



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 08:25 AM UTC

PDB ID : 1SPA / pdb_00001spa
Title : ROLE OF ASP222 IN THE CATALYTIC MECHANISM OF ESCHERICHIA COLI ASPARTATE AMINOTRANSFERASE: THE AMINO ACID RESIDUE WHICH ENHANCES THE FUNCTION OF THE ENZYME-BOUND COENZYME PYRIDOXAL 5'-PHOSPHATE
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Deposited on : 1993-01-26
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

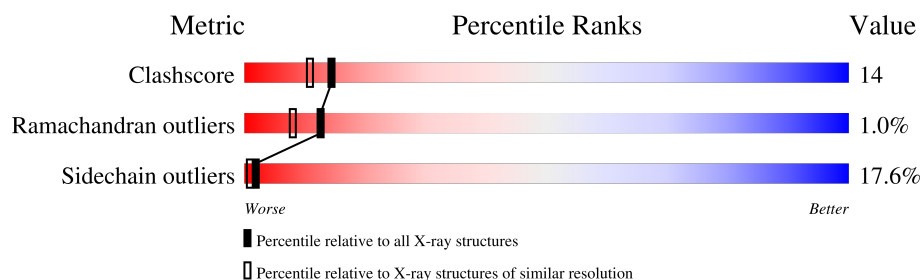
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	396	53% 32% 13% .

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3183 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.

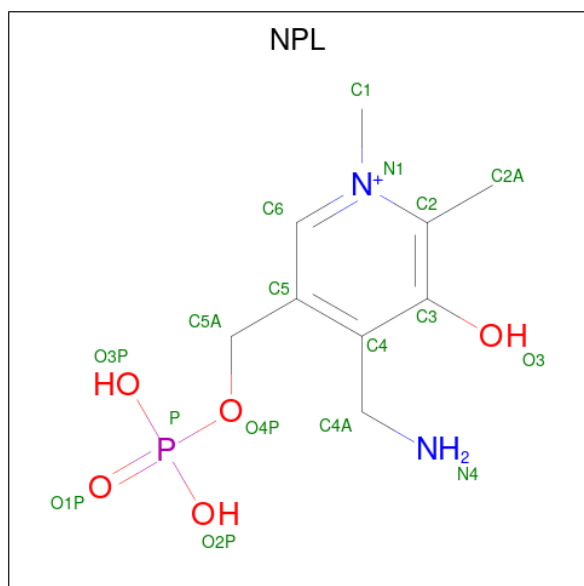
- Molecule 1 is a protein called ASPARTATE AMINOTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	396	3066	1935	536	582	13	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	222	ALA	ASP	conflict	UNP P00509

- Molecule 2 is N-METHYL-4-DEOXY-4-AMINO-PYRIDOXAL-5-PHOSPHATE (CCD ID: NPL) (formula: $C_9H_{16}N_2O_5P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	17	9	2	5	1	0	0

- Molecule 3 is water.

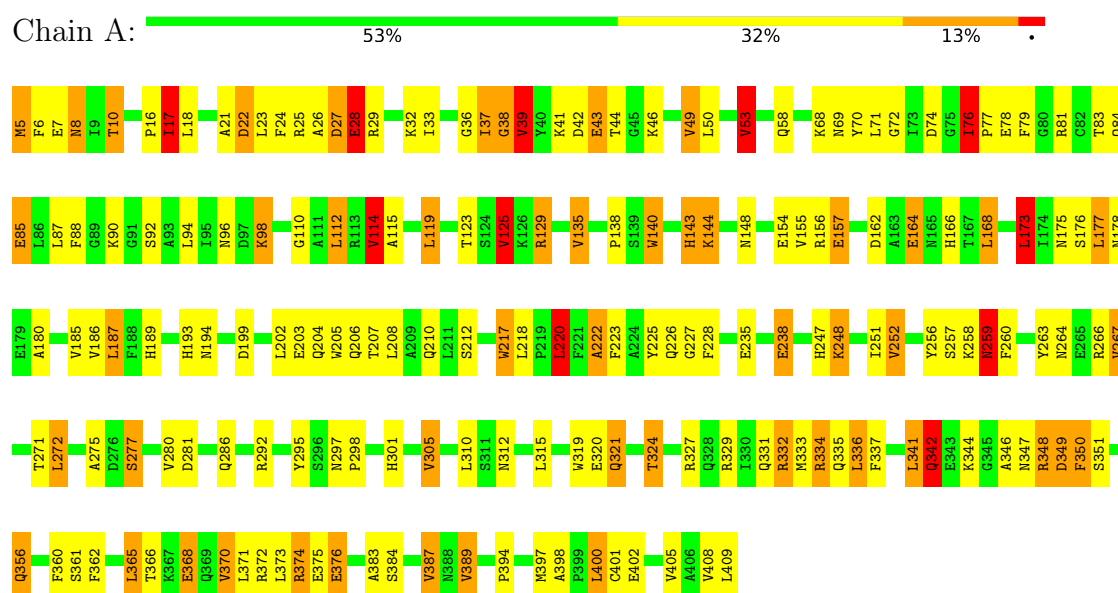
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	100	Total 100	O 100	0	0

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: ASPARTATE AMINOTRANSFERASE



4 Data and refinement statistics

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

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	154.80Å 86.20Å 79.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.209 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3183	wwPDB-VP
Average B, all atoms (Å ²)	28.0	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: NPL

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	1.28	17/3127 (0.5%)	2.07	123/4236 (2.9%)

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

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	368	GLU	CA-CB	9.16	1.67	1.53
1	A	17	ILE	CA-CB	8.56	1.64	1.54
1	A	143	HIS	CD2-NE2	-7.67	1.29	1.37
1	A	76	ILE	CA-CB	6.92	1.63	1.53
1	A	266	ARG	CA-C	-6.91	1.47	1.53
1	A	247	HIS	CD2-NE2	-6.48	1.30	1.37
1	A	49	VAL	CA-CB	6.46	1.62	1.54
1	A	166	HIS	CD2-NE2	-6.29	1.30	1.37
1	A	193	HIS	CG-ND1	-5.84	1.31	1.38
1	A	85	GLU	CA-CB	-5.76	1.44	1.53
1	A	189	HIS	CD2-NE2	-5.64	1.31	1.37
1	A	193	HIS	CD2-NE2	-5.61	1.31	1.37
1	A	114	VAL	CA-CB	5.60	1.61	1.54
1	A	301	HIS	CD2-NE2	-5.37	1.31	1.37
1	A	349	ASP	CA-CB	5.27	1.61	1.53
1	A	217	TRP	CG-CD2	-5.25	1.34	1.43
1	A	189	HIS	CG-ND1	-5.15	1.32	1.38

All (123) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	37	ILE	N-CA-C	13.67	124.57	110.62
1	A	259	ASN	OD1-CG-ND2	-11.20	111.40	122.60
1	A	349	ASP	N-CA-C	-11.07	90.36	108.41
1	A	349	ASP	CA-CB-CG	10.62	123.22	112.60
1	A	350	PHE	CA-CB-CG	-10.45	103.35	113.80
1	A	259	ASN	CA-CB-CG	9.62	122.22	112.60
1	A	387	VAL	N-CA-CB	-9.25	95.29	111.39
1	A	368	GLU	CA-CB-CG	9.19	132.48	114.10
1	A	115	ALA	N-CA-C	-9.15	101.25	111.14
1	A	321	GLN	CG-CD-NE2	8.88	129.71	116.40
1	A	342	GLN	OE1-CD-NE2	-8.73	113.87	122.60
1	A	58	GLN	OE1-CD-NE2	-8.67	113.93	122.60
1	A	387	VAL	CB-CA-C	8.59	123.41	110.62
1	A	222	ALA	N-CA-C	-8.47	94.92	108.73
1	A	10	THR	CA-CB-OG1	-8.36	97.06	109.60
1	A	199	ASP	CA-CB-CG	8.23	120.83	112.60
1	A	148	ASN	OD1-CG-ND2	-7.97	114.63	122.60
1	A	210	GLN	OE1-CD-NE2	-7.87	114.73	122.60
1	A	148	ASN	CB-CG-ND2	7.82	128.13	116.40
1	A	368	GLU	CB-CG-CD	7.75	125.77	112.60
1	A	42	ASP	CA-CB-CG	7.52	120.12	112.60
1	A	312	ASN	OD1-CG-ND2	-7.49	115.11	122.60
1	A	321	GLN	OE1-CD-NE2	-7.34	115.25	122.60
1	A	301	HIS	N-CA-C	7.30	118.93	110.97
1	A	368	GLU	N-CA-C	-7.26	103.36	111.28
1	A	286	GLN	OE1-CD-NE2	-7.23	115.37	122.60
1	A	27	ASP	CA-C-O	7.23	128.44	120.92
1	A	74	ASP	CA-CB-CG	7.20	119.80	112.60
1	A	374	ARG	CA-CB-CG	7.01	128.12	114.10
1	A	334	ARG	NE-CZ-NH2	-6.95	112.95	119.20
1	A	257	SER	CA-C-O	-6.93	113.54	120.82
1	A	259	ASN	CB-CG-ND2	6.87	126.70	116.40
1	A	228	PHE	N-CA-C	6.86	121.65	113.28
1	A	375	GLU	CA-CB-CG	6.86	127.83	114.10
1	A	204	GLN	OE1-CD-NE2	-6.86	115.74	122.60
1	A	228	PHE	CA-CB-CG	-6.66	107.14	113.80
1	A	39	VAL	N-CA-CB	-6.65	101.30	112.47
1	A	226	GLN	OE1-CD-NE2	-6.53	116.07	122.60
1	A	356	GLN	OE1-CD-NE2	-6.52	116.08	122.60
1	A	332	ARG	CB-CA-C	-6.52	99.77	110.85
1	A	125	VAL	O-C-N	6.51	130.38	122.83
1	A	72	GLY	N-CA-C	-6.46	104.27	112.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	264	ASN	CA-CB-CG	6.40	119.00	112.60
1	A	8	ASN	OD1-CG-ND2	-6.36	116.24	122.60
1	A	337	PHE	N-CA-C	-6.27	104.36	111.14
1	A	238	GLU	N-CA-CB	-6.25	100.59	110.22
1	A	10	THR	CA-CB-CG2	6.23	121.09	110.50
1	A	166	HIS	CB-CG-CD2	-6.23	123.10	131.20
1	A	7	GLU	CA-C-N	-6.20	113.93	123.17
1	A	7	GLU	C-N-CA	-6.20	113.93	123.17
1	A	53	VAL	N-CA-C	-6.19	103.15	111.44
1	A	58	GLN	N-CA-C	-6.17	104.64	111.36
1	A	217	TRP	CA-CB-CG	6.13	125.24	113.60
1	A	226	GLN	CG-CD-NE2	6.12	125.58	116.40
1	A	334	ARG	NE-CZ-NH1	6.09	127.59	121.50
1	A	49	VAL	CB-CA-C	6.08	119.10	110.84
1	A	175	ASN	CA-CB-CG	6.01	118.61	112.60
1	A	68	LYS	N-CA-C	-5.92	104.44	112.26
1	A	332	ARG	N-CA-CB	5.91	118.91	110.16
1	A	173	LEU	CA-C-O	-5.90	114.83	120.90
1	A	319	TRP	CG-CD2-CE3	5.89	139.79	133.90
1	A	389	VAL	N-CA-CB	-5.87	102.55	110.54
1	A	76	ILE	CB-CA-C	5.84	116.37	110.94
1	A	342	GLN	CG-CD-NE2	5.80	125.11	116.40
1	A	123	THR	CA-CB-OG1	-5.78	100.94	109.60
1	A	305	VAL	O-C-N	-5.74	116.28	121.91
1	A	342	GLN	CB-CG-CD	5.74	122.35	112.60
1	A	335	GLN	OE1-CD-NE2	-5.72	116.88	122.60
1	A	43	GLU	CB-CG-CD	5.68	122.25	112.60
1	A	297	ASN	N-CA-C	-5.66	103.12	108.22
1	A	154	GLU	CB-CG-CD	5.66	122.22	112.60
1	A	247	HIS	CA-CB-CG	-5.64	108.16	113.80
1	A	178	ASN	OD1-CG-ND2	-5.63	116.97	122.60
1	A	341	LEU	N-CA-C	-5.63	104.53	111.40
1	A	71	LEU	N-CA-C	-5.62	103.27	110.53
1	A	347	ASN	CB-CG-ND2	5.62	124.83	116.40
1	A	162	ASP	CA-CB-CG	5.62	118.22	112.60
1	A	376	GLU	N-CA-C	5.62	119.30	112.23
1	A	401	CYS	CA-CB-SG	-5.61	101.50	114.40
1	A	405	VAL	CA-CB-CG2	-5.60	100.88	110.40
1	A	38	GLY	O-C-N	-5.59	115.58	122.23
1	A	298	PRO	CB-CA-C	5.58	117.72	110.92
1	A	267	VAL	N-CA-CB	-5.57	104.39	111.64
1	A	6	PHE	CA-CB-CG	5.54	119.34	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	49	VAL	N-CA-C	-5.54	100.44	108.36
1	A	374	ARG	CB-CG-CD	-5.52	98.61	111.30
1	A	143	HIS	CB-CG-CD2	-5.51	124.04	131.20
1	A	84	GLN	OE1-CD-NE2	-5.49	117.11	122.60
1	A	173	LEU	N-CA-C	-5.46	104.96	111.03
1	A	39	VAL	O-C-N	-5.40	117.19	122.97
1	A	331	GLN	CA-C-O	-5.39	113.77	119.97
1	A	81	ARG	CA-CB-CG	5.38	124.87	114.10
1	A	119	LEU	N-CA-C	-5.36	105.52	111.36
1	A	260	PHE	N-CA-C	-5.35	105.58	113.61
1	A	37	ILE	CB-CA-C	-5.34	104.94	112.14
1	A	275	ALA	N-CA-C	5.32	117.83	111.71
1	A	96	ASN	CB-CG-ND2	5.31	124.37	116.40
1	A	301	HIS	CB-CG-CD2	-5.28	124.34	131.20
1	A	194	ASN	O-C-N	-5.25	116.52	121.35
1	A	187	LEU	CD1-CG-CD2	-5.25	99.25	110.80
1	A	272	LEU	CB-CA-C	-5.24	101.60	110.19
1	A	194	ASN	N-CA-CB	-5.24	103.37	110.23
1	A	204	GLN	CG-CD-NE2	5.23	124.25	116.40
1	A	112	LEU	O-C-N	-5.22	116.69	122.07
1	A	205	TRP	CE2-CD2-CG	-5.20	100.95	107.20
1	A	347	ASN	OD1-CG-ND2	-5.20	117.40	122.60
1	A	281	ASP	CA-CB-CG	5.18	117.78	112.60
1	A	347	ASN	O-C-N	5.18	129.96	123.23
1	A	334	ARG	CG-CD-NE	-5.16	100.65	112.00
1	A	248	LYS	N-CA-C	-5.16	105.35	111.69
1	A	220	LEU	O-C-N	-5.14	116.33	123.12
1	A	140	TRP	CG-CD2-CE3	5.10	139.00	133.90
1	A	389	VAL	O-C-N	-5.10	116.92	121.87
1	A	247	HIS	CB-CG-CD2	-5.09	124.58	131.20
1	A	350	PHE	N-CA-C	-5.09	106.76	113.17
1	A	310	LEU	N-CA-C	5.06	117.53	111.71
1	A	372	ARG	CA-CB-CG	-5.05	104.00	114.10
1	A	226	GLN	CA-CB-CG	-5.04	104.02	114.10
1	A	28	GLU	CB-CG-CD	5.03	121.15	112.60
1	A	115	ALA	CA-C-N	5.03	127.33	120.54
1	A	115	ALA	C-N-CA	5.03	127.33	120.54
1	A	7	GLU	CB-CG-CD	5.01	121.12	112.60
1	A	238	GLU	CA-CB-CG	5.01	124.11	114.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	295	TYR	Sidechain

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	3066	0	3019	86	2
2	A	17	0	14	1	0
3	A	100	0	0	3	0
All	All	3183	0	3033	86	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 14.

All (86) 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:348:ARG:HG3	1:A:350:PHE:CD2	2.06	0.88
1:A:28:GLU:HG3	1:A:29:ARG:HD2	1.58	0.84
1:A:258:LYS:HD3	1:A:263:TYR:HE1	1.43	0.83
1:A:50:LEU:HB2	1:A:53:VAL:HG13	1.61	0.82
1:A:43:GLU:HG2	1:A:329:ARG:HD2	1.62	0.80
1:A:22:ASP:HA	1:A:25:ARG:HG2	1.65	0.78
1:A:222:ALA:HB1	2:A:413:NPL:H23	1.66	0.75
1:A:79:PHE:O	1:A:83:THR:HG23	1.87	0.74
1:A:37:ILE:HD13	1:A:41:LYS:HD2	1.70	0.72
1:A:258:LYS:HD3	1:A:263:TYR:CE1	2.30	0.66
1:A:173:LEU:HD23	1:A:177:LEU:HD23	1.78	0.66
1:A:90:LYS:H	1:A:90:LYS:HD2	1.63	0.64
1:A:320:GLU:O	1:A:324:THR:HG23	1.98	0.64
1:A:125:VAL:HG21	1:A:185:VAL:HG23	1.79	0.64
1:A:17:ILE:HD13	1:A:36:GLY:O	1.98	0.63
1:A:348:ARG:HG3	1:A:350:PHE:HD2	1.61	0.63
1:A:22:ASP:HA	1:A:25:ARG:CG	2.30	0.61
1:A:203:GLU:O	1:A:207:THR:HG23	2.00	0.61
1:A:235:GLU:O	1:A:238:GLU:HB3	2.01	0.60
1:A:373:LEU:HD13	1:A:408:VAL:HG21	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:GLY:O	1:A:114:VAL:HG13	2.03	0.57
1:A:252:VAL:HG13	1:A:271:THR:HB	1.88	0.56
1:A:212:SER:HA	1:A:217:TRP:CD1	2.42	0.55
1:A:227:GLY:O	1:A:327:ARG:HD3	2.07	0.55
1:A:21:ALA:O	1:A:24:PHE:HB3	2.08	0.54
1:A:251:ILE:HG12	1:A:272:LEU:HD23	1.90	0.54
1:A:356:GLN:NE2	1:A:361:SER:HB2	2.22	0.54
1:A:374:ARG:HG2	1:A:374:ARG:HH11	1.73	0.53
1:A:362:PHE:CE1	1:A:384:SER:HB2	2.44	0.53
1:A:37:ILE:CD1	1:A:41:LYS:HD2	2.38	0.52
1:A:346:ALA:H	1:A:348:ARG:NH1	2.07	0.52
1:A:98:LYS:HA	1:A:98:LYS:HZ2	1.75	0.51
1:A:341:LEU:HD13	1:A:350:PHE:CD2	2.45	0.51
1:A:38:GLY:HA2	1:A:360:PHE:CZ	2.45	0.51
1:A:334:ARG:HH22	1:A:361:SER:HB3	1.75	0.51
1:A:356:GLN:HE22	1:A:361:SER:HB2	1.75	0.51
1:A:83:THR:HG22	1:A:256:TYR:OH	2.11	0.50
1:A:140:TRP:HB3	1:A:143:HIS:CE1	2.47	0.50
1:A:344:LYS:NZ	1:A:398:ALA:HB1	2.25	0.50
1:A:144:LYS:HB3	1:A:144:LYS:NZ	2.27	0.50
1:A:187:LEU:HD21	1:A:222:ALA:HB2	1.93	0.50
1:A:53:VAL:HB	1:A:305:VAL:HG11	1.94	0.50
1:A:370:VAL:HG21	1:A:383:ALA:HA	1.94	0.49
1:A:38:GLY:HA2	1:A:360:PHE:HZ	1.78	0.49
1:A:256:TYR:HA	1:A:259:ASN:ND2	2.27	0.49
1:A:27:ASP:HB3	1:A:32:LYS:HD3	1.93	0.49
1:A:129:ARG:NH2	1:A:156:ARG:HD3	2.28	0.48
1:A:185:VAL:HG11	1:A:220:LEU:HD12	1.95	0.48
1:A:256:TYR:HB2	1:A:267:VAL:HG13	1.96	0.48
1:A:256:TYR:HA	1:A:259:ASN:HD21	1.79	0.47
1:A:180:ALA:HB3	1:A:217:TRP:CH2	2.49	0.47
1:A:76:ILE:O	1:A:79:PHE:HB3	2.15	0.47
1:A:346:ALA:HB3	1:A:348:ARG:NH2	2.30	0.46
1:A:85:GLU:HB3	1:A:90:LYS:HE3	1.97	0.46
1:A:258:LYS:HE2	3:A:518:HOH:O	2.16	0.46
1:A:76:ILE:O	1:A:76:ILE:HG13	2.14	0.45
1:A:324:THR:HG22	3:A:421:HOH:O	2.16	0.45
1:A:129:ARG:HH21	1:A:156:ARG:HD3	1.81	0.45
1:A:135:VAL:O	1:A:157:GLU:HA	2.16	0.45
1:A:366:THR:O	1:A:370:VAL:HG13	2.17	0.45
1:A:225:TYR:CE1	1:A:258:LYS:HG3	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:LEU:HD13	1:A:143:HIS:HD2	1.82	0.45
1:A:202:LEU:HD13	1:A:238:GLU:HG2	1.99	0.45
1:A:168:LEU:HD21	1:A:173:LEU:HD12	1.99	0.44
1:A:277:SER:HA	1:A:280:VAL:HG12	1.98	0.44
1:A:140:TRP:HE3	1:A:143:HIS:NE2	2.15	0.44
1:A:177:LEU:HD13	1:A:177:LEU:HA	1.79	0.44
1:A:344:LYS:HD3	1:A:344:LYS:HA	1.62	0.44
1:A:39:VAL:HG22	1:A:263:TYR:CE1	2.53	0.44
1:A:336:LEU:HB3	1:A:397:MET:HE3	2.00	0.44
1:A:342:GLN:HE21	1:A:342:GLN:N	2.15	0.44
1:A:334:ARG:HG2	1:A:389:VAL:HG11	1.99	0.44
1:A:333:MET:HE1	1:A:394:PRO:HG3	2.00	0.43
1:A:76:ILE:HA	1:A:77:PRO:HD3	1.88	0.43
1:A:33:ILE:HD13	1:A:400:LEU:HB2	2.00	0.43
1:A:186:VAL:HG12	1:A:217:TRP:HB3	2.01	0.43
1:A:356:GLN:NE2	1:A:361:SER:CB	2.82	0.42
1:A:321:GLN:NE2	3:A:424:HOH:O	2.52	0.42
1:A:144:LYS:HG2	1:A:155:VAL:HG11	2.02	0.41
1:A:342:GLN:HA	1:A:348:ARG:NH1	2.36	0.41
1:A:69:ASN:ND2	1:A:70:TYR:H	2.18	0.41
1:A:365:LEU:HD12	1:A:365:LEU:HA	1.80	0.40
1:A:90:LYS:HD2	1:A:90:LYS:N	2.33	0.40
1:A:334:ARG:HH22	1:A:361:SER:CB	2.35	0.40
1:A:88:PHE:O	1:A:92:SER:HB2	2.22	0.40
1:A:346:ALA:HB3	1:A:348:ARG:CZ	2.52	0.40

All (2) 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:A:5:MET:SD	1:A:368:GLU:CB[6_555]	1.93	0.27
1:A:5:MET:N	1:A:368:GLU:OE1[6_555]	1.96	0.24

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	394/396 (100%)	369 (94%)	21 (5%)	4 (1%)	12	8

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	28	GLU
1	A	26	ALA
1	A	164	GLU
1	A	16	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	319/319 (100%)	263 (82%)	56 (18%)	2	1

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	MET
1	A	8	ASN
1	A	10	THR
1	A	17	ILE
1	A	18	LEU
1	A	22	ASP
1	A	23	LEU
1	A	39	VAL
1	A	44	THR
1	A	46	LYS
1	A	49	VAL
1	A	53	VAL
1	A	76	ILE
1	A	78	GLU
1	A	87	LEU

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Mol	Chain	Res	Type
1	A	94	LEU
1	A	98	LYS
1	A	114	VAL
1	A	119	LEU
1	A	125	VAL
1	A	129	ARG
1	A	135	VAL
1	A	138	PRO
1	A	144	LYS
1	A	157	GLU
1	A	164	GLU
1	A	168	LEU
1	A	173	LEU
1	A	176	SER
1	A	177	LEU
1	A	206	GLN
1	A	208	LEU
1	A	218	LEU
1	A	220	LEU
1	A	223	PHE
1	A	248	LYS
1	A	252	VAL
1	A	259	ASN
1	A	277	SER
1	A	292	ARG
1	A	315	LEU
1	A	324	THR
1	A	332	ARG
1	A	336	LEU
1	A	342	GLN
1	A	348	ARG
1	A	349	ASP
1	A	351	SER
1	A	365	LEU
1	A	370	VAL
1	A	371	LEU
1	A	376	GLU
1	A	387	VAL
1	A	400	LEU
1	A	402	GLU
1	A	409	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	63	ASN
1	A	69	ASN
1	A	84	GLN
1	A	166	HIS
1	A	193	HIS
1	A	194	ASN
1	A	247	HIS
1	A	259	ASN
1	A	321	GLN
1	A	356	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 [i](#)

1 ligand is 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	$\# Z > 2$
2	NPL	A	413	-	17,17,17	1.32	3 (17%)	23,25,25	2.21	5 (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
2	NPL	A	413	-	-	3/8/8/8	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	413	NPL	C2-N1	2.72	1.40	1.36
2	A	413	NPL	P-O4P	-2.63	1.52	1.60
2	A	413	NPL	P-O2P	-2.40	1.45	1.54

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	413	NPL	C3-C2-N1	5.68	121.35	118.35
2	A	413	NPL	O2P-P-O4P	-4.36	95.29	106.67
2	A	413	NPL	C2A-C2-C3	-3.97	116.83	120.82
2	A	413	NPL	C2-C3-C4	-3.33	118.96	120.52
2	A	413	NPL	O4P-P-O1P	2.73	113.81	106.44

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	413	NPL	C3-C4-C4A-N4
2	A	413	NPL	C5-C4-C4A-N4
2	A	413	NPL	C5A-O4P-P-O1P

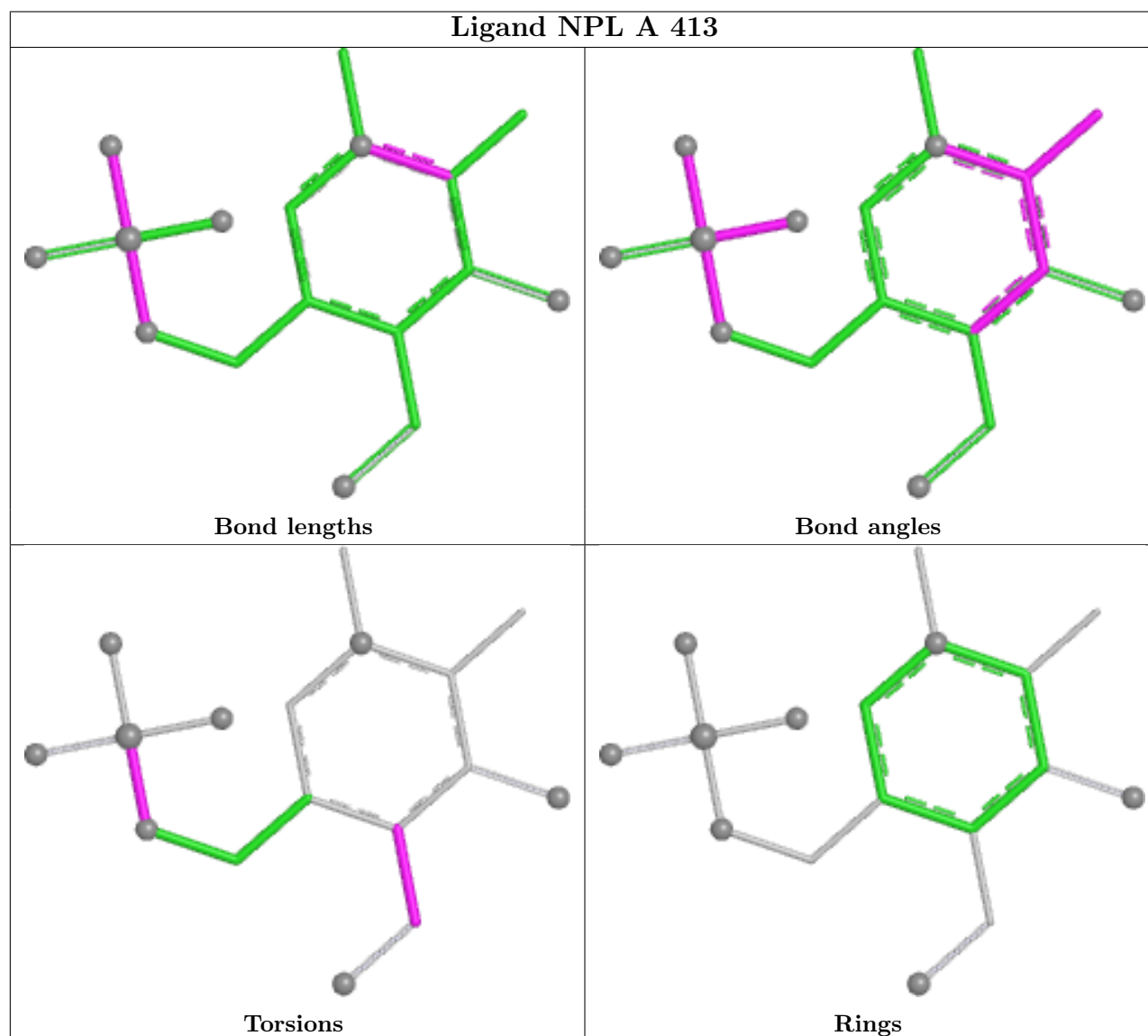
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	413	NPL	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be

highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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.