



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:05 PM UTC

PDB ID : 5NHF / pdb_00005nhf
Title : Human Erk2 with an Erk1/2 inhibitor
Authors : Debreczeni, J.E.; Ward, R.A.; Bethel, P.; Cook, C.; Davies, E.; Eckersley, K.; Fairley, G.; Feron, L.; Flemington, V.; Graham, M.A.; Greenwood, R.; Hopcroft, P.; Howard, T.D.; Hudson, J.; James, M.; Jones, C.D.; Jones, C.R.; Lamont, S.; Lewis, R.; Lindsay, N.; Roberts, K.; Simpson, I.; StGallay, S.; Swallow, S.; Tonge, M.
Deposited on : 2017-03-21
Resolution : 2.14 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)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)

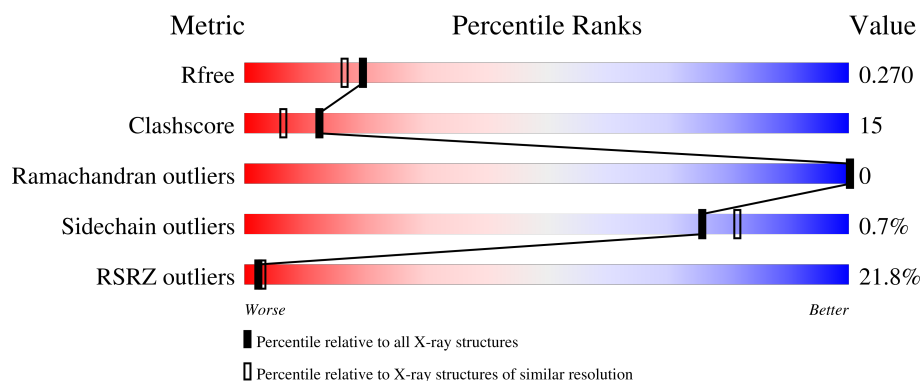
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.14 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}	180053	3689 (2.16-2.12)
Clashscore	190562	3812 (2.16-2.12)
Ramachandran outliers	187476	3773 (2.16-2.12)
Sidechain outliers	187428	3772 (2.16-2.12)
RSRZ outliers	180081	3691 (2.16-2.12)

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 $\leq 5\%$. 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	381	19% 62% 23% • 11%

Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2767 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 Mitogen-activated protein kinase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	340	Total	C	N	O	S	0	0	0
			2659	1715	446	484	14			

There are 22 discrepancies between the modelled and reference sequences:

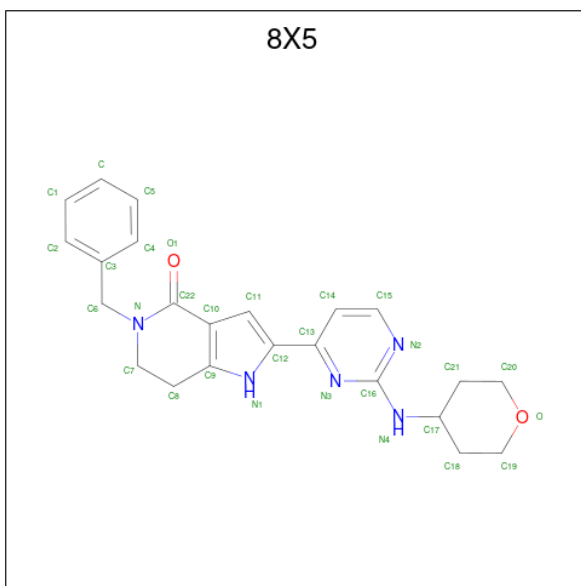
Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	MET	-	initiating methionine	UNP P28482
A	-17	HIS	-	expression tag	UNP P28482
A	-16	HIS	-	expression tag	UNP P28482
A	-15	HIS	-	expression tag	UNP P28482
A	-14	HIS	-	expression tag	UNP P28482
A	-13	HIS	-	expression tag	UNP P28482
A	-12	HIS	-	expression tag	UNP P28482
A	-11	GLY	-	expression tag	UNP P28482
A	-10	GLY	-	expression tag	UNP P28482
A	-9	GLY	-	expression tag	UNP P28482
A	-8	GLU	-	expression tag	UNP P28482
A	-7	ASN	-	expression tag	UNP P28482
A	-6	LEU	-	expression tag	UNP P28482
A	-5	TYR	-	expression tag	UNP P28482
A	-4	PHE	-	expression tag	UNP P28482
A	-3	GLN	-	expression tag	UNP P28482
A	-2	GLY	-	expression tag	UNP P28482
A	-1	SER	-	expression tag	UNP P28482
A	0	HIS	-	expression tag	UNP P28482
A	46	LEU	VAL	conflict	UNP P28482
A	361	LEU	-	expression tag	UNP P28482
A	362	GLU	-	expression tag	UNP P28482

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is 2-[2-(oxan-4-ylamino)pyrimidin-4-yl]-5-(phenylmethyl)-6,7-dihydro-1 {H}-pyrrolo[3,2-c]pyridin-4-one (CCD ID: 8X5) (formula: C₂₃H₂₅N₅O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			30	23	5	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	73	Total 73	O 73	0	0

- Molecule 1: Mitogen-activated protein kinase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	49.02Å 70.05Å 60.36Å 90.00° 109.31° 90.00°	Depositor
Resolution (Å)	44.20 – 2.14 44.20 – 2.14	Depositor EDS
% Data completeness (in resolution range)	98.0 (44.20-2.14) 98.0 (44.20-2.14)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 2.14Å)	Xtriage
Refinement program	BUSTER 2.11.6	Depositor
R, R_{free}	0.247 , 0.259 0.258 , 0.270	Depositor DCC
R_{free} test set	1312 reflections (6.27%)	wwPDB-VP
Wilson B-factor (Å ²)	40.7	Xtriage
Anisotropy	0.471	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 58.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2767	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.47% 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: CME, SO4, 8X5

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	0.69	20/2713 (0.7%)	1.07	31/3698 (0.8%)

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	92	ALA	C-N	7.55	1.41	1.33
1	A	279	PHE	C-N	6.55	1.41	1.34
1	A	225	ARG	C-N	6.26	1.41	1.33
1	A	322	GLU	C-N	6.15	1.41	1.33
1	A	64	TYR	CA-C	-6.07	1.44	1.52
1	A	267	LEU	C-N	5.89	1.41	1.33
1	A	246	SER	C-N	5.67	1.41	1.33
1	A	273	VAL	C-N	5.63	1.41	1.33
1	A	195	ALA	C-N	5.49	1.40	1.34
1	A	196	PRO	N-CD	5.48	1.55	1.47
1	A	280	PRO	N-CD	5.45	1.55	1.47
1	A	93	PRO	N-CD	5.43	1.55	1.47
1	A	339	PRO	N-CD	5.43	1.55	1.47
1	A	323	PRO	N-CD	5.40	1.55	1.47
1	A	247	PRO	N-CD	5.34	1.55	1.47
1	A	226	PRO	N-CD	5.30	1.55	1.47
1	A	268	PRO	N-CD	5.28	1.55	1.47
1	A	311	PRO	N-CD	5.27	1.55	1.47
1	A	274	PRO	N-CD	5.25	1.55	1.47
1	A	152	PRO	N-CD	5.21	1.55	1.47

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	179	ASP	N-CA-C	9.14	121.24	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	235	ASP	N-CA-C	-7.53	103.07	111.28
1	A	37	GLY	N-CA-C	7.36	124.64	110.66
1	A	322	GLU	CA-C-N	-6.92	113.17	120.03
1	A	322	GLU	C-N-CA	-6.92	113.17	120.03
1	A	246	SER	CA-C-N	-6.85	112.93	119.85
1	A	246	SER	C-N-CA	-6.85	112.93	119.85
1	A	64	TYR	N-CA-C	-6.78	104.82	113.23
1	A	267	LEU	CA-C-N	-6.78	113.00	119.85
1	A	267	LEU	C-N-CA	-6.78	113.00	119.85
1	A	92	ALA	CA-C-N	-6.67	113.05	120.45
1	A	92	ALA	C-N-CA	-6.67	113.05	120.45
1	A	279	PHE	CA-C-N	-6.65	112.86	120.04
1	A	279	PHE	C-N-CA	-6.65	112.86	120.04
1	A	225	ARG	CA-C-N	-6.54	113.03	119.90
1	A	225	ARG	C-N-CA	-6.54	113.03	119.90
1	A	48	LYS	N-CA-C	6.53	120.56	111.74
1	A	273	VAL	CA-C-N	-6.49	113.27	119.76
1	A	273	VAL	C-N-CA	-6.49	113.27	119.76
1	A	63	THR	O-C-N	-6.44	114.81	122.15
1	A	200	LEU	N-CA-C	6.42	121.38	111.56
1	A	191	ARG	N-CA-C	6.17	118.01	111.28
1	A	167	ASP	N-CA-C	6.05	118.00	110.91
1	A	28	LEU	N-CA-C	5.80	118.74	110.10
1	A	233	TYR	N-CA-C	5.65	120.70	111.37
1	A	195	ALA	CA-C-N	-5.62	113.23	119.19
1	A	195	ALA	C-N-CA	-5.62	113.23	119.19
1	A	59	PHE	N-CA-C	5.53	119.65	113.01
1	A	227	ILE	N-CA-C	5.33	116.10	110.72
1	A	82	ASN	N-CA-C	5.20	118.45	112.57
1	A	96	GLU	N-CA-C	5.20	116.94	111.28

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2659	0	2553	81	1
2	A	5	0	0	0	0
3	A	30	0	0	0	0
4	A	73	0	0	0	0
All	All	2767	0	2553	81	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 15.

All (81) 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:56:ILE:HG22	1:A:58:PRO:HD3	1.14	1.10
1:A:56:ILE:HG22	1:A:58:PRO:CD	1.99	0.93
1:A:54:LYS:HE3	1:A:56:ILE:HD11	1.64	0.79
1:A:56:ILE:CG2	1:A:58:PRO:HD3	2.07	0.77
1:A:179:ASP:OD1	1:A:205:TYR:N	2.18	0.76
1:A:98:MET:O	1:A:344:LYS:NZ	2.18	0.76
1:A:253:ASN:HA	1:A:261:ARG:HH22	1.54	0.71
1:A:89:ILE:HG21	1:A:351:THR:HG22	1.71	0.71
1:A:302:ILE:HD11	1:A:306:GLN:HG2	1.74	0.69
1:A:184:LEU:HD23	1:A:233:TYR:HE1	1.58	0.67
1:A:54:LYS:CE	1:A:56:ILE:HD11	2.25	0.66
1:A:253:ASN:HA	1:A:261:ARG:NH2	2.12	0.64
1:A:81:GLU:HB2	1:A:135:ARG:HH12	1.64	0.63
1:A:184:LEU:CD2	1:A:233:TYR:HE1	2.11	0.63
1:A:227:ILE:HG23	1:A:228:PHE:CD2	2.34	0.62
1:A:302:ILE:HG13	1:A:306:GLN:CD	2.23	0.62
1:A:89:ILE:HG21	1:A:351:THR:CG2	2.29	0.62
1:A:89:ILE:HD12	1:A:354:PHE:CE1	2.34	0.61
1:A:73:LYS:HE2	1:A:350:GLU:OE2	2.02	0.60
1:A:116:LEU:HD13	1:A:220:GLU:HG2	1.84	0.60
1:A:117:LYS:HG3	1:A:118:THR:HG23	1.85	0.59
1:A:187:TYR:CE2	1:A:189:ALA:HB3	2.37	0.59
1:A:289:LEU:HG	1:A:293:MET:HE2	1.85	0.59
1:A:111:ASP:OD1	1:A:113:TYR:HB3	2.02	0.59
1:A:75:LEU:CB	1:A:86:ILE:HD12	2.33	0.57
1:A:234:LEU:O	1:A:234:LEU:HG	2.04	0.57
1:A:42:ALA:HB3	1:A:53:ILE:CD1	2.35	0.56
1:A:114:LYS:O	1:A:117:LYS:HG2	2.07	0.55
1:A:179:ASP:OD1	1:A:205:TYR:HB2	2.07	0.55
1:A:75:LEU:HB2	1:A:86:ILE:CD1	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:21:VAL:HG12	1:A:102:TYR:CD2	2.42	0.54
1:A:42:ALA:HB3	1:A:53:ILE:HD12	1.89	0.54
1:A:95:ILE:HG22	1:A:96:GLU:OE1	2.07	0.54
1:A:75:LEU:HB3	1:A:86:ILE:HD12	1.89	0.54
1:A:243:ILE:HD12	1:A:278:LEU:HD11	1.89	0.53
1:A:161:CME:HA	1:A:161:CME:HE2	1.91	0.52
1:A:191:ARG:HA	1:A:194:ARG:HG3	1.92	0.52
1:A:36:TYR:CD1	1:A:36:TYR:N	2.78	0.52
1:A:36:TYR:CD2	1:A:58:PRO:HB3	2.45	0.52
1:A:28:LEU:HA	1:A:41:SER:O	2.09	0.51
1:A:43:TYR:HA	1:A:50:ARG:HD2	1.92	0.51
1:A:173:VAL:CG2	1:A:332:ASP:HA	2.42	0.50
1:A:62:GLN:NE2	1:A:338:LEU:O	2.31	0.49
1:A:89:ILE:HD13	1:A:351:THR:HG22	1.94	0.49
1:A:56:ILE:C	1:A:58:PRO:HD3	2.37	0.49
1:A:200:LEU:HD21	1:A:237:LEU:HD21	1.95	0.48
1:A:223:SER:HB3	1:A:279:PHE:HE1	1.78	0.48
1:A:19:PHE:CD2	1:A:28:LEU:HD21	2.49	0.48
1:A:187:TYR:HE2	1:A:189:ALA:HB3	1.78	0.48
1:A:36:TYR:HD2	1:A:58:PRO:HB3	1.78	0.48
1:A:75:LEU:HB2	1:A:86:ILE:HD12	1.95	0.48
1:A:95:ILE:HG23	1:A:96:GLU:HG3	1.96	0.48
1:A:184:LEU:CD2	1:A:233:TYR:CE1	2.95	0.47
1:A:223:SER:O	1:A:224:ASN:HB2	2.12	0.47
1:A:151:LYS:HE3	1:A:154:ASN:ND2	2.29	0.47
1:A:95:ILE:HG23	1:A:96:GLU:N	2.29	0.47
1:A:63:THR:HG23	1:A:64:TYR:CD1	2.50	0.46
1:A:115:LEU:HD12	1:A:119:GLN:HG3	1.98	0.45
1:A:179:ASP:OD2	1:A:202:SER:OG	2.31	0.45
1:A:21:VAL:CG1	1:A:102:TYR:CD2	2.99	0.45
1:A:36:TYR:H	1:A:36:TYR:HD1	1.64	0.45
1:A:223:SER:HB3	1:A:279:PHE:CE1	2.51	0.45
1:A:90:ILE:HB	1:A:102:TYR:HB2	1.97	0.45
1:A:118:THR:OG1	1:A:119:GLN:HG2	2.16	0.45
1:A:218:LEU:HD22	1:A:293:MET:HE1	1.98	0.44
1:A:75:LEU:CB	1:A:86:ILE:CD1	2.94	0.44
1:A:71:GLU:OE2	1:A:167:ASP:HA	2.18	0.43
1:A:184:LEU:HD22	1:A:233:TYR:CE1	2.52	0.43
1:A:232:HIS:CD2	1:A:235:ASP:HB2	2.53	0.42
1:A:36:TYR:HB3	1:A:64:TYR:CD2	2.54	0.42
1:A:41:SER:OG	1:A:50:ARG:NH2	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:327:ALA:HA	1:A:328:PRO:HD2	1.93	0.42
1:A:21:VAL:HG11	1:A:53:ILE:HG21	2.01	0.42
1:A:232:HIS:CE1	1:A:235:ASP:HB2	2.55	0.41
1:A:163:LEU:C	1:A:163:LEU:HD23	2.44	0.41
1:A:247:PRO:HG2	1:A:252:LEU:HD21	2.02	0.41
1:A:59:PHE:CZ	1:A:98:MET:HE3	2.55	0.41
1:A:146:LEU:HD22	1:A:207:LYS:HA	2.03	0.41
1:A:232:HIS:NE2	1:A:235:ASP:HB2	2.35	0.40
1:A:36:TYR:CD2	1:A:64:TYR:HB3	2.57	0.40
1:A:43:TYR:CD1	1:A:43:TYR:C	2.99	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:96:GLU:OE2	1:A:265:LEU:CD1[1_454]	2.05	0.15

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	335/381 (88%)	327 (98%)	8 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	276/334 (83%)	274 (99%)	2 (1%)	76	81

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	PHE
1	A	36	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	ASN
1	A	97	GLN
1	A	119	GLN
1	A	123	ASN
1	A	238	ASN
1	A	269	HIS
1	A	281	ASN
1	A	355	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue 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	$\# Z > 2$
1	CME	A	161	1	8,9,10	0.93	0	6,9,11	0.49	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
1	CME	A	161	1	-	0/5/8/10	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	161	CME	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	$\# Z > 2$
2	SO4	A	401	-	4,4,4	0.27	0	6,6,6	0.12	0
3	8X5	A	402	-	33,34,34	0.29	0	40,47,47	0.70	1 (2%)

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	8X5	A	402	-	-	2/12/33/33	0/5/5/5

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	402	8X5	C7-C8-C9	-2.26	104.44	108.52

There are no chirality outliers.

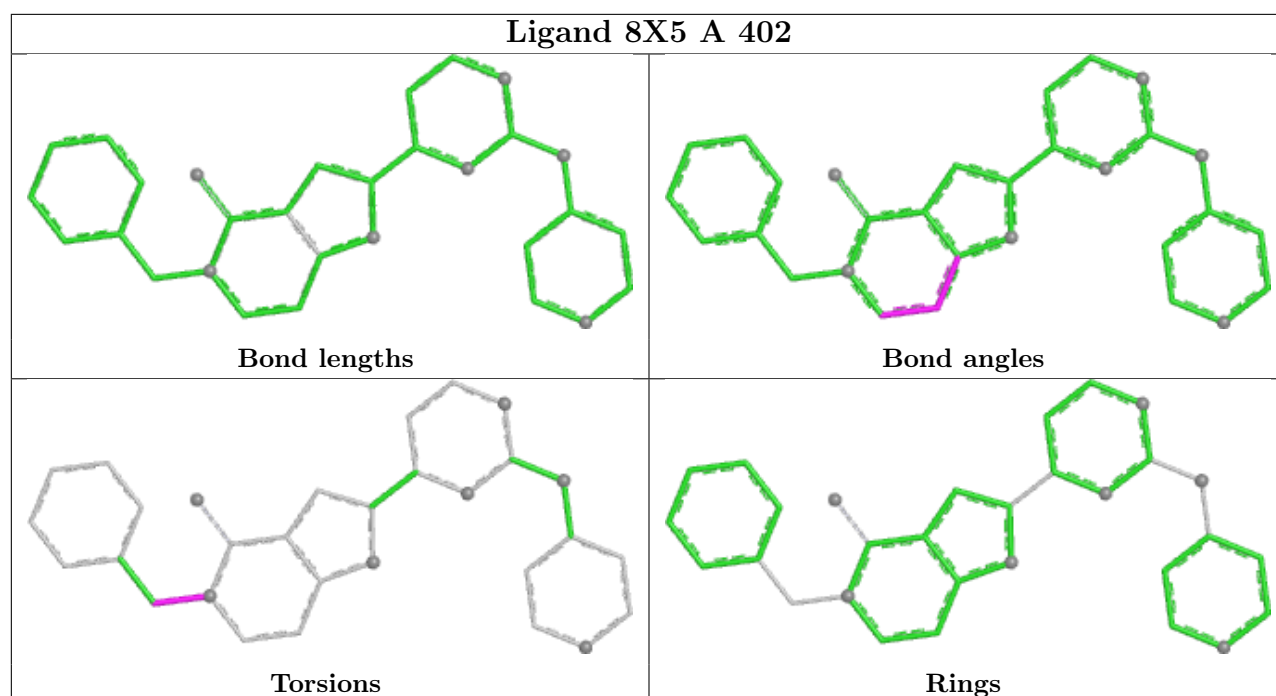
All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	402	8X5	C3-C6-N-C22
3	A	402	8X5	C3-C6-N-C7

There are no ring outliers.

No monomer is involved in short contacts.

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 ⓘ

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	339/381 (88%)	1.33	74 (21%) 2 3	27, 47, 79, 105	0

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	331	PHE	5.2
1	A	332	ASP	4.9
1	A	40	CYS	4.8
1	A	16	GLY	4.8
1	A	14	VAL	4.2
1	A	350	GLU	4.1
1	A	234	LEU	4.0
1	A	338	LEU	3.8
1	A	21	VAL	3.8
1	A	31	ILE	3.7
1	A	20	ASP	3.7
1	A	188	VAL	3.6
1	A	61	HIS	3.5
1	A	106	ASP	3.4
1	A	329	PHE	3.4
1	A	355	GLN	3.4
1	A	187	TYR	3.3
1	A	39	VAL	3.1
1	A	38	MET	3.1
1	A	95	ILE	3.1
1	A	56	ILE	3.0
1	A	209	ILE	3.0
1	A	354	PHE	3.0
1	A	339	PRO	2.9
1	A	34	GLY	2.9
1	A	59	PHE	2.8
1	A	179	ASP	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	112	LEU	2.7
1	A	22	GLY	2.7
1	A	254	CYS	2.7
1	A	337	ASP	2.7
1	A	45	ASN	2.6
1	A	347	ILE	2.6
1	A	88	ASP	2.6
1	A	33	GLU	2.5
1	A	26	THR	2.5
1	A	19	PHE	2.5
1	A	351	THR	2.5
1	A	37	GLY	2.5
1	A	30	TYR	2.4
1	A	43	TYR	2.4
1	A	102	TYR	2.4
1	A	94	THR	2.4
1	A	121	LEU	2.4
1	A	230	GLY	2.4
1	A	251	ASP	2.4
1	A	58	PRO	2.3
1	A	101	VAL	2.3
1	A	252	LEU	2.3
1	A	89	ILE	2.2
1	A	352	ALA	2.2
1	A	327	ALA	2.2
1	A	233	TYR	2.2
1	A	27	ASN	2.2
1	A	103	ILE	2.2
1	A	274	PRO	2.2
1	A	18	VAL	2.2
1	A	186	GLU	2.2
1	A	258	LEU	2.2
1	A	15	ARG	2.2
1	A	36	TYR	2.2
1	A	305	GLU	2.1
1	A	306	GLN	2.1
1	A	175	ASP	2.1
1	A	303	GLU	2.1
1	A	97	GLN	2.1
1	A	32	GLY	2.1
1	A	256	ILE	2.1
1	A	120	HIS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	206	THR	2.1
1	A	269	HIS	2.0
1	A	118	THR	2.0
1	A	42	ALA	2.0
1	A	261	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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. 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	B-factors(Å ²)	Q<0.9
1	CME	A	161	10/11	0.83	0.19	41,47,56,57	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

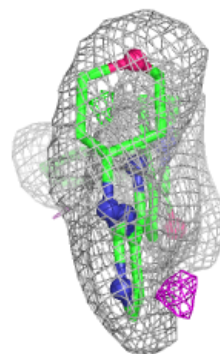
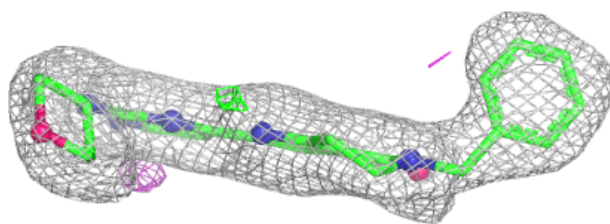
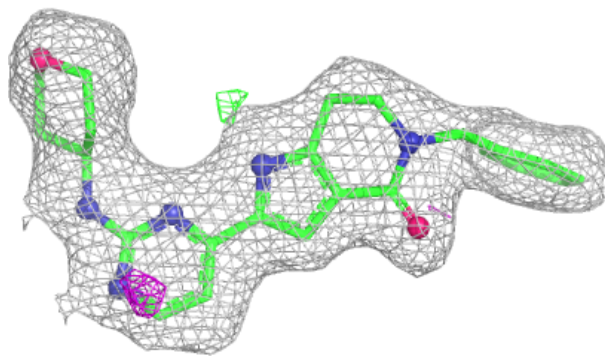
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SO4	A	401	5/5	0.82	0.12	82,83,83,84	0
3	8X5	A	402	30/30	0.89	0.12	40,45,51,51	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around 8X5 A 402:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.