



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 16, 2016 – 08:11 PM EST

PDB ID : 5K8O
Title : Mn²⁺/5NSA-bound 5-nitroanthranilate aminohydrolase
Authors : Kalyoncu, S.
Deposited on : 2016-05-30
Resolution : 2.89 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20028320
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) : rb-20028320

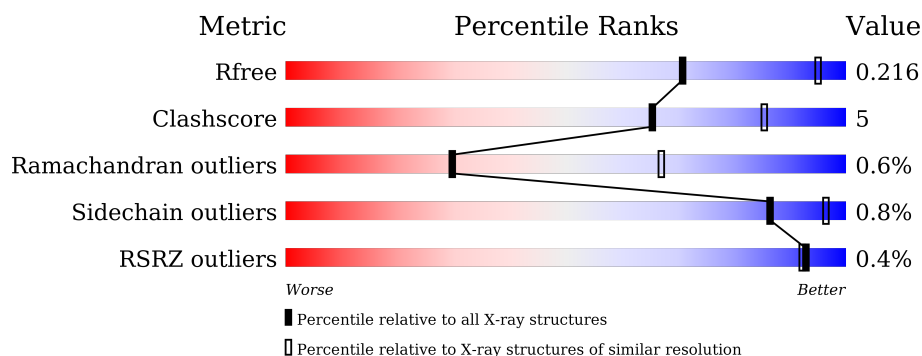
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.89 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	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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	425	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> % 86% 12% .. </div> </div>
1	B	425	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green, yellow, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 88% 11% . </div> </div>
1	C	425	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green, yellow, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 87% 12% . </div> </div>
1	D	425	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> % 84% 16% </div> </div>
1	E	425	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green, yellow, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 89% 10% . </div> </div>
1	F	425	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green, yellow, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 87% 11% .. </div> </div>

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Mol	Chain	Length	Quality of chain
1	G	425	 88% 11% .
1	H	425	 86% 13% .

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 26050 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 5-nitroanthranilic acid aminohydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	422	Total	C	N	O	S	0	0	0
			3241	2049	563	610	19			
1	B	425	Total	C	N	O	S	0	0	0
			3258	2059	566	613	20			
1	C	421	Total	C	N	O	S	0	0	0
			3235	2046	562	608	19			
1	D	425	Total	C	N	O	S	0	0	0
			3258	2059	566	613	20			
1	E	421	Total	C	N	O	S	0	0	0
			3235	2046	562	608	19			
1	F	421	Total	C	N	O	S	0	0	0
			3235	2046	562	608	19			
1	G	422	Total	C	N	O	S	0	0	0
			3241	2049	563	610	19			
1	H	421	Total	C	N	O	S	0	0	0
			3235	2046	562	608	19			

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

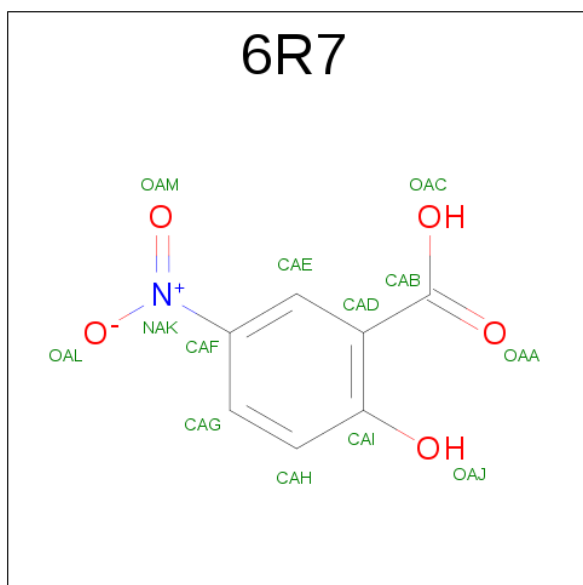
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	1	Total	Mn	0	0
			1	1		
2	D	1	Total	Mn	0	0
			1	1		
2	E	1	Total	Mn	0	0
			1	1		
2	H	1	Total	Mn	0	0
			1	1		
2	B	1	Total	Mn	0	0
			1	1		
2	C	1	Total	Mn	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mn	0	0
			1	1		
2	F	1	Total	Mn	0	0
			1	1		

- Molecule 3 is 5-nitrosalicylic acid (three-letter code: 6R7) (formula: $C_7H_5NO_5$).

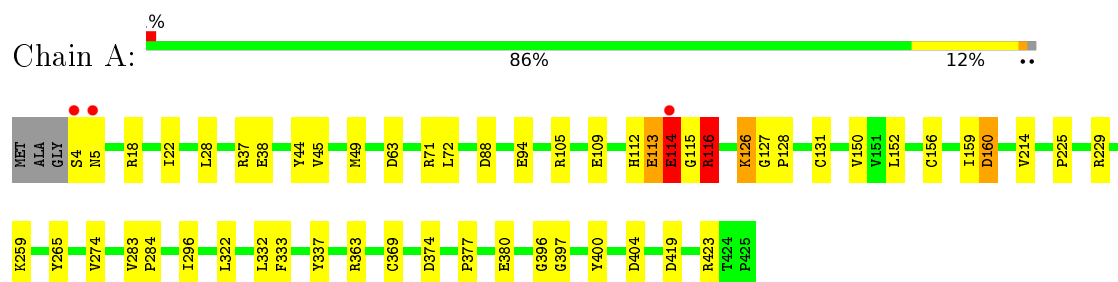


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			13	7	1	5		
3	B	1	Total	C	N	O	0	0
			13	7	1	5		
3	C	1	Total	C	N	O	0	0
			13	7	1	5		
3	D	1	Total	C	N	O	0	0
			13	7	1	5		
3	E	1	Total	C	N	O	0	0
			13	7	1	5		
3	F	1	Total	C	N	O	0	0
			13	7	1	5		
3	G	1	Total	C	N	O	1	0
			13	7	1	5		
3	H	1	Total	C	N	O	0	0
			13	7	1	5		

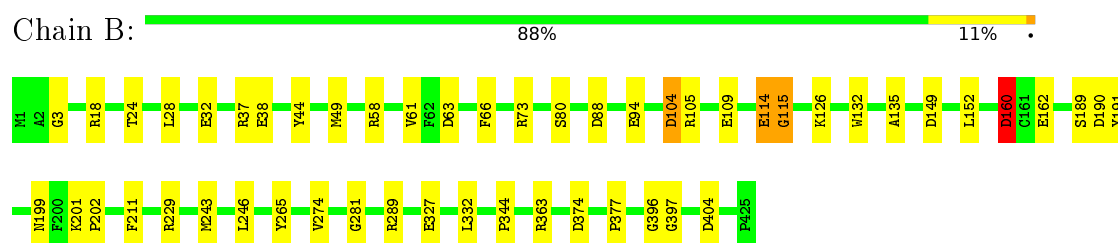
3 Residue-property plots [i](#)

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.

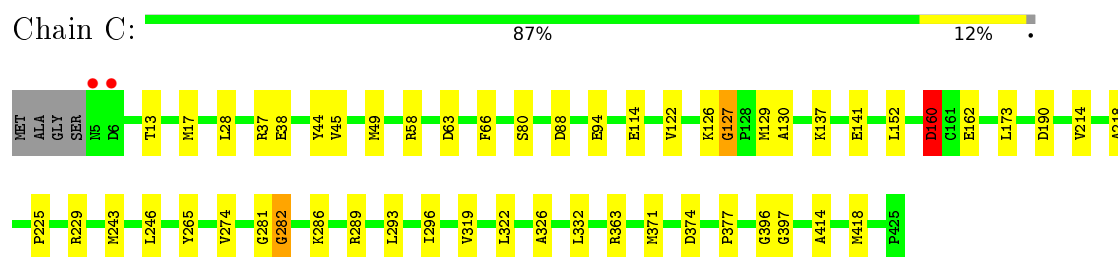
- Molecule 1: 5-nitroanthranilic acid aminohydrolase



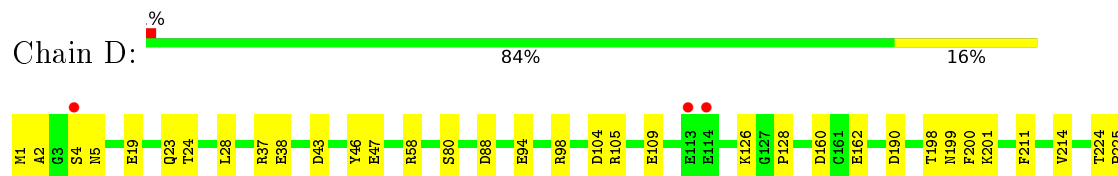
- Molecule 1: 5-nitroanthranilic acid aminohydrolase



- Molecule 1: 5-nitroanthranilic acid aminohydrolase



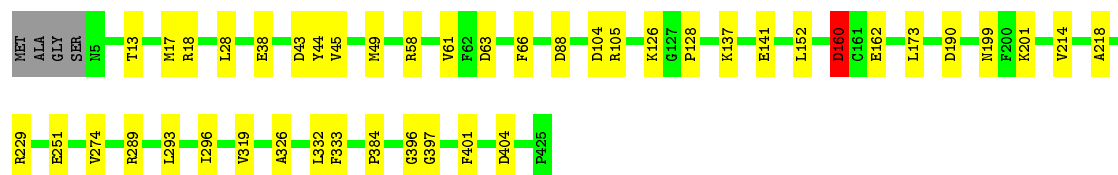
- Molecule 1: 5-nitroanthranilic acid aminohydrolase





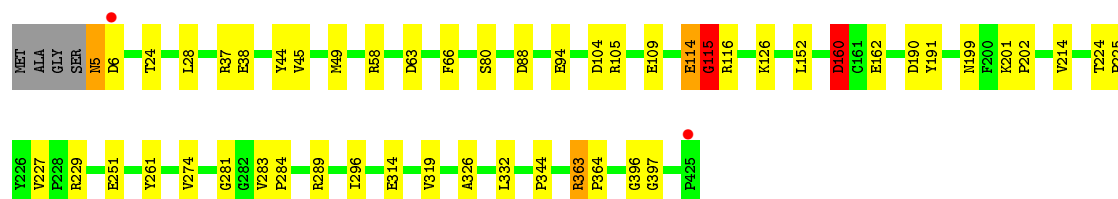
- Molecule 1: 5-nitroanthranilic acid aminohydrolase

Chain E: 89% 10%



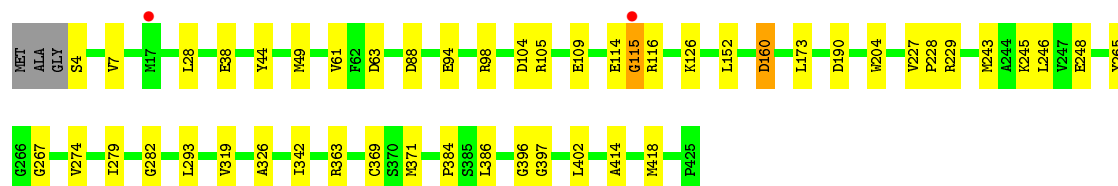
- Molecule 1: 5-nitroanthranilic acid aminohydrolase

Chain F: 87% 11%



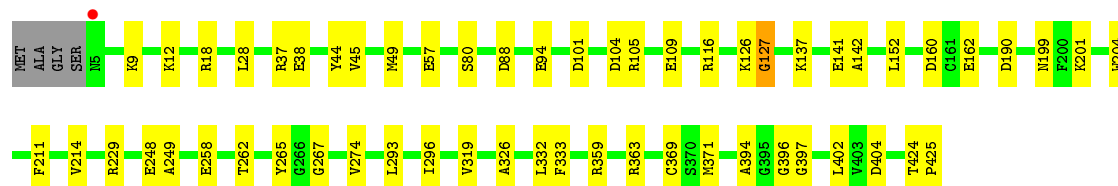
- Molecule 1: 5-nitroanthranilic acid aminohydrolase

Chain G: 88% 11%



- Molecule 1: 5-nitroanthranilic acid aminohydrolase

Chain H: 86% 13%



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	186.77Å 249.34Å 249.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.83 – 2.89 49.83 – 2.89	Depositor EDS
% Data completeness (in resolution range)	92.2 (49.83-2.89) 92.2 (49.83-2.89)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.57 (at 2.91Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.184 , 0.218 0.179 , 0.216	Depositor DCC
R_{free} test set	1998 reflections (1.68%)	DCC
Wilson B-factor (Å ²)	51.3	Xtriage
Anisotropy	0.163	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 16.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	26050	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.13% 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: MN, 6R7

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	0.65	1/3315 (0.0%)	0.69	5/4499 (0.1%)
1	B	0.66	0/3332	0.62	2/4521 (0.0%)
1	C	0.65	0/3309	0.60	1/4491 (0.0%)
1	D	0.66	0/3332	0.63	0/4521
1	E	0.66	0/3309	0.61	0/4491
1	F	0.66	0/3309	0.63	1/4491 (0.0%)
1	G	0.61	1/3315 (0.0%)	0.61	1/4499 (0.0%)
1	H	0.64	1/3309 (0.0%)	0.62	1/4491 (0.0%)
All	All	0.65	3/26530 (0.0%)	0.63	11/36004 (0.0%)

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	2
1	B	0	1
1	C	0	1
All	All	0	4

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	369	CYS	CB-SG	-5.74	1.72	1.81
1	H	369	CYS	CB-SG	-5.63	1.72	1.81
1	G	369	CYS	CB-SG	-5.12	1.73	1.81

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	115	GLY	N-CA-C	-9.24	90.00	113.10
1	A	127	GLY	N-CA-C	-6.66	96.45	113.10
1	A	116	ARG	CA-CB-CG	6.58	127.89	113.40
1	B	115	GLY	N-CA-C	-6.39	97.12	113.10
1	A	116	ARG	NE-CZ-NH1	6.00	123.30	120.30
1	A	126	LYS	N-CA-C	5.66	126.28	111.00
1	B	281	GLY	N-CA-C	-5.53	99.28	113.10
1	G	282	GLY	N-CA-C	-5.39	99.63	113.10
1	H	127	GLY	N-CA-C	-5.28	99.91	113.10
1	C	282	GLY	N-CA-C	-5.07	100.43	113.10
1	A	71	ARG	NE-CZ-NH2	-5.07	117.77	120.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	113	GLU	Peptide
1	A	114	GLU	Peptide
1	B	114	GLU	Peptide
1	C	127	GLY	Peptide

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	3241	0	3170	40	0
1	B	3258	0	3190	36	0
1	C	3235	0	3165	34	0
1	D	3258	0	3190	38	0
1	E	3235	0	3165	27	0
1	F	3235	0	3165	33	0
1	G	3241	0	3170	31	0
1	H	3235	0	3165	33	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
2	E	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
3	A	13	0	0	1	0
3	B	13	0	0	2	0
3	C	13	0	0	0	0
3	D	13	0	0	1	0
3	E	13	0	0	2	0
3	F	13	0	0	0	0
3	G	13	0	0	1	0
3	H	13	0	0	1	0
All	All	26050	0	25380	256	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 (256) 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:F:363:ARG:HD3	1:F:364:PRO:HD2	1.27	1.13
1:A:114:GLU:HG2	1:A:116:ARG:HG3	1.46	0.94
1:G:114:GLU:O	1:G:116:ARG:N	2.05	0.89
1:A:265:TYR:O	1:A:363:ARG:NH2	2.07	0.86
1:B:32:GLU:OE2	1:B:37:ARG:NH1	2.08	0.86
1:G:28:LEU:HG	1:G:126:LYS:HD2	1.56	0.85
1:H:265:TYR:O	1:H:363:ARG:NH2	2.12	0.83
1:A:28:LEU:HG	1:A:126:LYS:HD2	1.62	0.81
1:A:114:GLU:HG3	1:A:116:ARG:N	1.97	0.80
1:B:28:LEU:HG	1:B:126:LYS:HD2	1.62	0.79
1:B:18:ARG:NH2	1:B:404:ASP:OD1	2.15	0.79
1:D:28:LEU:HG	1:D:126:LYS:HD2	1.65	0.79
1:F:28:LEU:HG	1:F:126:LYS:HD2	1.62	0.79
1:C:28:LEU:HG	1:C:126:LYS:HD3	1.66	0.78
1:B:265:TYR:O	1:B:363:ARG:NH2	2.17	0.78
1:E:28:LEU:HG	1:E:126:LYS:HD2	1.64	0.77
1:H:28:LEU:HG	1:H:126:LYS:HD2	1.66	0.77
1:F:37:ARG:NH2	1:F:94:GLU:OE2	2.16	0.77
1:C:37:ARG:NH2	1:C:94:GLU:OE2	2.17	0.76
1:D:265:TYR:O	1:D:363:ARG:NH2	2.17	0.76
1:F:274:VAL:O	1:G:229:ARG:NH1	2.20	0.75
1:B:104:ASP:OD1	1:B:105:ARG:N	2.20	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:265:TYR:O	1:G:363:ARG:NH2	2.19	0.74
1:F:105:ARG:NH1	1:F:109:GLU:OE2	2.20	0.74
1:A:274:VAL:O	1:H:229:ARG:NH1	2.19	0.74
1:H:105:ARG:NH1	1:H:109:GLU:OE2	2.20	0.74
1:H:214:VAL:HB	1:H:296:ILE:HG22	1.69	0.73
1:C:229:ARG:NH1	1:D:274:VAL:O	2.23	0.72
1:B:105:ARG:NH1	1:B:109:GLU:OE2	2.22	0.72
1:D:214:VAL:HB	1:D:296:ILE:HG22	1.74	0.70
1:A:114:GLU:CG	1:A:116:ARG:HG3	2.22	0.70
1:E:214:VAL:HB	1:E:296:ILE:HG22	1.74	0.69
1:A:37:ARG:NH2	1:A:94:GLU:OE2	2.17	0.68
1:H:38:GLU:HG2	1:H:88:ASP:HB3	1.74	0.68
1:B:274:VAL:O	1:E:229:ARG:NH1	2.26	0.67
1:D:350:GLU:O	1:D:354:ARG:HG3	1.94	0.67
1:B:37:ARG:NH2	1:B:94:GLU:OE2	2.27	0.67
1:H:248:GLU:HG3	1:H:249:ALA:N	2.10	0.66
1:C:265:TYR:O	1:C:363:ARG:NH2	2.29	0.66
1:D:37:ARG:NH2	1:D:94:GLU:OE2	2.29	0.65
1:D:43:ASP:O	1:D:47:GLU:HG2	1.96	0.65
1:A:116:ARG:HH11	1:A:116:ARG:HG2	1.62	0.65
1:A:214:VAL:HB	1:A:296:ILE:HG22	1.79	0.64
1:G:105:ARG:NH1	1:G:109:GLU:OE2	2.31	0.64
1:G:114:GLU:C	1:G:116:ARG:H	2.02	0.63
1:H:37:ARG:NH2	1:H:94:GLU:OE2	2.25	0.63
1:F:114:GLU:OE1	1:F:116:ARG:NE	2.33	0.62
1:C:122:VAL:HA	1:C:126:LYS:O	1.99	0.62
1:D:105:ARG:NH1	1:D:109:GLU:OE1	2.33	0.62
1:F:214:VAL:HB	1:F:296:ILE:HG22	1.82	0.62
3:H:502:6R7:OAC	3:H:502:6R7:OAJ	2.15	0.61
1:G:38:GLU:HG2	1:G:88:ASP:HB3	1.83	0.61
1:A:113:GLU:O	1:A:115:GLY:N	2.32	0.60
1:A:114:GLU:HG3	1:A:116:ARG:H	1.67	0.60
1:A:4:SER:OG	1:A:5:ASN:N	2.34	0.59
1:B:88:ASP:N	1:B:88:ASP:OD1	2.35	0.59
1:B:229:ARG:NH1	1:E:274:VAL:O	2.34	0.59
1:A:18:ARG:NH1	1:A:404:ASP:OD1	2.36	0.58
1:A:229:ARG:NH1	1:H:274:VAL:O	2.37	0.58
1:A:423:ARG:HG3	1:A:423:ARG:HH11	1.69	0.58
1:F:289:ARG:NH1	3:G:502:6R7:OAL	2.37	0.58
1:D:43:ASP:OD1	1:D:58:ARG:NH1	2.38	0.57
3:B:502:6R7:OAL	1:E:289:ARG:NH1	2.39	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:319:VAL:HG11	1:D:326:ALA:HB3	1.87	0.56
1:D:80:SER:HB2	1:D:190:ASP:H	1.70	0.56
1:F:114:GLU:HG2	1:F:115:GLY:N	2.20	0.56
1:H:248:GLU:HG3	1:H:249:ALA:H	1.70	0.56
1:D:38:GLU:HG2	1:D:88:ASP:HB3	1.87	0.56
1:G:396:GLY:N	1:G:397:GLY:HA2	2.21	0.56
1:F:162:GLU:HG3	1:F:332:LEU:HD22	1.88	0.56
1:F:38:GLU:HG2	1:F:88:ASP:HB3	1.87	0.55
1:H:258:GLU:O	1:H:262:THR:OG1	2.23	0.55
1:A:116:ARG:HH11	1:A:116:ARG:CG	2.18	0.55
1:C:80:SER:HB2	1:C:190:ASP:H	1.72	0.55
1:C:274:VAL:O	1:D:229:ARG:NH1	2.40	0.54
1:G:114:GLU:HG3	1:G:115:GLY:N	2.22	0.54
1:B:396:GLY:N	1:B:397:GLY:HA2	2.22	0.54
1:C:214:VAL:HB	1:C:296:ILE:HG22	1.88	0.54
1:F:229:ARG:NH1	1:G:274:VAL:O	2.41	0.54
1:D:396:GLY:N	1:D:397:GLY:HA2	2.22	0.54
1:D:1:MET:HG3	1:D:2:ALA:N	2.22	0.54
1:E:38:GLU:HG2	1:E:88:ASP:HB3	1.90	0.54
1:B:80:SER:HB2	1:B:190:ASP:H	1.73	0.53
1:E:396:GLY:N	1:E:397:GLY:HA2	2.22	0.53
1:H:162:GLU:HG3	1:H:332:LEU:HD13	1.91	0.53
1:A:49:MET:HE1	1:A:152:LEU:HB2	1.91	0.52
1:C:38:GLU:HG2	1:C:88:ASP:HB3	1.90	0.52
1:E:58:ARG:HG2	1:E:66:PHE:CD1	2.44	0.52
1:E:128:PRO:HD3	1:E:401:PHE:HE2	1.75	0.52
1:G:245:LYS:O	1:G:248:GLU:HG2	2.10	0.52
1:C:28:LEU:HD13	1:C:44:TYR:CD1	2.45	0.51
1:A:114:GLU:HG2	1:A:116:ARG:CG	2.32	0.51
1:A:63:ASP:N	1:A:63:ASP:OD1	2.39	0.51
1:C:63:ASP:N	1:C:63:ASP:OD1	2.44	0.51
1:D:211:PHE:HB2	1:D:332:LEU:HB3	1.91	0.51
1:E:162:GLU:HG3	1:E:332:LEU:HD22	1.92	0.51
1:H:18:ARG:NH2	1:H:404:ASP:OD1	2.30	0.51
1:F:261:TYR:OH	1:F:314:GLU:OE2	2.26	0.51
1:D:5:ASN:OD1	1:D:5:ASN:N	2.38	0.51
1:A:225:PRO:HG2	1:H:371:MET:HB3	1.93	0.51
1:H:211:PHE:HB2	1:H:332:LEU:HB3	1.92	0.51
1:C:396:GLY:N	1:C:397:GLY:HA2	2.26	0.50
1:F:49:MET:HE1	1:F:152:LEU:HB2	1.93	0.50
1:B:38:GLU:HG2	1:B:88:ASP:HB3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:127:GLY:CA	1:C:130:ALA:H	2.25	0.50
1:C:371:MET:HB3	1:D:225:PRO:HG2	1.93	0.50
1:E:319:VAL:HG11	1:E:326:ALA:HB3	1.94	0.50
3:B:502:6R7:OAC	3:B:502:6R7:OAJ	2.27	0.50
1:D:23:GLN:HG3	1:D:24:THR:N	2.25	0.50
1:E:61:VAL:HG13	1:G:173:LEU:HD11	1.93	0.50
1:G:414:ALA:O	1:G:418:MET:HG3	2.12	0.50
1:H:45:VAL:O	1:H:49:MET:HG3	2.12	0.50
1:C:162:GLU:HG3	1:C:332:LEU:HD22	1.94	0.49
1:E:49:MET:HE1	1:E:152:LEU:HB2	1.94	0.49
1:D:342:ILE:HD11	1:D:345:LEU:HD23	1.94	0.49
1:C:286:LYS:HD3	1:D:98:ARG:NH2	2.28	0.49
1:D:199:ASN:O	1:D:201:LYS:HG3	2.13	0.49
1:C:289:ARG:NH1	3:D:502:6R7:OAL	2.46	0.49
1:A:22:ILE:HD13	1:A:112:HIS:CD2	2.48	0.49
1:F:319:VAL:HG11	1:F:326:ALA:HB3	1.94	0.48
3:A:502:6R7:OAJ	3:A:502:6R7:OAA	2.29	0.48
1:A:28:LEU:HD13	1:A:44:TYR:CD1	2.48	0.48
1:B:73:ARG:HG3	1:B:149:ASP:OD1	2.14	0.48
1:H:396:GLY:N	1:H:397:GLY:HA2	2.28	0.48
1:F:396:GLY:N	1:F:397:GLY:HA2	2.27	0.48
3:E:502:6R7:OAA	3:E:502:6R7:OAJ	2.27	0.48
1:F:224:THR:O	1:F:227:VAL:HG12	2.13	0.48
1:A:396:GLY:N	1:A:397:GLY:HA2	2.27	0.48
1:H:28:LEU:HD13	1:H:44:TYR:CD1	2.49	0.48
1:H:332:LEU:HG	1:H:333:PHE:N	2.28	0.48
1:F:28:LEU:HD13	1:F:44:TYR:CD1	2.49	0.47
1:A:116:ARG:HD3	1:A:400:TYR:CD1	2.49	0.47
1:A:45:VAL:O	1:A:49:MET:HG3	2.13	0.47
1:D:162:GLU:HG3	1:D:332:LEU:HD22	1.95	0.47
1:E:28:LEU:HD13	1:E:44:TYR:CD1	2.48	0.47
1:E:63:ASP:OD1	1:E:63:ASP:N	2.47	0.47
1:F:24:THR:O	1:F:28:LEU:HB2	2.14	0.47
1:C:126:LYS:HB3	1:C:129:MET:HE3	1.95	0.47
1:C:374:ASP:O	1:C:377:PRO:HD2	2.15	0.47
1:A:38:GLU:HG2	1:A:88:ASP:HB3	1.97	0.47
1:B:243:MET:HE3	1:B:246:LEU:HB3	1.96	0.47
1:G:204:TRP:CE2	1:G:267:GLY:HA2	2.50	0.47
1:A:38:GLU:OE2	1:A:156:CYS:HB2	2.14	0.46
1:B:211:PHE:HB2	1:B:332:LEU:HB3	1.96	0.46
1:B:49:MET:HE1	1:B:152:LEU:HB2	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:88:ASP:OD1	1:H:88:ASP:N	2.46	0.46
1:B:160:ASP:N	1:B:160:ASP:OD1	2.48	0.46
1:E:88:ASP:N	1:E:88:ASP:OD1	2.39	0.46
1:B:114:GLU:HG2	1:B:115:GLY:N	2.30	0.46
1:B:229:ARG:NH2	1:E:251:GLU:OE1	2.48	0.46
1:B:61:VAL:HG13	1:C:173:LEU:HD11	1.97	0.46
1:H:204:TRP:CE2	1:H:267:GLY:HA2	2.50	0.46
1:G:49:MET:HE1	1:G:152:LEU:HB2	1.97	0.46
1:H:126:LYS:N	1:H:127:GLY:HA3	2.30	0.46
1:D:224:THR:O	1:D:227:VAL:HG12	2.16	0.46
1:C:49:MET:HE1	1:C:152:LEU:HB2	1.98	0.46
1:F:58:ARG:HG2	1:F:66:PHE:CD1	2.51	0.46
1:B:162:GLU:HG3	1:B:332:LEU:HD22	1.98	0.45
1:C:414:ALA:O	1:C:418:MET:HG3	2.16	0.45
1:G:243:MET:HE3	1:G:246:LEU:HB3	1.97	0.45
1:F:191:TYR:OH	1:F:344:PRO:HG2	2.16	0.45
1:G:227:VAL:HA	1:G:228:PRO:HD3	1.85	0.45
1:B:63:ASP:OD1	1:B:63:ASP:N	2.45	0.45
1:F:225:PRO:HG2	1:G:371:MET:HB3	1.98	0.45
1:G:342:ILE:HD11	1:G:386:LEU:HD23	1.99	0.45
1:A:160:ASP:N	1:A:160:ASP:OD1	2.50	0.44
1:A:332:LEU:HG	1:A:333:PHE:N	2.31	0.44
1:C:126:LYS:HA	1:C:127:GLY:HA2	1.41	0.44
1:G:104:ASP:OD1	1:G:105:ARG:N	2.50	0.44
1:G:63:ASP:OD1	1:G:63:ASP:N	2.51	0.44
1:H:319:VAL:HG11	1:H:326:ALA:HB3	2.00	0.44
1:A:419:ASP:OD1	1:A:423:ARG:NH1	2.48	0.44
1:B:132:TRP:HA	1:B:135:ALA:HB3	1.99	0.44
1:H:116:ARG:HG2	1:H:402:LEU:HD23	2.00	0.44
1:H:49:MET:HE1	1:H:152:LEU:HB2	1.99	0.44
1:B:114:GLU:HG2	1:B:115:GLY:CA	2.48	0.44
1:E:18:ARG:NH2	1:E:404:ASP:OD1	2.37	0.44
1:C:127:GLY:HA3	1:C:130:ALA:H	1.82	0.44
1:G:319:VAL:HG11	1:G:326:ALA:HB3	2.00	0.44
1:H:80:SER:HB2	1:H:190:ASP:H	1.83	0.44
1:B:374:ASP:O	1:B:377:PRO:HD2	2.18	0.43
1:D:227:VAL:HA	1:D:228:PRO:HD3	1.82	0.43
1:G:4:SER:HB2	1:G:7:VAL:HG23	2.00	0.43
1:F:5:ASN:HB2	1:F:6:ASP:H	1.45	0.43
1:E:43:ASP:OD1	1:E:58:ARG:NH1	2.51	0.43
1:A:105:ARG:NH1	1:A:109:GLU:OE2	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:137:LYS:HE3	1:E:141:GLU:OE2	2.18	0.43
1:H:94:GLU:HG3	1:H:94:GLU:H	1.57	0.43
1:H:199:ASN:HB2	1:H:201:LYS:HE3	2.00	0.43
1:A:88:ASP:OD2	1:A:159:ILE:HG13	2.18	0.43
1:D:46:TYR:C	1:D:46:TYR:CD1	2.92	0.43
1:G:88:ASP:N	1:G:88:ASP:OD1	2.50	0.43
1:B:191:TYR:OH	1:B:344:PRO:HG2	2.19	0.43
1:C:58:ARG:HG2	1:C:66:PHE:CD1	2.53	0.43
1:F:45:VAL:O	1:F:49:MET:HG3	2.19	0.43
1:F:63:ASP:OD1	1:F:63:ASP:N	2.51	0.43
1:A:128:PRO:O	1:A:131:CYS:HB2	2.19	0.42
1:E:13:THR:O	1:E:17:MET:HG3	2.19	0.42
1:E:45:VAL:O	1:E:49:MET:HG3	2.18	0.42
1:F:283:VAL:HA	1:F:284:PRO:HD3	1.87	0.42
1:A:322:LEU:HA	1:A:322:LEU:HD23	1.88	0.42
1:F:199:ASN:HB2	1:F:201:LYS:HE3	2.01	0.42
1:F:80:SER:HB2	1:F:190:ASP:H	1.85	0.42
1:F:251:GLU:OE1	1:G:229:ARG:NH2	2.52	0.42
1:A:259:LYS:HB3	1:A:259:LYS:HE3	1.82	0.42
1:A:94:GLU:H	1:A:94:GLU:HG3	1.49	0.42
1:B:199:ASN:O	1:B:201:LYS:HG3	2.19	0.42
1:D:4:SER:HB2	1:D:5:ASN:OD1	2.19	0.42
1:B:289:ARG:NH1	3:E:502:6R7:OAL	2.52	0.42
1:B:374:ASP:C	1:B:377:PRO:HD2	2.40	0.42
1:G:116:ARG:HG2	1:G:402:LEU:HD23	2.02	0.42
1:H:9:LYS:HE3	1:H:142:ALA:O	2.20	0.42
1:C:322:LEU:HD23	1:C:322:LEU:HA	1.85	0.42
1:F:160:ASP:N	1:F:160:ASP:OD1	2.53	0.42
1:B:201:LYS:HA	1:B:202:PRO:HD3	1.90	0.42
1:A:72:LEU:HB3	1:A:150:VAL:HB	2.01	0.41
1:B:80:SER:HB2	1:B:189:SER:HB2	2.02	0.41
1:D:128:PRO:HD3	1:D:401:PHE:HE2	1.85	0.41
1:C:243:MET:HE3	1:C:246:LEU:HB3	2.02	0.41
1:D:424:THR:HA	1:D:425:PRO:HD3	1.94	0.41
1:H:424:THR:HA	1:H:425:PRO:HD3	1.95	0.41
1:B:94:GLU:HG3	1:B:94:GLU:H	1.60	0.41
1:D:198:THR:HG22	1:D:373:ARG:HH12	1.85	0.41
1:G:94:GLU:H	1:G:94:GLU:HG3	1.68	0.41
1:B:24:THR:O	1:B:28:LEU:HB2	2.20	0.41
1:C:319:VAL:HG11	1:C:326:ALA:HB3	2.02	0.41
1:D:211:PHE:HB3	1:D:331:PHE:CE1	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:398:ASN:OD1	1:D:399:THR:N	2.53	0.41
1:G:28:LEU:HD13	1:G:44:TYR:CD1	2.55	0.41
1:E:173:LEU:HD11	1:G:61:VAL:HG13	2.01	0.41
1:H:394:ALA:C	1:H:396:GLY:H	2.23	0.41
1:A:283:VAL:HA	1:A:284:PRO:HD3	1.87	0.41
1:D:281:GLY:HA2	1:D:282:GLY:HA2	1.90	0.41
1:F:201:LYS:HA	1:F:202:PRO:HD3	1.93	0.41
1:F:281:GLY:O	1:G:279:ILE:N	2.46	0.41
1:A:374:ASP:O	1:A:377:PRO:HD2	2.20	0.41
1:B:58:ARG:HG2	1:B:66:PHE:CD1	2.55	0.41
1:D:363:ARG:HG2	1:D:364:PRO:HD2	2.02	0.41
1:E:332:LEU:HG	1:E:333:PHE:N	2.34	0.41
1:A:337:TYR:CE1	1:A:380:GLU:HB2	2.56	0.41
1:C:13:THR:O	1:C:17:MET:HG3	2.21	0.41
1:C:225:PRO:HG2	1:D:371:MET:HB3	2.03	0.41
1:D:200:PHE:HB3	1:D:353:HIS:HE1	1.85	0.41
1:D:283:VAL:HA	1:D:284:PRO:HD3	1.82	0.41
1:B:28:LEU:HD13	1:B:44:TYR:CD1	2.56	0.41
1:C:137:LYS:HE3	1:C:141:GLU:OE2	2.21	0.41
1:C:160:ASP:OD1	1:C:160:ASP:N	2.54	0.41
1:E:190:ASP:O	1:E:384:PRO:HD2	2.21	0.41
1:D:329:LYS:NZ	1:H:57:GLU:OE2	2.32	0.41
1:C:281:GLY:HA2	1:C:282:GLY:HA2	1.88	0.40
1:E:160:ASP:OD1	1:E:160:ASP:N	2.53	0.40
1:E:199:ASN:O	1:E:201:LYS:HG3	2.21	0.40
1:F:88:ASP:N	1:F:88:ASP:OD1	2.54	0.40
1:C:45:VAL:O	1:C:49:MET:HG3	2.21	0.40
1:G:190:ASP:O	1:G:384:PRO:HD2	2.21	0.40
1:H:137:LYS:HE3	1:H:141:GLU:OE2	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	420/425 (99%)	398 (95%)	21 (5%)	1 (0%)	52	84
1	B	423/425 (100%)	400 (95%)	20 (5%)	3 (1%)	26	63
1	C	419/425 (99%)	396 (94%)	20 (5%)	3 (1%)	26	63
1	D	423/425 (100%)	400 (95%)	21 (5%)	2 (0%)	34	71
1	E	419/425 (99%)	399 (95%)	16 (4%)	4 (1%)	19	54
1	F	419/425 (99%)	397 (95%)	18 (4%)	4 (1%)	19	54
1	G	420/425 (99%)	400 (95%)	18 (4%)	2 (0%)	34	71
1	H	419/425 (99%)	396 (94%)	22 (5%)	1 (0%)	52	84
All	All	3362/3400 (99%)	3186 (95%)	156 (5%)	20 (1%)	30	67

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	114	GLU
1	F	114	GLU
1	G	115	GLY
1	A	114	GLU
1	D	104	ASP
1	E	104	ASP
1	F	104	ASP
1	H	104	ASP
1	B	104	ASP
1	D	160	ASP
1	E	105	ARG
1	F	160	ASP
1	B	160	ASP
1	C	160	ASP
1	C	218	ALA
1	E	160	ASP
1	E	218	ALA
1	G	160	ASP
1	B	3	GLY
1	F	115	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	330/331 (100%)	327 (99%)	3 (1%)	84	96
1	B	331/331 (100%)	329 (99%)	2 (1%)	90	97
1	C	329/331 (99%)	327 (99%)	2 (1%)	90	97
1	D	331/331 (100%)	329 (99%)	2 (1%)	90	97
1	E	329/331 (99%)	327 (99%)	2 (1%)	90	97
1	F	329/331 (99%)	326 (99%)	3 (1%)	84	96
1	G	330/331 (100%)	327 (99%)	3 (1%)	84	96
1	H	329/331 (99%)	324 (98%)	5 (2%)	72	92
All	All	2638/2648 (100%)	2616 (99%)	22 (1%)	86	96

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	114	GLU
1	A	116	ARG
1	A	160	ASP
1	B	160	ASP
1	B	327	GLU
1	C	160	ASP
1	C	293	LEU
1	D	19	GLU
1	D	293	LEU
1	E	160	ASP
1	E	293	LEU
1	F	5	ASN
1	F	160	ASP
1	F	363	ARG
1	G	98	ARG
1	G	160	ASP
1	G	293	LEU
1	H	12	LYS
1	H	101	ASP
1	H	160	ASP
1	H	293	LEU
1	H	359	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	256	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 16 ligands modelled in this entry, 8 are monoatomic - leaving 8 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	6R7	A	502	2	8,13,13	1.62	1 (12%)	13,18,18	1.31	3 (23%)
3	6R7	B	502	2	8,13,13	1.98	1 (12%)	13,18,18	1.35	3 (23%)
3	6R7	C	502	2	8,13,13	1.56	1 (12%)	13,18,18	1.31	1 (7%)
3	6R7	D	502	2	8,13,13	1.57	1 (12%)	13,18,18	0.96	1 (7%)
3	6R7	E	502	2	8,13,13	1.84	1 (12%)	13,18,18	1.77	5 (38%)
3	6R7	F	502	2	8,13,13	1.85	1 (12%)	13,18,18	1.59	2 (15%)
3	6R7	G	502	2	8,13,13	1.33	1 (12%)	13,18,18	1.36	1 (7%)
3	6R7	H	502	2	8,13,13	1.95	1 (12%)	13,18,18	2.45	4 (30%)

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	6R7	A	502	2	-	0/4/8/8	0/1/1/1
3	6R7	B	502	2	-	0/4/8/8	0/1/1/1
3	6R7	C	502	2	-	0/4/8/8	0/1/1/1
3	6R7	D	502	2	-	0/4/8/8	0/1/1/1
3	6R7	E	502	2	-	0/4/8/8	0/1/1/1
3	6R7	F	502	2	-	0/4/8/8	0/1/1/1
3	6R7	G	502	2	-	0/4/8/8	0/1/1/1
3	6R7	H	502	2	-	0/4/8/8	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	G	502	6R7	OAM-NAK	3.44	1.29	1.22
3	D	502	6R7	OAM-NAK	4.27	1.31	1.22
3	C	502	6R7	OAM-NAK	4.30	1.31	1.22
3	A	502	6R7	OAM-NAK	4.41	1.31	1.22
3	F	502	6R7	OAM-NAK	5.12	1.32	1.22
3	E	502	6R7	OAM-NAK	5.14	1.32	1.22
3	H	502	6R7	OAM-NAK	5.27	1.33	1.22
3	B	502	6R7	OAM-NAK	5.46	1.33	1.22

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	502	6R7	CAI-CAD-CAB	-4.03	117.54	121.64
3	H	502	6R7	CAI-CAD-CAB	-3.76	117.81	121.64
3	H	502	6R7	CAE-CAF-NAK	-3.56	115.61	118.74
3	H	502	6R7	CAG-CAH-CAI	-2.63	117.86	120.50
3	B	502	6R7	CAI-CAD-CAB	-2.23	119.37	121.64
3	E	502	6R7	CAG-CAH-CAI	-2.21	118.28	120.50
3	A	502	6R7	CAG-CAH-CAI	-2.16	118.34	120.50
3	E	502	6R7	CAI-CAD-CAB	-2.10	119.51	121.64
3	B	502	6R7	CAG-CAH-CAI	-2.07	118.43	120.50
3	E	502	6R7	CAE-CAD-CAI	2.02	120.25	117.63
3	D	502	6R7	CAG-CAF-NAK	2.05	120.98	119.51
3	A	502	6R7	CAE-CAD-CAI	2.06	120.30	117.63
3	F	502	6R7	CAE-CAD-CAI	2.26	120.56	117.63
3	B	502	6R7	CAG-CAF-NAK	2.47	121.29	119.51
3	A	502	6R7	CAG-CAF-NAK	2.60	121.38	119.51
3	C	502	6R7	CAG-CAF-NAK	3.13	121.76	119.51
3	G	502	6R7	CAG-CAF-NAK	3.26	121.85	119.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	502	6R7	OAM-NAK-CAF	3.32	121.06	118.67
3	E	502	6R7	CAG-CAF-NAK	3.43	121.97	119.51
3	H	502	6R7	CAG-CAF-NAK	5.84	123.70	119.51

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502	6R7	1	0
3	B	502	6R7	2	0
3	D	502	6R7	1	0
3	E	502	6R7	2	0
3	G	502	6R7	1	0
3	H	502	6R7	1	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	422/425 (99%)	-0.34	3 (0%) 89 88	26, 38, 60, 101	0
1	B	425/425 (100%)	-0.43	0 100 100	24, 34, 57, 92	0
1	C	421/425 (99%)	-0.50	2 (0%) 91 90	23, 37, 54, 101	0
1	D	425/425 (100%)	-0.34	4 (0%) 85 84	24, 34, 66, 92	0
1	E	421/425 (99%)	-0.54	0 100 100	24, 34, 53, 84	0
1	F	421/425 (99%)	-0.53	2 (0%) 91 90	24, 35, 52, 79	0
1	G	422/425 (99%)	-0.29	2 (0%) 91 90	26, 41, 62, 102	0
1	H	421/425 (99%)	-0.45	1 (0%) 95 95	25, 40, 62, 97	0
All	All	3378/3400 (99%)	-0.43	14 (0%) 93 92	23, 36, 61, 102	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	5	ASN	4.1
1	D	114	GLU	3.3
1	A	114	GLU	3.0
1	D	113	GLU	2.6
1	D	402	LEU	2.6
1	F	425	PRO	2.5
1	A	4	SER	2.4
1	C	5	ASN	2.3
1	D	4	SER	2.3
1	G	17	MET	2.1
1	F	6	ASP	2.1
1	G	115	GLY	2.1
1	C	6	ASP	2.1
1	H	5	ASN	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(Å ²)	Q<0.9
3	6R7	H	502	13/13	0.95	0.15	0.35	35,38,40,44	5
3	6R7	D	502	13/13	0.95	0.15	0.31	39,40,43,49	4
3	6R7	A	502	13/13	0.96	0.14	0.03	40,44,49,50	6
3	6R7	B	502	13/13	0.97	0.14	-0.04	37,40,46,46	6
3	6R7	E	502	13/13	0.96	0.13	-0.38	34,36,41,47	5
3	6R7	C	502	13/13	0.96	0.13	-0.47	34,39,42,51	4
3	6R7	G	502	13/13	0.96	0.12	-1.23	40,43,47,50	4
3	6R7	F	502	13/13	0.97	0.10	-1.50	32,33,34,38	4
2	MN	A	501	1/1	0.99	0.03	-2.91	37,37,37,37	0
2	MN	E	501	1/1	0.99	0.03	-3.12	38,38,38,38	0
2	MN	D	501	1/1	0.97	0.04	-3.15	38,38,38,38	0
2	MN	B	501	1/1	0.99	0.05	-3.99	41,41,41,41	0
2	MN	H	501	1/1	0.99	0.03	-4.14	39,39,39,39	0
2	MN	G	501	1/1	0.98	0.02	-5.14	41,41,41,41	0
2	MN	F	501	1/1	0.99	0.02	-5.35	33,33,33,33	0
2	MN	C	501	1/1	0.99	0.03	-6.00	38,38,38,38	0

6.5 Other polymers [i](#)

There are no such residues in this entry.