1.30	1.37
1	B	46	HIS	CD2-NE2	-6.28	1.30	1.37
1	B	117	HIS	CD2-NE2	-6.25	1.30	1.37
1	C	46	HIS	CD2-NE2	-6.23	1.30	1.37
1	C	35	HIS	CD2-NE2	-6.22	1.31	1.37
1	A	46	HIS	CD2-NE2	-6.11	1.31	1.37
1	A	117	HIS	CD2-NE2	-5.69	1.31	1.37
1	C	117	HIS	CG-ND1	-5.61	1.32	1.38
1	B	83	HIS	CG-ND1	-5.47	1.32	1.38
1	A	46	HIS	CG-ND1	-5.37	1.32	1.38
1	B	35	HIS	CD2-NE2	-5.24	1.32	1.37
1	C	46	HIS	CG-ND1	-5.18	1.32	1.38
1	C	80	VAL	CA-CB	5.14	1.59	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	83	HIS	CG-ND1	-5.14	1.32	1.38
1	D	46	HIS	CG-ND1	-5.13	1.32	1.38
1	B	117	HIS	CG-ND1	-5.12	1.32	1.38

All (96) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	35	HIS	CB-CG-CD2	-9.75	118.52	131.20
1	C	28	GLN	OE1-CD-NE2	-8.96	113.64	122.60
1	D	103	LYS	CA-CB-CG	8.84	131.78	114.10
1	C	110	PHE	CA-CB-CG	8.44	122.24	113.80
1	C	46	HIS	ND1-CG-CD2	8.26	114.36	106.10
1	C	107	GLN	OE1-CD-NE2	-8.16	114.44	122.60
1	A	10	ASN	OD1-CG-ND2	-8.16	114.44	122.60
1	B	47	ASN	CA-CB-CG	8.13	120.73	112.60
1	A	18	ASN	CA-CB-CG	7.93	120.53	112.60
1	C	77	ASP	CA-CB-CG	7.83	120.43	112.60
1	D	35	HIS	CB-CG-ND1	7.78	134.38	122.70
1	C	28	GLN	CG-CD-NE2	7.70	127.95	116.40
1	C	45	GLY	O-C-N	-7.63	116.06	122.77
1	B	98	ASP	CA-CB-CG	7.58	120.18	112.60
1	A	35	HIS	CB-CG-CD2	-7.57	121.36	131.20
1	C	47	ASN	CB-CG-ND2	7.48	127.61	116.40
1	A	10	ASN	CB-CG-ND2	7.46	127.58	116.40
1	C	47	ASN	OD1-CG-ND2	-7.20	115.40	122.60
1	B	42	ASN	CA-CB-CG	7.17	119.77	112.60
1	D	38	ASN	OD1-CG-ND2	-7.12	115.48	122.60
1	D	110	PHE	CA-CB-CG	7.01	120.81	113.80
1	A	117	HIS	ND1-CG-CD2	6.92	113.02	106.10
1	D	46	HIS	ND1-CG-CD2	6.92	113.02	106.10
1	D	117	HIS	ND1-CG-CD2	6.77	112.87	106.10
1	C	46	HIS	CB-CG-CD2	-6.71	122.48	131.20
1	D	103	LYS	CB-CA-C	-6.68	98.13	109.83
1	A	47	ASN	CA-CB-CG	6.64	119.24	112.60
1	C	126	THR	CA-CB-OG1	-6.60	99.70	109.60
1	A	46	HIS	ND1-CG-CD2	6.56	112.66	106.10
1	B	122	LYS	CG-CD-CE	-6.47	96.42	111.30
1	A	35	HIS	CB-CG-ND1	6.39	132.29	122.70
1	D	35	HIS	CA-CB-CG	6.39	120.19	113.80
1	C	117	HIS	ND1-CG-CD2	6.38	112.48	106.10
1	A	2	GLU	CA-CB-CG	-6.25	101.61	114.10
1	A	15	PHE	N-CA-C	-6.13	100.31	110.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	16	ASN	OD1-CG-ND2	-6.12	116.48	122.60
1	D	55	ASP	N-CA-C	6.11	120.90	113.38
1	D	103	LYS	CB-CG-CD	-6.11	97.24	111.30
1	D	68	LEU	N-CA-C	6.11	118.91	111.82
1	B	87	ILE	N-CA-C	6.09	118.15	108.89
1	A	77	ASP	N-CA-C	-6.08	99.95	109.25
1	B	35	HIS	CB-CG-ND1	6.06	131.79	122.70
1	B	46	HIS	ND1-CG-CD2	6.03	112.13	106.10
1	B	35	HIS	CB-CG-CD2	-5.98	123.42	131.20
1	C	38	ASN	OD1-CG-ND2	-5.98	116.62	122.60
1	B	117	HIS	ND1-CG-CD2	5.95	112.05	106.10
1	D	38	ASN	CB-CG-ND2	5.93	125.30	116.40
1	A	85	LYS	N-CA-C	-5.93	102.89	110.53
1	C	35	HIS	CB-CG-CD2	-5.91	123.51	131.20
1	A	117	HIS	CG-CD2-NE2	-5.91	101.30	107.20
1	B	126	THR	N-CA-CB	-5.81	101.07	110.71
1	C	16	ASN	OD1-CG-ND2	-5.79	116.81	122.60
1	B	114	PHE	CA-C-N	5.76	127.04	119.84
1	B	114	PHE	C-N-CA	5.76	127.04	119.84
1	A	15	PHE	CA-CB-CG	5.74	119.54	113.80
1	B	77	ASP	CA-CB-CG	5.73	118.33	112.60
1	A	23	ASP	CA-CB-CG	5.71	118.31	112.60
1	D	122	LYS	CA-CB-CG	5.70	125.50	114.10
1	D	114	PHE	CA-C-N	5.69	126.96	119.84
1	D	114	PHE	C-N-CA	5.69	126.96	119.84
1	D	10	ASN	CA-CB-CG	5.68	118.28	112.60
1	B	117	HIS	N-CA-C	5.57	118.28	111.82
1	D	64	MET	CA-CB-CG	-5.56	102.97	114.10
1	A	40	PRO	CA-C-N	5.55	127.66	120.44
1	A	40	PRO	C-N-CA	5.55	127.66	120.44
1	D	46	HIS	CB-CG-CD2	-5.53	124.02	131.20
1	C	82	ALA	O-C-N	-5.51	117.20	123.48
1	D	12	GLN	CA-CB-CG	5.50	125.09	114.10
1	A	46	HIS	CG-CD2-NE2	-5.45	101.75	107.20
1	D	77	ASP	N-CA-C	-5.42	101.83	109.96
1	A	45	GLY	O-C-N	-5.40	116.27	122.91
1	C	107	GLN	CG-CD-NE2	5.38	124.48	116.40
1	D	98	ASP	CA-CB-CG	5.38	117.98	112.60
1	B	27	LYS	N-CA-C	-5.36	105.09	111.69
1	B	69	ASP	CA-CB-CG	5.36	117.96	112.60
1	D	83	HIS	CB-CG-CD2	-5.31	124.30	131.20
1	C	18	ASN	CB-CG-ND2	5.30	124.35	116.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	22	VAL	N-CA-C	-5.27	100.86	108.71
1	C	94	SER	CA-CB-OG	5.24	121.58	111.10
1	C	47	ASN	CA-CB-CG	5.24	117.84	112.60
1	D	117	HIS	CA-C-O	-5.23	113.95	119.97
1	B	6	ASP	N-CA-C	-5.23	101.18	109.59
1	A	124	THR	CA-C-O	-5.22	115.12	121.28
1	C	46	HIS	CG-CD2-NE2	-5.14	102.06	107.20
1	C	84	THR	N-CA-C	-5.11	102.44	110.36
1	B	43	VAL	N-CA-C	-5.10	107.84	112.12
1	B	79	ARG	NE-CZ-NH1	5.10	126.60	121.50
1	C	114	PHE	CA-C-N	5.09	125.38	119.93
1	C	114	PHE	C-N-CA	5.09	125.38	119.93
1	B	71	ASP	CA-CB-CG	5.08	117.68	112.60
1	C	35	HIS	CB-CG-ND1	5.07	130.31	122.70
1	C	39	LEU	CA-C-N	5.03	124.92	119.85
1	C	39	LEU	C-N-CA	5.03	124.92	119.85
1	D	12	GLN	OE1-CD-NE2	-5.03	117.58	122.60
1	B	107	GLN	OE1-CD-NE2	-5.02	117.58	122.60
1	C	24	LYS	CA-CB-CG	5.00	124.10	114.10

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	967	216	951	16	0
1	B	967	216	951	11	1
1	C	967	216	951	11	0
1	D	967	216	951	16	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	4	0	0	0	0
4	A	89	178	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	92	184	0	2	0
4	C	78	156	0	1	0
4	D	78	156	0	0	0
All	All	4213	1538	3804	49	1

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 6.

All (49) 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:31:VAL:HG21	1:A:48:MET:HE3	1.77	0.67
1:C:47:ASN:HD21	1:C:113:THR:H	1.43	0.66
1:A:53:ALA:HA	1:A:109:MET:HG2	1.85	0.59
1:C:107:GLN:HE22	1:D:58:GLY:HA2	1.69	0.58
1:C:32:ASN:HA	1:C:94:SER:HB3	1.85	0.58
1:A:33:LEU:O	1:A:92:LYS:HD2	2.04	0.57
1:A:11:ASP:HA	1:A:44:MET:HG2	1.87	0.56
1:A:13:MET:HE3	1:A:120:LEU:HG	1.88	0.55
1:D:48:MET:O	1:D:83:HIS:HA	2.06	0.54
1:A:104:GLU:HG2	4:A:674:HOH:O	2.08	0.54
1:C:31:VAL:HG21	1:C:48:MET:HE3	1.89	0.54
1:A:13:MET:HE1	1:C:13:MET:HE1	1.89	0.54
1:C:122:LYS:NZ	1:D:57:GLN:HE22	2.06	0.54
1:D:23:ASP:HB3	1:D:26:CYS:SG	2.50	0.52
1:D:7:ILE:HG23	1:D:33:LEU:HD12	1.91	0.52
1:B:46:HIS:CE1	1:B:121:MET:SD	3.04	0.51
1:D:34:SER:HB3	1:D:92:LYS:HD3	1.95	0.49
1:B:31:VAL:HG11	1:B:48:MET:HG3	1.93	0.49
1:A:46:HIS:CE1	1:A:121:MET:SD	3.06	0.49
1:B:7:ILE:HG22	1:B:32:ASN:O	2.13	0.49
1:C:53:ALA:HA	1:C:109:MET:HG2	1.95	0.48
1:A:13:MET:HG2	1:A:44:MET:SD	2.55	0.46
1:B:46:HIS:HA	1:B:112:CYS:SG	2.55	0.46
1:A:55:ASP:O	1:A:59:VAL:HG23	2.15	0.46
1:A:4:SER:HA	1:A:30:THR:O	2.15	0.46
1:D:22:VAL:HG13	1:D:29:PHE:CD1	2.52	0.45
1:D:103:LYS:HD3	1:D:106:GLU:HB2	1.99	0.44
1:C:122:LYS:HZ3	1:D:57:GLN:HE22	1.65	0.44
1:B:10:ASN:HB2	4:B:716:HOH:O	2.16	0.44
1:D:35:HIS:O	1:D:89:SER:HA	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:85:LYS:HB3	1:C:85:LYS:HE2	1.81	0.44
1:C:107:GLN:NE2	1:D:58:GLY:HA2	2.31	0.43
1:A:20:ILE:HG13	1:A:110:PHE:CE1	2.53	0.43
1:B:21:THR:HG23	1:B:128:LYS:HB2	2.00	0.43
1:D:103:LYS:HE2	1:D:103:LYS:HB3	1.71	0.43
1:C:64:MET:HG3	1:C:115:PRO:HG3	2.01	0.43
1:A:56:MET:HE3	4:A:600:HOH:O	2.17	0.43
1:D:48:MET:HE2	1:D:110:PHE:CB	2.48	0.43
1:D:53:ALA:HA	1:D:109:MET:HG2	2.00	0.43
1:A:44:MET:HE1	4:C:781:HOH:O	2.19	0.42
1:D:8:GLN:O	1:D:15:PHE:HA	2.19	0.42
1:B:64:MET:HG3	1:B:115:PRO:HG3	2.01	0.42
1:B:64:MET:HB2	4:B:796:HOH:O	2.19	0.42
1:B:5:VAL:HG11	1:B:20:ILE:HD13	2.02	0.41
1:D:56:MET:HG3	1:D:111:PHE:CE1	2.55	0.41
1:B:66:SER:O	1:B:70:LYS:HG2	2.20	0.41
1:A:87:ILE:HB	1:A:91:GLU:HB2	2.02	0.41
1:A:4:SER:HB3	1:A:30:THR:HB	2.03	0.41
1:B:55:ASP:O	1:B:59:VAL:HG23	2.21	0.40

All (1) 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)
1:B:100:SER:HG	1:D:24:LYS:HZ2[2_564]	1.27	0.33

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	126/128 (98%)	123 (98%)	2 (2%)	1 (1%)	16 16

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	126/128 (98%)	119 (94%)	6 (5%)	1 (1%)	16	16
1	C	126/128 (98%)	119 (94%)	7 (6%)	0	100	100
1	D	126/128 (98%)	122 (97%)	4 (3%)	0	100	100
All	All	504/512 (98%)	483 (96%)	19 (4%)	2 (0%)	30	34

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	2	GLU
1	A	115	PRO

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	110/110 (100%)	102 (93%)	8 (7%)	13	15
1	B	110/110 (100%)	102 (93%)	8 (7%)	13	15
1	C	110/110 (100%)	104 (94%)	6 (6%)	19	24
1	D	110/110 (100%)	105 (96%)	5 (4%)	24	33
All	All	440/440 (100%)	413 (94%)	27 (6%)	17	20

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	ASN
1	A	21	THR
1	A	40	PRO
1	A	48	MET
1	A	94	SER
1	A	107	GLN
1	A	115	PRO
1	A	127	LEU
1	B	5	VAL

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Mol	Chain	Res	Type
1	B	12	GLN
1	B	25	SER
1	B	35	HIS
1	B	100	SER
1	B	115	PRO
1	B	126	THR
1	B	127	LEU
1	C	47	ASN
1	C	98	ASP
1	C	100	SER
1	C	115	PRO
1	C	126	THR
1	C	127	LEU
1	D	18	ASN
1	D	25	SER
1	D	35	HIS
1	D	68	LEU
1	D	100	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	GLN
1	C	47	ASN
1	C	107	GLN
1	D	57	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 4 are monoatomic - leaving 1 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	$\# Z > 2$
3	NO3	A	130	-	1,3,3	0.62	0	0,3,3	-	-

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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.