



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 03:30 AM GMT

PDB ID : 1NGP
Title : N1G9 (IGG1-LAMBDA) FAB FRAGMENT COMPLEXED WITH (4-HYDROXY-3-NITROPHENYL)ACETATE
Authors : Mizutani, R.; Satow, Y.
Deposited on : 1995-06-23
Resolution : 2.40 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

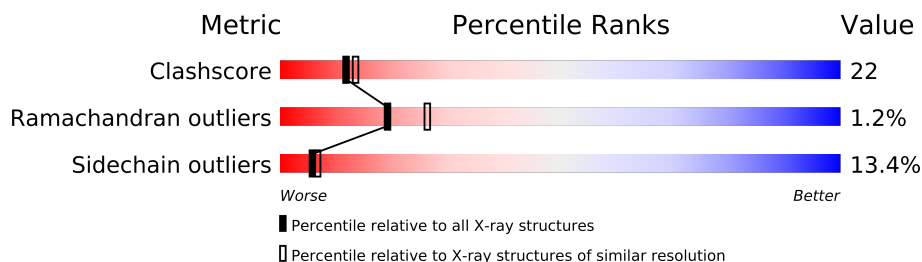
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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	79885	2789 (2.40-2.40)
Ramachandran outliers	78287	2736 (2.40-2.40)
Sidechain outliers	78261	2737 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	L	215	
2	H	222	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3380 atoms, of which 0 are hydrogen and 0 are deuterium.

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 N1G9 (IGG1-LAMBDA).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	211	Total	C	N	O	S	0	1	0
			1584	995	265	318	6			

- Molecule 2 is a protein called N1G9 (IGG1-LAMBDA).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	215	Total	C	N	O	S	0	1	0
			1646	1051	268	319	8			

There are 32 discrepancies between the modelled and reference sequences:

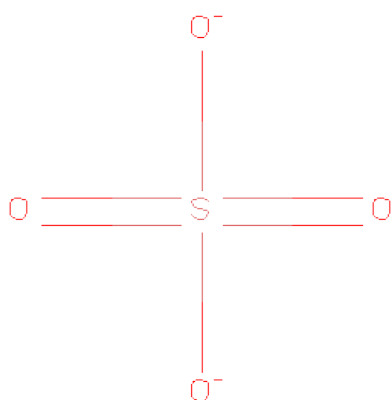
Chain	Residue	Modelled	Actual	Comment	Reference
H	1	GLN	GLU	CONFLICT	UNP P01751
H	3	GLN	LYS	CONFLICT	UNP P01751
H	5	GLN	HIS	CONFLICT	UNP P01751
H	6	GLN	GLU	CONFLICT	UNP P01751
H	7	PRO	SER	CONFLICT	UNP P01751
H	9	ALA	PRO	CONFLICT	UNP P01751
H	10	GLU	GLY	CONFLICT	UNP P01751
H	13	LYS	ARG	CONFLICT	UNP P01751
H	16	ALA	THR	CONFLICT	UNP P01751
H	20	LEU	ILE	CONFLICT	UNP P01751
H	31	SER	ASN	CONFLICT	UNP P01751
H	34	MET	LEU	CONFLICT	UNP P01751
H	35	HIS	GLY	CONFLICT	UNP P01751
H	43	ARG	HIS	CONFLICT	UNP P01751
H	50	ARG	ASP	CONFLICT	UNP P01751
H	52	ASP	TYR	CONFLICT	UNP P01751
H	54	ASN	GLY	CONFLICT	UNP P01751
H	55	SER	GLY	CONFLICT	UNP P01751
H	57	GLY	TYR	CONFLICT	UNP P01751
H	59	LYS	ASN	CONFLICT	UNP P01751

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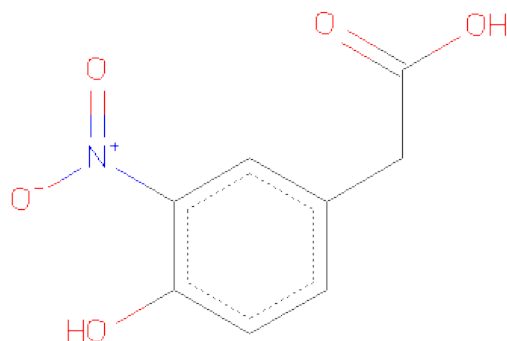
Chain	Residue	Modelled	Actual	Comment	Reference
H	66	SER	GLY	CONFLICT	UNP P01751
H	72	VAL	ALA	CONFLICT	UNP P01751
H	74	LYS	THR	CONFLICT	UNP P01751
H	75	PRO	SER	CONFLICT	UNP P01751
H	95	TYR	PHE	CONFLICT	UNP P01751
H	99	TYR	-	INSERTION	UNP P01751
H	100	ASP	PHE	CONFLICT	UNP P01751
H	?	-	TYR	DELETION	UNP P01751
H	?	-	TRP	DELETION	UNP P01751
H	109	TYR	VAL	CONFLICT	UNP P01751
H	116	LEU	VAL	CONFLICT	UNP P01751
H	194	PRO	THR	CONFLICT	UNP P01751

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	H	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is 2-(4-HYDROXY-3-NITROPHENYL)ACETICACID (three-letter code: NPA) (formula: C₈H₇NO₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	H	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 5 is water.

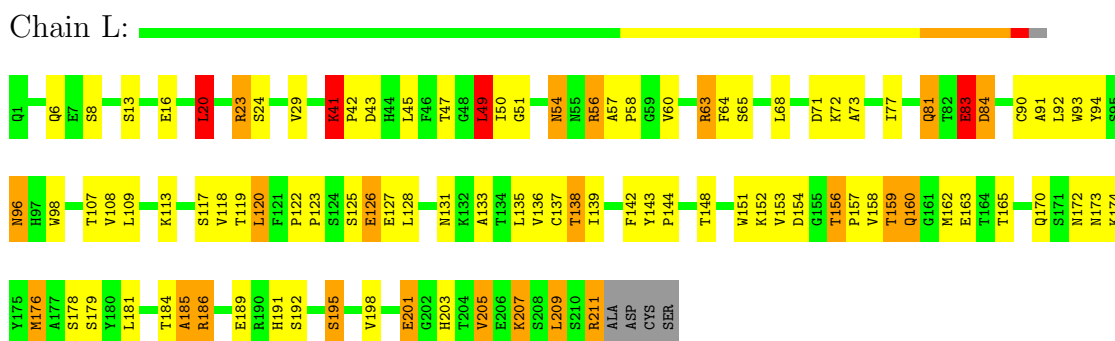
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	H	73	Total	O	0	0
			73	73		
5	L	58	Total	O	0	0
			58	58		

3 Residue-property plots

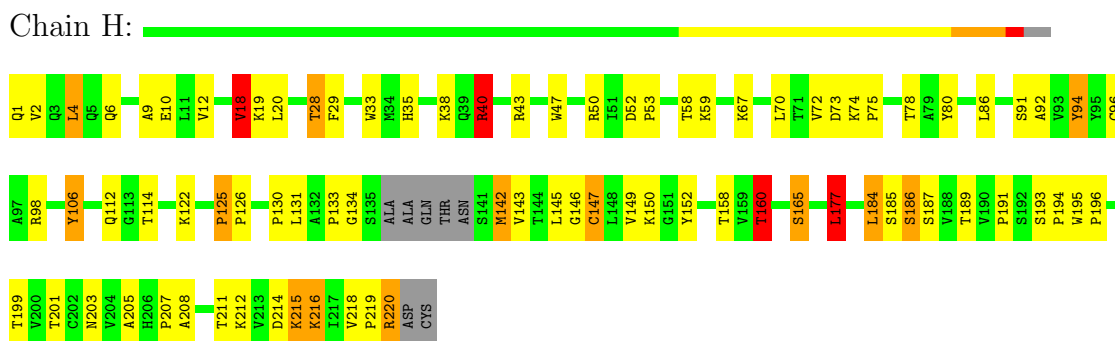
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

Note EDS was not executed.

• Molecule 1: N1G9 (IGG1-LAMBDA)



• Molecule 2: N1G9 (IGG1-LAMBDA)



4 Data and refinement statistics

Xtrriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	110.69Å 110.69Å 98.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.40	Depositor
% Data completeness (in resolution range)	89.3 (15.00-2.40)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ, X-PLOR 2.1	Depositor
R, R_{free}	0.196 , 0.250	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	3380	wwPDB-VP
Average B, all atoms (Å ²)	26.0	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: NPA, SO4

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	L	0.95	0/1627	1.65	21/2225 (0.9%)
2	H	0.94	0/1699	1.74	28/2320 (1.2%)
All	All	0.94	0/3326	1.70	49/4545 (1.1%)

There are no bond length outliers.

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	40	ARG	NE-CZ-NH1	16.16	128.38	120.30
2	H	94	TYR	CB-CG-CD2	11.04	127.63	121.00
1	L	56	ARG	NE-CZ-NH2	9.93	125.26	120.30
1	L	23	ARG	NE-CZ-NH1	9.74	125.17	120.30
1	L	143	TYR	CB-CG-CD1	-8.47	115.92	121.00
2	H	40	ARG	NE-CZ-NH2	-8.47	116.07	120.30
2	H	147	CYS	CB-CA-C	8.04	126.49	110.40
2	H	40	ARG	CD-NE-CZ	7.27	133.78	123.60
2	H	52	ASP	CB-CG-OD1	7.17	124.75	118.30
2	H	80	TYR	CB-CG-CD1	7.16	125.29	121.00
1	L	143	TYR	CB-CG-CD2	6.86	125.12	121.00
2	H	52	ASP	CB-CG-OD2	-6.86	112.13	118.30
1	L	209	LEU	CA-CB-CG	6.81	130.97	115.30
1	L	186	ARG	NE-CZ-NH1	6.58	123.59	120.30
1	L	20	LEU	CA-CB-CG	6.53	130.31	115.30
2	H	177	LEU	CA-CB-CG	6.51	130.28	115.30
1	L	84	ASP	CB-CG-OD1	6.44	124.10	118.30
1	L	98	TRP	N-CA-CB	6.33	121.99	110.60
2	H	214	ASP	CA-CB-CG	6.30	127.26	113.40
1	L	56	ARG	CD-NE-CZ	-6.27	114.82	123.60
1	L	71	ASP	CB-CG-OD1	6.14	123.83	118.30
2	H	94	TYR	CB-CG-CD1	-6.06	117.36	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	83	GLU	CG-CD-OE2	6.01	130.32	118.30
2	H	98	ARG	NE-CZ-NH1	-5.97	117.31	120.30
2	H	214	ASP	CB-CG-OD1	5.97	123.67	118.30
2	H	4	LEU	CA-CB-CG	5.95	128.97	115.30
1	L	41	LYS	CA-CB-CG	5.90	126.38	113.40
2	H	73	ASP	CB-CG-OD2	5.83	123.54	118.30
2	H	98	ARG	CD-NE-CZ	-5.83	115.44	123.60
1	L	126	GLU	OE1-CD-OE2	5.82	130.29	123.30
2	H	43	ARG	NE-CZ-NH1	5.80	123.20	120.30
1	L	65	SER	CB-CA-C	5.59	120.72	110.10
2	H	18	VAL	CB-CA-C	-5.57	100.82	111.40
1	L	49	LEU	CA-CB-CG	5.57	128.10	115.30
1	L	117	SER	O-C-N	5.51	131.52	122.70
2	H	9	ALA	N-CA-CB	-5.49	102.42	110.10
2	H	96	CYS	O-C-N	5.48	131.47	122.70
2	H	10	GLU	N-CA-CB	5.44	120.39	110.60
1	L	41	LYS	N-CA-CB	5.39	120.30	110.60
1	L	90	CYS	N-CA-CB	5.39	120.30	110.60
2	H	78	THR	CA-CB-OG1	-5.36	97.75	109.00
2	H	152	TYR	CB-CG-CD2	5.34	124.20	121.00
2	H	40	ARG	N-CA-C	-5.30	96.69	111.00
2	H	160	THR	N-CA-CB	-5.29	100.25	110.30
2	H	152	TYR	CB-CG-CD1	-5.28	117.83	121.00
2	H	147	CYS	CA-CB-SG	5.23	123.42	114.00
1	L	63	ARG	NE-CZ-NH1	-5.09	117.75	120.30
1	L	201	GLU	CB-CG-CD	5.05	127.83	114.20
2	H	165	SER	O-C-N	5.02	130.73	122.70

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	L	1584	0	1527	80	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	H	1646	0	1603	63	1
3	H	5	0	0	1	0
4	H	14	0	5	3	0
5	H	73	0	0	3	0
5	L	58	0	0	3	0
All	All	3380	0	3135	139	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 22.

All (139) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:L:211:ARG:HH11	1:L:211:ARG:HG2	1.14	1.06
1:L:153:VAL:HG12	1:L:158:VAL:HG22	1.40	1.02
2:H:35:HIS:CD2	2:H:47:TRP:HE1	1.87	0.91
2:H:35:HIS:HD2	2:H:47:TRP:HE1	0.94	0.89
1:L:51:GLY:HA3	2:H:106:TYR:HB3	1.56	0.88
1:L:189:GLU:O	1:L:211:ARG:NH2	2.06	0.88
1:L:131:ASN:HA	1:L:185:ALA:HB2	1.58	0.86
2:H:191:PRO:O	2:H:194:PRO:HD2	1.82	0.79
1:L:123:PRO:HD3	1:L:135:LEU:HD23	1.63	0.78
1:L:138:THR:HG22	5:L:234:HOH:O	1.84	0.77
1:L:123:PRO:HD3	1:L:135:LEU:CD2	2.15	0.76
2:H:35:HIS:HD2	2:H:47:TRP:NE1	1.78	0.74
1:L:153:VAL:HG12	1:L:158:VAL:CG2	2.17	0.73
1:L:131:ASN:HA	1:L:185:ALA:CB	2.18	0.73
2:H:59:LYS:NZ	4:H:302:NPA:O3	2.20	0.72
2:H:199:THR:HG21	2:H:216:LYS:HD3	1.71	0.71
1:L:211:ARG:NH1	1:L:211:ARG:HG2	1.89	0.71
1:L:56:ARG:NH1	1:L:64:PHE:O	2.24	0.70
1:L:41:LYS:HB3	1:L:42:PRO:HD2	1.73	0.70
1:L:122:PRO:HB3	1:L:209:LEU:HD21	1.74	0.69
2:H:191:PRO:C	2:H:194:PRO:HD2	2.14	0.68
2:H:130:PRO:O	2:H:131:LEU:HD23	1.94	0.68
1:L:172:ASN:O	1:L:173:ASN:HB2	1.95	0.67
1:L:120:LEU:HD13	1:L:209:LEU:HB2	1.77	0.67
2:H:193:SER:HB2	2:H:194:PRO:HD3	1.77	0.66
2:H:6:GLN:O	2:H:112:GLN:OE1	2.14	0.66
1:L:172:ASN:ND2	1:L:174:LYS:HD2	2.10	0.65
1:L:159:THR:O	1:L:160:GLN:HB2	1.97	0.65
1:L:122:PRO:HA	1:L:135:LEU:CD2	2.26	0.65
1:L:131:ASN:CA	1:L:185:ALA:HB2	2.26	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:L:51:GLY:CA	2:H:106:TYR:HB3	2.25	0.64
1:L:170:GLN:HG2	1:L:176:MET:HG2	1.79	0.63
2:H:18:VAL:HG22	2:H:86:LEU:HD11	1.80	0.63
1:L:185:ALA:O	1:L:189:GLU:HG2	2.00	0.61
2:H:18:VAL:CG2	2:H:86:LEU:HD11	2.31	0.61
2:H:184:LEU:HD23	2:H:185:SER:N	2.17	0.60
1:L:93:TRP:CH2	1:L:96:ASN:HA	2.36	0.60
2:H:184:LEU:HD23	2:H:184:LEU:C	2.21	0.60
2:H:218:VAL:HG23	2:H:219:PRO:HD2	1.83	0.60
2:H:146:GLY:HA2	2:H:187:SER:HA	1.83	0.59
1:L:81:GLN:HE21	1:L:81:GLN:HA	1.66	0.59
1:L:91:ALA:O	1:L:92:LEU:HD23	2.03	0.59
1:L:47:THR:HG23	5:H:320:HOH:O	2.01	0.59
1:L:64:PHE:CE1	1:L:77:ILE:HG12	2.39	0.58
1:L:24:SER:OG	1:L:94:TYR:OH	2.06	0.57
2:H:160:THR:HG22	2:H:203:ASN:HB2	1.86	0.57
2:H:199:THR:CG2	2:H:216:LYS:HD3	2.34	0.56
2:H:219:PRO:O	2:H:220:ARG:HB2	2.04	0.56
2:H:40:ARG:NH1	3:H:301:SO4:O4	2.38	0.56
1:L:107:THR:HG21	1:L:144:PRO:HB3	1.89	0.55
1:L:29:VAL:HG11	1:L:73:ALA:HB2	1.88	0.55
1:L:162:MET:HG3	1:L:162:MET:O	2.07	0.54
1:L:165:THR:HB	1:L:178:SER:H	1.72	0.54
1:L:184:THR:C	1:L:186:ARG:N	2.59	0.54
1:L:189:GLU:C	1:L:211:ARG:HH21	2.07	0.54
1:L:159:THR:O	1:L:159:THR:HG23	2.08	0.53
1:L:123:PRO:CG	1:L:133:ALA:HB1	2.37	0.53
2:H:35:HIS:CD2	2:H:50:ARG:HB3	2.44	0.53
1:L:83:GLU:H	1:L:83:GLU:CD	2.12	0.53
1:L:122:PRO:HA	1:L:135:LEU:HD23	1.89	0.53
2:H:196:PRO:HB3	2:H:219:PRO:HG3	1.91	0.52
1:L:96:ASN:N	1:L:96:ASN:ND2	2.55	0.52
2:H:70:LEU:N	2:H:70:LEU:HD23	2.23	0.52
1:L:184:THR:O	1:L:186:ARG:N	2.42	0.52
1:L:128:LEU:O	1:L:131:ASN:N	2.40	0.51
1:L:123:PRO:HD3	1:L:135:LEU:HD21	1.93	0.51
1:L:156:THR:O	1:L:157:PRO:C	2.49	0.50
1:L:96:ASN:N	1:L:96:ASN:HD22	2.10	0.50
1:L:93:TRP:CZ2	1:L:96:ASN:HA	2.46	0.50
2:H:142[B]:MET:HE3	2:H:189:THR:HG22	1.95	0.49
2:H:131:LEU:HB2	2:H:146:GLY:C	2.33	0.48
1:L:13:SER:H	1:L:16:GLU:HG3	1.77	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:L:118:VAL:HG21	1:L:205:VAL:HG22	1.96	0.48
1:L:119:THR:O	1:L:137:CYS:HA	2.13	0.48
1:L:20:LEU:HD23	1:L:20:LEU:N	2.29	0.48
1:L:81:GLN:CA	1:L:81:GLN:HE21	2.27	0.48
2:H:18:VAL:HG22	2:H:86:LEU:CD1	2.44	0.47
1:L:54:ASN:C	1:L:54:ASN:HD22	2.18	0.47
1:L:184:THR:OG1	1:L:186:ARG:HB2	2.14	0.47
1:L:152:LYS:O	1:L:195:SER:N	2.47	0.47
1:L:158:VAL:O	1:L:162:MET:HE3	2.15	0.47
2:H:12:VAL:HG11	2:H:18:VAL:HG13	1.97	0.47
2:H:201:THR:HG22	2:H:203:ASN:ND2	2.30	0.47
1:L:127:GLU:OE2	2:H:150:LYS:NZ	2.47	0.47
2:H:74:LYS:N	2:H:75:PRO:CD	2.78	0.47
2:H:220:ARG:HH11	2:H:220:ARG:HG2	1.79	0.46
1:L:81:GLN:HB3	1:L:83:GLU:HG2	1.97	0.46
2:H:130:PRO:HD3	2:H:215:LYS:HE2	1.97	0.46
1:L:49:LEU:HA	1:L:60:VAL:HG21	1.98	0.46
2:H:220:ARG:NH1	2:H:220:ARG:HG2	2.30	0.46
1:L:184:THR:O	1:L:185:ALA:C	2.53	0.46
1:L:68:LEU:HD23	1:L:73:ALA:HA	1.97	0.46
2:H:126:PRO:HB2	2:H:149:VAL:HG13	1.96	0.45
1:L:63:ARG:NH2	1:L:84:ASP:OD1	2.36	0.45
2:H:50:ARG:O	2:H:58:THR:HA	2.16	0.45
2:H:196:PRO:CB	2:H:219:PRO:HG3	2.47	0.45
2:H:53:PRO:HA	2:H:72:VAL:HG21	1.98	0.45
2:H:146:GLY:HA2	2:H:186:SER:O	2.17	0.45
1:L:120:LEU:HA	1:L:136:VAL:O	2.17	0.45
2:H:112:GLN:O	2:H:112:GLN:HG3	2.15	0.45
1:L:23:ARG:HG2	1:L:72:LYS:HD3	1.99	0.45
1:L:126:GLU:OE1	2:H:215:LYS:NZ	2.45	0.45
1:L:151:TRP:HA	1:L:195:SER:O	2.16	0.45
1:L:148:THR:O	1:L:198:VAL:HA	2.17	0.45
1:L:96:ASN:HD22	1:L:96:ASN:H	1.64	0.44
1:L:57:ALA:HB1	1:L:58:PRO:HD2	1.99	0.44
2:H:12:VAL:HG21	2:H:86:LEU:CD1	2.48	0.44
2:H:53:PRO:HB3	2:H:72:VAL:HG21	1.99	0.44
2:H:205:ALA:HB2	2:H:212:LYS:HD2	1.99	0.44
2:H:195:TRP:CG	2:H:196:PRO:HA	2.52	0.44
2:H:125:PRO:O	2:H:125:PRO:HG2	2.17	0.44
2:H:158:THR:HG22	5:H:323:HOH:O	2.16	0.44
2:H:207:PRO:O	2:H:208:ALA:C	2.56	0.44
1:L:50:ILE:HG22	1:L:51:GLY:N	2.33	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:145:LEU:O	2:H:187:SER:HA	2.18	0.43
1:L:131:ASN:C	1:L:185:ALA:HB2	2.38	0.43
1:L:139:ILE:HG22	1:L:142:PHE:CE1	2.53	0.43
1:L:154:ASP:OD2	1:L:191:HIS:HB3	2.18	0.43
2:H:50:ARG:HG2	4:H:302:NPA:O2	2.19	0.43
2:H:126:PRO:CB	2:H:149:VAL:HG13	2.48	0.43
1:L:184:THR:C	1:L:186:ARG:H	2.20	0.43
2:H:191:PRO:CB	2:H:194:PRO:CD	2.97	0.43
2:H:28:THR:O	2:H:29:PHE:C	2.57	0.42
2:H:177:LEU:HD11	5:H:321:HOH:O	2.18	0.42
1:L:158:VAL:O	1:L:162:MET:CE	2.68	0.42
2:H:195:TRP:CD1	2:H:196:PRO:HA	2.55	0.41
1:L:81:GLN:O	1:L:108:VAL:HG11	2.19	0.41
1:L:6:GLN:HB3	5:L:221:HOH:O	2.19	0.41
1:L:16:GLU:OE2	5:L:248:HOH:O	2.21	0.41
2:H:38:LYS:HG3	2:H:94:TYR:CE1	2.55	0.41
1:L:113:LYS:HG2	1:L:144:PRO:HD3	2.02	0.41
2:H:12:VAL:HG21	2:H:86:LEU:HD13	2.02	0.41
2:H:20:LEU:HD22	2:H:114:THR:HG21	2.03	0.41
1:L:207:LYS:HA	1:L:207:LYS:HD2	1.72	0.41
1:L:170:GLN:CG	1:L:176:MET:HG2	2.50	0.41
2:H:33:TRP:HZ3	4:H:302:NPA:C4	2.34	0.40
2:H:91:SER:O	2:H:92:ALA:HB2	2.22	0.40
2:H:130:PRO:C	2:H:131:LEU:HD23	2.42	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	Distance(Å)	Clash(Å)
2:H:19:LYS:NZ	2:H:19:LYS:NZ[8.665]	2.12	0.08

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	L	210/215 (98%)	190 (90%)	17 (8%)	3 (1%)	16	22
2	H	212/222 (96%)	191 (90%)	19 (9%)	2 (1%)	25	35
All	All	422/437 (97%)	381 (90%)	36 (8%)	5 (1%)	19	26

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	43	ASP
2	H	133	PRO
1	L	160	GLN
1	L	185	ALA
2	H	134	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	L	175/182 (96%)	149 (85%)	26 (15%)	4	5
2	H	186/192 (97%)	163 (88%)	23 (12%)	7	8
All	All	361/374 (96%)	312 (86%)	49 (14%)	6	6

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	8	SER
1	L	20	LEU
1	L	41	LYS
1	L	45	LEU
1	L	49	LEU
1	L	54	ASN
1	L	81	GLN
1	L	83	GLU
1	L	96	ASN
1	L	109	LEU
1	L	120	LEU
1	L	125	SER

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Mol	Chain	Res	Type
1	L	138	THR
1	L	156	THR
1	L	159	THR
1	L	163	GLU
1	L	176	MET
1	L	179	SER
1	L	181	LEU
1	L	192	SER
1	L	195	SER
1	L	201	GLU
1	L	203	HIS
1	L	205	VAL
1	L	207	LYS
1	L	211	ARG
2	H	1	GLN
2	H	2	VAL
2	H	4	LEU
2	H	18	VAL
2	H	28	THR
2	H	40	ARG
2	H	67	LYS
2	H	106	TYR
2	H	122	LYS
2	H	125	PRO
2	H	142[A]	MET
2	H	142[B]	MET
2	H	143	VAL
2	H	147	CYS
2	H	160	THR
2	H	165	SER
2	H	177	LEU
2	H	184	LEU
2	H	186	SER
2	H	211	THR
2	H	215	LYS
2	H	216	LYS
2	H	220	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	54	ASN

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Mol	Chain	Res	Type
1	L	81	GLN
1	L	96	ASN
2	H	1	GLN
2	H	35	HIS

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

2 ligands are 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	$\# Z > 2$
3	SO4	H	301	-	4,4,4	1.01	0	6,6,6	0.44	0
4	NPA	H	302	-	14,14,14	1.51	3 (21%)	19,19,19	1.48	4 (21%)

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	SO4	H	301	-	-	0/0/0/0	0/0/0/0
4	NPA	H	302	-	-	0/8/8/8	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	302	NPA	C3-N1	-3.52	1.41	1.46
4	H	302	NPA	O1-N1	-2.44	1.21	1.25
4	H	302	NPA	C2-C1	-2.42	1.34	1.39

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	302	NPA	O3-C4-C3	2.78	128.22	120.02
4	H	302	NPA	O4-C8-C7	-2.52	116.99	123.19
4	H	302	NPA	O3-C4-C5	-2.04	113.77	119.37
4	H	302	NPA	C2-C3-C4	2.00	123.67	121.38

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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 will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.