



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:04 AM UTC

PDB ID : 1G1Q / pdb_00001g1q
Title : Crystal structure of P-selectin lectin/EGF domains
Authors : Somers, W.S.; Camphausen, R.T.
Deposited on : 2000-10-13
Resolution : 2.40 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

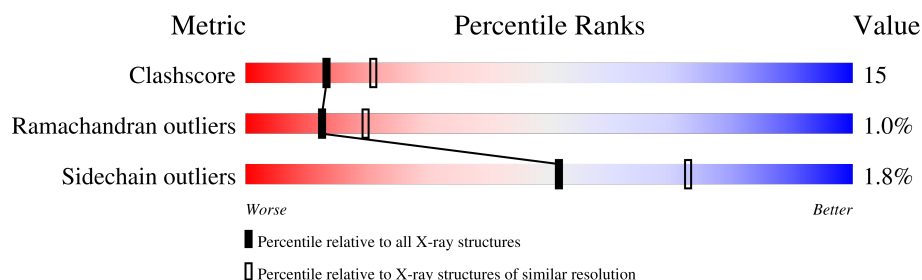
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)

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	162	
1	B	162	
1	C	162	
1	D	162	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5414 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 P-SELECTIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	160	Total	C	N	O	S	0	0	0
			1315	829	222	253	11			
1	B	160	Total	C	N	O	S	0	0	0
			1319	832	223	253	11			
1	C	159	Total	C	N	O	S	0	0	0
			1307	825	220	251	11			
1	D	160	Total	C	N	O	S	0	0	0
			1319	832	223	253	11			

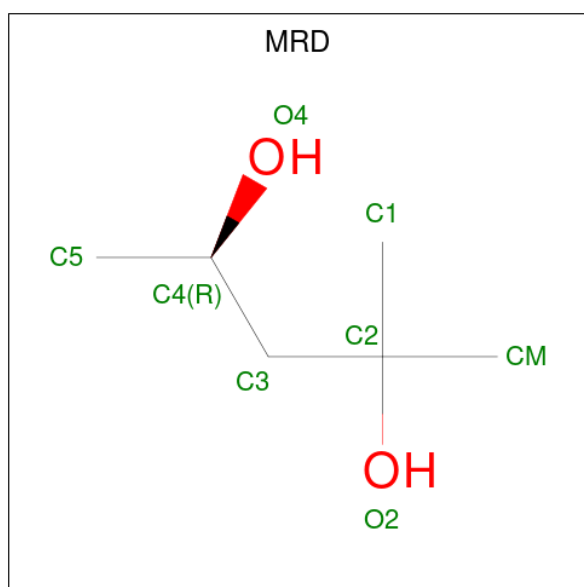
There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	158	ASP	GLU	conflict	UNP P16109
A	159	ASP	-	cloning artifact	UNP P16109
A	160	ASP	-	cloning artifact	UNP P16109
A	161	ASP	-	cloning artifact	UNP P16109
A	162	LYS	-	cloning artifact	UNP P16109
B	158	ASP	GLU	conflict	UNP P16109
B	159	ASP	-	cloning artifact	UNP P16109
B	160	ASP	-	cloning artifact	UNP P16109
B	161	ASP	-	cloning artifact	UNP P16109
B	162	LYS	-	cloning artifact	UNP P16109
C	158	ASP	GLU	conflict	UNP P16109
C	159	ASP	-	cloning artifact	UNP P16109
C	160	ASP	-	cloning artifact	UNP P16109
C	161	ASP	-	cloning artifact	UNP P16109
C	162	LYS	-	cloning artifact	UNP P16109
D	158	ASP	GLU	conflict	UNP P16109
D	159	ASP	-	cloning artifact	UNP P16109
D	160	ASP	-	cloning artifact	UNP P16109
D	161	ASP	-	cloning artifact	UNP P16109
D	162	LYS	-	cloning artifact	UNP P16109

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

- Molecule 3 is (4R)-2-METHYLPENTANE-2,4-DIOL (CCD ID: MRD) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 8 6 2	0	0
3	C	1	Total C O 8 6 2	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	53	Total O 53 53	0	0
4	B	39	Total O 39 39	0	0
4	C	30	Total O 30 30	0	0

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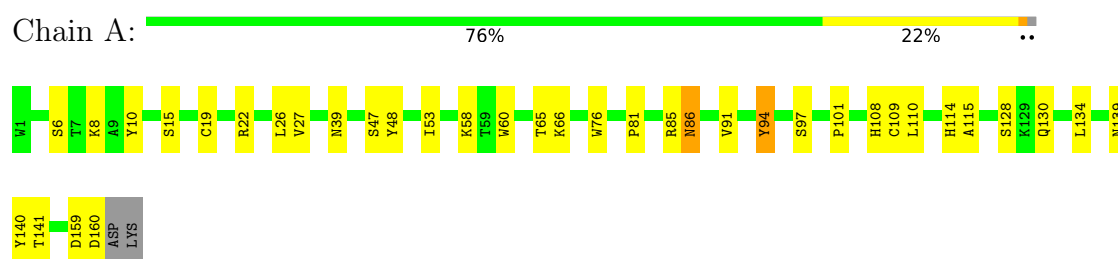
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	12	Total	O	0	0
			12	12		

3 Residue-property plots [i](#)

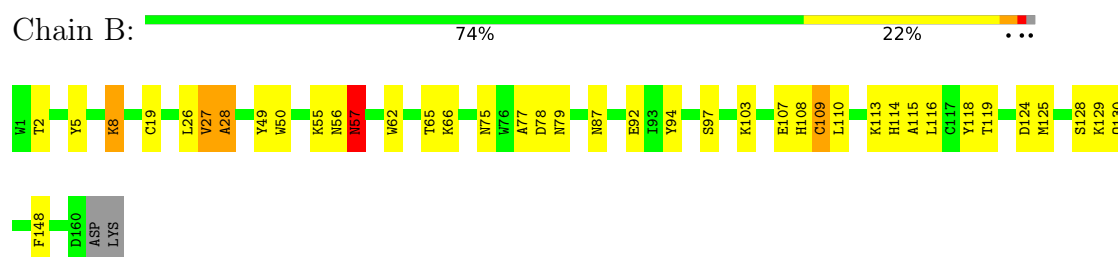
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

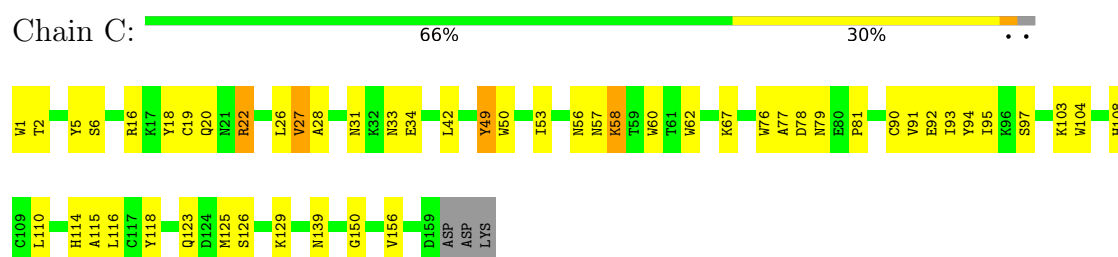
• Molecule 1: P-SELECTIN



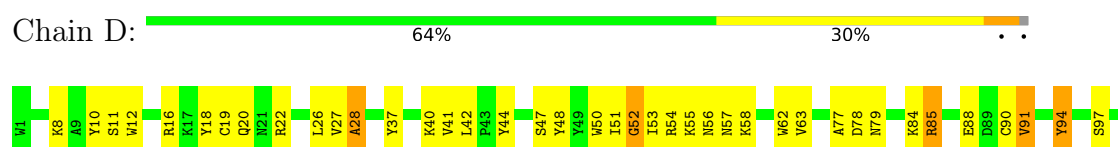
• Molecule 1: P-SELECTIN



• Molecule 1: P-SELECTIN



• Molecule 1: P-SELECTIN





4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	81.00Å 60.83Å 91.44Å 90.00° 103.58° 90.00°	Depositor
Resolution (Å)	14.00 – 2.40	Depositor
% Data completeness (in resolution range)	(Not available) (14.00-2.40)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.213 , 0.252	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5414	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	0.64	0/1354	1.00	4/1841 (0.2%)
1	B	0.53	0/1358	0.98	3/1845 (0.2%)
1	C	0.53	0/1346	0.91	4/1830 (0.2%)
1	D	0.51	0/1358	0.93	3/1845 (0.2%)
All	All	0.56	0/5416	0.96	14/7361 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	27	VAL	N-CA-C	6.89	119.55	109.63
1	A	91	VAL	N-CA-C	6.04	116.56	107.75
1	D	91	VAL	N-CA-C	6.00	118.09	108.85
1	D	44	TYR	N-CA-C	5.79	118.17	108.96
1	D	28	ALA	N-CA-C	-5.70	100.89	109.95
1	A	109	CYS	N-CA-C	5.62	119.61	112.87
1	C	27	VAL	N-CA-C	5.58	120.94	109.34
1	B	27	VAL	N-CA-C	5.54	117.43	109.51
1	C	150	GLY	CA-C-N	5.47	125.30	119.28
1	C	150	GLY	C-N-CA	5.47	125.30	119.28
1	B	109	CYS	N-CA-C	5.34	118.90	112.38
1	B	28	ALA	N-CA-C	-5.27	101.11	109.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	27	VAL	CB-CA-C	-5.17	102.89	111.32
1	C	49	TYR	N-CA-C	5.16	117.66	109.50

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	94	TYR	Sidechain

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	1315	0	1208	22	0
1	B	1319	0	1219	36	0
1	C	1307	0	1202	41	0
1	D	1319	0	1219	54	0
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	8	0	14	0	0
3	C	8	0	14	0	0
4	A	53	0	0	1	0
4	B	39	0	0	3	0
4	C	30	0	0	2	0
4	D	12	0	0	1	0
All	All	5414	0	4876	150	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 15.

All (150) 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:D:85:ARG:HH21	1:D:85:ARG:HB2	1.11	1.12

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:85:ARG:HB2	1:D:85:ARG:NH2	1.70	1.05
1:D:55:LYS:HZ3	1:D:58:LYS:HD3	1.26	0.98
1:D:55:LYS:NZ	1:D:58:LYS:HD3	1.92	0.84
1:D:55:LYS:HZ1	1:D:58:LYS:HA	1.43	0.83
1:B:125:MET:HG2	1:B:129:LYS:HZ1	1.42	0.83
1:B:125:MET:HE3	1:B:129:LYS:HZ2	1.50	0.76
1:B:125:MET:HE3	1:B:129:LYS:NZ	2.02	0.75
1:B:125:MET:HG2	1:B:129:LYS:HE3	1.69	0.73
1:D:52:GLY:O	1:D:63:VAL:HG22	1.90	0.72
1:C:27:VAL:HG11	1:C:116:LEU:HG	1.72	0.71
1:D:51:ILE:HG13	1:D:53:ILE:HG22	1.72	0.71
1:B:125:MET:HA	1:B:129:LYS:HE3	1.73	0.71
1:D:27:VAL:HG11	1:D:116:LEU:HG	1.73	0.69
1:D:77:ALA:HB2	1:D:103:LYS:HB3	1.73	0.69
1:D:27:VAL:HG21	1:D:116:LEU:HD23	1.75	0.69
1:B:125:MET:HG2	1:B:129:LYS:NZ	2.06	0.69
1:A:108:HIS:HD2	1:A:110:LEU:H	1.41	0.69
1:B:125:MET:HG2	1:B:129:LYS:CE	2.23	0.69
1:D:50:TRP:CD1	1:D:113:LYS:HB2	2.28	0.69
1:D:85:ARG:HH21	1:D:85:ARG:CB	1.99	0.66
1:D:107:GLU:HG3	1:D:111:LYS:HG3	1.78	0.66
1:C:57:ASN:C	1:C:58:LYS:HD3	2.21	0.65
1:A:85:ARG:HA	1:D:130:GLN:HE21	1.61	0.65
1:B:124:ASP:HB3	4:B:834:HOH:O	1.96	0.64
1:D:16:ARG:O	1:D:20:GLN:HG3	1.99	0.62
1:D:41:VAL:HG23	1:D:42:LEU:N	2.15	0.61
1:D:53:ILE:HB	1:D:62:TRP:CZ3	2.36	0.61
1:D:138:GLY:O	1:D:139:ASN:HB2	1.99	0.61
1:D:50:TRP:CE2	1:D:113:LYS:HD2	2.37	0.59
1:D:77:ALA:HA	1:D:103:LYS:HG2	1.84	0.58
1:B:19:CYS:SG	1:B:26:LEU:HD23	2.44	0.58
1:D:28:ALA:HA	1:D:62:TRP:CE3	2.39	0.58
1:A:76:TRP:CD2	1:A:81:PRO:HD3	2.39	0.57
1:D:128:SER:O	1:D:130:GLN:HG3	2.04	0.56
1:C:6:SER:HB3	1:C:18:TYR:CE1	2.41	0.56
1:A:22:ARG:HH11	1:A:22:ARG:HG3	1.71	0.56
1:B:94:TYR:HB3	1:B:97:SER:HB2	1.88	0.55
1:A:19:CYS:SG	1:A:26:LEU:HD23	2.47	0.55
1:C:18:TYR:O	1:C:22:ARG:HG3	2.06	0.55
1:D:16:ARG:NH2	1:D:63:VAL:HB	2.21	0.55
1:B:78:ASP:O	1:B:79:ASN:HB2	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:125:MET:CG	1:B:129:LYS:HZ1	2.17	0.54
1:D:50:TRP:CZ3	1:D:90:CYS:HB3	2.42	0.54
1:C:19:CYS:SG	1:C:26:LEU:HD23	2.48	0.54
1:D:139:ASN:CG	1:D:140:TYR:H	2.16	0.54
1:B:107:GLU:HG3	1:B:108:HIS:N	2.23	0.53
1:D:97:SER:O	1:D:101:PRO:HD3	2.08	0.53
1:D:78:ASP:O	1:D:79:ASN:HB2	2.09	0.53
1:C:22:ARG:HG3	1:C:22:ARG:HH11	1.72	0.53
1:C:27:VAL:CG1	1:C:116:LEU:HG	2.38	0.53
1:D:53:ILE:HB	1:D:62:TRP:CE3	2.42	0.53
1:C:91:VAL:CG1	1:C:104:TRP:HB3	2.39	0.53
1:C:49:TYR:O	1:C:92:GLU:HA	2.08	0.53
1:C:58:LYS:HD3	1:C:58:LYS:N	2.23	0.53
1:D:27:VAL:CG1	1:D:116:LEU:HG	2.38	0.53
1:B:50:TRP:CD1	1:B:113:LYS:HB2	2.44	0.53
1:D:151:PRO:HG2	4:D:807:HOH:O	2.09	0.52
1:D:26:LEU:HB3	1:D:52:GLY:N	2.25	0.52
1:B:2:THR:OG1	1:B:119:THR:HB	2.10	0.51
1:C:22:ARG:HH11	1:C:22:ARG:CG	2.24	0.51
1:C:94:TYR:HB3	1:C:97:SER:HB2	1.93	0.51
1:B:49:TYR:O	1:B:92:GLU:HA	2.10	0.51
1:B:56:ASN:C	1:B:57:ASN:CG	2.80	0.50
1:A:108:HIS:CD2	1:A:110:LEU:H	2.25	0.50
1:B:56:ASN:O	1:B:57:ASN:CG	2.54	0.50
1:A:22:ARG:HG3	1:A:22:ARG:NH1	2.26	0.50
1:A:39:ASN:HB3	4:A:851:HOH:O	2.12	0.50
1:B:128:SER:O	1:B:130:GLN:HG3	2.12	0.49
1:C:78:ASP:O	1:C:79:ASN:HB2	2.12	0.49
1:D:8:LYS:HB2	1:D:10:TYR:CZ	2.47	0.49
1:C:91:VAL:HG11	1:C:104:TRP:HB3	1.95	0.49
1:A:8:LYS:HE2	1:A:10:TYR:OH	2.12	0.49
1:C:108:HIS:HD2	1:C:110:LEU:HB2	1.78	0.49
1:C:50:TRP:CZ3	1:C:90:CYS:HB3	2.48	0.48
1:C:139:ASN:ND2	4:C:832:HOH:O	2.46	0.48
1:A:6:SER:O	1:A:114:HIS:HD2	1.96	0.48
1:D:12:TRP:HD1	1:D:50:TRP:CE3	2.31	0.48
1:D:18:TYR:C	1:D:18:TYR:CD1	2.92	0.48
1:C:33:ASN:HB3	1:D:149:TYR:CD2	2.48	0.48
1:B:65:THR:O	1:B:66:LYS:HB2	2.14	0.47
1:D:55:LYS:NZ	1:D:58:LYS:HA	2.22	0.47
1:C:94:TYR:HB3	1:C:97:SER:CB	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:67:LYS:HD2	4:C:829:HOH:O	2.14	0.47
1:D:77:ALA:CB	1:D:103:LYS:HB3	2.44	0.47
1:D:77:ALA:HA	1:D:103:LYS:CG	2.45	0.47
1:D:84:LYS:O	1:D:84:LYS:HG3	2.14	0.47
1:D:37:TYR:O	1:D:40:LYS:HG2	2.14	0.47
1:D:56:ASN:HB3	1:D:57:ASN:H	1.48	0.46
1:C:22:ARG:CG	1:C:22:ARG:NH1	2.79	0.46
1:C:77:ALA:HB2	1:C:103:LYS:HB3	1.97	0.46
1:D:18:TYR:CE1	1:D:22:ARG:HD3	2.50	0.46
1:A:139:ASN:CG	1:A:140:TYR:H	2.24	0.46
1:C:2:THR:O	1:C:118:TYR:HA	2.16	0.45
1:C:42:LEU:HB2	1:C:95:ILE:HG21	1.97	0.45
1:C:16:ARG:O	1:C:20:GLN:HG3	2.16	0.45
1:C:6:SER:HB3	1:C:18:TYR:CZ	2.51	0.45
1:D:41:VAL:HG23	1:D:42:LEU:H	1.78	0.45
1:B:2:THR:HG1	1:B:119:THR:HB	1.81	0.45
1:B:28:ALA:HA	1:B:62:TRP:CE3	2.51	0.45
1:C:93:ILE:HG23	1:C:95:ILE:HG12	1.98	0.45
1:C:125:MET:HA	1:C:129:LYS:HE3	1.98	0.45
1:D:125:MET:HA	1:D:125:MET:HE2	1.99	0.45
1:B:108:HIS:CD2	1:B:110:LEU:H	2.35	0.45
1:C:123:GLN:H	1:C:126:SER:HB2	1.82	0.45
1:C:123:GLN:NE2	1:C:123:GLN:HA	2.32	0.44
1:C:31:ASN:OD1	1:C:34:GLU:HG3	2.18	0.44
1:B:75:ASN:ND2	1:B:103:LYS:HA	2.32	0.44
1:B:77:ALA:HB2	1:B:103:LYS:HB3	2.00	0.44
1:C:123:GLN:HA	1:C:123:GLN:HE21	1.83	0.44
1:B:94:TYR:HB3	1:B:97:SER:CB	2.48	0.43
1:A:15:SER:HB3	1:A:115:ALA:HB2	1.99	0.43
1:A:134:LEU:HD23	1:A:141:THR:OG1	2.17	0.43
1:A:86:ASN:OD1	1:D:125:MET:HE1	2.19	0.43
1:C:125:MET:HG2	1:C:129:LYS:NZ	2.33	0.43
1:D:91:VAL:HG11	1:D:104:TRP:HB3	2.00	0.43
1:D:8:LYS:HB2	1:D:10:TYR:OH	2.19	0.42
1:D:41:VAL:CG2	1:D:42:LEU:N	2.80	0.42
1:A:97:SER:O	1:A:101:PRO:HD3	2.19	0.42
1:C:108:HIS:CD2	1:C:110:LEU:HB2	2.53	0.42
1:C:6:SER:O	1:C:114:HIS:HD2	2.02	0.42
1:B:107:GLU:CG	1:B:108:HIS:N	2.82	0.42
1:B:114:HIS:HE1	4:B:840:HOH:O	2.03	0.42
1:B:130:GLN:HB3	1:B:148:PHE:CG	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:ILE:HD11	1:A:60:TRP:HB3	2.01	0.41
1:B:27:VAL:HG11	1:B:116:LEU:HG	2.02	0.41
1:D:19:CYS:SG	1:D:26:LEU:HD23	2.60	0.41
1:A:159:ASP:O	1:A:160:ASP:CB	2.67	0.41
1:B:5:TYR:HA	1:B:115:ALA:O	2.20	0.41
1:B:8:LYS:HG2	4:B:831:HOH:O	2.19	0.41
1:D:94:TYR:HB3	1:D:97:SER:CB	2.51	0.41
1:A:128:SER:C	1:A:130:GLN:N	2.77	0.41
1:C:28:ALA:HA	1:C:62:TRP:CE3	2.56	0.41
1:C:53:ILE:HB	1:C:62:TRP:CZ3	2.55	0.41
1:A:47:SER:O	1:A:48:TYR:HB2	2.21	0.41
1:A:58:LYS:HG2	1:A:58:LYS:O	2.21	0.41
1:C:76:TRP:CD2	1:C:81:PRO:HD3	2.56	0.41
1:D:47:SER:O	1:D:48:TYR:HB2	2.20	0.41
1:C:5:TYR:HA	1:C:115:ALA:O	2.21	0.41
1:D:50:TRP:NE1	1:D:113:LYS:HB2	2.34	0.41
1:D:85:ARG:NH2	1:D:88:GLU:OE1	2.54	0.41
1:A:65:THR:O	1:A:66:LYS:HB2	2.21	0.40
1:B:2:THR:O	1:B:118:TYR:HA	2.20	0.40
1:D:12:TRP:CZ2	1:D:54:ARG:HD3	2.56	0.40
1:A:94:TYR:HB3	1:A:97:SER:HB2	2.03	0.40
1:B:87:ASN:ND2	1:B:109:CYS:HB2	2.36	0.40
1:C:27:VAL:HG13	1:C:116:LEU:C	2.46	0.40
1:C:53:ILE:HD11	1:C:60:TRP:HB3	2.04	0.40
1:B:55:LYS:C	1:B:56:ASN:HD22	2.29	0.40
1:B:125:MET:CE	1:B:129:LYS:NZ	2.80	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	158/162 (98%)	150 (95%)	8 (5%)	0	100	100
1	B	158/162 (98%)	147 (93%)	10 (6%)	1 (1%)	21	32
1	C	157/162 (97%)	142 (90%)	14 (9%)	1 (1%)	21	32
1	D	158/162 (98%)	137 (87%)	17 (11%)	4 (2%)	4	5
All	All	631/648 (97%)	576 (91%)	49 (8%)	6 (1%)	12	20

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	139	ASN
1	B	57	ASN
1	C	56	ASN
1	D	11	SER
1	D	52	GLY
1	D	94	TYR

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	142/147 (97%)	141 (99%)	1 (1%)	76	88
1	B	143/147 (97%)	141 (99%)	2 (1%)	59	79
1	C	141/147 (96%)	137 (97%)	4 (3%)	38	60
1	D	143/147 (97%)	140 (98%)	3 (2%)	47	69
All	All	569/588 (97%)	559 (98%)	10 (2%)	51	73

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	86	ASN
1	B	8	LYS
1	B	57	ASN
1	C	1	TRP
1	C	22	ARG

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Mol	Chain	Res	Type
1	C	58	LYS
1	C	156	VAL
1	D	85	ARG
1	D	121	SER
1	D	125	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	ASN
1	A	108	HIS
1	A	114	HIS
1	B	56	ASN
1	B	75	ASN
1	B	79	ASN
1	B	87	ASN
1	B	108	HIS
1	C	21	ASN
1	C	79	ASN
1	C	108	HIS
1	C	114	HIS
1	C	123	GLN
1	C	139	ASN
1	D	21	ASN
1	D	57	ASN
1	D	87	ASN
1	D	130	GLN

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 oligosaccharides 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	$\# Z > 2$
3	MRD	C	805	-	7,7,7	0.95	1 (14%)	9,10,10	0.46	0
3	MRD	A	806	-	7,7,7	0.76	0	9,10,10	0.41	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
3	MRD	C	805	-	-	0/5/5/5	-
3	MRD	A	806	-	-	0/5/5/5	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	805	MRD	O2-C2	-2.35	1.39	1.44

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

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.